



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 01:03 AM UTC

PDB ID : 6YFO / pdb_00006yfo
Title : Virus-like particle of bacteriophage GQ-907
Authors : Rumnieks, J.; Kalnins, G.; Sisovs, M.; Lieknina, I.; Tars, K.
Deposited on : 2020-03-26
Resolution : 3.48 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

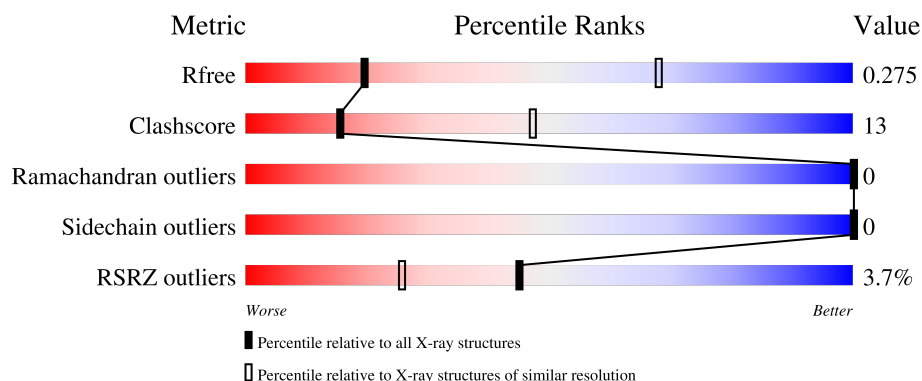
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.48 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1083 (3.52-3.44)
Clashscore	190562	1139 (3.52-3.44)
Ramachandran outliers	187476	1111 (3.52-3.44)
Sidechain outliers	187428	1112 (3.52-3.44)
RSRZ outliers	180081	1082 (3.52-3.44)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	AA	131	5% 79% 21%
1	AB	131	2% 68% 32%
1	AC	131	5% 73% 27%
1	AD	131	5% 80% 20%
1	AE	131	2% 68% 32%

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Mol	Chain	Length	Quality of chain
1	AF	131	5% 73% 27%
1	AG	131	4% 79% 21%
1	AH	131	5% 68% 32%
1	AI	131	8% 69% 31%
1	AJ	131	5% 79% 21%
1	AK	131	2% 69% 31%
1	AL	131	2% 76% 24%
1	AM	131	3% 79% 21%
1	AN	131	2% 69% 31%
1	AO	131	5% 69% 31%
1	AP	131	5% 82% 18%
1	AQ	131	5% 71% 29%
1	AR	131	3% 73% 27%
1	AS	131	3% 80% 20%
1	AT	131	2% 68% 32%
1	AU	131	11% 73% 27%
1	AV	131	2% 80% 20%
1	AW	131	2% 69% 31%
1	AX	131	8% 73% 27%
1	AY	131	3% 82% 18%
1	AZ	131	2% 71% 29%
1	BA	131	2% 75% 25%
1	BB	131	2% 79% 21%
1	BC	131	2% 70% 30%
1	BD	131	2% 76% 24%

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Mol	Chain	Length	Quality of chain	
1	BE	131	4%	79% 21%
1	BF	131	4%	70% 30%
1	BG	131	4%	74% 26%
1	BH	131	10%	82% 18%
1	BI	131	5%	70% 30%
1	BJ	131	2%	76% 24%
1	BK	131	6%	80% 20%
1	BL	131		69% 31%
1	BM	131	9%	69% 31%
1	BN	131	5%	83% 17%
1	BO	131	5%	69% 31%
1	BP	131	5%	73% 27%
1	BQ	131	5%	83% 17%
1	BR	131	4%	69% 31%
1	BS	131	2%	78% 22%
1	BT	131	%	81% 19%
1	BU	131	5%	69% 31%
1	BV	131	5%	73% 27%
1	BW	131	3%	79% 21%
1	BX	131	2%	67% 33%
1	BY	131	5%	69% 31%
1	BZ	131	5%	82% 18%
1	CA	131		68% 32%
1	CB	131	2%	76% 24%
1	CC	131	2%	77% 23%

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Mol	Chain	Length	Quality of chain	
1	CD	131	5%	66% 34%
1	CE	131	5%	69% 31%
1	CF	131	6%	80% 20%
1	CG	131	2%	69% 31%
1	CH	131	8%	79% 21%
1	CI	131	2%	77% 23%
1	CJ	131	4%	69% 31%
1	CK	131	7%	74% 26%
1	CL	131	2%	80% 20%
1	CM	131	2%	69% 31%
1	CN	131	2%	75% 25%
1	CO	131	2%	82% 18%
1	CP	131		70% 30%
1	CQ	131	2%	81% 19%
1	CR	131	5%	81% 19%
1	CS	131	4%	69% 31%
1	CT	131	8%	74% 26%
1	CU	131	2%	80% 20%
1	CV	131	2%	67% 33%
1	CW	131	2%	76% 24%
1	CX	131	4%	80% 20%
1	CY	131	2%	70% 30%
1	CZ	131	%	76% 24%
1	DA	131	2%	79% 21%
1	DB	131	2%	69% 31%

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Mol	Chain	Length	Quality of chain
1	DC	131	 2% 74% 26%
1	DD	131	 7% 82% 18%
1	DE	131	 6% 67% 33%
1	DF	131	 5% 77% 23%
1	DG	131	 2% 78% 22%
1	DH	131	 3% 67% 33%
1	DI	131	 8% 69% 31%
1	DJ	131	 0% 82% 18%
1	DK	131	 4% 71% 29%
1	DL	131	 1% 80% 20%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 87480 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called coat protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	AA	131	Total	C	N	O	S	0	0	0
			972	607	166	198	1			
1	AB	131	Total	C	N	O	S	0	0	0
			972	607	166	198	1			
1	AC	131	Total	C	N	O	S	0	0	0
			972	607	166	198	1			
1	AD	131	Total	C	N	O	S	0	0	0
			972	607	166	198	1			
1	AE	131	Total	C	N	O	S	0	0	0
			972	607	166	198	1			
1	AF	131	Total	C	N	O	S	0	0	0
			972	607	166	198	1			
1	AG	131	Total	C	N	O	S	0	0	0
			972	607	166	198	1			
1	AH	131	Total	C	N	O	S	0	0	0
			972	607	166	198	1			
1	AI	131	Total	C	N	O	S	0	0	0
			972	607	166	198	1			
1	AJ	131	Total	C	N	O	S	0	0	0
			972	607	166	198	1			
1	AK	131	Total	C	N	O	S	0	0	0
			972	607	166	198	1			
1	AL	131	Total	C	N	O	S	0	0	0
			972	607	166	198	1			
1	AM	131	Total	C	N	O	S	0	0	0
			972	607	166	198	1			
1	AN	131	Total	C	N	O	S	0	0	0
			972	607	166	198	1			
1	AO	131	Total	C	N	O	S	0	0	0
			972	607	166	198	1			
1	AP	131	Total	C	N	O	S	0	0	0
			972	607	166	198	1			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	AQ	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	AR	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	AS	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	AT	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	AU	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	AV	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	AW	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	AX	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	AY	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	AZ	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	BA	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	BB	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	BC	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	BD	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	BE	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	BF	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	BG	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	BH	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	BI	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	BJ	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	BK	131	Total 972	C 607	N 166	O 198	S 1	0	0	0

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	BL	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	BM	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	BN	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	BO	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	BP	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	BQ	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	BR	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	BS	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	BT	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	BU	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	BV	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	BW	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	BX	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	BY	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	BZ	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	CA	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	CB	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	CC	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	CD	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	CE	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	CF	131	Total 972	C 607	N 166	O 198	S 1	0	0	0

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	CG	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	CH	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	CI	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	CJ	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	CK	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	CL	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	CM	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	CN	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	CO	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	CP	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	CQ	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	CR	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	CS	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	CT	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	CU	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	CV	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	CW	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	CX	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	CY	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	CZ	131	Total 972	C 607	N 166	O 198	S 1	0	0	0
1	DA	131	Total 972	C 607	N 166	O 198	S 1	0	0	0

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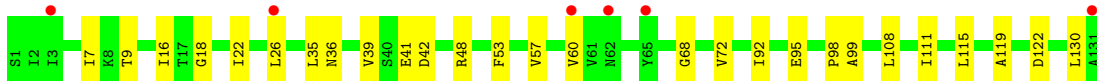
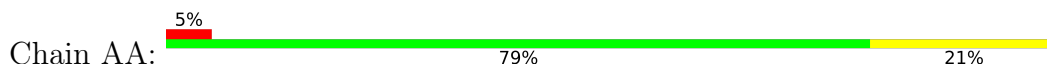
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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	DB	131	Total	C	N	O	S	0	0	0
			972	607	166	198	1			
1	DC	131	Total	C	N	O	S	0	0	0
			972	607	166	198	1			
1	DD	131	Total	C	N	O	S	0	0	0
			972	607	166	198	1			
1	DE	131	Total	C	N	O	S	0	0	0
			972	607	166	198	1			
1	DF	131	Total	C	N	O	S	0	0	0
			972	607	166	198	1			
1	DG	131	Total	C	N	O	S	0	0	0
			972	607	166	198	1			
1	DH	131	Total	C	N	O	S	0	0	0
			972	607	166	198	1			
1	DI	131	Total	C	N	O	S	0	0	0
			972	607	166	198	1			
1	DJ	131	Total	C	N	O	S	0	0	0
			972	607	166	198	1			
1	DK	131	Total	C	N	O	S	0	0	0
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1	DL	131	Total	C	N	O	S	0	0	0
			972	607	166	198	1			

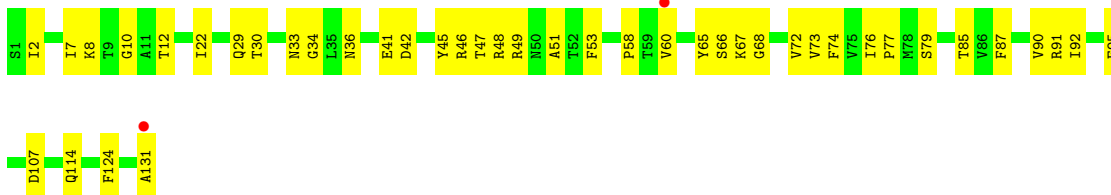
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

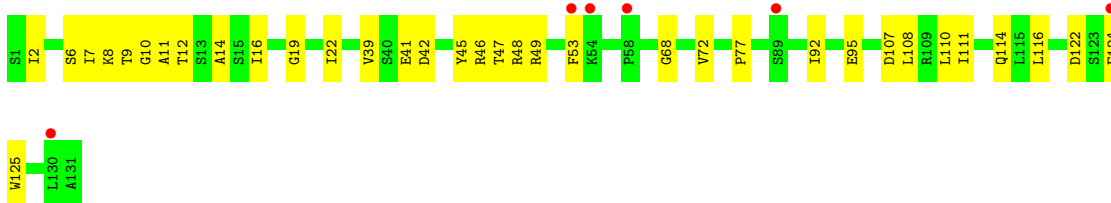
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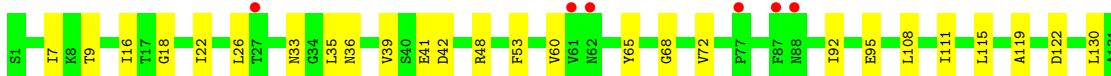
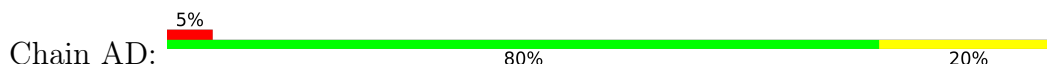
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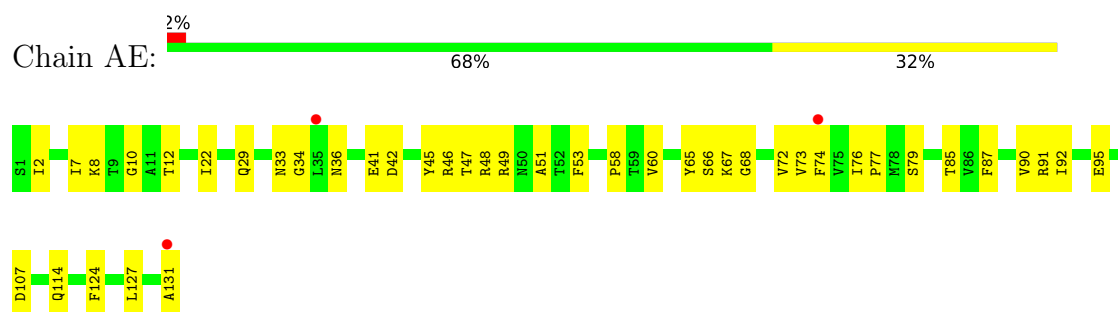
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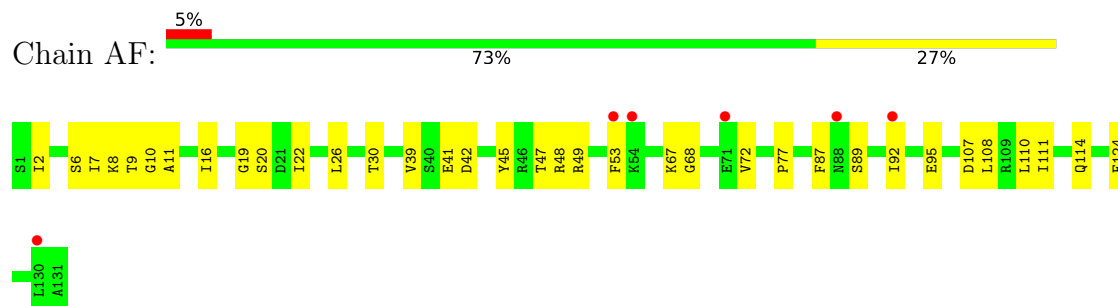
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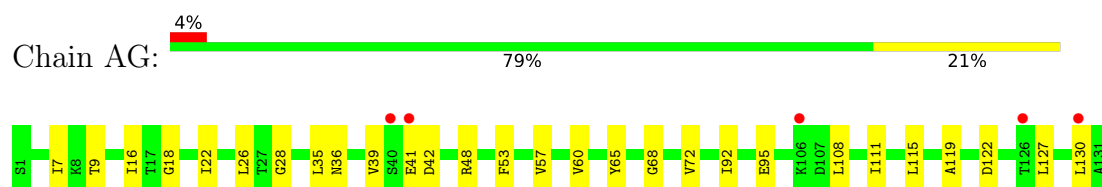
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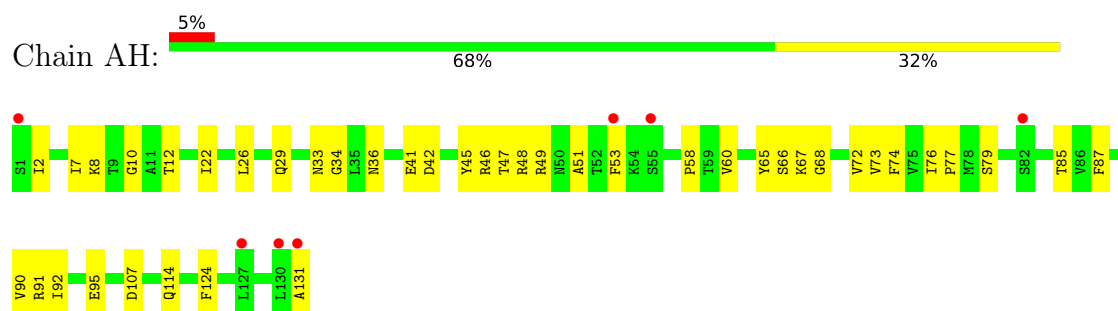
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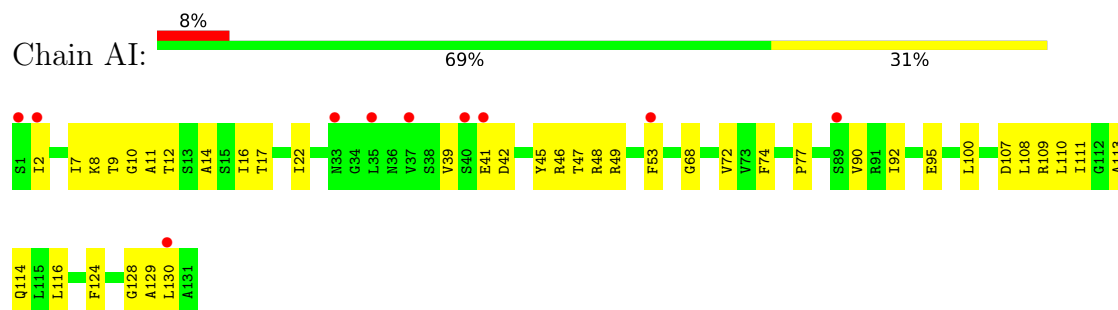
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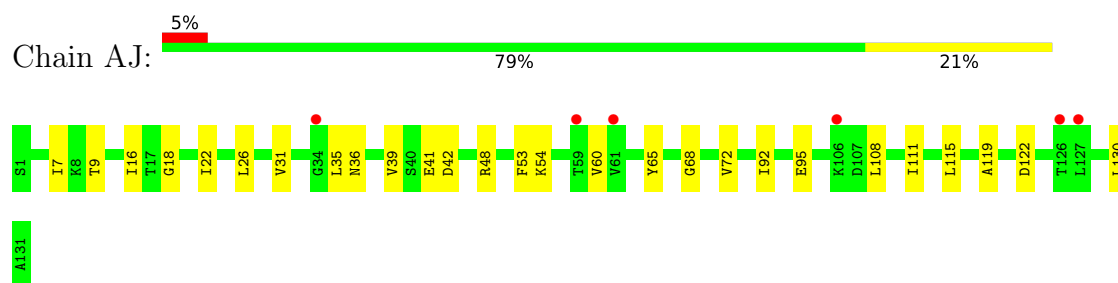
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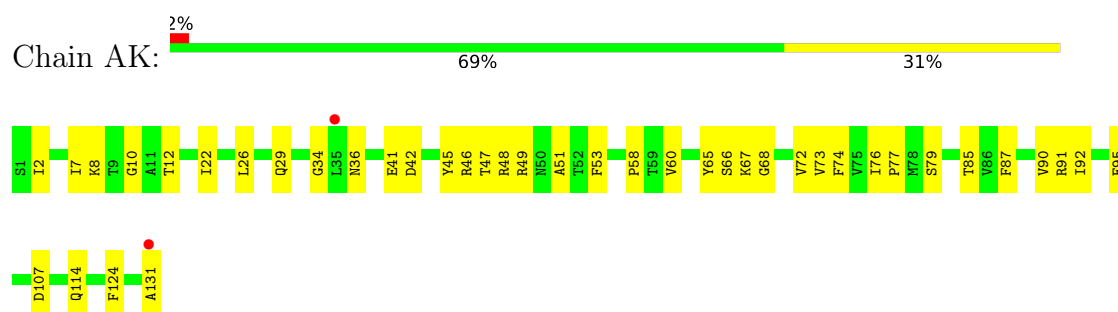
- Molecule 1: coat protein



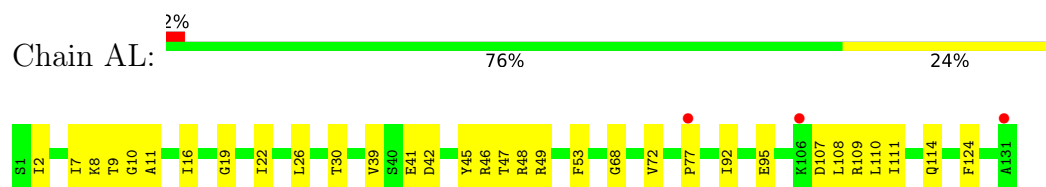
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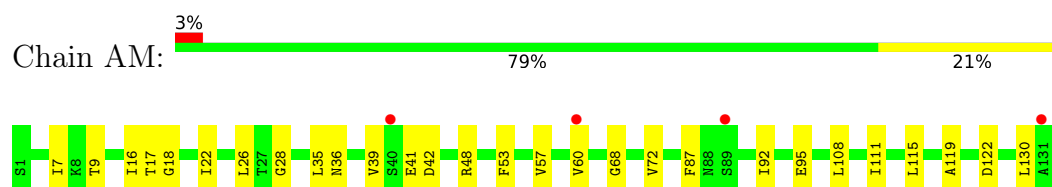
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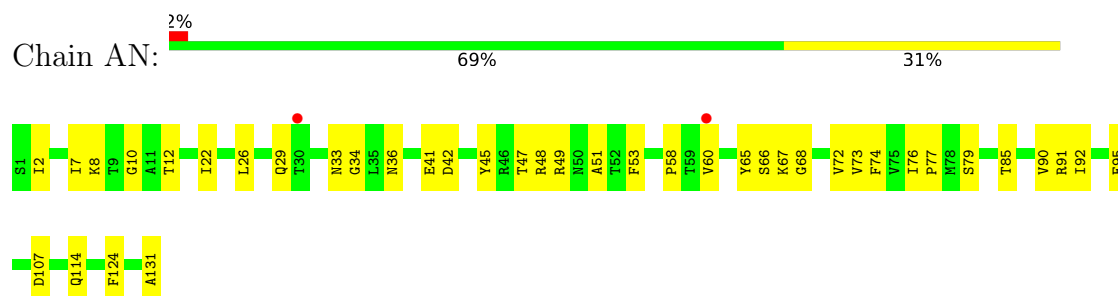
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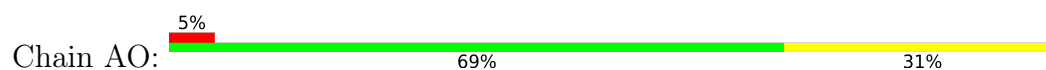
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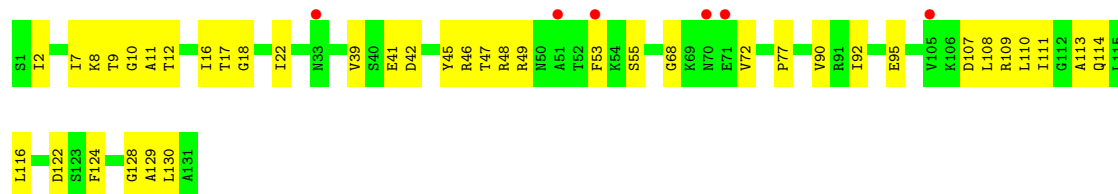


- Molecule 1: coat protein

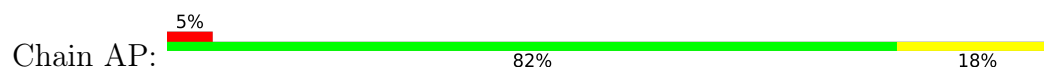


- Molecule 1: coat protein

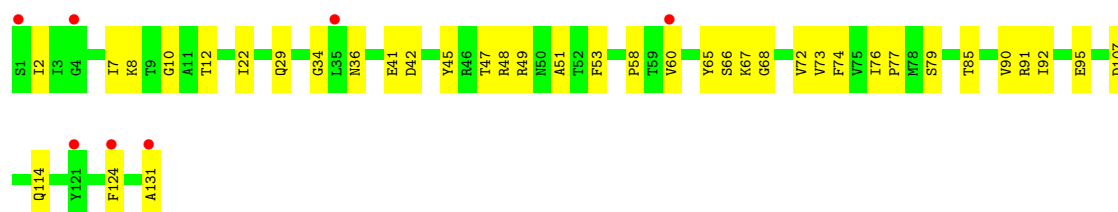
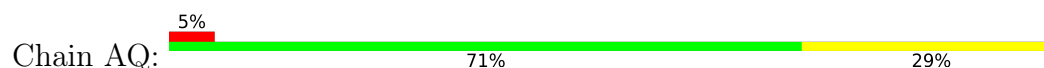




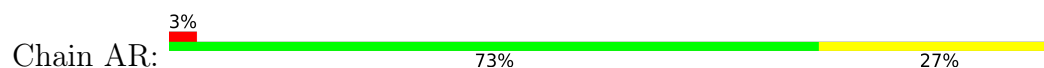
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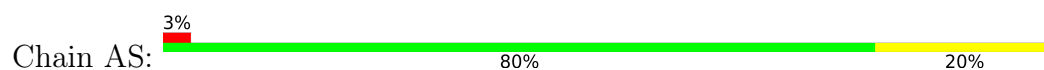
- Molecule 1: coat protein



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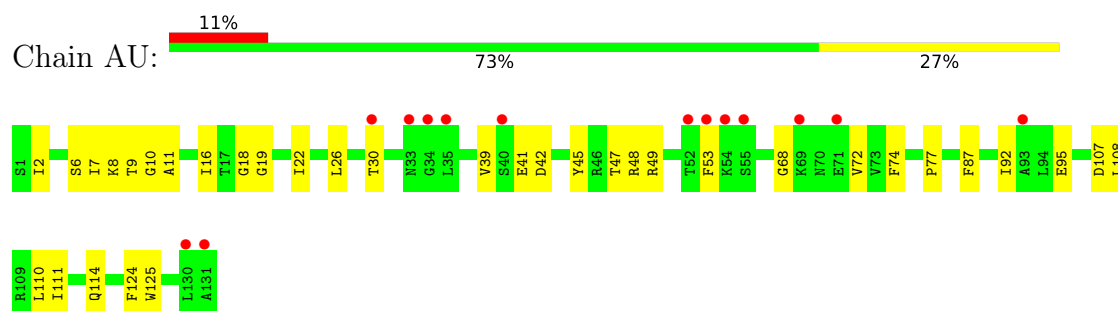
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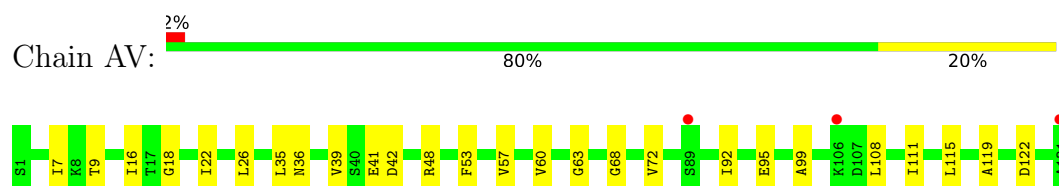
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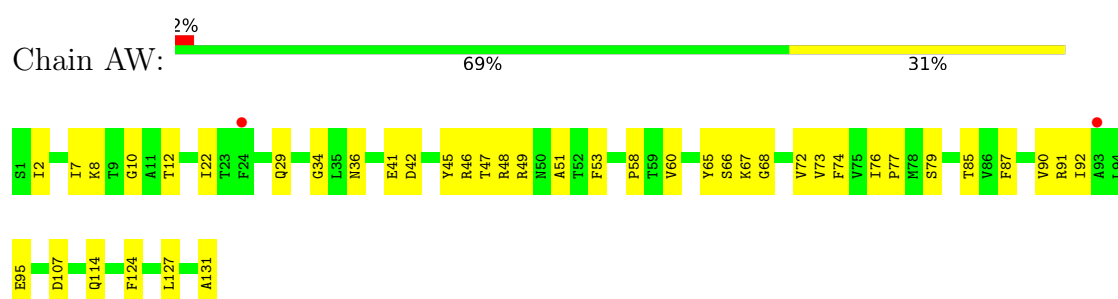
- Molecule 1: coat protein



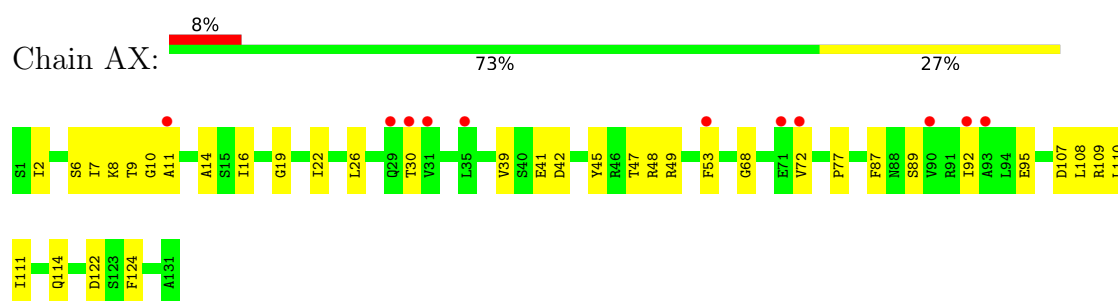
- Molecule 1: coat protein



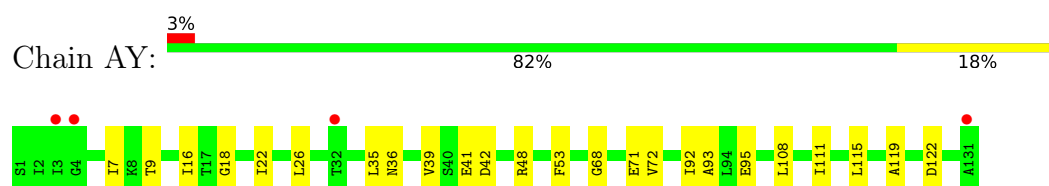
- Molecule 1: coat protein



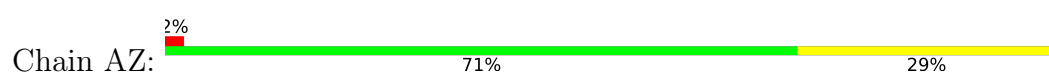
- Molecule 1: coat protein

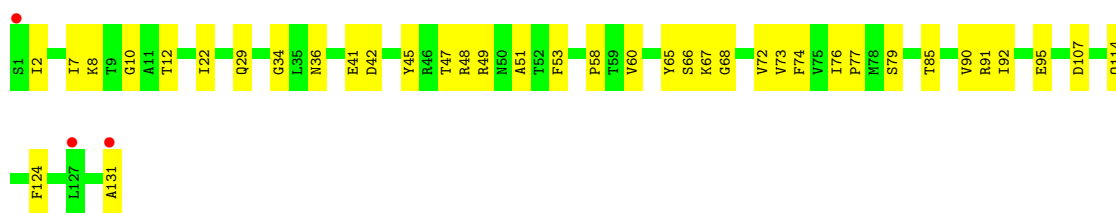


- Molecule 1: coat protein

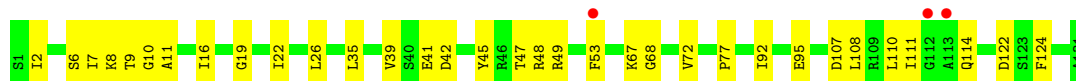
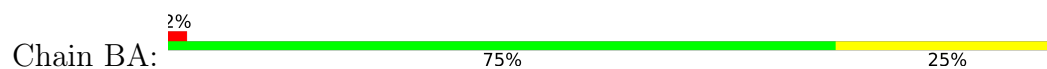


- Molecule 1: coat protein

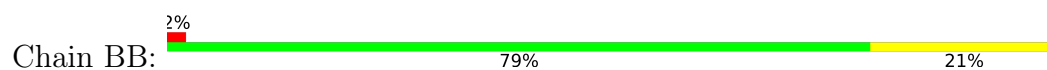




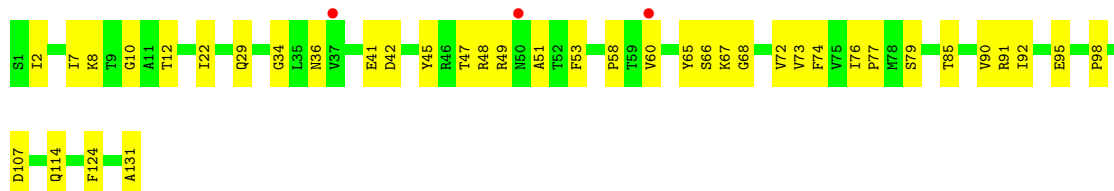
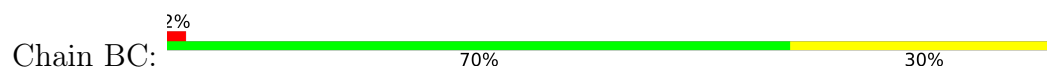
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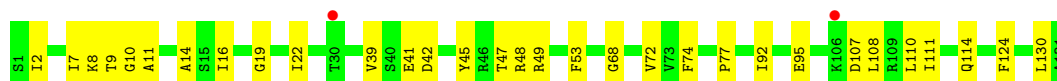
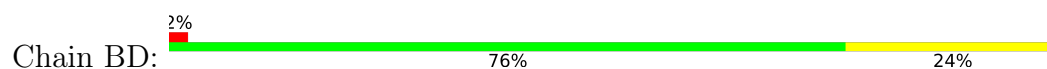
- Molecule 1: coat protein



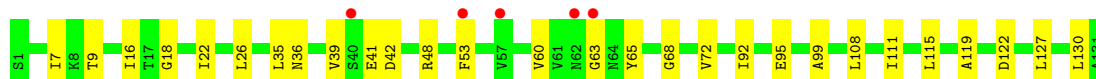
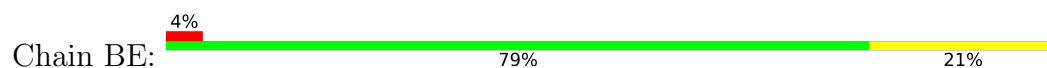
- Molecule 1: coat protein



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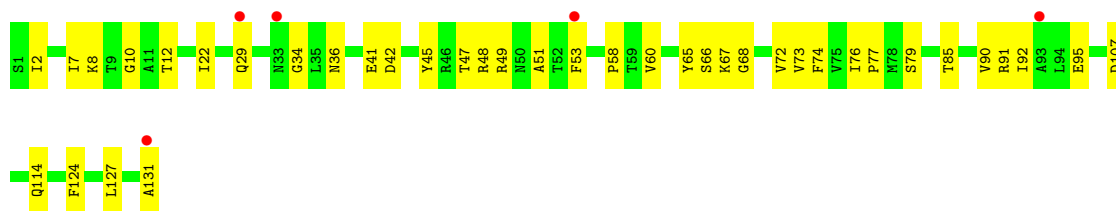


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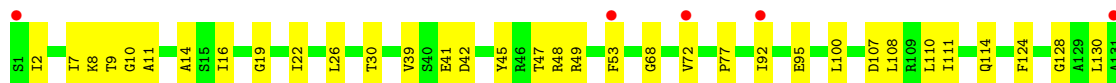
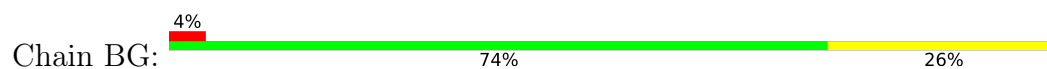


- Molecule 1: coat protein

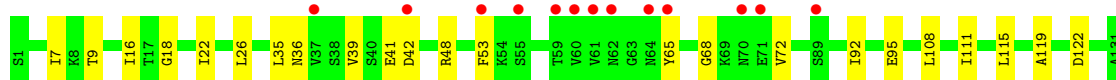
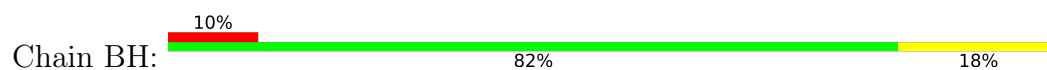




- Molecule 1: coat protein



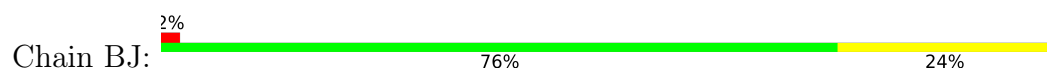
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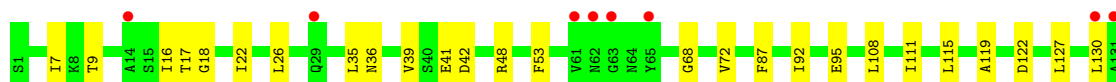
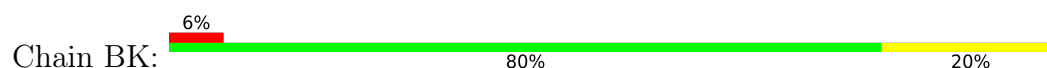
- Molecule 1: coat protein



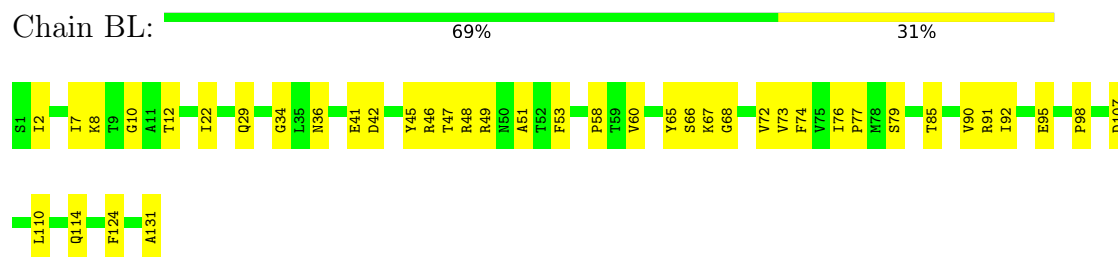
- Molecule 1: coat protein



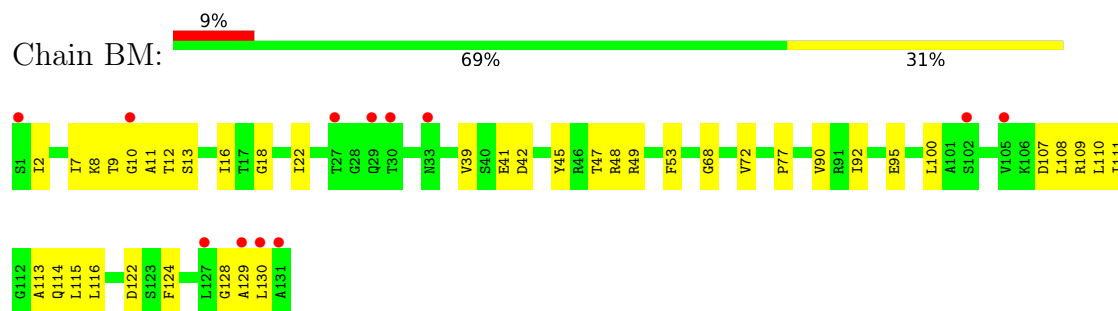
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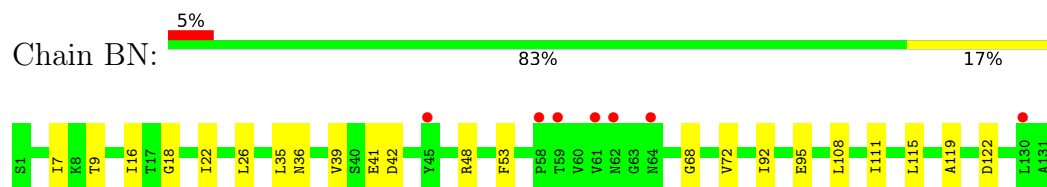
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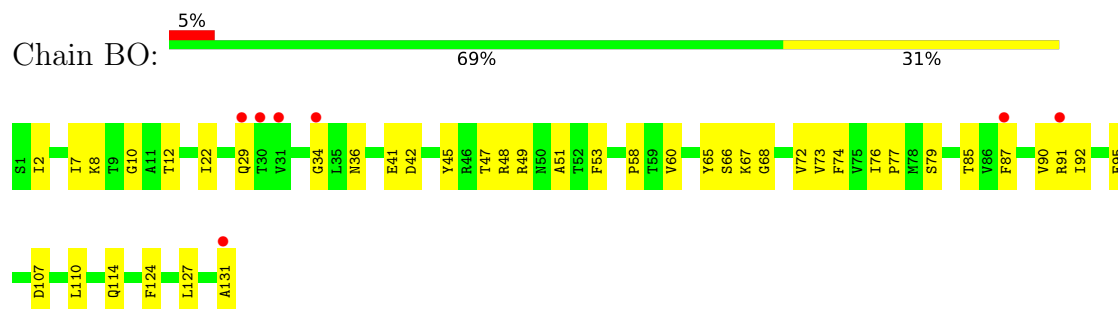
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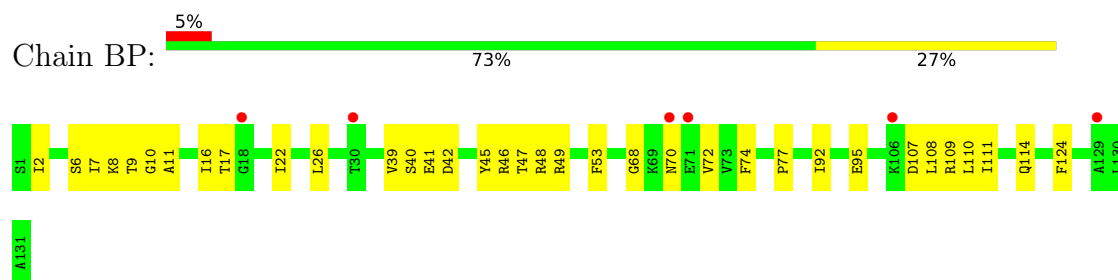
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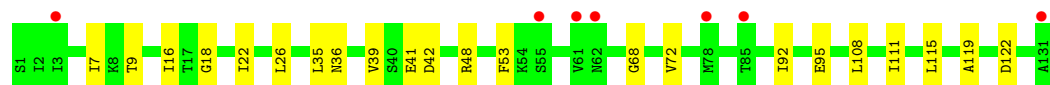
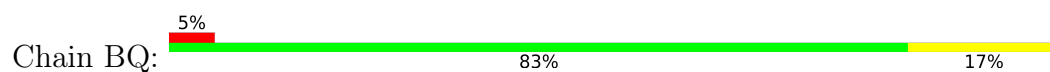
- Molecule 1: coat protein



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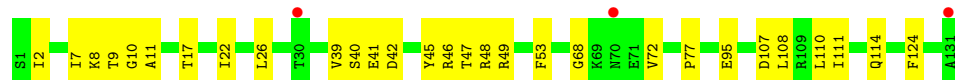
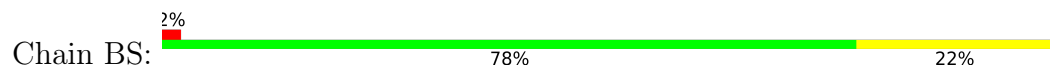
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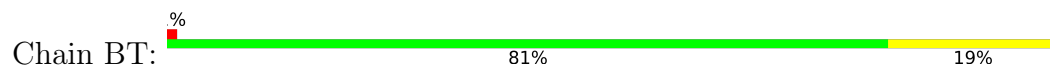
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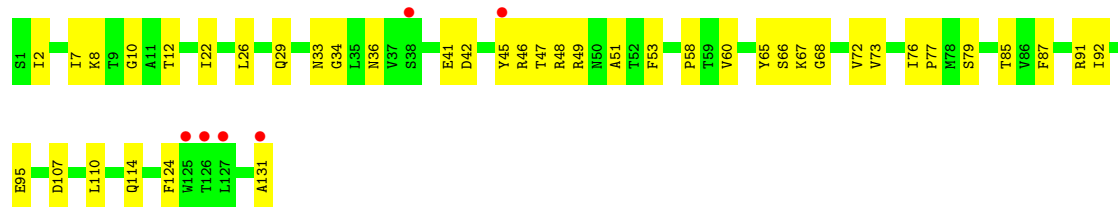
- Molecule 1: coat protein



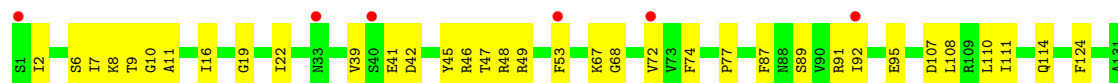
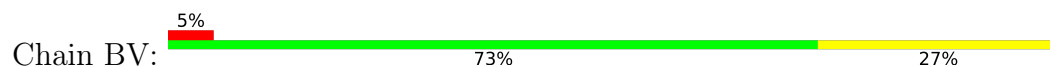
- Molecule 1: coat protein



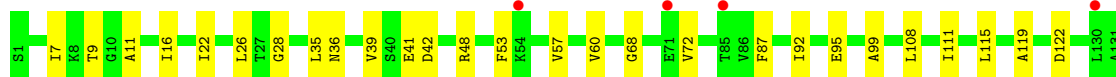
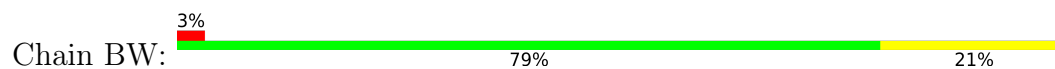
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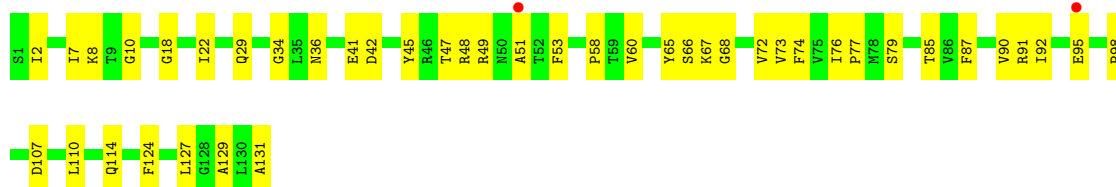
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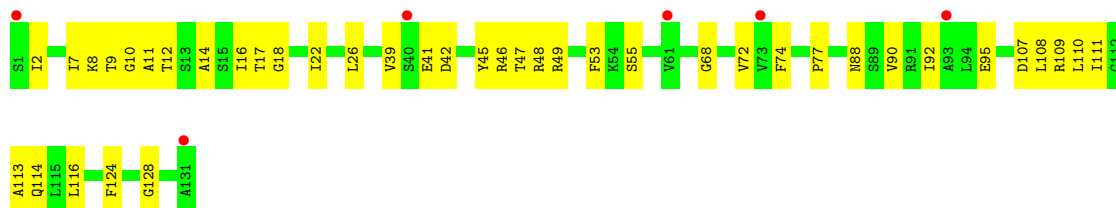
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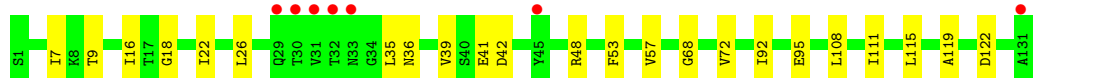
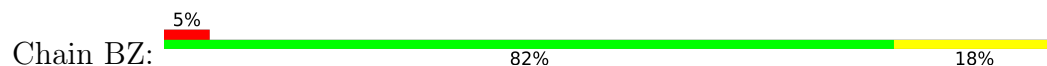
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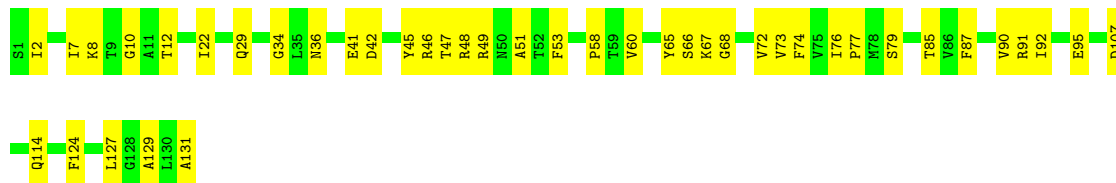
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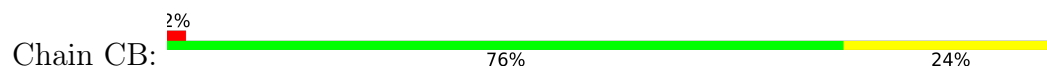
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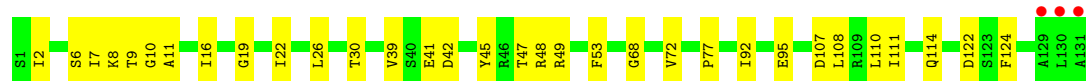


- Molecule 1: coat protein

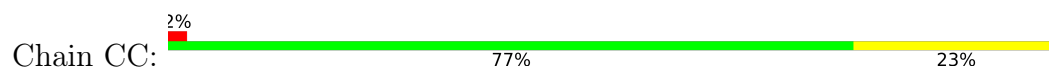


- Molecule 1: coat protein

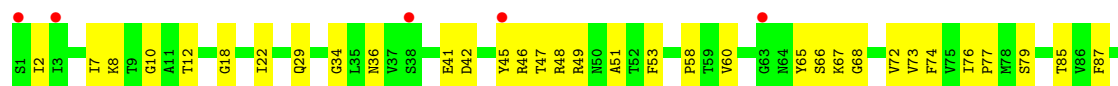




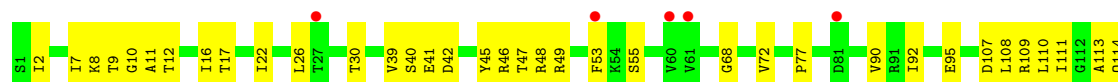
- Molecule 1: coat protein



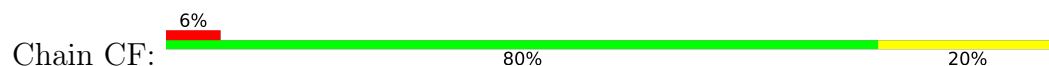
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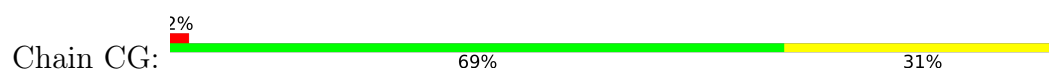
- Molecule 1: coat protein



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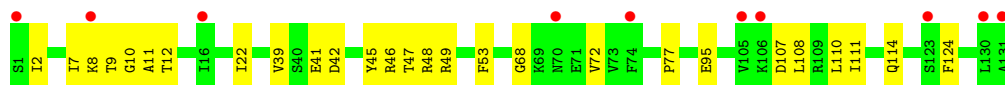
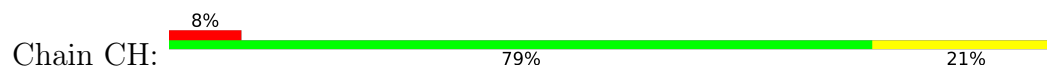


- Molecule 1: coat protein

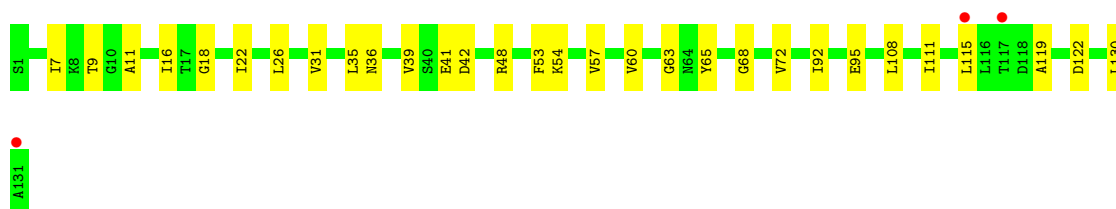
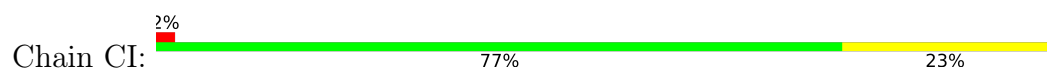




- Molecule 1: coat protein



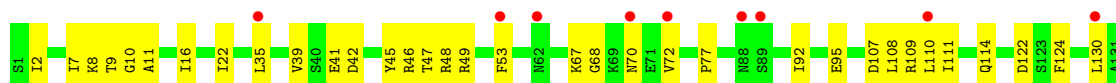
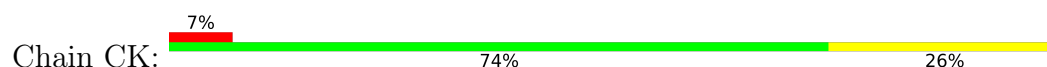
- Molecule 1: coat protein



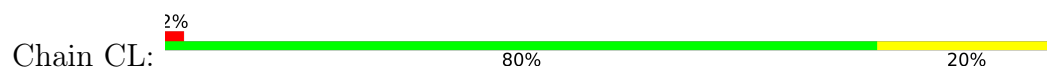
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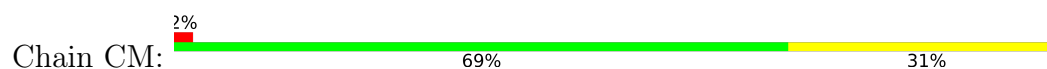
- Molecule 1: coat protein

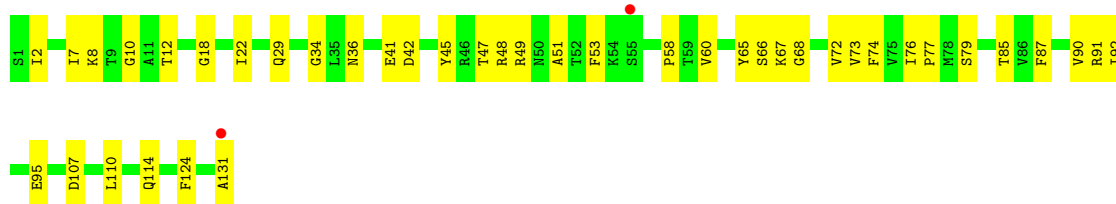


- Molecule 1: coat protein

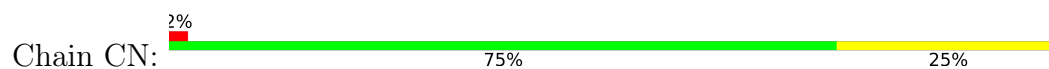


- Molecule 1: coat protein

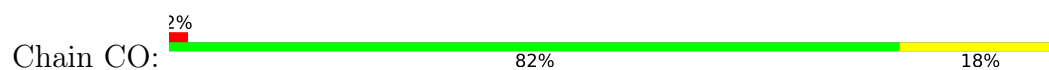




- Molecule 1: coat protein



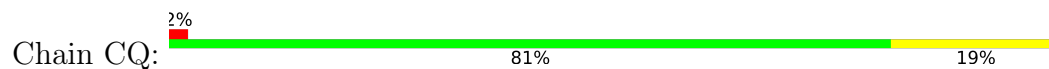
- Molecule 1: coat protein



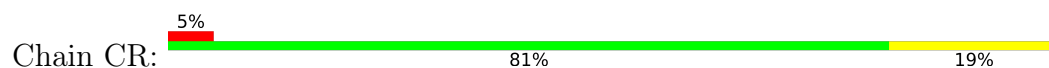
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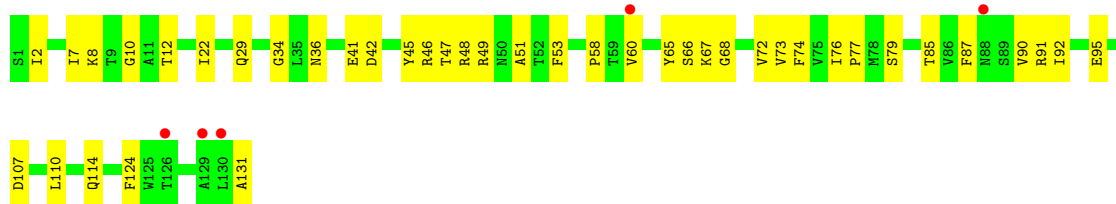


- Molecule 1: coat protein

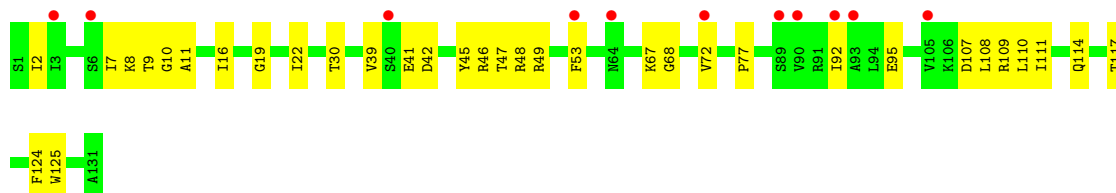
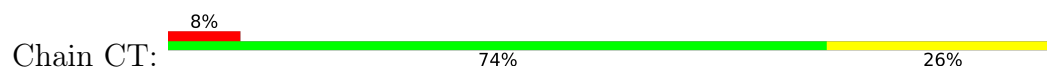


- Molecule 1: coat protein

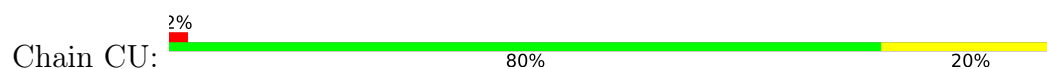




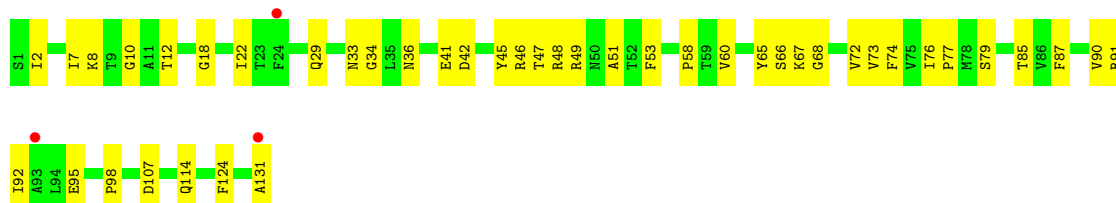
- Molecule 1: coat protein



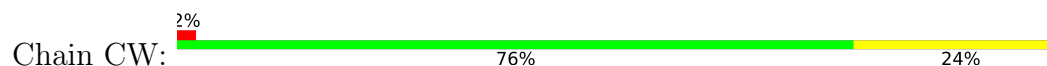
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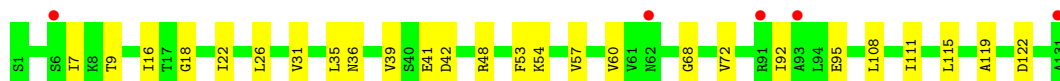
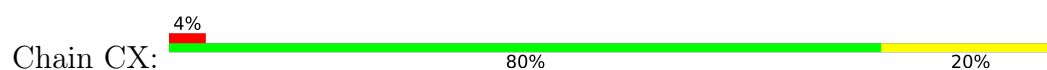
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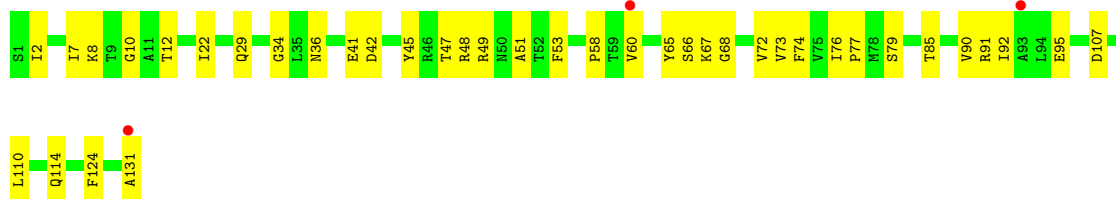
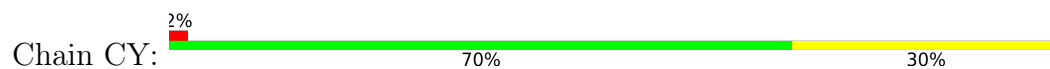
- Molecule 1: coat protein



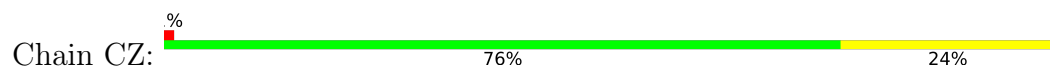
- Molecule 1: coat protein



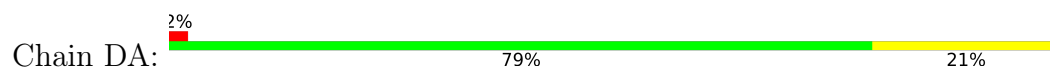
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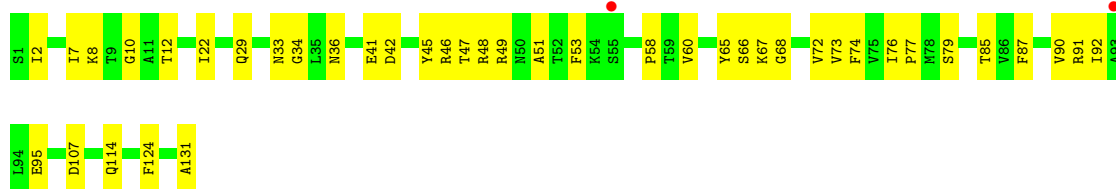
- Molecule 1: coat protein



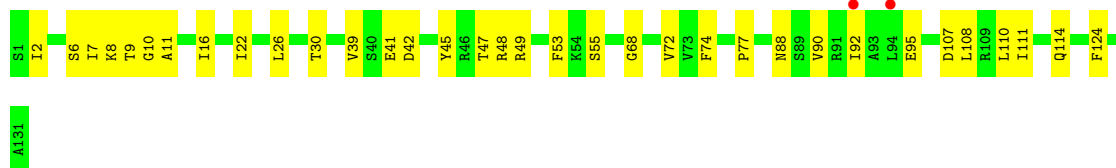
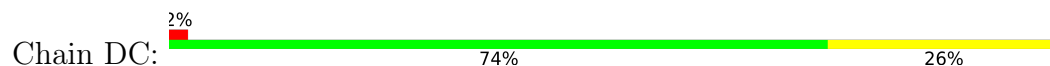
- Molecule 1: coat protein



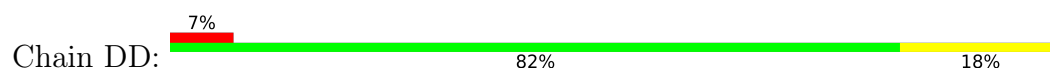
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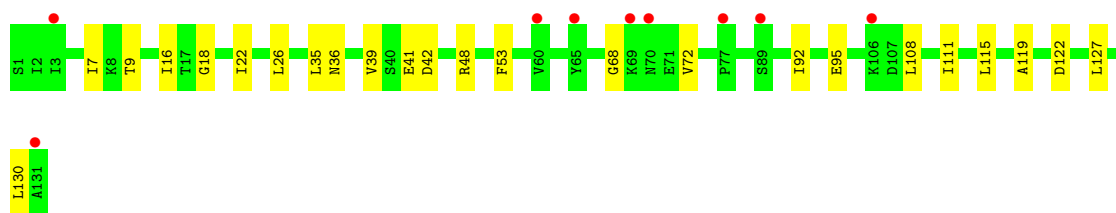


- Molecule 1: coat protein

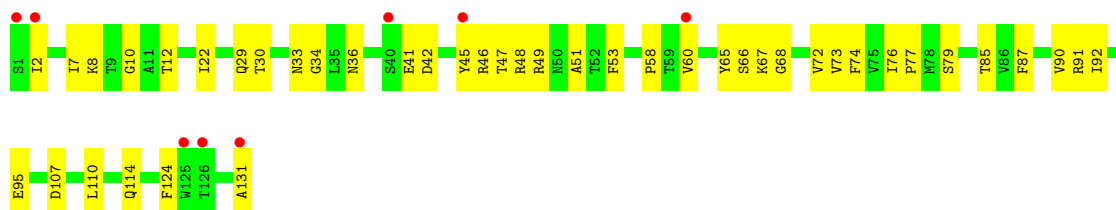


- Molecule 1: coat protein

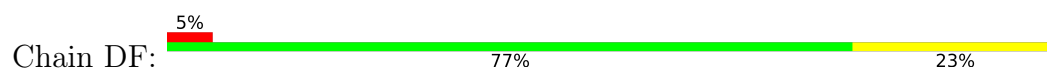




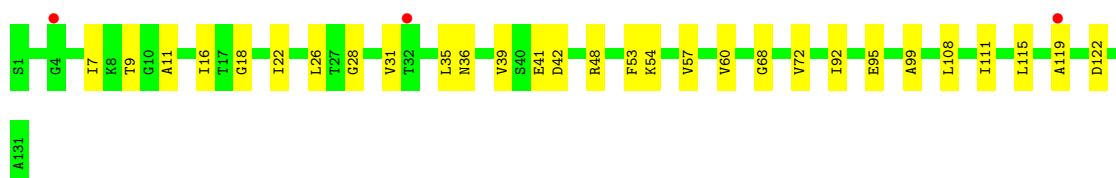
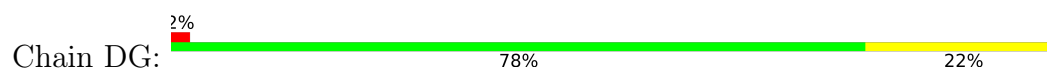
- Molecule 1: coat protein



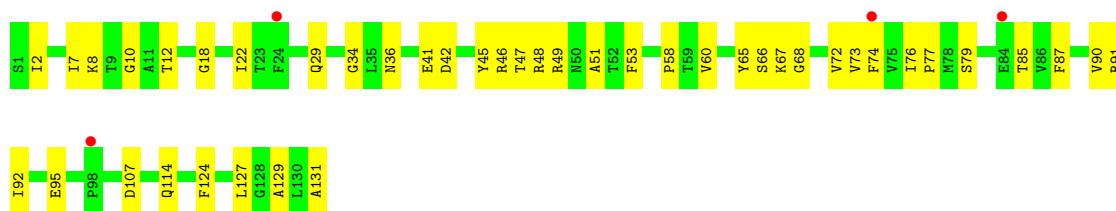
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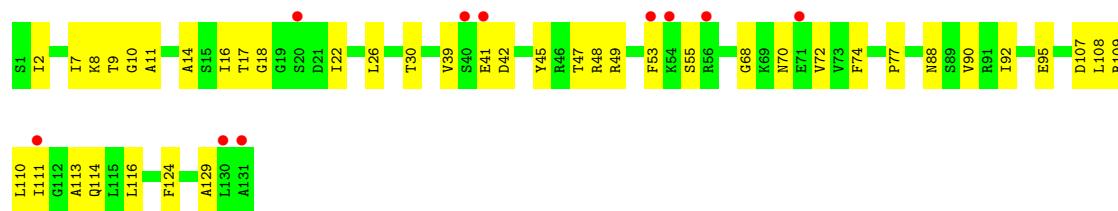


- Molecule 1: coat protein



- Molecule 1: coat protein





● Molecule 1: coat protein

Chain DJ: 82% 18%



● Molecule 1: coat protein

Chain DK: 4% 71% 29%



● Molecule 1: coat protein

Chain DL: 80% 20%



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	415.01Å 335.00Å 291.78Å 90.00° 134.66° 90.00°	Depositor
Resolution (Å)	49.14 – 3.48 49.14 – 3.48	Depositor EDS
% Data completeness (in resolution range)	96.5 (49.14-3.48) 98.1 (49.14-3.48)	Depositor EDS
R_{merge}	0.45	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.97 (at 3.48Å)	Xtriage
Refinement program	PHENIX 1.14 _3260	Depositor
R, R_{free}	0.269 , 0.271 0.273 , 0.275	Depositor DCC
R_{free} test set	10014 reflections (2.83%)	wwPDB-VP
Wilson B-factor (Å ²)	121.7	Xtriage
Anisotropy	0.231	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 86.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.37$, $\langle L^2 \rangle = 0.20$	Xtriage
Estimated twinning fraction	0.079 for h+2*k,-h-l 0.094 for h,-k,-h-l 0.095 for -h-2*k,-k,l	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	87480	wwPDB-VP
Average B, all atoms (Å ²)	110.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.44% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	AA	0.24	0/985	0.53	0/1338
1	AB	0.23	0/985	0.53	0/1338
1	AC	0.23	0/985	0.53	0/1338
1	AD	0.24	0/985	0.53	0/1338
1	AE	0.23	0/985	0.53	0/1338
1	AF	0.23	0/985	0.53	0/1338
1	AG	0.24	0/985	0.53	0/1338
1	AH	0.24	0/985	0.53	0/1338
1	AI	0.23	0/985	0.54	0/1338
1	AJ	0.24	0/985	0.53	0/1338
1	AK	0.24	0/985	0.53	0/1338
1	AL	0.23	0/985	0.53	0/1338
1	AM	0.24	0/985	0.53	0/1338
1	AN	0.23	0/985	0.53	0/1338
1	AO	0.23	0/985	0.53	0/1338
1	AP	0.24	0/985	0.53	0/1338
1	AQ	0.24	0/985	0.53	0/1338
1	AR	0.23	0/985	0.53	0/1338
1	AS	0.24	0/985	0.53	0/1338
1	AT	0.24	0/985	0.53	0/1338
1	AU	0.23	0/985	0.53	0/1338
1	AV	0.24	0/985	0.53	0/1338
1	AW	0.24	0/985	0.53	0/1338
1	AX	0.23	0/985	0.53	0/1338
1	AY	0.24	0/985	0.53	0/1338
1	AZ	0.24	0/985	0.53	0/1338
1	BA	0.23	0/985	0.54	0/1338
1	BB	0.24	0/985	0.53	0/1338
1	BC	0.24	0/985	0.53	0/1338
1	BD	0.23	0/985	0.53	0/1338
1	BE	0.24	0/985	0.53	0/1338
1	BF	0.24	0/985	0.53	0/1338
1	BG	0.23	0/985	0.53	0/1338
1	BH	0.24	0/985	0.52	0/1338

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	BI	0.23	0/985	0.53	0/1338
1	BJ	0.23	0/985	0.53	0/1338
1	BK	0.24	0/985	0.53	0/1338
1	BL	0.23	0/985	0.53	0/1338
1	BM	0.23	0/985	0.53	0/1338
1	BN	0.24	0/985	0.53	0/1338
1	BO	0.24	0/985	0.53	0/1338
1	BP	0.23	0/985	0.54	0/1338
1	BQ	0.24	0/985	0.53	0/1338
1	BR	0.24	0/985	0.53	0/1338
1	BS	0.23	0/985	0.53	0/1338
1	BT	0.24	0/985	0.53	0/1338
1	BU	0.24	0/985	0.53	0/1338
1	BV	0.23	0/985	0.53	0/1338
1	BW	0.24	0/985	0.52	0/1338
1	BX	0.24	0/985	0.53	0/1338
1	BY	0.23	0/985	0.53	0/1338
1	BZ	0.24	0/985	0.53	0/1338
1	CA	0.24	0/985	0.53	0/1338
1	CB	0.23	0/985	0.53	0/1338
1	CC	0.24	0/985	0.53	0/1338
1	CD	0.24	0/985	0.53	0/1338
1	CE	0.23	0/985	0.53	0/1338
1	CF	0.24	0/985	0.53	0/1338
1	CG	0.24	0/985	0.53	0/1338
1	CH	0.23	0/985	0.53	0/1338
1	CI	0.24	0/985	0.53	0/1338
1	CJ	0.24	0/985	0.53	0/1338
1	CK	0.23	0/985	0.53	0/1338
1	CL	0.24	0/985	0.53	0/1338
1	CM	0.24	0/985	0.53	0/1338
1	CN	0.23	0/985	0.53	0/1338
1	CO	0.24	0/985	0.53	0/1338
1	CP	0.24	0/985	0.53	0/1338
1	CQ	0.23	0/985	0.53	0/1338
1	CR	0.24	0/985	0.53	0/1338
1	CS	0.24	0/985	0.53	0/1338
1	CT	0.23	0/985	0.53	0/1338
1	CU	0.24	0/985	0.53	0/1338
1	CV	0.24	0/985	0.53	0/1338
1	CW	0.23	0/985	0.53	0/1338
1	CX	0.24	0/985	0.53	0/1338
1	CY	0.24	0/985	0.53	0/1338

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	CZ	0.23	0/985	0.53	0/1338
1	DA	0.24	0/985	0.53	0/1338
1	DB	0.24	0/985	0.53	0/1338
1	DC	0.23	0/985	0.53	0/1338
1	DD	0.24	0/985	0.53	0/1338
1	DE	0.24	0/985	0.53	0/1338
1	DF	0.23	0/985	0.54	0/1338
1	DG	0.24	0/985	0.53	0/1338
1	DH	0.24	0/985	0.53	0/1338
1	DI	0.23	0/985	0.53	0/1338
1	DJ	0.24	0/985	0.53	0/1338
1	DK	0.24	0/985	0.53	0/1338
1	DL	0.23	0/985	0.53	0/1338
All	All	0.23	0/88650	0.53	0/120420

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	AA	972	0	977	26	0
1	AB	972	0	977	39	0
1	AC	972	0	977	41	0
1	AD	972	0	977	25	0
1	AE	972	0	977	42	0
1	AF	972	0	977	41	0
1	AG	972	0	977	27	0
1	AH	972	0	977	38	0
1	AI	972	0	977	52	0
1	AJ	972	0	977	26	0
1	AK	972	0	977	40	0
1	AL	972	0	977	35	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	AM	972	0	977	28	0
1	AN	972	0	977	35	0
1	AO	972	0	977	51	0
1	AP	972	0	977	20	1
1	AQ	972	0	977	29	0
1	AR	972	0	977	37	1
1	AS	972	0	977	28	0
1	AT	972	0	977	39	0
1	AU	972	0	977	39	0
1	AV	972	0	977	25	0
1	AW	972	0	977	38	0
1	AX	972	0	977	37	0
1	AY	972	0	977	21	0
1	AZ	972	0	977	29	0
1	BA	972	0	977	32	0
1	BB	972	0	977	24	0
1	BC	972	0	977	39	0
1	BD	972	0	977	34	0
1	BE	972	0	977	29	0
1	BF	972	0	977	41	0
1	BG	972	0	977	40	0
1	BH	972	0	977	21	0
1	BI	972	0	977	31	0
1	BJ	972	0	977	43	0
1	BK	972	0	977	25	0
1	BL	972	0	977	43	0
1	BM	972	0	977	56	0
1	BN	972	0	977	20	0
1	BO	972	0	977	36	0
1	BP	972	0	977	46	0
1	BQ	972	0	977	21	0
1	BR	972	0	977	34	0
1	BS	972	0	977	30	1
1	BT	972	0	977	24	0
1	BU	972	0	977	36	0
1	BV	972	0	977	40	0
1	BW	972	0	977	27	0
1	BX	972	0	977	44	0
1	BY	972	0	977	54	0
1	BZ	972	0	977	21	0
1	CA	972	0	977	44	0
1	CB	972	0	977	32	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	CC	972	0	977	29	0
1	CD	972	0	977	46	0
1	CE	972	0	977	51	0
1	CF	972	0	977	27	0
1	CG	972	0	977	30	2
1	CH	972	0	977	27	1
1	CI	972	0	977	28	0
1	CJ	972	0	977	34	0
1	CK	972	0	977	46	0
1	CL	972	0	977	25	0
1	CM	972	0	977	35	0
1	CN	972	0	977	44	0
1	CO	972	0	977	21	0
1	CP	972	0	977	37	0
1	CQ	972	0	977	19	0
1	CR	972	0	977	24	0
1	CS	972	0	977	37	0
1	CT	972	0	977	37	0
1	CU	972	0	977	24	0
1	CV	972	0	977	39	0
1	CW	972	0	977	35	0
1	CX	972	0	977	24	0
1	CY	972	0	977	36	0
1	CZ	972	0	977	35	0
1	DA	972	0	977	26	0
1	DB	972	0	977	36	0
1	DC	972	0	977	39	0
1	DD	972	0	977	24	0
1	DE	972	0	977	38	0
1	DF	972	0	977	33	0
1	DG	972	0	977	29	0
1	DH	972	0	977	40	0
1	DI	972	0	977	55	0
1	DJ	972	0	977	23	0
1	DK	972	0	977	29	0
1	DL	972	0	977	22	0
All	All	87480	0	87930	2350	3

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 13.

All (2350) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BA:16:ILE:HD13	1:CW:114:GLN:HE21	1.10	1.14
1:AX:16:ILE:HD13	1:CT:114:GLN:HE21	1.12	1.12
1:BA:114:GLN:HE21	1:CW:16:ILE:HD13	1.11	1.10
1:AX:114:GLN:HE21	1:CT:16:ILE:HD13	1.18	1.08
1:AL:16:ILE:HD13	1:CB:114:GLN:HE21	1.19	1.07
1:AL:114:GLN:HE21	1:CB:16:ILE:HD13	1.17	1.03
1:BF:65:TYR:CD2	1:BS:46:ARG:HG2	1.98	0.97
1:AR:114:GLN:HE21	1:CK:16:ILE:HD13	1.28	0.96
1:AF:92:ILE:HG12	1:BV:92:ILE:HG12	1.53	0.91
1:AI:110:LEU:HD23	1:BY:16:ILE:HD11	1.50	0.91
1:BA:16:ILE:HD13	1:CW:114:GLN:NE2	1.86	0.90
1:BM:16:ILE:HD13	1:DI:114:GLN:HE21	1.34	0.90
1:AR:16:ILE:HD13	1:CK:114:GLN:HE21	1.37	0.89
1:AF:16:ILE:HD13	1:BV:114:GLN:HE21	1.37	0.89
1:AO:16:ILE:HD11	1:CE:110:LEU:HD23	1.54	0.89
1:AO:16:ILE:HD13	1:CE:114:GLN:HE21	1.36	0.88
1:BA:114:GLN:NE2	1:CW:16:ILE:HD13	1.87	0.88
1:BM:110:LEU:HD23	1:DI:16:ILE:HD11	1.55	0.88
1:BJ:16:ILE:HD13	1:DF:114:GLN:HE21	1.38	0.87
1:AF:26:LEU:HB3	1:DD:130:LEU:O	1.74	0.87
1:AX:114:GLN:NE2	1:CT:16:ILE:HD13	1.89	0.86
1:AX:16:ILE:HD13	1:CT:114:GLN:NE2	1.91	0.86
1:BG:16:ILE:HD13	1:DC:114:GLN:HE21	1.41	0.86
1:BM:108:LEU:HD22	1:DI:90:VAL:HG12	1.55	0.86
1:AO:114:GLN:HE21	1:CE:16:ILE:HD13	1.39	0.85
1:BM:16:ILE:HD11	1:DI:110:LEU:HD23	1.58	0.85
1:AC:16:ILE:HD13	1:BP:114:GLN:HE21	1.40	0.85
1:AF:114:GLN:HE21	1:BV:16:ILE:HD13	1.40	0.85
1:AO:110:LEU:HD23	1:CE:16:ILE:HD11	1.57	0.85
1:CD:65:TYR:HE2	1:CK:46:ARG:HA	1.40	0.84
1:BJ:114:GLN:HE21	1:DF:16:ILE:HD13	1.42	0.84
1:AL:16:ILE:HD13	1:CB:114:GLN:NE2	1.93	0.84
1:BF:65:TYR:HD2	1:BS:46:ARG:HG2	1.41	0.83
1:AE:87:PHE:CE2	1:BE:60:VAL:HG11	2.12	0.83
1:BX:65:TYR:HE2	1:CN:46:ARG:HA	1.43	0.83
1:AE:87:PHE:HE2	1:BE:60:VAL:HG11	1.43	0.82
1:AA:16:ILE:HD13	1:AB:114:GLN:HE21	1.45	0.82
1:CU:16:ILE:HD13	1:CV:114:GLN:HE21	1.45	0.82
1:AJ:16:ILE:HD13	1:AK:114:GLN:HE21	1.45	0.82
1:AU:114:GLN:HE21	1:CN:16:ILE:HD13	1.44	0.82
1:DD:16:ILE:HD13	1:DE:114:GLN:HE21	1.45	0.82
1:BZ:16:ILE:HD13	1:CA:114:GLN:HE21	1.45	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BJ:46:ARG:HG2	1:BL:65:TYR:CD2	2.14	0.82
1:BW:16:ILE:HD13	1:BX:114:GLN:HE21	1.45	0.82
1:DA:16:ILE:HD13	1:DB:114:GLN:HE21	1.45	0.82
1:AI:16:ILE:HD11	1:BY:110:LEU:HD23	1.62	0.82
1:BE:16:ILE:HD13	1:BF:114:GLN:HE21	1.44	0.82
1:BN:16:ILE:HD13	1:BO:114:GLN:HE21	1.45	0.82
1:BQ:16:ILE:HD13	1:BR:114:GLN:HE21	1.45	0.82
1:DJ:16:ILE:HD13	1:DK:114:GLN:HE21	1.45	0.82
1:AV:16:ILE:HD13	1:AW:114:GLN:HE21	1.45	0.82
1:BH:16:ILE:HD13	1:BI:114:GLN:HE21	1.45	0.82
1:BJ:46:ARG:HG2	1:BL:65:TYR:HD2	1.45	0.82
1:AL:114:GLN:NE2	1:CB:16:ILE:HD13	1.95	0.82
1:CI:16:ILE:HD13	1:CJ:114:GLN:HE21	1.45	0.82
1:AY:16:ILE:HD13	1:AZ:114:GLN:HE21	1.45	0.81
1:CX:16:ILE:HD13	1:CY:114:GLN:HE21	1.45	0.81
1:AD:16:ILE:HD13	1:AE:114:GLN:HE21	1.44	0.81
1:CC:16:ILE:HD13	1:CD:114:GLN:HE21	1.45	0.81
1:CF:16:ILE:HD13	1:CG:114:GLN:HE21	1.45	0.81
1:CR:16:ILE:HD13	1:CS:114:GLN:HE21	1.45	0.81
1:BD:16:ILE:HD13	1:CZ:114:GLN:HE21	1.43	0.81
1:DG:16:ILE:HD13	1:DH:114:GLN:HE21	1.45	0.81
1:AA:57:VAL:HG21	1:DH:29:GLN:HG3	1.62	0.80
1:AS:16:ILE:HD13	1:AT:114:GLN:HE21	1.45	0.80
1:AG:16:ILE:HD13	1:AH:114:GLN:HE21	1.45	0.80
1:CL:16:ILE:HD13	1:CM:114:GLN:HE21	1.45	0.80
1:CO:16:ILE:HD13	1:CP:114:GLN:HE21	1.45	0.80
1:AP:16:ILE:HD13	1:AQ:114:GLN:HE21	1.45	0.80
1:BT:16:ILE:HD13	1:BU:114:GLN:HE21	1.45	0.80
1:AI:16:ILE:HD13	1:BY:114:GLN:HE21	1.46	0.80
1:BB:16:ILE:HD13	1:BC:114:GLN:HE21	1.45	0.80
1:AM:16:ILE:HD13	1:AN:114:GLN:HE21	1.45	0.80
1:BK:16:ILE:HD13	1:BL:114:GLN:HE21	1.45	0.79
1:BM:130:LEU:HD21	1:DI:55:SER:HB3	1.65	0.78
1:BG:7:ILE:O	1:DC:22:ILE:HG22	1.84	0.77
1:AA:130:LEU:O	1:AU:26:LEU:HB3	1.85	0.77
1:CD:87:PHE:HE2	1:CI:60:VAL:HG11	1.48	0.76
1:AU:16:ILE:HD13	1:CN:114:GLN:HE21	1.49	0.76
1:BC:65:TYR:CD2	1:BP:46:ARG:HG2	2.21	0.75
1:CA:65:TYR:CD2	1:CH:46:ARG:HG2	2.22	0.75
1:BD:114:GLN:HE21	1:CZ:16:ILE:HD13	1.51	0.75
1:DG:16:ILE:HD13	1:DH:114:GLN:NE2	2.02	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CU:16:ILE:HD13	1:CV:114:GLN:NE2	2.02	0.75
1:AS:16:ILE:HD13	1:AT:114:GLN:NE2	2.02	0.74
1:AY:16:ILE:HD13	1:AZ:114:GLN:NE2	2.02	0.74
1:CI:16:ILE:HD13	1:CJ:114:GLN:NE2	2.02	0.74
1:BF:65:TYR:CE2	1:BS:46:ARG:HG2	2.22	0.74
1:BZ:16:ILE:HD13	1:CA:114:GLN:NE2	2.02	0.74
1:CD:87:PHE:CE2	1:CI:60:VAL:HG11	2.22	0.74
1:CO:16:ILE:HD13	1:CP:114:GLN:NE2	2.02	0.74
1:BQ:16:ILE:HD13	1:BR:114:GLN:NE2	2.02	0.74
1:CX:16:ILE:HD13	1:CY:114:GLN:NE2	2.02	0.74
1:AD:16:ILE:HD13	1:AE:114:GLN:NE2	2.02	0.74
1:AJ:16:ILE:HD13	1:AK:114:GLN:NE2	2.02	0.74
1:BH:16:ILE:HD13	1:BI:114:GLN:NE2	2.02	0.74
1:CA:46:ARG:HG3	1:CF:65:TYR:CE2	2.21	0.74
1:CA:87:PHE:CE2	1:CF:60:VAL:HG11	2.21	0.74
1:BM:12:THR:OG1	1:DI:17:THR:O	2.06	0.74
1:DA:16:ILE:HD13	1:DB:114:GLN:NE2	2.02	0.74
1:AT:87:PHE:HE2	1:AV:60:VAL:HG11	1.53	0.74
1:BE:16:ILE:HD13	1:BF:114:GLN:NE2	2.02	0.74
1:BM:100:LEU:HD13	1:DI:88:ASN:ND2	2.02	0.74
1:BW:16:ILE:HD13	1:BX:114:GLN:NE2	2.02	0.74
1:DJ:16:ILE:HD13	1:DK:114:GLN:NE2	2.02	0.74
1:AA:16:ILE:HD13	1:AB:114:GLN:NE2	2.02	0.74
1:AM:16:ILE:HD13	1:AN:114:GLN:NE2	2.02	0.74
1:BB:16:ILE:HD13	1:BC:114:GLN:NE2	2.02	0.74
1:CC:16:ILE:HD13	1:CD:114:GLN:NE2	2.02	0.73
1:CR:16:ILE:HD13	1:CS:114:GLN:NE2	2.02	0.73
1:DD:16:ILE:HD13	1:DE:114:GLN:NE2	2.02	0.73
1:AP:16:ILE:HD13	1:AQ:114:GLN:NE2	2.02	0.73
1:BK:16:ILE:HD13	1:BL:114:GLN:NE2	2.02	0.73
1:BT:16:ILE:HD13	1:BU:114:GLN:NE2	2.02	0.73
1:CW:46:ARG:HG2	1:CY:65:TYR:CD2	2.23	0.73
1:AG:16:ILE:HD13	1:AH:114:GLN:NE2	2.02	0.73
1:AU:92:ILE:HG12	1:CN:92:ILE:HG12	1.70	0.73
1:CF:16:ILE:HD13	1:CG:114:GLN:NE2	2.02	0.73
1:AV:16:ILE:HD13	1:AW:114:GLN:NE2	2.02	0.72
1:BN:16:ILE:HD13	1:BO:114:GLN:NE2	2.02	0.72
1:BX:65:TYR:CD2	1:CN:46:ARG:HG2	2.24	0.72
1:CL:16:ILE:HD13	1:CM:114:GLN:NE2	2.02	0.72
1:BM:109:ARG:NH2	1:DI:116:LEU:O	2.22	0.72
1:CD:65:TYR:CE2	1:CK:46:ARG:HA	2.22	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AC:7:ILE:O	1:BP:22:ILE:HG22	1.89	0.72
1:AE:85:THR:N	1:BE:63:GLY:O	2.22	0.72
1:AF:30:THR:OG1	1:DE:33:ASN:OD1	2.08	0.71
1:AT:87:PHE:CE2	1:AV:60:VAL:HG11	2.25	0.71
1:CJ:8:LYS:HB2	1:CJ:114:GLN:HE22	1.56	0.71
1:AE:8:LYS:HB2	1:AE:114:GLN:HE22	1.56	0.71
1:AU:16:ILE:HD11	1:CN:110:LEU:HD23	1.73	0.71
1:DE:8:LYS:HB2	1:DE:114:GLN:HE22	1.56	0.71
1:AA:60:VAL:HG11	1:DH:87:PHE:HE2	1.55	0.71
1:CM:8:LYS:HB2	1:CM:114:GLN:HE22	1.56	0.71
1:CG:8:LYS:HB2	1:CG:114:GLN:HE22	1.56	0.71
1:AK:8:LYS:HB2	1:AK:114:GLN:HE22	1.56	0.71
1:AZ:8:LYS:HB2	1:AZ:114:GLN:HE22	1.56	0.71
1:BC:8:LYS:HB2	1:BC:114:GLN:HE22	1.56	0.71
1:BF:8:LYS:HB2	1:BF:114:GLN:HE22	1.56	0.71
1:BM:116:LEU:O	1:DI:109:ARG:NE	2.24	0.71
1:AN:8:LYS:HB2	1:AN:114:GLN:HE22	1.56	0.71
1:BM:109:ARG:HG2	1:DI:116:LEU:HD22	1.73	0.71
1:BR:8:LYS:HB2	1:BR:114:GLN:HE22	1.56	0.71
1:AC:92:ILE:HG12	1:BP:92:ILE:HG12	1.73	0.70
1:BG:111:ILE:HD13	1:DC:74:PHE:CZ	2.26	0.70
1:AB:8:LYS:HB2	1:AB:114:GLN:HE22	1.56	0.70
1:AQ:8:LYS:HB2	1:AQ:114:GLN:HE22	1.56	0.70
1:BX:8:LYS:HB2	1:BX:114:GLN:HE22	1.56	0.70
1:CR:60:VAL:HG11	1:DB:87:PHE:HE2	1.56	0.70
1:AH:8:LYS:HB2	1:AH:114:GLN:HE22	1.56	0.70
1:BM:129:ALA:HB2	1:DG:26:LEU:O	1.91	0.70
1:CV:8:LYS:HB2	1:CV:114:GLN:HE22	1.56	0.70
1:DH:8:LYS:HB2	1:DH:114:GLN:HE22	1.56	0.70
1:AW:8:LYS:HB2	1:AW:114:GLN:HE22	1.56	0.70
1:CA:8:LYS:HB2	1:CA:114:GLN:HE22	1.56	0.70
1:AI:130:LEU:HD21	1:BY:55:SER:HB3	1.73	0.70
1:BI:8:LYS:HB2	1:BI:114:GLN:HE22	1.56	0.70
1:BO:8:LYS:HB2	1:BO:114:GLN:HE22	1.56	0.70
1:CD:8:LYS:HB2	1:CD:114:GLN:HE22	1.56	0.70
1:CY:8:LYS:HB2	1:CY:114:GLN:HE22	1.56	0.70
1:BF:65:TYR:HE2	1:BS:46:ARG:HA	1.56	0.70
1:BL:8:LYS:HB2	1:BL:114:GLN:HE22	1.56	0.70
1:BU:8:LYS:HB2	1:BU:114:GLN:HE22	1.56	0.70
1:AC:114:GLN:HE21	1:BP:16:ILE:HD13	1.57	0.70
1:CA:85:THR:N	1:CF:63:GLY:O	2.19	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CD:65:TYR:CD2	1:CK:46:ARG:HG2	2.26	0.69
1:BX:65:TYR:CE2	1:CN:46:ARG:HA	2.26	0.69
1:DB:8:LYS:HB2	1:DB:114:GLN:HE22	1.56	0.69
1:DK:8:LYS:HB2	1:DK:114:GLN:HE22	1.56	0.69
1:CP:8:LYS:HB2	1:CP:114:GLN:HE22	1.56	0.69
1:CS:8:LYS:HB2	1:CS:114:GLN:HE22	1.56	0.69
1:AQ:2:ILE:HD11	1:AQ:124:PHE:HB2	1.75	0.69
1:AT:2:ILE:HD11	1:AT:124:PHE:HB2	1.75	0.69
1:CP:2:ILE:HD11	1:CP:124:PHE:HB2	1.75	0.69
1:AT:8:LYS:HB2	1:AT:114:GLN:HE22	1.56	0.69
1:AX:6:SER:HA	1:CT:22:ILE:O	1.93	0.69
1:BD:39:VAL:HG12	1:BD:41:GLU:H	1.58	0.69
1:BG:92:ILE:HG12	1:DC:92:ILE:HG12	1.74	0.69
1:CT:39:VAL:HG12	1:CT:41:GLU:H	1.58	0.69
1:AF:87:PHE:O	1:BV:67:LYS:NZ	2.20	0.68
1:AX:39:VAL:HG12	1:AX:41:GLU:H	1.58	0.68
1:AK:2:ILE:HD11	1:AK:124:PHE:HB2	1.75	0.68
1:CD:65:TYR:HD2	1:CK:46:ARG:HG2	1.56	0.68
1:CJ:2:ILE:HD11	1:CJ:124:PHE:HB2	1.75	0.68
1:DL:39:VAL:HG12	1:DL:41:GLU:H	1.59	0.68
1:BS:39:VAL:HG12	1:BS:41:GLU:H	1.58	0.68
1:DI:39:VAL:HG12	1:DI:41:GLU:H	1.59	0.68
1:AB:2:ILE:HD11	1:AB:124:PHE:HB2	1.75	0.68
1:AE:2:ILE:HD11	1:AE:124:PHE:HB2	1.75	0.68
1:AE:42:ASP:HB2	1:AE:48:ARG:HG2	1.76	0.68
1:AK:42:ASP:HB2	1:AK:48:ARG:HG2	1.76	0.68
1:BG:39:VAL:HG12	1:BG:41:GLU:H	1.58	0.68
1:BR:42:ASP:HB2	1:BR:48:ARG:HG2	1.76	0.68
1:BX:42:ASP:HB2	1:BX:48:ARG:HG2	1.76	0.68
1:CN:39:VAL:HG12	1:CN:41:GLU:H	1.59	0.68
1:DB:2:ILE:HD11	1:DB:124:PHE:HB2	1.75	0.68
1:DF:39:VAL:HG12	1:DF:41:GLU:H	1.58	0.68
1:BY:39:VAL:HG12	1:BY:41:GLU:H	1.58	0.68
1:CB:39:VAL:HG12	1:CB:41:GLU:H	1.58	0.68
1:AH:29:GLN:HG3	1:CC:57:VAL:HG21	1.75	0.68
1:AI:114:GLN:HE21	1:BY:16:ILE:HD13	1.57	0.68
1:AQ:42:ASP:HB2	1:AQ:48:ARG:HG2	1.76	0.68
1:AT:42:ASP:HB2	1:AT:48:ARG:HG2	1.76	0.68
1:BO:2:ILE:HD11	1:BO:124:PHE:HB2	1.76	0.68
1:BX:65:TYR:HD2	1:CN:46:ARG:HG2	1.57	0.68
1:CH:39:VAL:HG12	1:CH:41:GLU:H	1.58	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CJ:42:ASP:HB2	1:CJ:48:ARG:HG2	1.76	0.68
1:DA:42:ASP:HB2	1:DA:48:ARG:HG3	1.76	0.68
1:AI:39:VAL:HG12	1:AI:41:GLU:H	1.58	0.68
1:CY:2:ILE:HD11	1:CY:124:PHE:HB2	1.75	0.68
1:AF:39:VAL:HG12	1:AF:41:GLU:H	1.59	0.68
1:AW:42:ASP:HB2	1:AW:48:ARG:HG2	1.76	0.68
1:BM:39:VAL:HG12	1:BM:41:GLU:H	1.58	0.68
1:BU:42:ASP:HB2	1:BU:48:ARG:HG2	1.76	0.68
1:BX:2:ILE:HD11	1:BX:124:PHE:HB2	1.75	0.68
1:DG:42:ASP:HB2	1:DG:48:ARG:HG3	1.76	0.68
1:AM:42:ASP:HB2	1:AM:48:ARG:HG3	1.76	0.68
1:BK:42:ASP:HB2	1:BK:48:ARG:HG3	1.76	0.68
1:BP:39:VAL:HG12	1:BP:41:GLU:H	1.59	0.68
1:CO:42:ASP:HB2	1:CO:48:ARG:HG3	1.76	0.68
1:CQ:39:VAL:HG12	1:CQ:41:GLU:H	1.58	0.68
1:DK:42:ASP:HB2	1:DK:48:ARG:HG2	1.76	0.68
1:AD:42:ASP:HB2	1:AD:48:ARG:HG3	1.76	0.68
1:AP:92:ILE:HG12	1:AQ:92:ILE:HG12	1.76	0.68
1:CA:65:TYR:HD2	1:CH:46:ARG:HG2	1.58	0.68
1:CO:92:ILE:HG12	1:CP:92:ILE:HG12	1.76	0.68
1:CW:39:VAL:HG12	1:CW:41:GLU:H	1.59	0.68
1:DB:42:ASP:HB2	1:DB:48:ARG:HG2	1.76	0.68
1:DJ:42:ASP:HB2	1:DJ:48:ARG:HG3	1.76	0.68
1:AO:39:VAL:HG12	1:AO:41:GLU:H	1.59	0.67
1:BJ:16:ILE:HD11	1:DF:110:LEU:HD23	1.74	0.67
1:BK:92:ILE:HG12	1:BL:92:ILE:HG12	1.76	0.67
1:AE:29:GLN:OE1	1:AE:36:ASN:ND2	2.28	0.67
1:AH:42:ASP:HB2	1:AH:48:ARG:HG2	1.76	0.67
1:AU:39:VAL:HG12	1:AU:41:GLU:H	1.58	0.67
1:AZ:2:ILE:HD11	1:AZ:124:PHE:HB2	1.75	0.67
1:BB:42:ASP:HB2	1:BB:48:ARG:HG3	1.76	0.67
1:BC:42:ASP:HB2	1:BC:48:ARG:HG2	1.76	0.67
1:BJ:39:VAL:HG12	1:BJ:41:GLU:H	1.58	0.67
1:BO:42:ASP:HB2	1:BO:48:ARG:HG2	1.76	0.67
1:CV:29:GLN:OE1	1:CV:36:ASN:ND2	2.28	0.67
1:AJ:92:ILE:HG12	1:AK:92:ILE:HG12	1.76	0.67
1:AN:29:GLN:OE1	1:AN:36:ASN:ND2	2.28	0.67
1:AS:42:ASP:HB2	1:AS:48:ARG:HG3	1.76	0.67
1:BC:65:TYR:HD2	1:BP:46:ARG:HG2	1.57	0.67
1:BF:42:ASP:HB2	1:BF:48:ARG:HG2	1.76	0.67
1:BL:2:ILE:HD11	1:BL:124:PHE:HB2	1.75	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BL:42:ASP:HB2	1:BL:48:ARG:HG2	1.76	0.67
1:CP:29:GLN:OE1	1:CP:36:ASN:ND2	2.28	0.67
1:AB:29:GLN:OE1	1:AB:36:ASN:ND2	2.28	0.67
1:AB:42:ASP:HB2	1:AB:48:ARG:HG2	1.76	0.67
1:AV:92:ILE:HG12	1:AW:92:ILE:HG12	1.76	0.67
1:BI:29:GLN:OE1	1:BI:36:ASN:ND2	2.28	0.67
1:BM:128:GLY:O	1:DG:28:GLY:HA3	1.95	0.67
1:CD:42:ASP:HB2	1:CD:48:ARG:HG2	1.76	0.67
1:CI:92:ILE:HG12	1:CJ:92:ILE:HG12	1.76	0.67
1:CP:42:ASP:HB2	1:CP:48:ARG:HG2	1.76	0.67
1:CU:42:ASP:HB2	1:CU:48:ARG:HG3	1.76	0.67
1:DH:2:ILE:HD11	1:DH:124:PHE:HB2	1.75	0.67
1:AG:92:ILE:HG12	1:AH:92:ILE:HG12	1.76	0.67
1:AK:29:GLN:OE1	1:AK:36:ASN:ND2	2.28	0.67
1:BO:29:GLN:OE1	1:BO:36:ASN:ND2	2.28	0.67
1:CE:39:VAL:HG12	1:CE:41:GLU:H	1.58	0.67
1:CG:29:GLN:OE1	1:CG:36:ASN:ND2	2.28	0.67
1:CU:92:ILE:HG12	1:CV:92:ILE:HG12	1.76	0.67
1:AM:92:ILE:HG12	1:AN:92:ILE:HG12	1.76	0.67
1:AT:29:GLN:OE1	1:AT:36:ASN:ND2	2.28	0.67
1:BR:29:GLN:OE1	1:BR:36:ASN:ND2	2.28	0.67
1:BV:39:VAL:HG12	1:BV:41:GLU:H	1.58	0.67
1:BX:29:GLN:OE1	1:BX:36:ASN:ND2	2.28	0.67
1:BZ:92:ILE:HG12	1:CA:92:ILE:HG12	1.76	0.67
1:CC:42:ASP:HB2	1:CC:48:ARG:HG3	1.76	0.67
1:CK:39:VAL:HG12	1:CK:41:GLU:H	1.59	0.67
1:CV:2:ILE:HD11	1:CV:124:PHE:HB2	1.75	0.67
1:CY:29:GLN:OE1	1:CY:36:ASN:ND2	2.28	0.67
1:DD:92:ILE:HG12	1:DE:92:ILE:HG12	1.76	0.67
1:AA:92:ILE:HG12	1:AB:92:ILE:HG12	1.76	0.67
1:AX:92:ILE:HG12	1:CT:92:ILE:HG12	1.76	0.67
1:BF:29:GLN:OE1	1:BF:36:ASN:ND2	2.28	0.67
1:BH:42:ASP:HB2	1:BH:48:ARG:HG3	1.76	0.67
1:BL:29:GLN:OE1	1:BL:36:ASN:ND2	2.28	0.67
1:BN:92:ILE:HG12	1:BO:92:ILE:HG12	1.76	0.67
1:BU:2:ILE:HD11	1:BU:124:PHE:HB2	1.75	0.67
1:CS:29:GLN:OE1	1:CS:36:ASN:ND2	2.28	0.67
1:DK:2:ILE:HD11	1:DK:124:PHE:HB2	1.75	0.67
1:AG:42:ASP:HB2	1:AG:48:ARG:HG3	1.76	0.67
1:AQ:29:GLN:OE1	1:AQ:36:ASN:ND2	2.28	0.67
1:AR:39:VAL:HG12	1:AR:41:GLU:H	1.59	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AW:29:GLN:OE1	1:AW:36:ASN:ND2	2.28	0.67
1:BC:2:ILE:HD11	1:BC:124:PHE:HB2	1.75	0.67
1:BT:92:ILE:HG12	1:BU:92:ILE:HG12	1.76	0.67
1:BU:29:GLN:OE1	1:BU:36:ASN:ND2	2.28	0.67
1:CL:92:ILE:HG12	1:CM:92:ILE:HG12	1.76	0.67
1:CM:42:ASP:HB2	1:CM:48:ARG:HG2	1.76	0.67
1:CR:42:ASP:HB2	1:CR:48:ARG:HG3	1.76	0.67
1:CX:92:ILE:HG12	1:CY:92:ILE:HG12	1.76	0.67
1:DE:29:GLN:OE1	1:DE:36:ASN:ND2	2.28	0.67
1:AL:39:VAL:HG12	1:AL:41:GLU:H	1.58	0.67
1:AR:114:GLN:NE2	1:CK:16:ILE:HD13	2.07	0.67
1:AW:2:ILE:HD11	1:AW:124:PHE:HB2	1.75	0.67
1:AY:92:ILE:HG12	1:AZ:92:ILE:HG12	1.76	0.67
1:BW:42:ASP:HB2	1:BW:48:ARG:HG3	1.76	0.67
1:CA:87:PHE:HE2	1:CF:60:VAL:HG11	1.56	0.67
1:CD:29:GLN:OE1	1:CD:36:ASN:ND2	2.28	0.67
1:CJ:29:GLN:OE1	1:CJ:36:ASN:ND2	2.28	0.67
1:DD:42:ASP:HB2	1:DD:48:ARG:HG3	1.76	0.67
1:DE:2:ILE:HD11	1:DE:124:PHE:HB2	1.75	0.67
1:DH:42:ASP:HB2	1:DH:48:ARG:HG2	1.76	0.67
1:AN:2:ILE:HD11	1:AN:124:PHE:HB2	1.75	0.67
1:CS:2:ILE:HD11	1:CS:124:PHE:HB2	1.75	0.67
1:CX:42:ASP:HB2	1:CX:48:ARG:HG3	1.76	0.67
1:AH:2:ILE:HD11	1:AH:124:PHE:HB2	1.75	0.66
1:AP:42:ASP:HB2	1:AP:48:ARG:HG3	1.76	0.66
1:AZ:29:GLN:OE1	1:AZ:36:ASN:ND2	2.28	0.66
1:BI:2:ILE:HD11	1:BI:124:PHE:HB2	1.75	0.66
1:BR:2:ILE:HD11	1:BR:124:PHE:HB2	1.75	0.66
1:BX:45:TYR:HD1	1:BX:48:ARG:HH21	1.44	0.66
1:CG:2:ILE:HD11	1:CG:124:PHE:HB2	1.75	0.66
1:CI:42:ASP:HB2	1:CI:48:ARG:HG3	1.76	0.66
1:CL:42:ASP:HB2	1:CL:48:ARG:HG3	1.76	0.66
1:CM:2:ILE:HD11	1:CM:124:PHE:HB2	1.75	0.66
1:DA:92:ILE:HG12	1:DB:92:ILE:HG12	1.76	0.66
1:DB:29:GLN:OE1	1:DB:36:ASN:ND2	2.28	0.66
1:DG:92:ILE:HG12	1:DH:92:ILE:HG12	1.76	0.66
1:DH:29:GLN:OE1	1:DH:36:ASN:ND2	2.28	0.66
1:DK:29:GLN:OE1	1:DK:36:ASN:ND2	2.28	0.66
1:AA:42:ASP:HB2	1:AA:48:ARG:HG3	1.76	0.66
1:AN:42:ASP:HB2	1:AN:48:ARG:HG2	1.76	0.66
1:AZ:42:ASP:HB2	1:AZ:48:ARG:HG2	1.76	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BE:42:ASP:HB2	1:BE:48:ARG:HG3	1.76	0.66
1:BF:2:ILE:HD11	1:BF:124:PHE:HB2	1.75	0.66
1:BL:45:TYR:HD1	1:BL:48:ARG:HH21	1.44	0.66
1:CA:45:TYR:HD1	1:CA:48:ARG:HH21	1.43	0.66
1:DJ:92:ILE:HG12	1:DK:92:ILE:HG12	1.76	0.66
1:AH:29:GLN:OE1	1:AH:36:ASN:ND2	2.28	0.66
1:AJ:42:ASP:HB2	1:AJ:48:ARG:HG3	1.76	0.66
1:BC:29:GLN:OE1	1:BC:36:ASN:ND2	2.28	0.66
1:BH:92:ILE:HG12	1:BI:92:ILE:HG12	1.76	0.66
1:BQ:42:ASP:HB2	1:BQ:48:ARG:HG3	1.76	0.66
1:BW:92:ILE:HG12	1:BX:92:ILE:HG12	1.76	0.66
1:CA:42:ASP:HB2	1:CA:48:ARG:HG2	1.76	0.66
1:CC:92:ILE:HG12	1:CD:92:ILE:HG12	1.76	0.66
1:CF:42:ASP:HB2	1:CF:48:ARG:HG3	1.76	0.66
1:CV:42:ASP:HB2	1:CV:48:ARG:HG2	1.76	0.66
1:AC:39:VAL:HG12	1:AC:41:GLU:H	1.59	0.66
1:AV:42:ASP:HB2	1:AV:48:ARG:HG3	1.76	0.66
1:BZ:42:ASP:HB2	1:BZ:48:ARG:HG3	1.76	0.66
1:CD:2:ILE:HD11	1:CD:124:PHE:HB2	1.75	0.66
1:CD:45:TYR:HD1	1:CD:48:ARG:HH21	1.43	0.66
1:CP:45:TYR:HD1	1:CP:48:ARG:HH21	1.44	0.66
1:DB:45:TYR:HD1	1:DB:48:ARG:HH21	1.44	0.66
1:DE:42:ASP:HB2	1:DE:48:ARG:HG2	1.76	0.66
1:AT:45:TYR:HD1	1:AT:48:ARG:HH21	1.43	0.66
1:CA:2:ILE:HD11	1:CA:124:PHE:HB2	1.75	0.66
1:CS:42:ASP:HB2	1:CS:48:ARG:HG2	1.76	0.66
1:AS:92:ILE:HG12	1:AT:92:ILE:HG12	1.76	0.66
1:AZ:45:TYR:HD1	1:AZ:48:ARG:HH21	1.43	0.66
1:BA:39:VAL:HG12	1:BA:41:GLU:H	1.58	0.66
1:BE:92:ILE:HG12	1:BF:92:ILE:HG12	1.76	0.66
1:CY:42:ASP:HB2	1:CY:48:ARG:HG2	1.76	0.66
1:BM:16:ILE:HG21	1:DI:114:GLN:NE2	2.10	0.66
1:CG:42:ASP:HB2	1:CG:48:ARG:HG2	1.76	0.66
1:AY:42:ASP:HB2	1:AY:48:ARG:HG3	1.76	0.66
1:BB:92:ILE:HG12	1:BC:92:ILE:HG12	1.76	0.66
1:BC:45:TYR:HD1	1:BC:48:ARG:HH21	1.44	0.66
1:BN:42:ASP:HB2	1:BN:48:ARG:HG3	1.76	0.66
1:BT:42:ASP:HB2	1:BT:48:ARG:HG3	1.76	0.66
1:CM:29:GLN:OE1	1:CM:36:ASN:ND2	2.28	0.66
1:DK:45:TYR:HD1	1:DK:48:ARG:HH21	1.43	0.66
1:AH:87:PHE:HE2	1:CC:60:VAL:HG11	1.58	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BI:42:ASP:HB2	1:BI:48:ARG:HG2	1.76	0.66
1:CA:29:GLN:OE1	1:CA:36:ASN:ND2	2.28	0.66
1:CA:46:ARG:HG3	1:CF:65:TYR:CD2	2.31	0.66
1:DH:45:TYR:HD1	1:DH:48:ARG:HH21	1.44	0.66
1:AQ:45:TYR:HD1	1:AQ:48:ARG:HH21	1.43	0.66
1:BJ:110:LEU:HD23	1:DF:16:ILE:HD11	1.78	0.66
1:CY:45:TYR:HD1	1:CY:48:ARG:HH21	1.44	0.66
1:CZ:39:VAL:HG12	1:CZ:41:GLU:H	1.59	0.66
1:AB:33:ASN:OD1	1:AU:30:THR:OG1	2.13	0.65
1:AI:108:LEU:HD22	1:BY:90:VAL:HG12	1.77	0.65
1:AW:45:TYR:HD1	1:AW:48:ARG:HH21	1.44	0.65
1:BD:7:ILE:O	1:CZ:22:ILE:HG22	1.96	0.65
1:BQ:92:ILE:HG12	1:BR:92:ILE:HG12	1.76	0.65
1:CM:45:TYR:HD1	1:CM:48:ARG:HH21	1.44	0.65
1:CV:45:TYR:HD1	1:CV:48:ARG:HH21	1.43	0.65
1:DC:39:VAL:HG12	1:DC:41:GLU:H	1.59	0.65
1:AD:92:ILE:HG12	1:AE:92:ILE:HG12	1.76	0.65
1:CR:92:ILE:HG12	1:CS:92:ILE:HG12	1.76	0.65
1:DE:45:TYR:HD1	1:DE:48:ARG:HH21	1.43	0.65
1:AK:45:TYR:HD1	1:AK:48:ARG:HH21	1.44	0.65
1:BR:45:TYR:HD1	1:BR:48:ARG:HH21	1.43	0.65
1:CF:92:ILE:HG12	1:CG:92:ILE:HG12	1.76	0.65
1:CJ:45:TYR:HD1	1:CJ:48:ARG:HH21	1.44	0.65
1:BG:100:LEU:HD13	1:DC:88:ASN:ND2	2.12	0.65
1:BI:45:TYR:HD1	1:BI:48:ARG:HH21	1.43	0.65
1:BF:45:TYR:HD1	1:BF:48:ARG:HH21	1.44	0.64
1:BM:116:LEU:HD22	1:DI:109:ARG:HG2	1.79	0.64
1:CG:45:TYR:HD1	1:CG:48:ARG:HH21	1.44	0.64
1:AE:45:TYR:HD1	1:AE:48:ARG:HH21	1.44	0.64
1:AN:45:TYR:HD1	1:AN:48:ARG:HH21	1.43	0.64
1:CS:45:TYR:HD1	1:CS:48:ARG:HH21	1.44	0.64
1:BG:114:GLN:HE21	1:DC:16:ILE:HD13	1.63	0.64
1:BO:87:PHE:HE2	1:DJ:60:VAL:HG11	1.62	0.64
1:AL:92:ILE:HG12	1:CB:92:ILE:HG12	1.79	0.64
1:AC:46:ARG:HA	1:DH:65:TYR:HE2	1.61	0.64
1:AH:45:TYR:HD1	1:AH:48:ARG:HH21	1.43	0.64
1:AI:12:THR:OG1	1:BY:17:THR:O	2.11	0.64
1:AH:26:LEU:O	1:CD:129:ALA:HB2	1.97	0.64
1:AO:16:ILE:HG21	1:CE:114:GLN:NE2	2.13	0.64
1:BO:45:TYR:HD1	1:BO:48:ARG:HH21	1.44	0.64
1:BU:45:TYR:HD1	1:BU:48:ARG:HH21	1.44	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BA:92:ILE:HG12	1:CW:92:ILE:HG12	1.80	0.63
1:AE:46:ARG:HG3	1:BE:65:TYR:CE2	2.33	0.63
1:AB:45:TYR:HD1	1:AB:48:ARG:HH21	1.44	0.63
1:BE:127:LEU:HD13	1:BS:40:SER:HB2	1.79	0.63
1:BM:7:ILE:O	1:DI:22:ILE:HG22	1.98	0.63
1:CR:57:VAL:HG21	1:DB:29:GLN:HG3	1.80	0.63
1:AR:16:ILE:HD11	1:CK:110:LEU:HD23	1.81	0.62
1:AF:6:SER:HA	1:BV:22:ILE:O	1.98	0.62
1:BD:92:ILE:HG12	1:CZ:92:ILE:HG12	1.80	0.62
1:CR:60:VAL:HG11	1:DB:87:PHE:CE2	2.33	0.62
1:AA:60:VAL:HG11	1:DH:87:PHE:CE2	2.33	0.62
1:AH:87:PHE:CE2	1:CC:60:VAL:HG11	2.33	0.62
1:AB:30:THR:OG1	1:AS:33:ASN:OD1	2.16	0.62
1:AE:127:LEU:HD21	1:DE:48:ARG:HH12	1.63	0.62
1:CB:47:THR:HA	1:CB:77:PRO:HG2	1.82	0.62
1:DC:47:THR:HA	1:DC:77:PRO:HG2	1.82	0.62
1:AL:22:ILE:O	1:CB:6:SER:HA	1.99	0.62
1:BC:65:TYR:HE2	1:BP:46:ARG:HA	1.63	0.62
1:BS:47:THR:HA	1:BS:77:PRO:HG2	1.82	0.62
1:CE:47:THR:HA	1:CE:77:PRO:HG2	1.82	0.62
1:CQ:47:THR:HA	1:CQ:77:PRO:HG2	1.82	0.62
1:AW:65:TYR:CD2	1:BV:46:ARG:HG2	2.35	0.62
1:BJ:47:THR:HA	1:BJ:77:PRO:HG2	1.82	0.62
1:AF:8:LYS:HB2	1:AF:114:GLN:HE22	1.65	0.62
1:BJ:46:ARG:HA	1:BL:65:TYR:HE2	1.64	0.62
1:CK:47:THR:HA	1:CK:77:PRO:HG2	1.82	0.62
1:CW:47:THR:HA	1:CW:77:PRO:HG2	1.82	0.62
1:CZ:47:THR:HA	1:CZ:77:PRO:HG2	1.82	0.62
1:AI:8:LYS:HB2	1:AI:114:GLN:HE22	1.65	0.62
1:AL:8:LYS:HB2	1:AL:114:GLN:HE22	1.65	0.62
1:BJ:8:LYS:HB2	1:BJ:114:GLN:HE22	1.65	0.62
1:BJ:92:ILE:HG12	1:DF:92:ILE:HG12	1.82	0.62
1:BV:8:LYS:HB2	1:BV:114:GLN:HE22	1.65	0.62
1:DL:8:LYS:HB2	1:DL:114:GLN:HE22	1.65	0.62
1:BD:8:LYS:HB2	1:BD:114:GLN:HE22	1.65	0.61
1:BG:47:THR:HA	1:BG:77:PRO:HG2	1.82	0.61
1:AO:8:LYS:HB2	1:AO:114:GLN:HE22	1.65	0.61
1:BM:47:THR:HA	1:BM:77:PRO:HG2	1.82	0.61
1:BV:47:THR:HA	1:BV:77:PRO:HG2	1.82	0.61
1:CT:42:ASP:HB2	1:CT:48:ARG:HG2	1.83	0.61
1:DI:8:LYS:HB2	1:DI:114:GLN:HE22	1.65	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AI:47:THR:HA	1:AI:77:PRO:HG2	1.82	0.61
1:AO:46:ARG:HG2	1:CP:65:TYR:CD2	2.34	0.61
1:AU:8:LYS:HB2	1:AU:114:GLN:HE22	1.65	0.61
1:BA:8:LYS:HB2	1:BA:114:GLN:HE22	1.65	0.61
1:BD:22:ILE:HG22	1:CZ:7:ILE:O	2.01	0.61
1:BP:47:THR:HA	1:BP:77:PRO:HG2	1.82	0.61
1:BY:42:ASP:HB2	1:BY:48:ARG:HG2	1.83	0.61
1:CW:8:LYS:HB2	1:CW:114:GLN:HE22	1.65	0.61
1:AO:42:ASP:HB2	1:AO:48:ARG:HG2	1.83	0.61
1:BG:8:LYS:HB2	1:BG:114:GLN:HE22	1.65	0.61
1:BM:8:LYS:HB2	1:BM:114:GLN:HE22	1.65	0.61
1:BM:111:ILE:HD13	1:DI:74:PHE:CZ	2.34	0.61
1:BS:42:ASP:HB2	1:BS:48:ARG:HG2	1.83	0.61
1:CK:8:LYS:HB2	1:CK:114:GLN:HE22	1.65	0.61
1:CZ:8:LYS:HB2	1:CZ:114:GLN:HE22	1.65	0.61
1:CN:42:ASP:HB2	1:CN:48:ARG:HG2	1.83	0.61
1:AC:8:LYS:HB2	1:AC:114:GLN:HE22	1.65	0.61
1:AO:47:THR:HA	1:AO:77:PRO:HG2	1.82	0.61
1:AX:42:ASP:HB2	1:AX:48:ARG:HG2	1.83	0.61
1:DI:47:THR:HA	1:DI:77:PRO:HG2	1.82	0.61
1:AC:47:THR:HA	1:AC:77:PRO:HG2	1.82	0.61
1:AF:19:GLY:O	1:BV:8:LYS:HD2	2.00	0.61
1:AF:42:ASP:HB2	1:AF:48:ARG:HG2	1.83	0.61
1:BY:8:LYS:HB2	1:BY:114:GLN:HE22	1.65	0.61
1:CN:47:THR:HA	1:CN:77:PRO:HG2	1.82	0.61
1:CQ:8:LYS:HB2	1:CQ:114:GLN:HE22	1.65	0.61
1:AU:47:THR:HA	1:AU:77:PRO:HG2	1.82	0.61
1:AX:47:THR:HA	1:AX:77:PRO:HG2	1.82	0.61
1:BG:42:ASP:HB2	1:BG:48:ARG:HG2	1.83	0.61
1:BM:114:GLN:HE21	1:DI:16:ILE:HD13	1.65	0.61
1:CK:42:ASP:HB2	1:CK:48:ARG:HG2	1.83	0.61
1:AI:42:ASP:HB2	1:AI:48:ARG:HG2	1.83	0.61
1:AR:8:LYS:HB2	1:AR:114:GLN:HE22	1.65	0.61
1:AR:42:ASP:HB2	1:AR:48:ARG:HG2	1.83	0.61
1:AU:42:ASP:HB2	1:AU:48:ARG:HG2	1.83	0.61
1:BM:42:ASP:HB2	1:BM:48:ARG:HG2	1.83	0.61
1:BP:8:LYS:HB2	1:BP:114:GLN:HE22	1.65	0.61
1:BP:42:ASP:HB2	1:BP:48:ARG:HG2	1.83	0.61
1:CE:8:LYS:HB2	1:CE:114:GLN:HE22	1.65	0.61
1:DF:47:THR:HA	1:DF:77:PRO:HG2	1.82	0.61
1:AJ:65:TYR:CE2	1:CV:46:ARG:HG3	2.36	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AL:47:THR:HA	1:AL:77:PRO:HG2	1.82	0.60
1:BF:65:TYR:CE2	1:BS:46:ARG:HA	2.36	0.60
1:BS:8:LYS:HB2	1:BS:114:GLN:HE22	1.65	0.60
1:BY:47:THR:HA	1:BY:77:PRO:HG2	1.82	0.60
1:CH:8:LYS:HB2	1:CH:114:GLN:HE22	1.65	0.60
1:CZ:42:ASP:HB2	1:CZ:48:ARG:HG2	1.83	0.60
1:AF:47:THR:HA	1:AF:77:PRO:HG2	1.82	0.60
1:AU:110:LEU:HD23	1:CN:16:ILE:HD11	1.83	0.60
1:CT:47:THR:HA	1:CT:77:PRO:HG2	1.82	0.60
1:DC:42:ASP:HB2	1:DC:48:ARG:HG2	1.83	0.60
1:BD:42:ASP:HB2	1:BD:48:ARG:HG2	1.83	0.60
1:BD:47:THR:HA	1:BD:77:PRO:HG2	1.82	0.60
1:CQ:42:ASP:HB2	1:CQ:48:ARG:HG2	1.83	0.60
1:AC:110:LEU:HD23	1:BP:16:ILE:HD11	1.83	0.60
1:AX:8:LYS:HB2	1:AX:114:GLN:HE22	1.65	0.60
1:BV:42:ASP:HB2	1:BV:48:ARG:HG2	1.83	0.60
1:AB:87:PHE:CE2	1:AS:60:VAL:HG11	2.37	0.60
1:AN:29:GLN:HG3	1:BZ:57:VAL:HG21	1.84	0.60
1:CN:8:LYS:HB2	1:CN:114:GLN:HE22	1.65	0.60
1:AC:42:ASP:HB2	1:AC:48:ARG:HG2	1.83	0.60
1:CT:8:LYS:HB2	1:CT:114:GLN:HE22	1.65	0.60
1:CW:46:ARG:HG2	1:CY:65:TYR:CE2	2.36	0.60
1:AC:16:ILE:HD11	1:BP:110:LEU:HD23	1.84	0.60
1:CB:42:ASP:HB2	1:CB:48:ARG:HG2	1.83	0.60
1:CH:47:THR:HA	1:CH:77:PRO:HG2	1.82	0.60
1:CB:8:LYS:HB2	1:CB:114:GLN:HE22	1.65	0.60
1:AR:47:THR:HA	1:AR:77:PRO:HG2	1.82	0.60
1:BJ:42:ASP:HB2	1:BJ:48:ARG:HG2	1.83	0.60
1:CH:42:ASP:HB2	1:CH:48:ARG:HG2	1.83	0.60
1:AF:22:ILE:O	1:BV:6:SER:HA	2.02	0.60
1:BD:110:LEU:HD23	1:CZ:16:ILE:HD11	1.83	0.60
1:DL:47:THR:HA	1:DL:77:PRO:HG2	1.82	0.60
1:CE:42:ASP:HB2	1:CE:48:ARG:HG2	1.83	0.59
1:CW:42:ASP:HB2	1:CW:48:ARG:HG2	1.83	0.59
1:AF:8:LYS:HD2	1:BV:19:GLY:O	2.02	0.59
1:AF:114:GLN:NE2	1:BV:16:ILE:HD13	2.15	0.59
1:AO:109:ARG:HG2	1:CE:116:LEU:HD22	1.84	0.59
1:BA:42:ASP:HB2	1:BA:48:ARG:HG2	1.83	0.59
1:DC:8:LYS:HB2	1:DC:114:GLN:HE22	1.65	0.59
1:DF:8:LYS:HB2	1:DF:114:GLN:HE22	1.65	0.59
1:DL:42:ASP:HB2	1:DL:48:ARG:HG2	1.83	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AC:111:ILE:HD13	1:BP:74:PHE:CZ	2.37	0.59
1:AM:130:LEU:O	1:CB:26:LEU:HB3	2.02	0.59
1:BA:47:THR:HA	1:BA:77:PRO:HG2	1.82	0.59
1:CW:10:GLY:N	1:CW:107:ASP:OD1	2.34	0.59
1:DF:42:ASP:HB2	1:DF:48:ARG:HG2	1.83	0.59
1:AB:46:ARG:HG3	1:AS:65:TYR:CE2	2.38	0.59
1:DI:42:ASP:HB2	1:DI:48:ARG:HG2	1.83	0.59
1:DF:10:GLY:N	1:DF:107:ASP:OD1	2.34	0.59
1:AH:46:ARG:NH2	1:CC:99:ALA:HA	2.18	0.59
1:AI:116:LEU:HD22	1:BY:109:ARG:HG2	1.85	0.59
1:BO:87:PHE:CE2	1:DJ:60:VAL:HG11	2.38	0.59
1:CD:65:TYR:HE2	1:CK:46:ARG:CA	2.14	0.59
1:AG:130:LEU:O	1:CE:26:LEU:HB3	2.03	0.59
1:AL:42:ASP:HB2	1:AL:48:ARG:HG2	1.83	0.59
1:BU:29:GLN:HG3	1:DG:57:VAL:HG21	1.85	0.59
1:AE:33:ASN:OD1	1:BG:30:THR:OG1	2.20	0.59
1:BI:127:LEU:HD21	1:BL:48:ARG:HH12	1.68	0.59
1:BM:115:LEU:HD13	1:DI:72:VAL:HG11	1.84	0.59
1:AH:33:ASN:OD1	1:CE:30:THR:OG1	2.20	0.59
1:AO:108:LEU:HD22	1:CE:90:VAL:HG12	1.85	0.58
1:BG:110:LEU:HD23	1:DC:16:ILE:HD11	1.85	0.58
1:BM:116:LEU:O	1:DI:109:ARG:NH2	2.35	0.58
1:AF:16:ILE:HD13	1:BV:114:GLN:NE2	2.13	0.58
1:BG:16:ILE:HD11	1:DC:110:LEU:HD23	1.85	0.58
1:CW:46:ARG:HA	1:CY:65:TYR:HE2	1.67	0.58
1:AO:109:ARG:NH2	1:CE:116:LEU:O	2.37	0.58
1:CZ:10:GLY:N	1:CZ:107:ASP:OD1	2.34	0.58
1:AK:65:TYR:CD2	1:BY:46:ARG:HG2	2.39	0.58
1:BK:39:VAL:O	1:BK:48:ARG:HD2	2.04	0.58
1:AG:39:VAL:O	1:AG:48:ARG:HD2	2.04	0.58
1:BZ:39:VAL:O	1:BZ:48:ARG:HD2	2.04	0.58
1:DL:10:GLY:N	1:DL:107:ASP:OD1	2.34	0.58
1:AI:116:LEU:O	1:BY:109:ARG:NH2	2.36	0.58
1:AM:39:VAL:O	1:AM:48:ARG:HD2	2.04	0.58
1:AO:46:ARG:HG2	1:CP:65:TYR:CE2	2.39	0.58
1:AT:46:ARG:NH2	1:AV:99:ALA:HA	2.19	0.58
1:CO:39:VAL:O	1:CO:48:ARG:HD2	2.04	0.58
1:CR:39:VAL:O	1:CR:48:ARG:HD2	2.04	0.58
1:AA:39:VAL:O	1:AA:48:ARG:HD2	2.04	0.58
1:AY:39:VAL:O	1:AY:48:ARG:HD2	2.04	0.58
1:BA:67:LYS:NZ	1:CW:87:PHE:O	2.34	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BE:39:VAL:O	1:BE:48:ARG:HD2	2.04	0.58
1:AJ:39:VAL:O	1:AJ:48:ARG:HD2	2.04	0.57
1:AO:113:ALA:HB1	1:CE:113:ALA:HB1	1.85	0.57
1:BT:39:VAL:O	1:BT:48:ARG:HD2	2.04	0.57
1:AS:39:VAL:O	1:AS:48:ARG:HD2	2.04	0.57
1:CK:10:GLY:N	1:CK:107:ASP:OD1	2.34	0.57
1:CX:39:VAL:O	1:CX:48:ARG:HD2	2.04	0.57
1:DA:39:VAL:O	1:DA:48:ARG:HD2	2.04	0.57
1:AC:10:GLY:N	1:AC:107:ASP:OD1	2.34	0.57
1:DJ:39:VAL:O	1:DJ:48:ARG:HD2	2.04	0.57
1:BH:39:VAL:O	1:BH:48:ARG:HD2	2.04	0.57
1:BJ:7:ILE:O	1:DF:22:ILE:HG22	2.04	0.57
1:BQ:39:VAL:O	1:BQ:48:ARG:HD2	2.04	0.57
1:CI:39:VAL:O	1:CI:48:ARG:HD2	2.04	0.57
1:CT:10:GLY:N	1:CT:107:ASP:OD1	2.34	0.57
1:DC:10:GLY:N	1:DC:107:ASP:OD1	2.34	0.57
1:AB:47:THR:HA	1:AB:77:PRO:HG2	1.87	0.57
1:AJ:60:VAL:HG11	1:CV:87:PHE:CE2	2.39	0.57
1:AQ:47:THR:HA	1:AQ:77:PRO:HG2	1.87	0.57
1:AX:10:GLY:N	1:AX:107:ASP:OD1	2.34	0.57
1:DG:39:VAL:O	1:DG:48:ARG:HD2	2.04	0.57
1:AP:39:VAL:O	1:AP:48:ARG:HD2	2.04	0.57
1:BA:10:GLY:N	1:BA:107:ASP:OD1	2.34	0.57
1:BB:39:VAL:O	1:BB:48:ARG:HD2	2.04	0.57
1:BU:47:THR:HA	1:BU:77:PRO:HG2	1.87	0.57
1:CA:60:VAL:HG13	1:CA:65:TYR:HE1	1.70	0.57
1:CC:39:VAL:O	1:CC:48:ARG:HD2	2.04	0.57
1:DD:39:VAL:O	1:DD:48:ARG:HD2	2.04	0.57
1:DE:47:THR:HA	1:DE:77:PRO:HG2	1.87	0.57
1:AO:116:LEU:HD22	1:CE:109:ARG:HG2	1.87	0.57
1:AZ:47:THR:HA	1:AZ:77:PRO:HG2	1.87	0.57
1:AZ:60:VAL:HG13	1:AZ:65:TYR:HE1	1.70	0.57
1:BG:10:GLY:N	1:BG:107:ASP:OD1	2.34	0.57
1:AO:116:LEU:O	1:CE:109:ARG:NE	2.36	0.57
1:AQ:60:VAL:HG13	1:AQ:65:TYR:HE1	1.70	0.57
1:AU:10:GLY:N	1:AU:107:ASP:OD1	2.34	0.57
1:BC:60:VAL:HG13	1:BC:65:TYR:HE1	1.70	0.57
1:BL:47:THR:HA	1:BL:77:PRO:HG2	1.87	0.57
1:BM:90:VAL:HG12	1:DI:108:LEU:HD22	1.86	0.57
1:BW:39:VAL:O	1:BW:48:ARG:HD2	2.04	0.57
1:CF:39:VAL:O	1:CF:48:ARG:HD2	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CY:47:THR:HA	1:CY:77:PRO:HG2	1.87	0.57
1:DB:47:THR:HA	1:DB:77:PRO:HG2	1.87	0.57
1:AR:6:SER:HA	1:CK:22:ILE:O	2.04	0.57
1:AT:47:THR:HA	1:AT:77:PRO:HG2	1.87	0.57
1:BP:10:GLY:N	1:BP:107:ASP:OD1	2.34	0.57
1:BU:60:VAL:HG13	1:BU:65:TYR:HE1	1.70	0.57
1:CG:60:VAL:HG13	1:CG:65:TYR:HE1	1.70	0.57
1:DE:60:VAL:HG13	1:DE:65:TYR:HE1	1.70	0.57
1:AI:129:ALA:HB2	1:BW:26:LEU:O	2.05	0.56
1:AK:60:VAL:HG13	1:AK:65:TYR:HE1	1.70	0.56
1:BM:10:GLY:N	1:BM:107:ASP:OD1	2.34	0.56
1:BR:47:THR:HA	1:BR:77:PRO:HG2	1.87	0.56
1:BR:60:VAL:HG13	1:BR:65:TYR:HE1	1.70	0.56
1:BX:47:THR:HA	1:BX:77:PRO:HG2	1.87	0.56
1:CD:47:THR:HA	1:CD:77:PRO:HG2	1.87	0.56
1:CP:47:THR:HA	1:CP:77:PRO:HG2	1.87	0.56
1:AI:109:ARG:NH2	1:BY:116:LEU:O	2.37	0.56
1:AT:60:VAL:HG13	1:AT:65:TYR:HE1	1.70	0.56
1:BN:39:VAL:O	1:BN:48:ARG:HD2	2.04	0.56
1:BX:60:VAL:HG13	1:BX:65:TYR:HE1	1.70	0.56
1:CD:60:VAL:HG13	1:CD:65:TYR:HE1	1.70	0.56
1:CN:10:GLY:N	1:CN:107:ASP:OD1	2.34	0.56
1:DK:60:VAL:HG13	1:DK:65:TYR:HE1	1.70	0.56
1:AD:39:VAL:O	1:AD:48:ARG:HD2	2.04	0.56
1:AN:47:THR:HA	1:AN:77:PRO:HG2	1.87	0.56
1:BF:47:THR:HA	1:BF:77:PRO:HG2	1.87	0.56
1:CP:60:VAL:HG13	1:CP:65:TYR:HE1	1.70	0.56
1:CU:39:VAL:O	1:CU:48:ARG:HD2	2.04	0.56
1:CV:60:VAL:HG13	1:CV:65:TYR:HE1	1.70	0.56
1:DH:47:THR:HA	1:DH:77:PRO:HG2	1.87	0.56
1:AB:60:VAL:HG13	1:AB:65:TYR:HE1	1.70	0.56
1:AU:7:ILE:O	1:CN:22:ILE:HG22	2.05	0.56
1:BU:87:PHE:HE2	1:DG:60:VAL:HG11	1.69	0.56
1:DB:60:VAL:HG13	1:DB:65:TYR:HE1	1.70	0.56
1:AI:10:GLY:N	1:AI:107:ASP:OD1	2.34	0.56
1:AK:87:PHE:HE2	1:BW:60:VAL:HG11	1.71	0.56
1:AO:109:ARG:NE	1:CE:116:LEU:O	2.39	0.56
1:CH:10:GLY:N	1:CH:107:ASP:OD1	2.34	0.56
1:CJ:47:THR:HA	1:CJ:77:PRO:HG2	1.87	0.56
1:AI:22:ILE:HG22	1:BY:7:ILE:O	2.05	0.56
1:BD:22:ILE:O	1:CZ:6:SER:HA	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BO:47:THR:HA	1:BO:77:PRO:HG2	1.87	0.56
1:CL:39:VAL:O	1:CL:48:ARG:HD2	2.04	0.56
1:CL:119:ALA:HA	1:CL:122:ASP:HB2	1.88	0.56
1:AE:60:VAL:HG13	1:AE:65:TYR:HE1	1.70	0.56
1:AR:92:ILE:HG12	1:CK:92:ILE:HG12	1.86	0.56
1:AU:6:SER:HA	1:CN:22:ILE:O	2.06	0.56
1:AV:39:VAL:O	1:AV:48:ARG:HD2	2.04	0.56
1:BC:47:THR:HA	1:BC:77:PRO:HG2	1.87	0.56
1:CA:65:TYR:HE2	1:CH:46:ARG:HA	1.71	0.56
1:CI:119:ALA:HA	1:CI:122:ASP:HB2	1.88	0.56
1:CJ:60:VAL:HG13	1:CJ:65:TYR:HE1	1.70	0.56
1:CX:119:ALA:HA	1:CX:122:ASP:HB2	1.88	0.56
1:CY:60:VAL:HG13	1:CY:65:TYR:HE1	1.70	0.56
1:DH:60:VAL:HG13	1:DH:65:TYR:HE1	1.70	0.56
1:AA:119:ALA:HA	1:AA:122:ASP:HB2	1.88	0.56
1:AM:119:ALA:HA	1:AM:122:ASP:HB2	1.88	0.56
1:AW:60:VAL:HG13	1:AW:65:TYR:HE1	1.70	0.56
1:BH:119:ALA:HA	1:BH:122:ASP:HB2	1.88	0.56
1:BI:60:VAL:HG13	1:BI:65:TYR:HE1	1.70	0.56
1:BU:22:ILE:HA	1:BU:41:GLU:OE1	2.06	0.56
1:BZ:119:ALA:HA	1:BZ:122:ASP:HB2	1.88	0.56
1:CJ:34:GLY:HA3	1:CJ:53:PHE:O	2.06	0.56
1:CW:46:ARG:HG2	1:CY:65:TYR:HD2	1.67	0.56
1:AO:12:THR:OG1	1:CE:17:THR:O	2.20	0.56
1:AW:22:ILE:HA	1:AW:41:GLU:OE1	2.06	0.56
1:BB:119:ALA:HA	1:BB:122:ASP:HB2	1.88	0.56
1:BF:60:VAL:HG13	1:BF:65:TYR:HE1	1.70	0.56
1:BR:22:ILE:HA	1:BR:41:GLU:OE1	2.06	0.56
1:CO:119:ALA:HA	1:CO:122:ASP:HB2	1.88	0.56
1:CP:22:ILE:HA	1:CP:41:GLU:OE1	2.06	0.56
1:CU:119:ALA:HA	1:CU:122:ASP:HB2	1.88	0.56
1:CV:22:ILE:HA	1:CV:41:GLU:OE1	2.06	0.56
1:DB:22:ILE:HA	1:DB:41:GLU:OE1	2.06	0.56
1:DG:119:ALA:HA	1:DG:122:ASP:HB2	1.88	0.56
1:AK:47:THR:HA	1:AK:77:PRO:HG2	1.87	0.56
1:AW:34:GLY:HA3	1:AW:53:PHE:O	2.06	0.56
1:BI:22:ILE:HA	1:BI:41:GLU:OE1	2.06	0.56
1:BL:22:ILE:HA	1:BL:41:GLU:OE1	2.06	0.56
1:BM:108:LEU:HD22	1:DI:90:VAL:CG1	2.33	0.56
1:BY:10:GLY:N	1:BY:107:ASP:OD1	2.34	0.56
1:CR:119:ALA:HA	1:CR:122:ASP:HB2	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CV:47:THR:HA	1:CV:77:PRO:HG2	1.87	0.56
1:AD:119:ALA:HA	1:AD:122:ASP:HB2	1.88	0.55
1:AK:34:GLY:HA3	1:AK:53:PHE:O	2.06	0.55
1:AN:22:ILE:HA	1:AN:41:GLU:OE1	2.06	0.55
1:AW:47:THR:HA	1:AW:77:PRO:HG2	1.87	0.55
1:BC:22:ILE:HA	1:BC:41:GLU:OE1	2.06	0.55
1:CA:47:THR:HA	1:CA:77:PRO:HG2	1.87	0.55
1:CD:22:ILE:HA	1:CD:41:GLU:OE1	2.06	0.55
1:CD:98:PRO:HB2	1:CK:46:ARG:HG3	1.87	0.55
1:AE:47:THR:HA	1:AE:77:PRO:HG2	1.87	0.55
1:AH:60:VAL:HG13	1:AH:65:TYR:HE1	1.70	0.55
1:AN:34:GLY:HA3	1:AN:53:PHE:O	2.06	0.55
1:AS:119:ALA:HA	1:AS:122:ASP:HB2	1.88	0.55
1:BF:65:TYR:HE2	1:BS:46:ARG:CA	2.19	0.55
1:BK:119:ALA:HA	1:BK:122:ASP:HB2	1.88	0.55
1:BO:22:ILE:HA	1:BO:41:GLU:OE1	2.06	0.55
1:CF:119:ALA:HA	1:CF:122:ASP:HB2	1.88	0.55
1:CG:22:ILE:HA	1:CG:41:GLU:OE1	2.06	0.55
1:CP:34:GLY:HA3	1:CP:53:PHE:O	2.06	0.55
1:CS:22:ILE:HA	1:CS:41:GLU:OE1	2.06	0.55
1:CV:34:GLY:HA3	1:CV:53:PHE:O	2.06	0.55
1:DH:34:GLY:HA3	1:DH:53:PHE:O	2.06	0.55
1:AF:10:GLY:N	1:AF:107:ASP:OD1	2.34	0.55
1:AI:109:ARG:HG2	1:BY:116:LEU:HD22	1.88	0.55
1:AP:119:ALA:HA	1:AP:122:ASP:HB2	1.88	0.55
1:BC:34:GLY:HA3	1:BC:53:PHE:O	2.06	0.55
1:BO:60:VAL:HG13	1:BO:65:TYR:HE1	1.70	0.55
1:CA:22:ILE:HA	1:CA:41:GLU:OE1	2.06	0.55
1:CC:119:ALA:HA	1:CC:122:ASP:HB2	1.88	0.55
1:CG:34:GLY:HA3	1:CG:53:PHE:O	2.06	0.55
1:CM:47:THR:HA	1:CM:77:PRO:HG2	1.87	0.55
1:CM:60:VAL:HG13	1:CM:65:TYR:HE1	1.70	0.55
1:DJ:119:ALA:HA	1:DJ:122:ASP:HB2	1.88	0.55
1:AR:10:GLY:N	1:AR:107:ASP:OD1	2.34	0.55
1:BI:47:THR:HA	1:BI:77:PRO:HG2	1.87	0.55
1:BN:119:ALA:HA	1:BN:122:ASP:HB2	1.88	0.55
1:BX:22:ILE:HA	1:BX:41:GLU:OE1	2.06	0.55
1:CS:34:GLY:HA3	1:CS:53:PHE:O	2.06	0.55
1:CS:47:THR:HA	1:CS:77:PRO:HG2	1.87	0.55
1:CY:22:ILE:HA	1:CY:41:GLU:OE1	2.06	0.55
1:DE:34:GLY:HA3	1:DE:53:PHE:O	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:DH:22:ILE:HA	1:DH:41:GLU:OE1	2.06	0.55
1:DK:34:GLY:HA3	1:DK:53:PHE:O	2.06	0.55
1:AB:22:ILE:HA	1:AB:41:GLU:OE1	2.06	0.55
1:AE:34:GLY:HA3	1:AE:53:PHE:O	2.06	0.55
1:AH:34:GLY:HA3	1:AH:53:PHE:O	2.06	0.55
1:AZ:34:GLY:HA3	1:AZ:53:PHE:O	2.06	0.55
1:BQ:119:ALA:HA	1:BQ:122:ASP:HB2	1.88	0.55
1:BT:119:ALA:HA	1:BT:122:ASP:HB2	1.88	0.55
1:CM:22:ILE:HA	1:CM:41:GLU:OE1	2.06	0.55
1:CS:60:VAL:HG13	1:CS:65:TYR:HE1	1.70	0.55
1:DK:22:ILE:HA	1:DK:41:GLU:OE1	2.06	0.55
1:DK:47:THR:HA	1:DK:77:PRO:HG2	1.87	0.55
1:AH:22:ILE:HA	1:AH:41:GLU:OE1	2.06	0.55
1:AT:34:GLY:HA3	1:AT:53:PHE:O	2.06	0.55
1:BE:119:ALA:HA	1:BE:122:ASP:HB2	1.88	0.55
1:CA:34:GLY:HA3	1:CA:53:PHE:O	2.06	0.55
1:AV:119:ALA:HA	1:AV:122:ASP:HB2	1.88	0.55
1:BO:34:GLY:HA3	1:BO:53:PHE:O	2.06	0.55
1:BR:34:GLY:HA3	1:BR:53:PHE:O	2.06	0.55
1:CG:47:THR:HA	1:CG:77:PRO:HG2	1.87	0.55
1:DB:34:GLY:HA3	1:DB:53:PHE:O	2.06	0.55
1:AF:48:ARG:HH12	1:DD:127:LEU:HB3	1.72	0.55
1:AH:47:THR:HA	1:AH:77:PRO:HG2	1.87	0.55
1:AT:22:ILE:HA	1:AT:41:GLU:OE1	2.06	0.55
1:BX:34:GLY:HA3	1:BX:53:PHE:O	2.06	0.55
1:CE:10:GLY:N	1:CE:107:ASP:OD1	2.34	0.55
1:AZ:22:ILE:HA	1:AZ:41:GLU:OE1	2.06	0.55
1:BF:22:ILE:HA	1:BF:41:GLU:OE1	2.06	0.55
1:BL:60:VAL:HG13	1:BL:65:TYR:HE1	1.70	0.55
1:CD:34:GLY:HA3	1:CD:53:PHE:O	2.06	0.55
1:AI:12:THR:H	1:BY:18:GLY:HA2	1.71	0.55
1:AK:22:ILE:HA	1:AK:41:GLU:OE1	2.06	0.55
1:BI:34:GLY:HA3	1:BI:53:PHE:O	2.06	0.55
1:BS:10:GLY:N	1:BS:107:ASP:OD1	2.34	0.55
1:AI:90:VAL:HG12	1:BY:108:LEU:HD22	1.87	0.54
1:BF:34:GLY:HA3	1:BF:53:PHE:O	2.06	0.54
1:BU:34:GLY:HA3	1:BU:53:PHE:O	2.06	0.54
1:CQ:10:GLY:N	1:CQ:107:ASP:OD1	2.34	0.54
1:DD:119:ALA:HA	1:DD:122:ASP:HB2	1.88	0.54
1:AL:10:GLY:N	1:AL:107:ASP:OD1	2.34	0.54
1:AN:60:VAL:HG13	1:AN:65:TYR:HE1	1.70	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BU:87:PHE:CE2	1:DG:60:VAL:HG11	2.43	0.54
1:DA:119:ALA:HA	1:DA:122:ASP:HB2	1.88	0.54
1:AS:22:ILE:HD11	1:AS:39:VAL:HG11	1.90	0.54
1:AY:119:ALA:HA	1:AY:122:ASP:HB2	1.88	0.54
1:BJ:46:ARG:HG3	1:BL:98:PRO:HB2	1.88	0.54
1:BL:34:GLY:HA3	1:BL:53:PHE:O	2.06	0.54
1:CD:29:GLN:HG3	1:CI:57:VAL:HG21	1.90	0.54
1:CU:22:ILE:HD11	1:CU:39:VAL:HG11	1.90	0.54
1:CY:34:GLY:HA3	1:CY:53:PHE:O	2.06	0.54
1:DI:10:GLY:N	1:DI:107:ASP:OD1	2.34	0.54
1:DJ:22:ILE:HD11	1:DJ:39:VAL:HG11	1.90	0.54
1:AE:22:ILE:HA	1:AE:41:GLU:OE1	2.06	0.54
1:AQ:34:GLY:HA3	1:AQ:53:PHE:O	2.06	0.54
1:AR:109:ARG:NH1	1:CK:122:ASP:OD1	2.35	0.54
1:BB:22:ILE:HD11	1:BB:39:VAL:HG11	1.90	0.54
1:BH:22:ILE:HD11	1:BH:39:VAL:HG11	1.90	0.54
1:BT:22:ILE:HD11	1:BT:39:VAL:HG11	1.90	0.54
1:BW:119:ALA:HA	1:BW:122:ASP:HB2	1.88	0.54
1:DE:22:ILE:HA	1:DE:41:GLU:OE1	2.06	0.54
1:AB:34:GLY:HA3	1:AB:53:PHE:O	2.06	0.54
1:AE:46:ARG:HG3	1:BE:65:TYR:CD2	2.42	0.54
1:AG:119:ALA:HA	1:AG:122:ASP:HB2	1.88	0.54
1:AI:116:LEU:O	1:BY:109:ARG:NE	2.38	0.54
1:AJ:119:ALA:HA	1:AJ:122:ASP:HB2	1.88	0.54
1:AM:22:ILE:HD11	1:AM:39:VAL:HG11	1.90	0.54
1:AQ:22:ILE:HA	1:AQ:41:GLU:OE1	2.06	0.54
1:BD:16:ILE:HD11	1:CZ:110:LEU:HD23	1.90	0.54
1:BJ:46:ARG:HA	1:BL:65:TYR:CE2	2.41	0.54
1:CJ:22:ILE:HA	1:CJ:41:GLU:OE1	2.06	0.54
1:CM:34:GLY:HA3	1:CM:53:PHE:O	2.06	0.54
1:AM:60:VAL:HG11	1:CP:87:PHE:CE2	2.43	0.54
1:CD:58:PRO:HB2	1:CK:45:TYR:CE2	2.43	0.54
1:AA:22:ILE:HD11	1:AA:39:VAL:HG11	1.90	0.54
1:AD:22:ILE:HD11	1:AD:39:VAL:HG11	1.90	0.54
1:BI:127:LEU:HD23	1:BL:48:ARG:HH22	1.72	0.54
1:AP:22:ILE:HD11	1:AP:39:VAL:HG11	1.90	0.54
1:DD:22:ILE:HD11	1:DD:39:VAL:HG11	1.90	0.54
1:AD:33:ASN:OD1	1:DE:30:THR:OG1	2.20	0.53
1:AI:130:LEU:HD21	1:BY:55:SER:CB	2.37	0.53
1:BO:29:GLN:HG3	1:DJ:57:VAL:HG21	1.90	0.53
1:BD:111:ILE:HD13	1:CZ:74:PHE:CZ	2.42	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CC:22:ILE:HD11	1:CC:39:VAL:HG11	1.90	0.53
1:AJ:60:VAL:HG11	1:CV:87:PHE:HE2	1.73	0.53
1:AO:10:GLY:N	1:AO:107:ASP:OD1	2.34	0.53
1:AY:22:ILE:HD11	1:AY:39:VAL:HG11	1.90	0.53
1:CR:22:ILE:HD11	1:CR:39:VAL:HG11	1.90	0.53
1:DG:22:ILE:HD11	1:DG:39:VAL:HG11	1.90	0.53
1:AJ:22:ILE:HD11	1:AJ:39:VAL:HG11	1.90	0.53
1:AO:116:LEU:O	1:CE:109:ARG:NH2	2.40	0.53
1:AG:22:ILE:HD11	1:AG:39:VAL:HG11	1.90	0.53
1:AO:90:VAL:HG12	1:CE:108:LEU:HD22	1.89	0.53
1:BN:39:VAL:HG23	1:BN:42:ASP:H	1.74	0.53
1:BW:22:ILE:HD11	1:BW:39:VAL:HG11	1.90	0.53
1:CI:22:ILE:HD11	1:CI:39:VAL:HG11	1.90	0.53
1:CJ:29:GLN:HG3	1:DA:57:VAL:HG21	1.91	0.53
1:DJ:39:VAL:HG23	1:DJ:42:ASP:H	1.74	0.53
1:AV:22:ILE:HD11	1:AV:39:VAL:HG11	1.90	0.53
1:BJ:10:GLY:N	1:BJ:107:ASP:OD1	2.34	0.53
1:BW:39:VAL:HG23	1:BW:42:ASP:H	1.74	0.53
1:BX:87:PHE:HE2	1:CL:60:VAL:HG11	1.74	0.53
1:CF:39:VAL:HG23	1:CF:42:ASP:H	1.74	0.53
1:CX:22:ILE:HD11	1:CX:39:VAL:HG11	1.90	0.53
1:AA:39:VAL:HG23	1:AA:42:ASP:H	1.74	0.53
1:AV:39:VAL:HG23	1:AV:42:ASP:H	1.74	0.53
1:AG:39:VAL:HG23	1:AG:42:ASP:H	1.74	0.53
1:BK:22:ILE:HD11	1:BK:39:VAL:HG11	1.90	0.53
1:BQ:22:ILE:HD11	1:BQ:39:VAL:HG11	1.90	0.53
1:CB:10:GLY:N	1:CB:107:ASP:OD1	2.34	0.53
1:CI:39:VAL:HG23	1:CI:42:ASP:H	1.74	0.53
1:CM:29:GLN:HG3	1:CX:57:VAL:HG21	1.90	0.53
1:DA:22:ILE:HD11	1:DA:39:VAL:HG11	1.90	0.53
1:AF:22:ILE:HD12	1:AF:49:ARG:HH21	1.74	0.53
1:AK:48:ARG:HH12	1:BX:127:LEU:HD21	1.73	0.53
1:BL:49:ARG:HG2	1:BL:76:ILE:HG12	1.91	0.53
1:CF:22:ILE:HD11	1:CF:39:VAL:HG11	1.90	0.53
1:DG:39:VAL:HG23	1:DG:42:ASP:H	1.74	0.53
1:DH:49:ARG:HG2	1:DH:76:ILE:HG12	1.91	0.53
1:AO:22:ILE:HD12	1:AO:49:ARG:HH21	1.74	0.53
1:BA:22:ILE:HD12	1:BA:49:ARG:HH21	1.74	0.53
1:BD:22:ILE:HD12	1:BD:49:ARG:HH21	1.74	0.53
1:BM:130:LEU:HD21	1:DI:55:SER:CB	2.35	0.53
1:CE:22:ILE:HD12	1:CE:49:ARG:HH21	1.74	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AG:60:VAL:HG11	1:CS:87:PHE:CE2	2.44	0.52
1:AX:2:ILE:HD11	1:AX:124:PHE:HB2	1.92	0.52
1:BG:2:ILE:HD11	1:BG:124:PHE:HB2	1.91	0.52
1:BH:39:VAL:HG23	1:BH:42:ASP:H	1.74	0.52
1:BN:22:ILE:HD11	1:BN:39:VAL:HG11	1.90	0.52
1:BO:49:ARG:HG2	1:BO:76:ILE:HG12	1.91	0.52
1:CO:39:VAL:HG23	1:CO:42:ASP:H	1.74	0.52
1:DI:22:ILE:HD12	1:DI:49:ARG:HH21	1.74	0.52
1:DL:22:ILE:HD12	1:DL:49:ARG:HH21	1.74	0.52
1:AD:39:VAL:HG23	1:AD:42:ASP:H	1.74	0.52
1:AL:2:ILE:HD11	1:AL:124:PHE:HB2	1.92	0.52
1:AT:49:ARG:HG2	1:AT:76:ILE:HG12	1.91	0.52
1:BJ:2:ILE:HD11	1:BJ:124:PHE:HB2	1.92	0.52
1:BT:39:VAL:HG23	1:BT:42:ASP:H	1.74	0.52
1:BX:49:ARG:HG2	1:BX:76:ILE:HG12	1.91	0.52
1:BY:22:ILE:HD12	1:BY:49:ARG:HH21	1.74	0.52
1:CT:2:ILE:HD11	1:CT:124:PHE:HB2	1.92	0.52
1:DE:49:ARG:HG2	1:DE:76:ILE:HG12	1.91	0.52
1:AE:49:ARG:HG2	1:AE:76:ILE:HG12	1.91	0.52
1:AI:7:ILE:O	1:BY:22:ILE:HG22	2.08	0.52
1:AN:49:ARG:HG2	1:AN:76:ILE:HG12	1.91	0.52
1:AR:22:ILE:HD12	1:AR:49:ARG:HH21	1.74	0.52
1:AZ:49:ARG:HG2	1:AZ:76:ILE:HG12	1.91	0.52
1:BD:10:GLY:N	1:BD:107:ASP:OD1	2.34	0.52
1:CD:85:THR:N	1:CI:63:GLY:O	2.35	0.52
1:CE:2:ILE:HD11	1:CE:124:PHE:HB2	1.92	0.52
1:CN:22:ILE:HD12	1:CN:49:ARG:HH21	1.74	0.52
1:CO:22:ILE:HD11	1:CO:39:VAL:HG11	1.90	0.52
1:CQ:22:ILE:HD12	1:CQ:49:ARG:HH21	1.74	0.52
1:CR:39:VAL:HG23	1:CR:42:ASP:H	1.74	0.52
1:CZ:45:TYR:HA	1:CZ:48:ARG:HE	1.75	0.52
1:DC:22:ILE:HD12	1:DC:49:ARG:HH21	1.74	0.52
1:AC:45:TYR:HA	1:AC:48:ARG:HE	1.75	0.52
1:AX:22:ILE:HD12	1:AX:49:ARG:HH21	1.74	0.52
1:BA:2:ILE:HD11	1:BA:124:PHE:HB2	1.92	0.52
1:BC:49:ARG:HG2	1:BC:76:ILE:HG12	1.91	0.52
1:BM:113:ALA:HB1	1:DI:113:ALA:HB1	1.90	0.52
1:BQ:39:VAL:HG23	1:BQ:42:ASP:H	1.74	0.52
1:CJ:49:ARG:HG2	1:CJ:76:ILE:HG12	1.92	0.52
1:CL:22:ILE:HD11	1:CL:39:VAL:HG11	1.90	0.52
1:AI:2:ILE:HD11	1:AI:124:PHE:HB2	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AO:2:ILE:HD11	1:AO:124:PHE:HB2	1.92	0.52
1:AR:45:TYR:HA	1:AR:48:ARG:HE	1.75	0.52
1:BD:45:TYR:HA	1:BD:48:ARG:HE	1.75	0.52
1:BE:39:VAL:HG23	1:BE:42:ASP:H	1.74	0.52
1:BG:108:LEU:HD22	1:DC:90:VAL:HG12	1.92	0.52
1:BJ:45:TYR:HA	1:BJ:48:ARG:HE	1.75	0.52
1:BP:45:TYR:HA	1:BP:48:ARG:HE	1.75	0.52
1:CH:45:TYR:HB2	1:CH:48:ARG:HH21	1.75	0.52
1:CQ:2:ILE:HD11	1:CQ:124:PHE:HB2	1.92	0.52
1:CS:49:ARG:HG2	1:CS:76:ILE:HG12	1.91	0.52
1:CW:2:ILE:HD11	1:CW:124:PHE:HB2	1.92	0.52
1:CW:22:ILE:HD12	1:CW:49:ARG:HH21	1.74	0.52
1:CX:39:VAL:HG23	1:CX:42:ASP:H	1.74	0.52
1:DF:2:ILE:HD11	1:DF:124:PHE:HB2	1.92	0.52
1:DF:45:TYR:HA	1:DF:48:ARG:HE	1.75	0.52
1:AL:22:ILE:HD12	1:AL:49:ARG:HH21	1.74	0.52
1:AL:45:TYR:HB2	1:AL:48:ARG:HH21	1.75	0.52
1:AU:45:TYR:HB2	1:AU:48:ARG:HH21	1.75	0.52
1:BE:22:ILE:HD11	1:BE:39:VAL:HG11	1.90	0.52
1:CG:49:ARG:HG2	1:CG:76:ILE:HG12	1.92	0.52
1:CH:2:ILE:HD11	1:CH:124:PHE:HB2	1.92	0.52
1:CH:22:ILE:HD12	1:CH:49:ARG:HH21	1.74	0.52
1:CK:45:TYR:HA	1:CK:48:ARG:HE	1.75	0.52
1:DF:22:ILE:HD12	1:DF:49:ARG:HH21	1.74	0.52
1:AF:45:TYR:HA	1:AF:48:ARG:HE	1.75	0.52
1:AJ:39:VAL:HG23	1:AJ:42:ASP:H	1.74	0.52
1:AK:87:PHE:CE2	1:BW:60:VAL:HG11	2.45	0.52
1:AS:39:VAL:HG23	1:AS:42:ASP:H	1.74	0.52
1:AX:45:TYR:HA	1:AX:48:ARG:HE	1.75	0.52
1:BB:39:VAL:HG23	1:BB:42:ASP:H	1.74	0.52
1:BM:22:ILE:HD12	1:BM:49:ARG:HH21	1.74	0.52
1:BS:22:ILE:HD12	1:BS:49:ARG:HH21	1.74	0.52
1:BV:45:TYR:HA	1:BV:48:ARG:HE	1.75	0.52
1:BX:65:TYR:HE2	1:CN:46:ARG:CA	2.17	0.52
1:CK:22:ILE:HD12	1:CK:49:ARG:HH21	1.74	0.52
1:CL:39:VAL:HG23	1:CL:42:ASP:H	1.74	0.52
1:AC:2:ILE:HD11	1:AC:124:PHE:HB2	1.92	0.52
1:AO:114:GLN:NE2	1:CE:16:ILE:HG21	2.24	0.52
1:BS:2:ILE:HD11	1:BS:124:PHE:HB2	1.92	0.52
1:BV:45:TYR:HB2	1:BV:48:ARG:HH21	1.75	0.52
1:BZ:22:ILE:HD11	1:BZ:39:VAL:HG11	1.90	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CT:22:ILE:HD12	1:CT:49:ARG:HH21	1.74	0.52
1:DC:2:ILE:HD11	1:DC:124:PHE:HB2	1.92	0.52
1:DJ:68:GLY:O	1:DJ:95:GLU:HA	2.10	0.52
1:AB:49:ARG:HG2	1:AB:76:ILE:HG12	1.91	0.52
1:AI:45:TYR:HA	1:AI:48:ARG:HE	1.75	0.52
1:AM:39:VAL:HG23	1:AM:42:ASP:H	1.74	0.52
1:AQ:49:ARG:HG2	1:AQ:76:ILE:HG12	1.91	0.52
1:BA:22:ILE:HA	1:BA:41:GLU:OE1	2.10	0.52
1:BJ:22:ILE:HD12	1:BJ:49:ARG:HH21	1.74	0.52
1:BW:68:GLY:O	1:BW:95:GLU:HA	2.10	0.52
1:BX:29:GLN:HG3	1:CL:57:VAL:HG21	1.90	0.52
1:BZ:68:GLY:O	1:BZ:95:GLU:HA	2.10	0.52
1:CB:45:TYR:HA	1:CB:48:ARG:HE	1.75	0.52
1:CE:45:TYR:HA	1:CE:48:ARG:HE	1.75	0.52
1:DD:39:VAL:HG23	1:DD:42:ASP:H	1.74	0.52
1:DF:45:TYR:HB2	1:DF:48:ARG:HH21	1.75	0.52
1:AG:60:VAL:HG11	1:CS:87:PHE:HE2	1.75	0.52
1:AI:22:ILE:HA	1:AI:41:GLU:OE1	2.10	0.52
1:AJ:68:GLY:O	1:AJ:95:GLU:HA	2.10	0.52
1:BF:49:ARG:HG2	1:BF:76:ILE:HG12	1.91	0.52
1:BG:45:TYR:HB2	1:BG:48:ARG:HH21	1.75	0.52
1:BI:49:ARG:HG2	1:BI:76:ILE:HG12	1.91	0.52
1:BK:39:VAL:HG23	1:BK:42:ASP:H	1.74	0.52
1:BN:68:GLY:O	1:BN:95:GLU:HA	2.10	0.52
1:BP:22:ILE:HA	1:BP:41:GLU:OE1	2.10	0.52
1:BY:45:TYR:HB2	1:BY:48:ARG:HH21	1.75	0.52
1:CB:22:ILE:HD12	1:CB:49:ARG:HH21	1.74	0.52
1:CK:2:ILE:HD11	1:CK:124:PHE:HB2	1.92	0.52
1:CM:49:ARG:HG2	1:CM:76:ILE:HG12	1.91	0.52
1:DB:49:ARG:HG2	1:DB:76:ILE:HG12	1.92	0.52
1:AC:22:ILE:HA	1:AC:41:GLU:OE1	2.10	0.51
1:AO:45:TYR:HB2	1:AO:48:ARG:HH21	1.75	0.51
1:AP:39:VAL:HG23	1:AP:42:ASP:H	1.74	0.51
1:AP:68:GLY:O	1:AP:95:GLU:HA	2.10	0.51
1:AT:48:ARG:HH12	1:AW:127:LEU:HD21	1.74	0.51
1:AU:22:ILE:HD12	1:AU:49:ARG:HH21	1.74	0.51
1:AX:45:TYR:HB2	1:AX:48:ARG:HH21	1.75	0.51
1:BB:68:GLY:O	1:BB:95:GLU:HA	2.10	0.51
1:BM:45:TYR:HB2	1:BM:48:ARG:HH21	1.75	0.51
1:BP:2:ILE:HD11	1:BP:124:PHE:HB2	1.92	0.51
1:BV:22:ILE:HD12	1:BV:49:ARG:HH21	1.74	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CB:22:ILE:HA	1:CB:41:GLU:OE1	2.10	0.51
1:CB:45:TYR:HB2	1:CB:48:ARG:HH21	1.75	0.51
1:CC:39:VAL:HG23	1:CC:42:ASP:H	1.74	0.51
1:CE:22:ILE:HA	1:CE:41:GLU:OE1	2.10	0.51
1:CT:22:ILE:HA	1:CT:41:GLU:OE1	2.10	0.51
1:CU:68:GLY:O	1:CU:95:GLU:HA	2.10	0.51
1:CZ:22:ILE:HA	1:CZ:41:GLU:OE1	2.10	0.51
1:DD:68:GLY:O	1:DD:95:GLU:HA	2.10	0.51
1:AD:68:GLY:O	1:AD:95:GLU:HA	2.10	0.51
1:AF:45:TYR:HB2	1:AF:48:ARG:HH21	1.75	0.51
1:AI:22:ILE:HD12	1:AI:49:ARG:HH21	1.74	0.51
1:AL:45:TYR:HA	1:AL:48:ARG:HE	1.75	0.51
1:AR:22:ILE:HA	1:AR:41:GLU:OE1	2.10	0.51
1:AU:2:ILE:HD11	1:AU:124:PHE:HB2	1.91	0.51
1:AV:68:GLY:O	1:AV:95:GLU:HA	2.10	0.51
1:BA:45:TYR:HB2	1:BA:48:ARG:HH21	1.75	0.51
1:BD:45:TYR:HB2	1:BD:48:ARG:HH21	1.75	0.51
1:CD:49:ARG:HG2	1:CD:76:ILE:HG12	1.91	0.51
1:CE:45:TYR:HB2	1:CE:48:ARG:HH21	1.75	0.51
1:CN:45:TYR:HB2	1:CN:48:ARG:HH21	1.75	0.51
1:CR:68:GLY:O	1:CR:95:GLU:HA	2.10	0.51
1:CU:39:VAL:HG23	1:CU:42:ASP:H	1.74	0.51
1:CZ:22:ILE:HD12	1:CZ:49:ARG:HH21	1.74	0.51
1:DA:39:VAL:HG23	1:DA:42:ASP:H	1.74	0.51
1:AD:65:TYR:CE2	1:DE:46:ARG:HG3	2.45	0.51
1:AL:22:ILE:HA	1:AL:41:GLU:OE1	2.10	0.51
1:AS:68:GLY:O	1:AS:95:GLU:HA	2.10	0.51
1:AU:45:TYR:HA	1:AU:48:ARG:HE	1.75	0.51
1:AY:68:GLY:O	1:AY:95:GLU:HA	2.10	0.51
1:BA:45:TYR:HA	1:BA:48:ARG:HE	1.75	0.51
1:BD:16:ILE:HD13	1:CZ:114:GLN:NE2	2.22	0.51
1:BG:45:TYR:HA	1:BG:48:ARG:HE	1.75	0.51
1:BJ:45:TYR:HB2	1:BJ:48:ARG:HH21	1.75	0.51
1:BS:45:TYR:HA	1:BS:48:ARG:HE	1.75	0.51
1:BT:68:GLY:O	1:BT:95:GLU:HA	2.10	0.51
1:BY:22:ILE:HA	1:BY:41:GLU:OE1	2.10	0.51
1:BZ:39:VAL:HG23	1:BZ:42:ASP:H	1.74	0.51
1:CB:2:ILE:HD11	1:CB:124:PHE:HB2	1.92	0.51
1:CH:45:TYR:HA	1:CH:48:ARG:HE	1.75	0.51
1:CQ:22:ILE:HA	1:CQ:41:GLU:OE1	2.10	0.51
1:CQ:45:TYR:HA	1:CQ:48:ARG:HE	1.75	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CT:45:TYR:HB2	1:CT:48:ARG:HH21	1.75	0.51
1:CV:49:ARG:HG2	1:CV:76:ILE:HG12	1.92	0.51
1:CY:49:ARG:HG2	1:CY:76:ILE:HG12	1.91	0.51
1:BD:22:ILE:HA	1:BD:41:GLU:OE1	2.10	0.51
1:BH:68:GLY:O	1:BH:95:GLU:HA	2.10	0.51
1:BM:2:ILE:HD11	1:BM:124:PHE:HB2	1.92	0.51
1:BM:22:ILE:HA	1:BM:41:GLU:OE1	2.10	0.51
1:BM:45:TYR:HA	1:BM:48:ARG:HE	1.75	0.51
1:BR:49:ARG:HG2	1:BR:76:ILE:HG12	1.91	0.51
1:BS:45:TYR:HB2	1:BS:48:ARG:HH21	1.75	0.51
1:BY:2:ILE:HD11	1:BY:124:PHE:HB2	1.92	0.51
1:DL:45:TYR:HB2	1:DL:48:ARG:HH21	1.75	0.51
1:AI:45:TYR:HB2	1:AI:48:ARG:HH21	1.75	0.51
1:AM:60:VAL:HG11	1:CP:87:PHE:HE2	1.74	0.51
1:AO:22:ILE:HA	1:AO:41:GLU:OE1	2.10	0.51
1:BF:65:TYR:HE2	1:BS:46:ARG:CB	2.23	0.51
1:CN:22:ILE:HA	1:CN:41:GLU:OE1	2.10	0.51
1:CN:45:TYR:HA	1:CN:48:ARG:HE	1.75	0.51
1:CP:49:ARG:HG2	1:CP:76:ILE:HG12	1.92	0.51
1:AA:7:ILE:HD12	1:AA:111:ILE:HG13	1.93	0.51
1:AC:22:ILE:HD12	1:AC:49:ARG:HH21	1.74	0.51
1:AF:2:ILE:HD11	1:AF:124:PHE:HB2	1.92	0.51
1:AO:130:LEU:HD21	1:CE:55:SER:HB3	1.92	0.51
1:AR:2:ILE:HD11	1:AR:124:PHE:HB2	1.92	0.51
1:BK:68:GLY:O	1:BK:95:GLU:HA	2.10	0.51
1:BV:22:ILE:HA	1:BV:41:GLU:OE1	2.10	0.51
1:CC:9:THR:HG22	1:CC:111:ILE:HD12	1.93	0.51
1:CN:2:ILE:HD11	1:CN:124:PHE:HB2	1.92	0.51
1:DF:22:ILE:HA	1:DF:41:GLU:OE1	2.10	0.51
1:DI:45:TYR:HA	1:DI:48:ARG:HE	1.75	0.51
1:AD:9:THR:HG22	1:AD:111:ILE:HD12	1.93	0.51
1:AF:67:LYS:NZ	1:BV:87:PHE:O	2.27	0.51
1:AO:45:TYR:HA	1:AO:48:ARG:HE	1.75	0.51
1:BH:7:ILE:HD12	1:BH:111:ILE:HG13	1.93	0.51
1:BN:7:ILE:HD12	1:BN:111:ILE:HG13	1.93	0.51
1:BQ:68:GLY:O	1:BQ:95:GLU:HA	2.10	0.51
1:BU:46:ARG:NH2	1:DG:99:ALA:HA	2.25	0.51
1:BU:49:ARG:HG2	1:BU:76:ILE:HG12	1.91	0.51
1:CF:9:THR:HG22	1:CF:111:ILE:HD12	1.93	0.51
1:CL:68:GLY:O	1:CL:95:GLU:HA	2.10	0.51
1:CO:68:GLY:O	1:CO:95:GLU:HA	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CR:9:THR:HG22	1:CR:111:ILE:HD12	1.93	0.51
1:CU:9:THR:HG22	1:CU:111:ILE:HD12	1.93	0.51
1:CX:7:ILE:HD12	1:CX:111:ILE:HG13	1.93	0.51
1:DC:45:TYR:HB2	1:DC:48:ARG:HH21	1.75	0.51
1:DG:7:ILE:HD12	1:DG:111:ILE:HG13	1.93	0.51
1:DL:2:ILE:HD11	1:DL:124:PHE:HB2	1.92	0.51
1:DL:22:ILE:HA	1:DL:41:GLU:OE1	2.10	0.51
1:DL:45:TYR:HA	1:DL:48:ARG:HE	1.75	0.51
1:AG:68:GLY:O	1:AG:95:GLU:HA	2.10	0.51
1:AH:49:ARG:HG2	1:AH:76:ILE:HG12	1.91	0.51
1:AW:49:ARG:HG2	1:AW:76:ILE:HG12	1.91	0.51
1:AY:7:ILE:HD12	1:AY:111:ILE:HG13	1.93	0.51
1:AY:39:VAL:HG23	1:AY:42:ASP:H	1.74	0.51
1:BC:65:TYR:CE2	1:BP:46:ARG:HA	2.45	0.51
1:BE:9:THR:HG22	1:BE:111:ILE:HD12	1.93	0.51
1:BG:22:ILE:HA	1:BG:41:GLU:OE1	2.10	0.51
1:BJ:22:ILE:HA	1:BJ:41:GLU:OE1	2.10	0.51
1:BV:2:ILE:HD11	1:BV:124:PHE:HB2	1.91	0.51
1:BV:10:GLY:N	1:BV:107:ASP:OD1	2.34	0.51
1:DA:68:GLY:O	1:DA:95:GLU:HA	2.10	0.51
1:DC:22:ILE:HA	1:DC:41:GLU:OE1	2.10	0.51
1:DG:68:GLY:O	1:DG:95:GLU:HA	2.10	0.51
1:DI:22:ILE:HA	1:DI:41:GLU:OE1	2.10	0.51
1:DK:49:ARG:HG2	1:DK:76:ILE:HG12	1.91	0.51
1:AA:68:GLY:O	1:AA:95:GLU:HA	2.10	0.51
1:AI:17:THR:O	1:BY:12:THR:OG1	2.25	0.51
1:AR:110:LEU:HD23	1:CK:16:ILE:HD11	1.92	0.51
1:BB:9:THR:HG22	1:BB:111:ILE:HD12	1.93	0.51
1:BE:68:GLY:O	1:BE:95:GLU:HA	2.10	0.51
1:CA:49:ARG:HG2	1:CA:76:ILE:HG12	1.91	0.51
1:CU:7:ILE:HD12	1:CU:111:ILE:HG13	1.93	0.51
1:CW:45:TYR:HA	1:CW:48:ARG:HE	1.75	0.51
1:CZ:2:ILE:HD11	1:CZ:124:PHE:HB2	1.92	0.51
1:AE:48:ARG:HH12	1:BF:127:LEU:HD21	1.76	0.51
1:AM:7:ILE:HD12	1:AM:111:ILE:HG13	1.93	0.51
1:AM:68:GLY:O	1:AM:95:GLU:HA	2.10	0.51
1:AS:7:ILE:HD12	1:AS:111:ILE:HG13	1.93	0.51
1:AV:9:THR:HG22	1:AV:111:ILE:HD12	1.93	0.51
1:AX:22:ILE:HA	1:AX:41:GLU:OE1	2.10	0.51
1:BJ:26:LEU:HB3	1:BK:130:LEU:O	2.11	0.51
1:BK:7:ILE:HD12	1:BK:111:ILE:HG13	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:65:TYR:CE2	1:CH:46:ARG:HG2	2.46	0.51
1:CI:68:GLY:O	1:CI:95:GLU:HA	2.10	0.51
1:CM:87:PHE:HE2	1:CX:60:VAL:HG11	1.75	0.51
1:CW:45:TYR:HB2	1:CW:48:ARG:HH21	1.75	0.51
1:AF:22:ILE:HA	1:AF:41:GLU:OE1	2.10	0.50
1:AK:49:ARG:HG2	1:AK:76:ILE:HG12	1.91	0.50
1:BA:22:ILE:O	1:CW:6:SER:HA	2.11	0.50
1:BN:9:THR:HG22	1:BN:111:ILE:HD12	1.93	0.50
1:BT:7:ILE:HD12	1:BT:111:ILE:HG13	1.93	0.50
1:BW:7:ILE:HD12	1:BW:111:ILE:HG13	1.93	0.50
1:BZ:7:ILE:HD12	1:BZ:111:ILE:HG13	1.93	0.50
1:BZ:9:THR:HG22	1:BZ:111:ILE:HD12	1.93	0.50
1:CC:7:ILE:HD12	1:CC:111:ILE:HG13	1.93	0.50
1:CH:22:ILE:HA	1:CH:41:GLU:OE1	2.10	0.50
1:CQ:45:TYR:HB2	1:CQ:48:ARG:HH21	1.75	0.50
1:CT:45:TYR:HA	1:CT:48:ARG:HE	1.75	0.50
1:CZ:45:TYR:HB2	1:CZ:48:ARG:HH21	1.75	0.50
1:DD:7:ILE:HD12	1:DD:111:ILE:HG13	1.93	0.50
1:DG:9:THR:HG22	1:DG:111:ILE:HD12	1.93	0.50
1:AG:7:ILE:HD12	1:AG:111:ILE:HG13	1.93	0.50
1:AK:46:ARG:NH2	1:BW:99:ALA:HA	2.26	0.50
1:AY:9:THR:HG22	1:AY:111:ILE:HD12	1.93	0.50
1:BH:9:THR:HG22	1:BH:111:ILE:HD12	1.93	0.50
1:BP:22:ILE:HD12	1:BP:49:ARG:HH21	1.74	0.50
1:BS:22:ILE:HA	1:BS:41:GLU:OE1	2.10	0.50
1:CO:9:THR:HG22	1:CO:111:ILE:HD12	1.93	0.50
1:DJ:7:ILE:HD12	1:DJ:111:ILE:HG13	1.93	0.50
1:AA:9:THR:HG22	1:AA:111:ILE:HD12	1.93	0.50
1:AC:45:TYR:HB2	1:AC:48:ARG:HH21	1.75	0.50
1:AH:48:ARG:HH12	1:CD:127:LEU:HD21	1.76	0.50
1:AI:128:GLY:O	1:BW:28:GLY:HA3	2.11	0.50
1:AM:22:ILE:HA	1:AM:41:GLU:OE1	2.12	0.50
1:AR:45:TYR:HB2	1:AR:48:ARG:HH21	1.75	0.50
1:AU:22:ILE:HA	1:AU:41:GLU:OE1	2.10	0.50
1:AX:16:ILE:HG21	1:CT:114:GLN:NE2	2.26	0.50
1:BE:7:ILE:HD12	1:BE:111:ILE:HG13	1.93	0.50
1:BQ:7:ILE:HD12	1:BQ:111:ILE:HG13	1.93	0.50
1:BW:9:THR:HG22	1:BW:111:ILE:HD12	1.93	0.50
1:BY:45:TYR:HA	1:BY:48:ARG:HE	1.75	0.50
1:CF:68:GLY:O	1:CF:95:GLU:HA	2.10	0.50
1:CL:7:ILE:HD12	1:CL:111:ILE:HG13	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CX:68:GLY:O	1:CX:95:GLU:HA	2.10	0.50
1:DA:7:ILE:HD12	1:DA:111:ILE:HG13	1.93	0.50
1:DA:9:THR:HG22	1:DA:111:ILE:HD12	1.93	0.50
1:AB:46:ARG:HG3	1:AS:65:TYR:CD2	2.46	0.50
1:AY:22:ILE:HA	1:AY:41:GLU:OE1	2.12	0.50
1:BB:7:ILE:HD12	1:BB:111:ILE:HG13	1.93	0.50
1:BD:2:ILE:HD11	1:BD:124:PHE:HB2	1.92	0.50
1:BG:22:ILE:HD12	1:BG:49:ARG:HH21	1.74	0.50
1:CC:68:GLY:O	1:CC:95:GLU:HA	2.10	0.50
1:CK:22:ILE:HA	1:CK:41:GLU:OE1	2.10	0.50
1:CL:22:ILE:HA	1:CL:41:GLU:OE1	2.12	0.50
1:CR:22:ILE:HA	1:CR:41:GLU:OE1	2.12	0.50
1:AC:125:TRP:O	1:BP:70:ASN:ND2	2.44	0.50
1:BH:22:ILE:HA	1:BH:41:GLU:OE1	2.12	0.50
1:BK:9:THR:HG22	1:BK:111:ILE:HD12	1.93	0.50
1:BP:45:TYR:HB2	1:BP:48:ARG:HH21	1.75	0.50
1:CK:45:TYR:HB2	1:CK:48:ARG:HH21	1.75	0.50
1:CX:22:ILE:HA	1:CX:41:GLU:OE1	2.12	0.50
1:AB:48:ARG:HH12	1:AT:127:LEU:HD21	1.76	0.50
1:AD:7:ILE:HD12	1:AD:111:ILE:HG13	1.93	0.50
1:AE:87:PHE:CZ	1:BE:60:VAL:HG11	2.46	0.50
1:AJ:22:ILE:HA	1:AJ:41:GLU:OE1	2.12	0.50
1:AN:33:ASN:OD1	1:CB:30:THR:OG1	2.30	0.50
1:BJ:40:SER:HB2	1:BK:127:LEU:HD13	1.93	0.50
1:BN:22:ILE:HA	1:BN:41:GLU:OE1	2.12	0.50
1:CL:9:THR:HG22	1:CL:111:ILE:HD12	1.93	0.50
1:CW:22:ILE:HA	1:CW:41:GLU:OE1	2.10	0.50
1:DD:22:ILE:HA	1:DD:41:GLU:OE1	2.12	0.50
1:DI:2:ILE:HD11	1:DI:124:PHE:HB2	1.92	0.50
1:AV:22:ILE:HA	1:AV:41:GLU:OE1	2.12	0.50
1:BE:22:ILE:HA	1:BE:41:GLU:OE1	2.12	0.50
1:BT:9:THR:HG22	1:BT:111:ILE:HD12	1.93	0.50
1:CF:22:ILE:HA	1:CF:41:GLU:OE1	2.12	0.50
1:DC:45:TYR:HA	1:DC:48:ARG:HE	1.75	0.50
1:DI:45:TYR:HB2	1:DI:48:ARG:HH21	1.75	0.50
1:BF:48:ARG:HH12	1:BR:127:LEU:HD21	1.77	0.50
1:BX:98:PRO:HB2	1:CN:46:ARG:HG3	1.94	0.50
1:CC:22:ILE:HA	1:CC:41:GLU:OE1	2.12	0.50
1:CF:7:ILE:HD12	1:CF:111:ILE:HG13	1.93	0.50
1:CU:22:ILE:HA	1:CU:41:GLU:OE1	2.12	0.50
1:AG:9:THR:HG22	1:AG:111:ILE:HD12	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AM:9:THR:HG22	1:AM:111:ILE:HD12	1.93	0.50
1:BH:65:TYR:CE2	1:BL:46:ARG:HG3	2.47	0.50
1:BK:22:ILE:HA	1:BK:41:GLU:OE1	2.12	0.50
1:BW:22:ILE:HA	1:BW:41:GLU:OE1	2.12	0.50
1:CI:22:ILE:HA	1:CI:41:GLU:OE1	2.12	0.50
1:DG:22:ILE:HA	1:DG:41:GLU:OE1	2.12	0.50
1:DJ:22:ILE:HA	1:DJ:41:GLU:OE1	2.12	0.50
1:AI:46:ARG:HG2	1:CS:65:TYR:CD2	2.47	0.49
1:AW:65:TYR:HD2	1:BV:46:ARG:HG2	1.74	0.49
1:CI:7:ILE:HD12	1:CI:111:ILE:HG13	1.93	0.49
1:AP:9:THR:HG22	1:AP:111:ILE:HD12	1.93	0.49
1:CR:7:ILE:HD12	1:CR:111:ILE:HG13	1.93	0.49
1:DJ:9:THR:HG22	1:DJ:111:ILE:HD12	1.93	0.49
1:AO:55:SER:HB3	1:CE:130:LEU:HD21	1.94	0.49
1:AP:7:ILE:HD12	1:AP:111:ILE:HG13	1.93	0.49
1:AS:22:ILE:HA	1:AS:41:GLU:OE1	2.12	0.49
1:AV:7:ILE:HD12	1:AV:111:ILE:HG13	1.93	0.49
1:BQ:9:THR:HG22	1:BQ:111:ILE:HD12	1.93	0.49
1:AD:60:VAL:HG11	1:DE:87:PHE:CE2	2.48	0.49
1:AJ:9:THR:HG22	1:AJ:111:ILE:HD12	1.93	0.49
1:AO:108:LEU:HA	1:AO:111:ILE:HG22	1.95	0.49
1:BG:19:GLY:O	1:DC:8:LYS:HD2	2.13	0.49
1:BQ:22:ILE:HA	1:BQ:41:GLU:OE1	2.12	0.49
1:CI:9:THR:HG22	1:CI:111:ILE:HD12	1.93	0.49
1:CO:7:ILE:HD12	1:CO:111:ILE:HG13	1.93	0.49
1:CX:9:THR:HG22	1:CX:111:ILE:HD12	1.93	0.49
1:AD:22:ILE:HA	1:AD:41:GLU:OE1	2.12	0.49
1:AO:17:THR:O	1:CE:12:THR:OG1	2.23	0.49
1:AO:18:GLY:CA	1:CE:11:ALA:HA	2.42	0.49
1:BP:108:LEU:HA	1:BP:111:ILE:HG22	1.95	0.49
1:DA:22:ILE:HA	1:DA:41:GLU:OE1	2.12	0.49
1:AA:22:ILE:HA	1:AA:41:GLU:OE1	2.12	0.49
1:AR:108:LEU:HA	1:AR:111:ILE:HG22	1.95	0.49
1:BG:14:ALA:HB2	1:DC:16:ILE:HD12	1.95	0.49
1:BT:22:ILE:HA	1:BT:41:GLU:OE1	2.12	0.49
1:BY:9:THR:HG22	1:BY:111:ILE:HD12	1.95	0.49
1:BZ:22:ILE:HA	1:BZ:41:GLU:OE1	2.12	0.49
1:CO:22:ILE:HA	1:CO:41:GLU:OE1	2.12	0.49
1:DD:9:THR:HG22	1:DD:111:ILE:HD12	1.93	0.49
1:AB:87:PHE:HE2	1:AS:60:VAL:HG11	1.78	0.49
1:AC:108:LEU:HA	1:AC:111:ILE:HG22	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AG:22:ILE:HA	1:AG:41:GLU:OE1	2.12	0.49
1:AJ:7:ILE:HD12	1:AJ:111:ILE:HG13	1.93	0.49
1:BG:108:LEU:HA	1:BG:111:ILE:HG22	1.95	0.49
1:BM:108:LEU:HA	1:BM:111:ILE:HG22	1.95	0.49
1:BS:9:THR:HG22	1:BS:111:ILE:HD12	1.95	0.49
1:BS:108:LEU:HA	1:BS:111:ILE:HG22	1.95	0.49
1:CK:108:LEU:HA	1:CK:111:ILE:HG22	1.95	0.49
1:CQ:108:LEU:HA	1:CQ:111:ILE:HG22	1.95	0.49
1:AW:65:TYR:HE2	1:BV:46:ARG:HA	1.77	0.49
1:CN:108:LEU:HA	1:CN:111:ILE:HG22	1.95	0.49
1:CR:99:ALA:HA	1:DB:46:ARG:NH2	2.28	0.49
1:CZ:108:LEU:HA	1:CZ:111:ILE:HG22	1.95	0.49
1:DF:108:LEU:HA	1:DF:111:ILE:HG22	1.95	0.49
1:BJ:125:TRP:O	1:DF:70:ASN:ND2	2.46	0.49
1:BM:9:THR:HG22	1:BM:111:ILE:HD12	1.95	0.49
1:CI:130:LEU:O	1:DC:26:LEU:HB3	2.13	0.49
1:CK:9:THR:HG22	1:CK:111:ILE:HD12	1.95	0.49
1:AL:108:LEU:HA	1:AL:111:ILE:HG22	1.95	0.49
1:CU:57:VAL:HG21	1:CY:29:GLN:HG3	1.94	0.49
1:AC:22:ILE:O	1:BP:6:SER:HA	2.12	0.48
1:BD:108:LEU:HA	1:BD:111:ILE:HG22	1.95	0.48
1:CA:87:PHE:CZ	1:CF:60:VAL:HG11	2.47	0.48
1:CQ:9:THR:HG22	1:CQ:111:ILE:HD12	1.95	0.48
1:CT:9:THR:HG22	1:CT:111:ILE:HD12	1.95	0.48
1:DC:9:THR:HG22	1:DC:111:ILE:HD12	1.95	0.48
1:AD:130:LEU:O	1:BG:26:LEU:HB3	2.13	0.48
1:AW:87:PHE:CE2	1:BT:60:VAL:HG11	2.48	0.48
1:BB:22:ILE:HA	1:BB:41:GLU:OE1	2.12	0.48
1:BG:22:ILE:O	1:DC:6:SER:HA	2.13	0.48
1:BM:68:GLY:O	1:BM:95:GLU:HA	2.14	0.48
1:BX:79:SER:HA	1:BX:85:THR:HG22	1.96	0.48
1:DL:68:GLY:O	1:DL:95:GLU:HA	2.14	0.48
1:AL:9:THR:HG22	1:AL:111:ILE:HD12	1.95	0.48
1:AO:68:GLY:O	1:AO:95:GLU:HA	2.14	0.48
1:AR:9:THR:HG22	1:AR:111:ILE:HD12	1.95	0.48
1:AS:9:THR:HG22	1:AS:111:ILE:HD12	1.93	0.48
1:AU:9:THR:HG22	1:AU:111:ILE:HD12	1.95	0.48
1:AW:79:SER:HA	1:AW:85:THR:HG22	1.96	0.48
1:BA:108:LEU:HA	1:BA:111:ILE:HG22	1.95	0.48
1:CD:46:ARG:HG3	1:CI:65:TYR:CE2	2.48	0.48
1:CD:79:SER:HA	1:CD:85:THR:HG22	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CH:68:GLY:O	1:CH:95:GLU:HA	2.14	0.48
1:DI:9:THR:HG22	1:DI:111:ILE:HD12	1.95	0.48
1:DI:108:LEU:HA	1:DI:111:ILE:HG22	1.95	0.48
1:AE:79:SER:HA	1:AE:85:THR:HG22	1.96	0.48
1:AF:9:THR:HG22	1:AF:111:ILE:HD12	1.95	0.48
1:AO:18:GLY:HA2	1:CE:11:ALA:HA	1.96	0.48
1:AX:9:THR:HG22	1:AX:111:ILE:HD12	1.95	0.48
1:BG:68:GLY:O	1:BG:95:GLU:HA	2.14	0.48
1:BJ:68:GLY:O	1:BJ:95:GLU:HA	2.14	0.48
1:CA:65:TYR:CE2	1:CH:46:ARG:HA	2.48	0.48
1:CG:79:SER:HA	1:CG:85:THR:HG22	1.96	0.48
1:CN:68:GLY:O	1:CN:95:GLU:HA	2.14	0.48
1:CT:68:GLY:O	1:CT:95:GLU:HA	2.14	0.48
1:CZ:68:GLY:O	1:CZ:95:GLU:HA	2.14	0.48
1:DC:68:GLY:O	1:DC:95:GLU:HA	2.14	0.48
1:DH:79:SER:HA	1:DH:85:THR:HG22	1.96	0.48
1:AB:79:SER:HA	1:AB:85:THR:HG22	1.96	0.48
1:AI:111:ILE:HD13	1:BY:74:PHE:CZ	2.48	0.48
1:AK:79:SER:HA	1:AK:85:THR:HG22	1.96	0.48
1:AU:22:ILE:HG22	1:CN:7:ILE:O	2.14	0.48
1:BJ:46:ARG:HG2	1:BL:65:TYR:CE2	2.48	0.48
1:BT:130:LEU:O	1:DI:26:LEU:HB3	2.14	0.48
1:CA:79:SER:HA	1:CA:85:THR:HG22	1.96	0.48
1:DK:79:SER:HA	1:DK:85:THR:HG22	1.96	0.48
1:AA:7:ILE:O	1:AB:22:ILE:HG22	2.14	0.48
1:AI:108:LEU:HA	1:AI:111:ILE:HG22	1.95	0.48
1:AT:79:SER:HA	1:AT:85:THR:HG22	1.96	0.48
1:BA:9:THR:HG22	1:BA:111:ILE:HD12	1.95	0.48
1:BB:7:ILE:O	1:BC:22:ILE:HG22	2.14	0.48
1:BE:7:ILE:O	1:BF:22:ILE:HG22	2.14	0.48
1:BK:7:ILE:O	1:BL:22:ILE:HG22	2.14	0.48
1:BO:79:SER:HA	1:BO:85:THR:HG22	1.96	0.48
1:BP:68:GLY:O	1:BP:95:GLU:HA	2.14	0.48
1:BS:68:GLY:O	1:BS:95:GLU:HA	2.14	0.48
1:BV:68:GLY:O	1:BV:95:GLU:HA	2.14	0.48
1:BY:68:GLY:O	1:BY:95:GLU:HA	2.14	0.48
1:BZ:7:ILE:O	1:CA:22:ILE:HG22	2.14	0.48
1:CE:108:LEU:HA	1:CE:111:ILE:HG22	1.95	0.48
1:DF:9:THR:HG22	1:DF:111:ILE:HD12	1.95	0.48
1:AC:68:GLY:O	1:AC:95:GLU:HA	2.14	0.48
1:AD:7:ILE:O	1:AE:22:ILE:HG22	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AL:68:GLY:O	1:AL:95:GLU:HA	2.14	0.48
1:BD:68:GLY:O	1:BD:95:GLU:HA	2.14	0.48
1:BJ:122:ASP:OD1	1:DF:109:ARG:NH1	2.44	0.48
1:BX:87:PHE:CE2	1:CL:60:VAL:HG11	2.49	0.48
1:CE:68:GLY:O	1:CE:95:GLU:HA	2.14	0.48
1:CH:9:THR:HG22	1:CH:111:ILE:HD12	1.95	0.48
1:CH:108:LEU:HA	1:CH:111:ILE:HG22	1.95	0.48
1:CM:87:PHE:CE2	1:CX:60:VAL:HG11	2.48	0.48
1:DD:7:ILE:O	1:DE:22:ILE:HG22	2.14	0.48
1:DF:68:GLY:O	1:DF:95:GLU:HA	2.14	0.48
1:AP:22:ILE:HA	1:AP:41:GLU:OE1	2.12	0.48
1:AT:29:GLN:HG3	1:AV:57:VAL:HG21	1.96	0.48
1:AU:87:PHE:O	1:CN:67:LYS:NZ	2.29	0.48
1:AX:108:LEU:HA	1:AX:111:ILE:HG22	1.95	0.48
1:BC:65:TYR:CE2	1:BP:46:ARG:HG2	2.48	0.48
1:BC:79:SER:HA	1:BC:85:THR:HG22	1.95	0.48
1:BR:79:SER:HA	1:BR:85:THR:HG22	1.96	0.48
1:CK:68:GLY:O	1:CK:95:GLU:HA	2.14	0.48
1:CN:9:THR:HG22	1:CN:111:ILE:HD12	1.95	0.48
1:CU:7:ILE:O	1:CV:22:ILE:HG22	2.14	0.48
1:DJ:7:ILE:O	1:DK:22:ILE:HG22	2.14	0.48
1:DL:9:THR:HG22	1:DL:111:ILE:HD12	1.95	0.48
1:AQ:79:SER:HA	1:AQ:85:THR:HG22	1.96	0.48
1:BD:9:THR:HG22	1:BD:111:ILE:HD12	1.95	0.48
1:CD:58:PRO:HG2	1:CK:45:TYR:CD2	2.49	0.48
1:CE:9:THR:HG22	1:CE:111:ILE:HD12	1.95	0.48
1:CJ:79:SER:HA	1:CJ:85:THR:HG22	1.96	0.48
1:DC:108:LEU:HA	1:DC:111:ILE:HG22	1.95	0.48
1:AI:68:GLY:O	1:AI:95:GLU:HA	2.14	0.48
1:AM:57:VAL:HG21	1:CP:29:GLN:HG3	1.95	0.48
1:AU:68:GLY:O	1:AU:95:GLU:HA	2.14	0.48
1:AX:8:LYS:HD2	1:CT:19:GLY:O	2.13	0.48
1:BV:108:LEU:HA	1:BV:111:ILE:HG22	1.95	0.48
1:CC:7:ILE:O	1:CD:22:ILE:HG22	2.14	0.48
1:CW:9:THR:HG22	1:CW:111:ILE:HD12	1.95	0.48
1:CX:7:ILE:O	1:CY:22:ILE:HG22	2.14	0.48
1:AK:29:GLN:HB3	1:AK:36:ASN:HD22	1.80	0.47
1:AP:7:ILE:O	1:AQ:22:ILE:HG22	2.14	0.47
1:AP:22:ILE:HG22	1:AQ:7:ILE:O	2.14	0.47
1:BA:68:GLY:O	1:BA:95:GLU:HA	2.14	0.47
1:BG:9:THR:HG22	1:BG:111:ILE:HD12	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BK:22:ILE:HG22	1:BL:7:ILE:O	2.14	0.47
1:BR:29:GLN:HB3	1:BR:36:ASN:HD22	1.79	0.47
1:BW:7:ILE:O	1:BX:22:ILE:HG22	2.14	0.47
1:CW:108:LEU:HA	1:CW:111:ILE:HG22	1.95	0.47
1:CY:29:GLN:HB3	1:CY:36:ASN:HD22	1.79	0.47
1:DG:7:ILE:O	1:DH:22:ILE:HG22	2.14	0.47
1:DI:68:GLY:O	1:DI:95:GLU:HA	2.14	0.47
1:AE:29:GLN:HB3	1:AE:36:ASN:HD22	1.80	0.47
1:AJ:7:ILE:O	1:AK:22:ILE:HG22	2.14	0.47
1:AN:29:GLN:HB3	1:AN:36:ASN:HD22	1.79	0.47
1:AU:108:LEU:HA	1:AU:111:ILE:HG22	1.95	0.47
1:AV:7:ILE:O	1:AW:22:ILE:HG22	2.14	0.47
1:BH:22:ILE:HG22	1:BI:7:ILE:O	2.14	0.47
1:BL:79:SER:HA	1:BL:85:THR:HG22	1.95	0.47
1:BU:79:SER:HA	1:BU:85:THR:HG22	1.96	0.47
1:CB:9:THR:HG22	1:CB:111:ILE:HD12	1.95	0.47
1:CF:22:ILE:HG22	1:CG:7:ILE:O	2.14	0.47
1:CM:29:GLN:HB3	1:CM:36:ASN:HD22	1.80	0.47
1:CM:79:SER:HA	1:CM:85:THR:HG22	1.96	0.47
1:CP:29:GLN:HB3	1:CP:36:ASN:HD22	1.80	0.47
1:CP:79:SER:HA	1:CP:85:THR:HG22	1.95	0.47
1:CT:108:LEU:HA	1:CT:111:ILE:HG22	1.95	0.47
1:CW:68:GLY:O	1:CW:95:GLU:HA	2.14	0.47
1:AA:22:ILE:HG22	1:AB:7:ILE:O	2.14	0.47
1:AD:22:ILE:HG22	1:AE:7:ILE:O	2.14	0.47
1:AE:46:ARG:NH2	1:BE:99:ALA:HA	2.28	0.47
1:AO:9:THR:HG22	1:AO:111:ILE:HD12	1.95	0.47
1:AS:130:LEU:O	1:AX:26:LEU:HB3	2.15	0.47
1:AT:33:ASN:OD1	1:AX:30:THR:OG1	2.32	0.47
1:BB:22:ILE:HG22	1:BC:7:ILE:O	2.14	0.47
1:BC:29:GLN:HB3	1:BC:36:ASN:HD22	1.80	0.47
1:BF:79:SER:HA	1:BF:85:THR:HG22	1.96	0.47
1:BL:29:GLN:HB3	1:BL:36:ASN:HD22	1.80	0.47
1:BQ:7:ILE:O	1:BR:22:ILE:HG22	2.14	0.47
1:BQ:22:ILE:HG22	1:BR:7:ILE:O	2.14	0.47
1:BY:108:LEU:HA	1:BY:111:ILE:HG22	1.95	0.47
1:CL:22:ILE:HG22	1:CM:7:ILE:O	2.14	0.47
1:CV:29:GLN:HB3	1:CV:36:ASN:HD22	1.80	0.47
1:AC:9:THR:HG22	1:AC:111:ILE:HD12	1.95	0.47
1:AI:92:ILE:HG12	1:BY:92:ILE:HG12	1.94	0.47
1:AT:29:GLN:HB3	1:AT:36:ASN:HD22	1.80	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AY:7:ILE:O	1:AZ:22:ILE:HG22	2.14	0.47
1:BH:7:ILE:O	1:BI:22:ILE:HG22	2.14	0.47
1:BN:7:ILE:O	1:BO:22:ILE:HG22	2.14	0.47
1:BP:9:THR:HG22	1:BP:111:ILE:HD12	1.95	0.47
1:CF:7:ILE:O	1:CG:22:ILE:HG22	2.14	0.47
1:CO:7:ILE:O	1:CP:22:ILE:HG22	2.14	0.47
1:CS:79:SER:HA	1:CS:85:THR:HG22	1.96	0.47
1:DA:22:ILE:HG22	1:DB:7:ILE:O	2.14	0.47
1:DJ:22:ILE:HG22	1:DK:7:ILE:O	2.14	0.47
1:AC:122:ASP:OD1	1:BP:109:ARG:NH1	2.44	0.47
1:AF:68:GLY:O	1:AF:95:GLU:HA	2.14	0.47
1:AF:108:LEU:HA	1:AF:111:ILE:HG22	1.95	0.47
1:AG:127:LEU:HD13	1:CE:40:SER:HB2	1.95	0.47
1:AH:29:GLN:HB3	1:AH:36:ASN:HD22	1.80	0.47
1:AH:79:SER:HA	1:AH:85:THR:HG22	1.95	0.47
1:AI:9:THR:HG22	1:AI:111:ILE:HD12	1.95	0.47
1:AM:7:ILE:O	1:AN:22:ILE:HG22	2.14	0.47
1:AW:87:PHE:HE2	1:BT:60:VAL:HG11	1.79	0.47
1:BC:48:ARG:HH12	1:BO:127:LEU:HD21	1.79	0.47
1:BE:22:ILE:HG22	1:BF:7:ILE:O	2.14	0.47
1:BJ:45:TYR:CE2	1:BL:58:PRO:HD2	2.50	0.47
1:CI:7:ILE:O	1:CJ:22:ILE:HG22	2.14	0.47
1:CP:53:PHE:CD1	1:CP:72:VAL:HG22	2.50	0.47
1:CZ:9:THR:HG22	1:CZ:111:ILE:HD12	1.95	0.47
1:DE:29:GLN:HB3	1:DE:36:ASN:HD22	1.80	0.47
1:AJ:130:LEU:O	1:BY:26:LEU:HB3	2.14	0.47
1:AO:128:GLY:O	1:CC:28:GLY:HA3	2.14	0.47
1:AZ:53:PHE:CD1	1:AZ:72:VAL:HG22	2.50	0.47
1:AZ:79:SER:HA	1:AZ:85:THR:HG22	1.96	0.47
1:BI:79:SER:HA	1:BI:85:THR:HG22	1.95	0.47
1:BU:29:GLN:HB3	1:BU:36:ASN:HD22	1.79	0.47
1:CB:108:LEU:HA	1:CB:111:ILE:HG22	1.95	0.47
1:CQ:68:GLY:O	1:CQ:95:GLU:HA	2.14	0.47
1:CS:29:GLN:HB3	1:CS:36:ASN:HD22	1.80	0.47
1:CU:22:ILE:HG22	1:CV:7:ILE:O	2.14	0.47
1:AB:53:PHE:CD1	1:AB:72:VAL:HG22	2.50	0.47
1:AE:53:PHE:CD1	1:AE:72:VAL:HG22	2.50	0.47
1:AG:7:ILE:O	1:AH:22:ILE:HG22	2.14	0.47
1:AI:113:ALA:HB1	1:BY:113:ALA:HB1	1.97	0.47
1:AN:48:ARG:HH12	1:CA:127:LEU:HD21	1.80	0.47
1:AQ:53:PHE:CD1	1:AQ:72:VAL:HG22	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AR:22:ILE:HG22	1:CK:7:ILE:O	2.13	0.47
1:AR:68:GLY:O	1:AR:95:GLU:HA	2.14	0.47
1:AS:7:ILE:O	1:AT:22:ILE:HG22	2.14	0.47
1:AW:46:ARG:HG3	1:BT:65:TYR:CE2	2.50	0.47
1:AX:68:GLY:O	1:AX:95:GLU:HA	2.14	0.47
1:BB:127:LEU:HD13	1:BP:40:SER:HB2	1.97	0.47
1:BJ:108:LEU:HA	1:BJ:111:ILE:HG22	1.95	0.47
1:BO:53:PHE:CD1	1:BO:72:VAL:HG22	2.50	0.47
1:BT:7:ILE:O	1:BU:22:ILE:HG22	2.14	0.47
1:BX:29:GLN:HB3	1:BX:36:ASN:HD22	1.80	0.47
1:CA:29:GLN:HB3	1:CA:36:ASN:HD22	1.79	0.47
1:CB:68:GLY:O	1:CB:95:GLU:HA	2.14	0.47
1:CG:53:PHE:CD1	1:CG:72:VAL:HG22	2.50	0.47
1:CJ:87:PHE:HE2	1:DA:60:VAL:HG11	1.79	0.47
1:CL:7:ILE:O	1:CM:22:ILE:HG22	2.14	0.47
1:CR:7:ILE:O	1:CS:22:ILE:HG22	2.14	0.47
1:CS:53:PHE:CD1	1:CS:72:VAL:HG22	2.50	0.47
1:CV:79:SER:HA	1:CV:85:THR:HG22	1.95	0.47
1:CX:22:ILE:HG22	1:CY:7:ILE:O	2.14	0.47
1:CY:79:SER:HA	1:CY:85:THR:HG22	1.95	0.47
1:DB:29:GLN:HB3	1:DB:36:ASN:HD22	1.79	0.47
1:DE:79:SER:HA	1:DE:85:THR:HG22	1.96	0.47
1:AB:29:GLN:HB3	1:AB:36:ASN:HD22	1.80	0.47
1:AR:16:ILE:HD13	1:CK:114:GLN:NE2	2.17	0.47
1:AS:22:ILE:HG22	1:AT:7:ILE:O	2.14	0.47
1:BC:53:PHE:CD1	1:BC:72:VAL:HG22	2.50	0.47
1:BF:53:PHE:CD1	1:BF:72:VAL:HG22	2.50	0.47
1:BG:128:GLY:O	1:DA:28:GLY:HA3	2.15	0.47
1:BR:53:PHE:CD1	1:BR:72:VAL:HG22	2.50	0.47
1:CC:22:ILE:HG22	1:CD:7:ILE:O	2.14	0.47
1:CR:22:ILE:HG22	1:CS:7:ILE:O	2.14	0.47
1:CY:53:PHE:CD1	1:CY:72:VAL:HG22	2.50	0.47
1:DE:53:PHE:CD1	1:DE:72:VAL:HG22	2.50	0.47
1:DK:29:GLN:HB3	1:DK:36:ASN:HD22	1.80	0.47
1:DL:108:LEU:HA	1:DL:111:ILE:HG22	1.95	0.47
1:AH:53:PHE:CD1	1:AH:72:VAL:HG22	2.50	0.47
1:AI:110:LEU:O	1:AI:114:GLN:HG3	2.15	0.47
1:AJ:22:ILE:HG22	1:AK:7:ILE:O	2.14	0.47
1:BJ:9:THR:HG22	1:BJ:111:ILE:HD12	1.95	0.47
1:BZ:22:ILE:HG22	1:CA:7:ILE:O	2.14	0.47
1:CG:29:GLN:HB3	1:CG:36:ASN:HD22	1.80	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:DA:7:ILE:O	1:DB:22:ILE:HG22	2.14	0.47
1:DB:53:PHE:CD1	1:DB:72:VAL:HG22	2.50	0.47
1:DB:79:SER:HA	1:DB:85:THR:HG22	1.95	0.47
1:DK:53:PHE:CD1	1:DK:72:VAL:HG22	2.50	0.47
1:AM:22:ILE:HG22	1:AN:7:ILE:O	2.14	0.47
1:AN:53:PHE:CD1	1:AN:72:VAL:HG22	2.50	0.47
1:BP:110:LEU:O	1:BP:114:GLN:HG3	2.15	0.47
1:BV:9:THR:HG22	1:BV:111:ILE:HD12	1.95	0.47
1:BX:73:VAL:HG22	1:BX:91:ARG:HE	1.80	0.47
1:CJ:53:PHE:CD1	1:CJ:72:VAL:HG22	2.50	0.47
1:CJ:87:PHE:CE2	1:DA:60:VAL:HG11	2.50	0.47
1:AC:110:LEU:O	1:AC:114:GLN:HG3	2.15	0.46
1:AE:73:VAL:HG22	1:AE:91:ARG:HE	1.81	0.46
1:AK:53:PHE:CD1	1:AK:72:VAL:HG22	2.50	0.46
1:AL:110:LEU:O	1:AL:114:GLN:HG3	2.16	0.46
1:AO:110:LEU:O	1:AO:114:GLN:HG3	2.15	0.46
1:AT:53:PHE:CD1	1:AT:72:VAL:HG22	2.50	0.46
1:AY:22:ILE:HG22	1:AZ:7:ILE:O	2.14	0.46
1:BA:110:LEU:O	1:BA:114:GLN:HG3	2.15	0.46
1:BF:29:GLN:HB3	1:BF:36:ASN:HD22	1.79	0.46
1:BF:73:VAL:HG22	1:BF:91:ARG:HE	1.81	0.46
1:BR:73:VAL:HG22	1:BR:91:ARG:HE	1.80	0.46
1:BU:48:ARG:HH12	1:DH:127:LEU:HD21	1.80	0.46
1:BW:22:ILE:HG22	1:BX:7:ILE:O	2.14	0.46
1:CA:73:VAL:HG22	1:CA:91:ARG:HE	1.81	0.46
1:CD:29:GLN:HB3	1:CD:36:ASN:HD22	1.80	0.46
1:CT:110:LEU:O	1:CT:114:GLN:HG3	2.15	0.46
1:CV:53:PHE:CD1	1:CV:72:VAL:HG22	2.50	0.46
1:DL:110:LEU:O	1:DL:114:GLN:HG3	2.15	0.46
1:AG:22:ILE:HG22	1:AH:7:ILE:O	2.14	0.46
1:AG:28:GLY:HA3	1:BY:128:GLY:O	2.16	0.46
1:AN:79:SER:HA	1:AN:85:THR:HG22	1.96	0.46
1:AU:7:ILE:HB	1:AU:114:GLN:OE1	2.16	0.46
1:BL:53:PHE:CD1	1:BL:72:VAL:HG22	2.50	0.46
1:BT:22:ILE:HG22	1:BU:7:ILE:O	2.14	0.46
1:BU:53:PHE:CD1	1:BU:72:VAL:HG22	2.50	0.46
1:BV:110:LEU:O	1:BV:114:GLN:HG3	2.15	0.46
1:BX:53:PHE:CD1	1:BX:72:VAL:HG22	2.50	0.46
1:CH:110:LEU:O	1:CH:114:GLN:HG3	2.16	0.46
1:CI:22:ILE:HG22	1:CJ:7:ILE:O	2.14	0.46
1:CP:73:VAL:HG22	1:CP:91:ARG:HE	1.80	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CQ:110:LEU:O	1:CQ:114:GLN:HG3	2.15	0.46
1:CS:73:VAL:HG22	1:CS:91:ARG:HE	1.80	0.46
1:DD:22:ILE:HG22	1:DE:7:ILE:O	2.14	0.46
1:AK:65:TYR:HE2	1:BY:46:ARG:HA	1.80	0.46
1:AK:73:VAL:HG22	1:AK:91:ARG:HE	1.81	0.46
1:AR:7:ILE:O	1:CK:22:ILE:HG22	2.15	0.46
1:AW:65:TYR:CE2	1:BV:46:ARG:HG2	2.50	0.46
1:AZ:29:GLN:HB3	1:AZ:36:ASN:HD22	1.79	0.46
1:BD:110:LEU:O	1:BD:114:GLN:HG3	2.15	0.46
1:BI:73:VAL:HG22	1:BI:91:ARG:HE	1.81	0.46
1:BN:22:ILE:HG22	1:BO:7:ILE:O	2.14	0.46
1:CB:110:LEU:O	1:CB:114:GLN:HG3	2.16	0.46
1:CD:53:PHE:CD1	1:CD:72:VAL:HG22	2.50	0.46
1:CD:73:VAL:HG22	1:CD:91:ARG:HE	1.81	0.46
1:CJ:33:ASN:OD1	1:DC:30:THR:OG1	2.33	0.46
1:CT:7:ILE:HB	1:CT:114:GLN:OE1	2.16	0.46
1:DH:53:PHE:CD1	1:DH:72:VAL:HG22	2.50	0.46
1:DK:73:VAL:HG22	1:DK:91:ARG:HE	1.81	0.46
1:AE:127:LEU:CD2	1:DE:48:ARG:HH12	2.27	0.46
1:AV:22:ILE:HG22	1:AW:7:ILE:O	2.14	0.46
1:AX:7:ILE:HB	1:AX:114:GLN:OE1	2.16	0.46
1:BA:7:ILE:HB	1:BA:114:GLN:OE1	2.16	0.46
1:BU:73:VAL:HG22	1:BU:91:ARG:HE	1.80	0.46
1:CA:53:PHE:CD1	1:CA:72:VAL:HG22	2.50	0.46
1:CM:53:PHE:CD1	1:CM:72:VAL:HG22	2.50	0.46
1:CW:7:ILE:HB	1:CW:114:GLN:OE1	2.16	0.46
1:DF:110:LEU:O	1:DF:114:GLN:HG3	2.16	0.46
1:DH:29:GLN:HB3	1:DH:36:ASN:HD22	1.80	0.46
1:DL:7:ILE:HB	1:DL:114:GLN:OE1	2.16	0.46
1:AK:29:GLN:HG3	1:BW:57:VAL:HG21	1.98	0.46
1:AK:65:TYR:HD2	1:BY:46:ARG:HG2	1.79	0.46
1:BI:53:PHE:CD1	1:BI:72:VAL:HG22	2.50	0.46
1:BM:18:GLY:HA2	1:DI:11:ALA:HA	1.97	0.46
1:BO:29:GLN:HB3	1:BO:36:ASN:HD22	1.79	0.46
1:CQ:7:ILE:HB	1:CQ:114:GLN:OE1	2.16	0.46
1:DC:7:ILE:HB	1:DC:114:GLN:OE1	2.16	0.46
1:DE:73:VAL:HG22	1:DE:91:ARG:HE	1.81	0.46
1:DG:22:ILE:HG22	1:DH:7:ILE:O	2.14	0.46
1:AR:110:LEU:O	1:AR:114:GLN:HG3	2.15	0.46
1:AT:73:VAL:HG22	1:AT:91:ARG:HE	1.80	0.46
1:AU:110:LEU:O	1:AU:114:GLN:HG3	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AU:125:TRP:O	1:CN:70:ASN:ND2	2.49	0.46
1:AW:53:PHE:CD1	1:AW:72:VAL:HG22	2.50	0.46
1:AX:110:LEU:O	1:AX:114:GLN:HG3	2.15	0.46
1:BJ:7:ILE:HB	1:BJ:114:GLN:OE1	2.16	0.46
1:BM:7:ILE:HB	1:BM:114:GLN:OE1	2.16	0.46
1:CB:7:ILE:HB	1:CB:114:GLN:OE1	2.16	0.46
1:CM:73:VAL:HG22	1:CM:91:ARG:HE	1.80	0.46
1:CY:73:VAL:HG22	1:CY:91:ARG:HE	1.80	0.46
1:DC:110:LEU:O	1:DC:114:GLN:HG3	2.15	0.46
1:DF:7:ILE:HB	1:DF:114:GLN:OE1	2.16	0.46
1:AB:10:GLY:H	1:AB:107:ASP:CG	2.24	0.46
1:AW:29:GLN:HB3	1:AW:36:ASN:HD22	1.79	0.46
1:BC:10:GLY:H	1:BC:107:ASP:CG	2.24	0.46
1:BG:7:ILE:HB	1:BG:114:GLN:OE1	2.16	0.46
1:BJ:16:ILE:HG21	1:DF:114:GLN:NE2	2.31	0.46
1:BJ:46:ARG:CA	1:BL:65:TYR:HE2	2.27	0.46
1:BM:110:LEU:O	1:BM:114:GLN:HG3	2.15	0.46
1:BS:7:ILE:HB	1:BS:114:GLN:OE1	2.16	0.46
1:CK:7:ILE:HB	1:CK:114:GLN:OE1	2.16	0.46
1:CN:110:LEU:O	1:CN:114:GLN:HG3	2.16	0.46
1:CO:22:ILE:HG22	1:CP:7:ILE:O	2.14	0.46
1:CV:73:VAL:HG22	1:CV:91:ARG:HE	1.81	0.46
1:DH:73:VAL:HG22	1:DH:91:ARG:HE	1.81	0.46
1:AO:46:ARG:HA	1:CP:65:TYR:HE2	1.80	0.46
1:AZ:10:GLY:H	1:AZ:107:ASP:CG	2.24	0.46
1:AZ:73:VAL:HG22	1:AZ:91:ARG:HE	1.81	0.46
1:BI:29:GLN:HB3	1:BI:36:ASN:HD22	1.79	0.46
1:BO:10:GLY:H	1:BO:107:ASP:CG	2.24	0.46
1:BU:10:GLY:H	1:BU:107:ASP:CG	2.24	0.46
1:BY:110:LEU:O	1:BY:114:GLN:HG3	2.15	0.46
1:CJ:73:VAL:HG22	1:CJ:91:ARG:HE	1.81	0.46
1:CZ:7:ILE:HB	1:CZ:114:GLN:OE1	2.16	0.46
1:AK:10:GLY:H	1:AK:107:ASP:CG	2.24	0.46
1:AL:30:THR:OG1	1:CV:33:ASN:OD1	2.32	0.46
1:AT:10:GLY:H	1:AT:107:ASP:CG	2.24	0.46
1:BG:110:LEU:O	1:BG:114:GLN:HG3	2.15	0.46
1:BJ:11:ALA:HB3	1:BJ:110:LEU:HD23	1.98	0.46
1:CA:10:GLY:H	1:CA:107:ASP:CG	2.24	0.46
1:CE:110:LEU:O	1:CE:114:GLN:HG3	2.15	0.46
1:CH:7:ILE:HB	1:CH:114:GLN:OE1	2.16	0.46
1:CJ:10:GLY:H	1:CJ:107:ASP:CG	2.24	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CP:10:GLY:H	1:CP:107:ASP:CG	2.24	0.46
1:DI:7:ILE:HB	1:DI:114:GLN:OE1	2.16	0.46
1:AG:65:TYR:CE2	1:CS:46:ARG:HG3	2.50	0.46
1:AH:10:GLY:H	1:AH:107:ASP:CG	2.24	0.46
1:AX:89:SER:HB2	1:CT:95:GLU:HB2	1.97	0.46
1:BA:11:ALA:HB3	1:BA:110:LEU:HD23	1.98	0.46
1:BC:73:VAL:HG22	1:BC:91:ARG:HE	1.81	0.46
1:CB:11:ALA:HB3	1:CB:110:LEU:HD23	1.98	0.46
1:CE:7:ILE:HB	1:CE:114:GLN:OE1	2.16	0.46
1:CT:11:ALA:HB3	1:CT:110:LEU:HD23	1.98	0.46
1:DE:10:GLY:H	1:DE:107:ASP:CG	2.24	0.46
1:AE:48:ARG:HH22	1:BF:127:LEU:HD23	1.81	0.45
1:AI:14:ALA:HA	1:BY:16:ILE:HA	1.96	0.45
1:AO:114:GLN:NE2	1:CE:16:ILE:HD13	2.20	0.45
1:AQ:29:GLN:HB3	1:AQ:36:ASN:HD22	1.79	0.45
1:AW:73:VAL:HG22	1:AW:91:ARG:HE	1.81	0.45
1:BA:6:SER:HA	1:CW:22:ILE:O	2.16	0.45
1:BF:10:GLY:H	1:BF:107:ASP:CG	2.24	0.45
1:BF:65:TYR:CE2	1:BS:46:ARG:CG	2.97	0.45
1:BP:7:ILE:HB	1:BP:114:GLN:OE1	2.16	0.45
1:BS:11:ALA:HB3	1:BS:110:LEU:HD23	1.98	0.45
1:BS:110:LEU:O	1:BS:114:GLN:HG3	2.15	0.45
1:CG:73:VAL:HG22	1:CG:91:ARG:HE	1.80	0.45
1:CN:11:ALA:HB3	1:CN:110:LEU:HD23	1.98	0.45
1:AC:7:ILE:HB	1:AC:114:GLN:OE1	2.16	0.45
1:AE:10:GLY:H	1:AE:107:ASP:CG	2.24	0.45
1:AH:73:VAL:HG22	1:AH:91:ARG:HE	1.80	0.45
1:AL:46:ARG:HG2	1:CV:65:TYR:CD2	2.51	0.45
1:AO:7:ILE:HB	1:AO:114:GLN:OE1	2.16	0.45
1:BD:7:ILE:HB	1:BD:114:GLN:OE1	2.16	0.45
1:BI:42:ASP:O	1:BI:48:ARG:HD3	2.17	0.45
1:BJ:114:GLN:NE2	1:DF:16:ILE:HD13	2.22	0.45
1:BP:11:ALA:HB3	1:BP:110:LEU:HD23	1.98	0.45
1:CD:10:GLY:H	1:CD:107:ASP:CG	2.24	0.45
1:CJ:42:ASP:O	1:CJ:48:ARG:HD3	2.17	0.45
1:DB:10:GLY:H	1:DB:107:ASP:CG	2.24	0.45
1:DF:11:ALA:HB3	1:DF:110:LEU:HD23	1.98	0.45
1:DH:10:GLY:H	1:DH:107:ASP:CG	2.24	0.45
1:AL:7:ILE:HB	1:AL:114:GLN:OE1	2.16	0.45
1:AN:73:VAL:HG22	1:AN:91:ARG:HE	1.80	0.45
1:AX:19:GLY:O	1:CT:8:LYS:HD2	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BG:111:ILE:HD13	1:DC:74:PHE:CE2	2.51	0.45
1:BM:18:GLY:CA	1:DI:11:ALA:HA	2.47	0.45
1:BY:7:ILE:HB	1:BY:114:GLN:OE1	2.16	0.45
1:CJ:29:GLN:HB3	1:CJ:36:ASN:HD22	1.79	0.45
1:CM:42:ASP:O	1:CM:48:ARG:HD3	2.17	0.45
1:CV:42:ASP:O	1:CV:48:ARG:HD3	2.17	0.45
1:CY:42:ASP:O	1:CY:48:ARG:HD3	2.17	0.45
1:DI:110:LEU:O	1:DI:114:GLN:HG3	2.16	0.45
1:AE:42:ASP:O	1:AE:48:ARG:HD3	2.17	0.45
1:AI:11:ALA:HB3	1:AI:110:LEU:HD23	1.98	0.45
1:BC:42:ASP:O	1:BC:48:ARG:HD3	2.17	0.45
1:BG:11:ALA:HB3	1:BG:110:LEU:HD23	1.98	0.45
1:BI:60:VAL:HG22	1:BI:65:TYR:CE1	2.52	0.45
1:BR:42:ASP:O	1:BR:48:ARG:HD3	2.17	0.45
1:BX:60:VAL:HG22	1:BX:65:TYR:CE1	2.52	0.45
1:CA:42:ASP:O	1:CA:48:ARG:HD3	2.17	0.45
1:CM:60:VAL:HG22	1:CM:65:TYR:CE1	2.52	0.45
1:CN:7:ILE:HB	1:CN:114:GLN:OE1	2.16	0.45
1:CS:42:ASP:O	1:CS:48:ARG:HD3	2.17	0.45
1:CZ:110:LEU:O	1:CZ:114:GLN:HG3	2.15	0.45
1:AF:110:LEU:O	1:AF:114:GLN:HG3	2.15	0.45
1:AH:60:VAL:HG22	1:AH:65:TYR:CE1	2.52	0.45
1:AH:65:TYR:HE2	1:CE:46:ARG:HA	1.82	0.45
1:AI:7:ILE:HB	1:AI:114:GLN:OE1	2.16	0.45
1:AR:7:ILE:HB	1:AR:114:GLN:OE1	2.16	0.45
1:AX:11:ALA:HB3	1:AX:110:LEU:HD23	1.98	0.45
1:AZ:60:VAL:HG22	1:AZ:65:TYR:CE1	2.52	0.45
1:BL:10:GLY:H	1:BL:107:ASP:CG	2.24	0.45
1:BL:73:VAL:HG22	1:BL:91:ARG:HE	1.81	0.45
1:BO:73:VAL:HG22	1:BO:91:ARG:HE	1.80	0.45
1:CD:42:ASP:O	1:CD:48:ARG:HD3	2.17	0.45
1:CS:60:VAL:HG22	1:CS:65:TYR:CE1	2.52	0.45
1:CW:110:LEU:O	1:CW:114:GLN:HG3	2.15	0.45
1:DB:42:ASP:O	1:DB:48:ARG:HD3	2.17	0.45
1:DH:60:VAL:HG22	1:DH:65:TYR:CE1	2.52	0.45
1:AB:42:ASP:O	1:AB:48:ARG:HD3	2.17	0.45
1:AC:11:ALA:HB3	1:AC:110:LEU:HD23	1.98	0.45
1:AF:7:ILE:HB	1:AF:114:GLN:OE1	2.16	0.45
1:AF:89:SER:HB2	1:BV:95:GLU:HB2	1.98	0.45
1:AN:42:ASP:O	1:AN:48:ARG:HD3	2.17	0.45
1:AW:10:GLY:H	1:AW:107:ASP:CG	2.24	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BI:10:GLY:H	1:BI:107:ASP:CG	2.24	0.45
1:BL:42:ASP:O	1:BL:48:ARG:HD3	2.17	0.45
1:BR:10:GLY:H	1:BR:107:ASP:CG	2.24	0.45
1:BV:7:ILE:HB	1:BV:114:GLN:OE1	2.16	0.45
1:BW:87:PHE:O	1:BX:67:LYS:NZ	2.41	0.45
1:BX:10:GLY:H	1:BX:107:ASP:CG	2.24	0.45
1:CD:60:VAL:HG22	1:CD:65:TYR:CE1	2.52	0.45
1:CG:10:GLY:H	1:CG:107:ASP:CG	2.24	0.45
1:CK:110:LEU:O	1:CK:114:GLN:HG3	2.16	0.45
1:DK:42:ASP:O	1:DK:48:ARG:HD3	2.17	0.45
1:AK:42:ASP:O	1:AK:48:ARG:HD3	2.17	0.45
1:AL:19:GLY:O	1:CB:8:LYS:HD2	2.15	0.45
1:AQ:10:GLY:H	1:AQ:107:ASP:CG	2.24	0.45
1:AQ:73:VAL:HG22	1:AQ:91:ARG:HE	1.80	0.45
1:AW:60:VAL:HG22	1:AW:65:TYR:CE1	2.52	0.45
1:BU:26:LEU:O	1:DH:129:ALA:HB2	2.17	0.45
1:CG:42:ASP:O	1:CG:48:ARG:HD3	2.17	0.45
1:CG:60:VAL:HG22	1:CG:65:TYR:CE1	2.52	0.45
1:CH:11:ALA:HB3	1:CH:110:LEU:HD23	1.98	0.45
1:CY:60:VAL:HG22	1:CY:65:TYR:CE1	2.52	0.45
1:DC:11:ALA:HB3	1:DC:110:LEU:HD23	1.98	0.45
1:AO:92:ILE:HG12	1:CE:92:ILE:HG12	1.98	0.45
1:AQ:42:ASP:O	1:AQ:48:ARG:HD3	2.17	0.45
1:AZ:42:ASP:O	1:AZ:48:ARG:HD3	2.17	0.45
1:BJ:110:LEU:O	1:BJ:114:GLN:HG3	2.15	0.45
1:BL:60:VAL:HG22	1:BL:65:TYR:CE1	2.52	0.45
1:CV:60:VAL:HG22	1:CV:65:TYR:CE1	2.52	0.45
1:CW:46:ARG:HA	1:CY:65:TYR:CE2	2.50	0.45
1:DB:60:VAL:HG22	1:DB:65:TYR:CE1	2.52	0.45
1:DK:60:VAL:HG22	1:DK:65:TYR:CE1	2.52	0.45
1:AB:60:VAL:HG22	1:AB:65:TYR:CE1	2.52	0.45
1:AI:39:VAL:HG12	1:AI:42:ASP:H	1.82	0.45
1:AK:60:VAL:HG22	1:AK:65:TYR:CE1	2.52	0.45
1:AT:42:ASP:O	1:AT:48:ARG:HD3	2.17	0.45
1:BB:130:LEU:O	1:BP:26:LEU:HB3	2.16	0.45
1:BF:60:VAL:HG22	1:BF:65:TYR:CE1	2.52	0.45
1:BO:60:VAL:HG22	1:BO:65:TYR:CE1	2.52	0.45
1:CM:10:GLY:H	1:CM:107:ASP:CG	2.24	0.45
1:CQ:39:VAL:HG12	1:CQ:42:ASP:H	1.82	0.45
1:CS:10:GLY:H	1:CS:107:ASP:CG	2.24	0.45
1:DC:39:VAL:HG12	1:DC:42:ASP:H	1.82	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AB:48:ARG:HH12	1:AT:127:LEU:CD2	2.29	0.45
1:AL:11:ALA:HB3	1:AL:110:LEU:HD23	1.98	0.45
1:AP:26:LEU:HD12	1:AP:36:ASN:O	2.17	0.45
1:BD:11:ALA:HB3	1:BD:110:LEU:HD23	1.98	0.45
1:BM:13:SER:O	1:DI:16:ILE:HG13	2.16	0.45
1:BX:42:ASP:O	1:BX:48:ARG:HD3	2.17	0.45
1:BY:11:ALA:HB3	1:BY:110:LEU:HD23	1.98	0.45
1:CY:10:GLY:H	1:CY:107:ASP:CG	2.24	0.45
1:CZ:39:VAL:HG12	1:CZ:42:ASP:H	1.82	0.45
1:DB:73:VAL:HG22	1:DB:91:ARG:HE	1.81	0.45
1:DH:42:ASP:O	1:DH:48:ARG:HD3	2.17	0.45
1:DL:39:VAL:HG12	1:DL:42:ASP:H	1.82	0.45
1:AF:11:ALA:HB3	1:AF:110:LEU:HD23	1.98	0.44
1:AJ:26:LEU:HD12	1:AJ:36:ASN:O	2.17	0.44
1:AN:60:VAL:HG22	1:AN:65:TYR:CE1	2.52	0.44
1:AT:60:VAL:HG22	1:AT:65:TYR:CE1	2.52	0.44
1:BC:60:VAL:HG22	1:BC:65:TYR:CE1	2.52	0.44
1:BQ:108:LEU:HA	1:BQ:111:ILE:HG22	2.00	0.44
1:CA:60:VAL:HG22	1:CA:65:TYR:CE1	2.52	0.44
1:CO:26:LEU:HD12	1:CO:36:ASN:O	2.18	0.44
1:CV:10:GLY:H	1:CV:107:ASP:CG	2.24	0.44
1:AC:39:VAL:HG12	1:AC:42:ASP:H	1.82	0.44
1:AG:26:LEU:HD12	1:AG:36:ASN:O	2.17	0.44
1:AH:42:ASP:O	1:AH:48:ARG:HD3	2.17	0.44
1:AR:11:ALA:HB3	1:AR:110:LEU:HD23	1.98	0.44
1:AU:11:ALA:HB3	1:AU:110:LEU:HD23	1.99	0.44
1:BM:16:ILE:HA	1:DI:14:ALA:HA	1.99	0.44
1:BM:116:LEU:O	1:DI:109:ARG:CZ	2.65	0.44
1:BO:42:ASP:O	1:BO:48:ARG:HD3	2.17	0.44
1:BS:39:VAL:HG12	1:BS:42:ASP:H	1.83	0.44
1:BU:42:ASP:O	1:BU:48:ARG:HD3	2.17	0.44
1:BY:39:VAL:HG12	1:BY:42:ASP:H	1.82	0.44
1:BZ:108:LEU:HA	1:BZ:111:ILE:HG22	2.00	0.44
1:CD:58:PRO:HD2	1:CK:45:TYR:CZ	2.52	0.44
1:CH:39:VAL:HG12	1:CH:42:ASP:H	1.82	0.44
1:CJ:60:VAL:HG22	1:CJ:65:TYR:CE1	2.52	0.44
1:CX:26:LEU:HD12	1:CX:36:ASN:O	2.18	0.44
1:DA:108:LEU:HA	1:DA:111:ILE:HG22	1.99	0.44
1:DE:42:ASP:O	1:DE:48:ARG:HD3	2.17	0.44
1:DI:11:ALA:HB3	1:DI:110:LEU:HD23	1.98	0.44
1:DJ:108:LEU:HA	1:DJ:111:ILE:HG22	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:108:LEU:HA	1:AA:111:ILE:HG22	2.00	0.44
1:AB:73:VAL:HG22	1:AB:91:ARG:HE	1.80	0.44
1:AO:39:VAL:HG12	1:AO:42:ASP:H	1.82	0.44
1:AQ:60:VAL:HG22	1:AQ:65:TYR:CE1	2.52	0.44
1:AR:39:VAL:HG12	1:AR:42:ASP:H	1.82	0.44
1:AS:26:LEU:HD12	1:AS:36:ASN:O	2.17	0.44
1:AS:108:LEU:HA	1:AS:111:ILE:HG22	2.00	0.44
1:BF:42:ASP:O	1:BF:48:ARG:HD3	2.17	0.44
1:BR:60:VAL:HG22	1:BR:65:TYR:CE1	2.52	0.44
1:CE:11:ALA:HB3	1:CE:110:LEU:HD23	1.98	0.44
1:CJ:68:GLY:O	1:CJ:95:GLU:HA	2.18	0.44
1:CO:108:LEU:HA	1:CO:111:ILE:HG22	1.99	0.44
1:DK:10:GLY:H	1:DK:107:ASP:CG	2.24	0.44
1:AB:87:PHE:CZ	1:AS:60:VAL:HG11	2.51	0.44
1:AE:68:GLY:O	1:AE:95:GLU:HA	2.18	0.44
1:AG:108:LEU:HA	1:AG:111:ILE:HG22	1.99	0.44
1:AL:114:GLN:NE2	1:CB:16:ILE:HG21	2.33	0.44
1:AO:11:ALA:HB3	1:AO:110:LEU:HD23	1.98	0.44
1:AR:125:TRP:O	1:CK:70:ASN:ND2	2.51	0.44
1:AV:108:LEU:HA	1:AV:111:ILE:HG22	2.00	0.44
1:AY:26:LEU:HD12	1:AY:36:ASN:O	2.18	0.44
1:BD:39:VAL:HG12	1:BD:42:ASP:H	1.82	0.44
1:BF:68:GLY:O	1:BF:95:GLU:HA	2.18	0.44
1:BN:108:LEU:HA	1:BN:111:ILE:HG22	2.00	0.44
1:BO:65:TYR:CD2	1:DL:46:ARG:HG2	2.52	0.44
1:BZ:26:LEU:HD12	1:BZ:36:ASN:O	2.18	0.44
1:CK:11:ALA:HB3	1:CK:110:LEU:HD23	1.98	0.44
1:CL:35:LEU:HD22	1:CM:131:ALA:H	1.83	0.44
1:CL:108:LEU:HA	1:CL:111:ILE:HG22	2.00	0.44
1:CP:42:ASP:O	1:CP:48:ARG:HD3	2.17	0.44
1:CW:11:ALA:HB3	1:CW:110:LEU:HD23	1.98	0.44
1:DD:26:LEU:HD12	1:DD:36:ASN:O	2.18	0.44
1:DJ:35:LEU:HD22	1:DK:131:ALA:H	1.83	0.44
1:AC:14:ALA:HB2	1:BP:16:ILE:HD12	1.98	0.44
1:AF:48:ARG:NH1	1:DD:127:LEU:HD22	2.32	0.44
1:AP:108:LEU:HA	1:AP:111:ILE:HG22	1.99	0.44
1:AV:26:LEU:HD12	1:AV:36:ASN:O	2.18	0.44
1:AW:42:ASP:O	1:AW:48:ARG:HD3	2.17	0.44
1:AX:39:VAL:HG12	1:AX:42:ASP:H	1.82	0.44
1:BD:74:PHE:CZ	1:CZ:111:ILE:HD13	2.52	0.44
1:BK:87:PHE:O	1:BL:67:LYS:NZ	2.41	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BL:58:PRO:HB3	1:BL:66:SER:O	2.18	0.44
1:BP:39:VAL:HG12	1:BP:42:ASP:H	1.82	0.44
1:BU:60:VAL:HG22	1:BU:65:TYR:CE1	2.52	0.44
1:BZ:35:LEU:HD22	1:CA:131:ALA:H	1.83	0.44
1:CC:26:LEU:HD12	1:CC:36:ASN:O	2.17	0.44
1:CF:35:LEU:HD22	1:CG:131:ALA:H	1.83	0.44
1:CF:108:LEU:HA	1:CF:111:ILE:HG22	2.00	0.44
1:CG:68:GLY:O	1:CG:95:GLU:HA	2.18	0.44
1:CI:108:LEU:HA	1:CI:111:ILE:HG22	1.99	0.44
1:DI:39:VAL:HG12	1:DI:42:ASP:H	1.82	0.44
1:AD:26:LEU:HD12	1:AD:36:ASN:O	2.17	0.44
1:AG:35:LEU:HD22	1:AH:131:ALA:H	1.83	0.44
1:AM:26:LEU:HD12	1:AM:36:ASN:O	2.17	0.44
1:AS:35:LEU:HD22	1:AT:131:ALA:H	1.83	0.44
1:AT:67:LYS:HD3	1:AT:95:GLU:HB3	2.00	0.44
1:AU:39:VAL:HG12	1:AU:42:ASP:H	1.82	0.44
1:AZ:67:LYS:HD3	1:AZ:95:GLU:HB3	2.00	0.44
1:BK:26:LEU:HD12	1:BK:36:ASN:O	2.17	0.44
1:BL:68:GLY:O	1:BL:95:GLU:HA	2.18	0.44
1:BT:35:LEU:HD22	1:BU:131:ALA:H	1.83	0.44
1:BU:68:GLY:O	1:BU:95:GLU:HA	2.18	0.44
1:BV:11:ALA:HB3	1:BV:110:LEU:HD23	1.98	0.44
1:CF:26:LEU:HD12	1:CF:36:ASN:O	2.17	0.44
1:CN:39:VAL:HG12	1:CN:42:ASP:H	1.82	0.44
1:CW:39:VAL:HG12	1:CW:42:ASP:H	1.82	0.44
1:CZ:11:ALA:HB3	1:CZ:110:LEU:HD23	1.98	0.44
1:DA:26:LEU:HD12	1:DA:36:ASN:O	2.18	0.44
1:DD:35:LEU:HD22	1:DE:131:ALA:H	1.83	0.44
1:DE:60:VAL:HG22	1:DE:65:TYR:CE1	2.52	0.44
1:DE:68:GLY:O	1:DE:95:GLU:HA	2.18	0.44
1:AA:35:LEU:HD22	1:AB:131:ALA:H	1.83	0.44
1:AK:68:GLY:O	1:AK:95:GLU:HA	2.18	0.44
1:AL:39:VAL:HG12	1:AL:42:ASP:H	1.82	0.44
1:AN:10:GLY:H	1:AN:107:ASP:CG	2.24	0.44
1:AT:68:GLY:O	1:AT:95:GLU:HA	2.18	0.44
1:BC:98:PRO:HB2	1:BP:46:ARG:HG3	2.00	0.44
1:BG:111:ILE:CD1	1:DC:74:PHE:CZ	2.99	0.44
1:BJ:22:ILE:HG22	1:DF:7:ILE:O	2.17	0.44
1:BO:58:PRO:HB3	1:BO:66:SER:O	2.18	0.44
1:BQ:26:LEU:HD12	1:BQ:36:ASN:O	2.17	0.44
1:BX:58:PRO:HB2	1:CN:45:TYR:CE2	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:58:PRO:HB3	1:CA:66:SER:O	2.18	0.44
1:CB:39:VAL:HG12	1:CB:42:ASP:H	1.82	0.44
1:CD:58:PRO:HB3	1:CD:66:SER:O	2.18	0.44
1:CL:26:LEU:HD12	1:CL:36:ASN:O	2.17	0.44
1:CR:35:LEU:HD22	1:CS:131:ALA:H	1.83	0.44
1:CR:108:LEU:HA	1:CR:111:ILE:HG22	2.00	0.44
1:CS:67:LYS:HD3	1:CS:95:GLU:HB3	2.00	0.44
1:CS:68:GLY:O	1:CS:95:GLU:HA	2.18	0.44
1:CT:30:THR:OG1	1:DB:33:ASN:OD1	2.33	0.44
1:CT:39:VAL:HG12	1:CT:42:ASP:H	1.82	0.44
1:AE:60:VAL:HG22	1:AE:65:TYR:CE1	2.52	0.44
1:AN:68:GLY:O	1:AN:95:GLU:HA	2.18	0.44
1:AU:18:GLY:HA2	1:CN:12:THR:H	1.83	0.44
1:AW:67:LYS:HD3	1:AW:95:GLU:HB3	2.00	0.44
1:BE:108:LEU:HA	1:BE:111:ILE:HG22	2.00	0.44
1:BG:22:ILE:HG22	1:DC:7:ILE:O	2.18	0.44
1:BI:68:GLY:O	1:BI:95:GLU:HA	2.18	0.44
1:BM:11:ALA:HB3	1:BM:110:LEU:HD23	1.98	0.44
1:BT:26:LEU:HD12	1:BT:36:ASN:O	2.17	0.44
1:BT:108:LEU:HA	1:BT:111:ILE:HG22	2.00	0.44
1:BX:67:LYS:HD3	1:BX:95:GLU:HB3	2.00	0.44
1:CG:58:PRO:HB3	1:CG:66:SER:O	2.18	0.44
1:CK:39:VAL:HG12	1:CK:42:ASP:H	1.82	0.44
1:CM:58:PRO:HB3	1:CM:66:SER:O	2.18	0.44
1:CP:60:VAL:HG22	1:CP:65:TYR:CE1	2.52	0.44
1:CR:26:LEU:HD12	1:CR:36:ASN:O	2.18	0.44
1:CV:67:LYS:HD3	1:CV:95:GLU:HB3	2.00	0.44
1:CV:68:GLY:O	1:CV:95:GLU:HA	2.18	0.44
1:DH:58:PRO:HB3	1:DH:66:SER:O	2.18	0.44
1:AA:26:LEU:HD12	1:AA:36:ASN:O	2.17	0.44
1:AN:26:LEU:O	1:CA:129:ALA:HB2	2.17	0.44
1:AO:46:ARG:HG2	1:CP:65:TYR:HD2	1.82	0.44
1:AW:58:PRO:HB3	1:AW:66:SER:O	2.18	0.44
1:BD:130:LEU:HD21	1:CZ:55:SER:HB3	1.99	0.44
1:BI:58:PRO:HB3	1:BI:66:SER:O	2.18	0.44
1:BJ:39:VAL:HG12	1:BJ:42:ASP:H	1.83	0.44
1:BN:26:LEU:HD12	1:BN:36:ASN:O	2.18	0.44
1:BR:58:PRO:HB3	1:BR:66:SER:O	2.18	0.44
1:BW:26:LEU:HD12	1:BW:36:ASN:O	2.17	0.44
1:CA:67:LYS:HD3	1:CA:95:GLU:HB3	2.00	0.44
1:CC:108:LEU:HA	1:CC:111:ILE:HG22	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CI:26:LEU:HD12	1:CI:36:ASN:O	2.18	0.44
1:CJ:58:PRO:HB3	1:CJ:66:SER:O	2.18	0.44
1:CS:58:PRO:HB3	1:CS:66:SER:O	2.18	0.44
1:DJ:26:LEU:HD12	1:DJ:36:ASN:O	2.18	0.44
1:AH:58:PRO:HB3	1:AH:66:SER:O	2.18	0.43
1:AK:58:PRO:HB3	1:AK:66:SER:O	2.18	0.43
1:AW:68:GLY:O	1:AW:95:GLU:HA	2.18	0.43
1:AZ:68:GLY:O	1:AZ:95:GLU:HA	2.18	0.43
1:BG:39:VAL:HG12	1:BG:42:ASP:H	1.82	0.43
1:BX:58:PRO:HB3	1:BX:66:SER:O	2.18	0.43
1:CO:35:LEU:HD22	1:CP:131:ALA:H	1.83	0.43
1:CY:58:PRO:HB3	1:CY:66:SER:O	2.18	0.43
1:DB:67:LYS:HD3	1:DB:95:GLU:HB3	2.00	0.43
1:DK:58:PRO:HB3	1:DK:66:SER:O	2.18	0.43
1:AA:99:ALA:HA	1:DH:46:ARG:NH2	2.33	0.43
1:AD:35:LEU:HD22	1:AE:131:ALA:H	1.83	0.43
1:AD:60:VAL:HG11	1:DE:87:PHE:HE2	1.82	0.43
1:AD:108:LEU:HA	1:AD:111:ILE:HG22	2.00	0.43
1:AF:39:VAL:HG12	1:AF:42:ASP:H	1.82	0.43
1:AN:58:PRO:HB3	1:AN:66:SER:O	2.18	0.43
1:AQ:58:PRO:HB3	1:AQ:66:SER:O	2.18	0.43
1:BB:108:LEU:HA	1:BB:111:ILE:HG22	2.00	0.43
1:BJ:16:ILE:HD13	1:DF:114:GLN:NE2	2.20	0.43
1:BM:129:ALA:HB1	1:DG:26:LEU:HB3	2.00	0.43
1:BR:68:GLY:O	1:BR:95:GLU:HA	2.18	0.43
1:BW:108:LEU:HA	1:BW:111:ILE:HG22	2.00	0.43
1:CA:58:PRO:HD2	1:CH:45:TYR:CE2	2.54	0.43
1:CQ:11:ALA:HB3	1:CQ:110:LEU:HD23	1.98	0.43
1:CU:26:LEU:HD12	1:CU:36:ASN:O	2.17	0.43
1:CU:35:LEU:HD22	1:CV:131:ALA:H	1.83	0.43
1:DB:68:GLY:O	1:DB:95:GLU:HA	2.18	0.43
1:DG:26:LEU:HD12	1:DG:36:ASN:O	2.18	0.43
1:DL:11:ALA:HB3	1:DL:110:LEU:HD23	1.98	0.43
1:AJ:35:LEU:HD22	1:AK:131:ALA:H	1.83	0.43
1:AO:16:ILE:HD13	1:CE:114:GLN:NE2	2.18	0.43
1:BE:26:LEU:HD12	1:BE:36:ASN:O	2.18	0.43
1:BH:26:LEU:HD12	1:BH:36:ASN:O	2.17	0.43
1:BK:35:LEU:HD22	1:BL:131:ALA:H	1.83	0.43
1:BM:109:ARG:O	1:DI:116:LEU:HD13	2.18	0.43
1:BQ:35:LEU:HD22	1:BR:131:ALA:H	1.83	0.43
1:CC:111:ILE:O	1:CC:115:LEU:HG	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CN:7:ILE:HD12	1:CN:111:ILE:HG13	2.01	0.43
1:CP:68:GLY:O	1:CP:95:GLU:HA	2.18	0.43
1:CT:7:ILE:HD12	1:CT:111:ILE:HG13	2.01	0.43
1:CX:35:LEU:HD22	1:CY:131:ALA:H	1.83	0.43
1:DB:58:PRO:HB3	1:DB:66:SER:O	2.18	0.43
1:DH:68:GLY:O	1:DH:95:GLU:HA	2.18	0.43
1:AC:16:ILE:HD13	1:BP:114:GLN:NE2	2.21	0.43
1:AJ:108:LEU:HA	1:AJ:111:ILE:HG22	2.00	0.43
1:AM:53:PHE:CD1	1:AM:72:VAL:HG22	2.54	0.43
1:BA:39:VAL:HG12	1:BA:42:ASP:H	1.82	0.43
1:BE:53:PHE:CD1	1:BE:72:VAL:HG22	2.54	0.43
1:BH:111:ILE:O	1:BH:115:LEU:HG	2.19	0.43
1:BI:67:LYS:HD3	1:BI:95:GLU:HB3	2.00	0.43
1:BK:108:LEU:HA	1:BK:111:ILE:HG22	2.00	0.43
1:BM:90:VAL:CG1	1:DI:108:LEU:HD22	2.48	0.43
1:BO:65:TYR:HE2	1:DL:46:ARG:HA	1.83	0.43
1:BU:58:PRO:HB3	1:BU:66:SER:O	2.18	0.43
1:BU:67:LYS:HD3	1:BU:95:GLU:HB3	2.00	0.43
1:CM:68:GLY:O	1:CM:95:GLU:HA	2.18	0.43
1:CP:67:LYS:HD3	1:CP:95:GLU:HB3	2.00	0.43
1:DD:111:ILE:O	1:DD:115:LEU:HG	2.19	0.43
1:DG:53:PHE:CD1	1:DG:72:VAL:HG22	2.54	0.43
1:DG:108:LEU:HA	1:DG:111:ILE:HG22	1.99	0.43
1:AF:48:ARG:HH12	1:DD:127:LEU:CB	2.31	0.43
1:AV:53:PHE:CD1	1:AV:72:VAL:HG22	2.54	0.43
1:AY:53:PHE:CD1	1:AY:72:VAL:HG22	2.54	0.43
1:BB:35:LEU:HD22	1:BC:131:ALA:H	1.83	0.43
1:BB:87:PHE:O	1:BC:67:LYS:NZ	2.41	0.43
1:BC:58:PRO:HB3	1:BC:66:SER:O	2.18	0.43
1:BC:67:LYS:HD3	1:BC:95:GLU:HB3	2.00	0.43
1:BC:68:GLY:O	1:BC:95:GLU:HA	2.18	0.43
1:BD:7:ILE:HD12	1:BD:111:ILE:HG13	2.01	0.43
1:BF:67:LYS:HD3	1:BF:95:GLU:HB3	2.00	0.43
1:BS:7:ILE:HD12	1:BS:111:ILE:HG13	2.01	0.43
1:BV:39:VAL:HG12	1:BV:42:ASP:H	1.82	0.43
1:BX:68:GLY:O	1:BX:95:GLU:HA	2.18	0.43
1:CA:68:GLY:O	1:CA:95:GLU:HA	2.18	0.43
1:CP:58:PRO:HB3	1:CP:66:SER:O	2.18	0.43
1:CU:111:ILE:O	1:CU:115:LEU:HG	2.19	0.43
1:CY:68:GLY:O	1:CY:95:GLU:HA	2.18	0.43
1:DE:58:PRO:HB3	1:DE:66:SER:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:DH:67:LYS:HD3	1:DH:95:GLU:HB3	2.00	0.43
1:AB:68:GLY:O	1:AB:95:GLU:HA	2.18	0.43
1:AK:67:LYS:HD3	1:AK:95:GLU:HB3	2.00	0.43
1:AY:111:ILE:O	1:AY:115:LEU:HG	2.19	0.43
1:BB:53:PHE:CD1	1:BB:72:VAL:HG22	2.54	0.43
1:BK:53:PHE:CD1	1:BK:72:VAL:HG22	2.54	0.43
1:BL:67:LYS:HD3	1:BL:95:GLU:HB3	2.00	0.43
1:BM:124:PHE:CE2	1:DI:70:ASN:HB3	2.53	0.43
1:BW:35:LEU:HD22	1:BX:131:ALA:H	1.83	0.43
1:CE:7:ILE:HD12	1:CE:111:ILE:HG13	2.01	0.43
1:CG:67:LYS:HD3	1:CG:95:GLU:HB3	2.00	0.43
1:CK:7:ILE:HD12	1:CK:111:ILE:HG13	2.01	0.43
1:DA:53:PHE:CD1	1:DA:72:VAL:HG22	2.54	0.43
1:DI:7:ILE:HD12	1:DI:111:ILE:HG13	2.01	0.43
1:AC:22:ILE:HG22	1:BP:7:ILE:O	2.18	0.43
1:AM:35:LEU:HD22	1:AN:131:ALA:H	1.83	0.43
1:AM:111:ILE:O	1:AM:115:LEU:HG	2.19	0.43
1:AO:18:GLY:HA2	1:CE:12:THR:H	1.83	0.43
1:AP:53:PHE:CD1	1:AP:72:VAL:HG22	2.54	0.43
1:AQ:68:GLY:O	1:AQ:95:GLU:HA	2.18	0.43
1:BE:111:ILE:O	1:BE:115:LEU:HG	2.19	0.43
1:BF:58:PRO:HB3	1:BF:66:SER:O	2.18	0.43
1:BH:108:LEU:HA	1:BH:111:ILE:HG22	2.00	0.43
1:BJ:7:ILE:HD12	1:BJ:111:ILE:HG13	2.01	0.43
1:BK:111:ILE:O	1:BK:115:LEU:HG	2.19	0.43
1:BM:39:VAL:HG12	1:BM:42:ASP:H	1.82	0.43
1:BM:92:ILE:HG12	1:DI:92:ILE:HG12	2.00	0.43
1:BM:122:ASP:OD1	1:DI:109:ARG:NH1	2.40	0.43
1:BN:53:PHE:CD1	1:BN:72:VAL:HG22	2.54	0.43
1:BR:67:LYS:HD3	1:BR:95:GLU:HB3	2.00	0.43
1:BT:111:ILE:O	1:BT:115:LEU:HG	2.19	0.43
1:BZ:111:ILE:O	1:BZ:115:LEU:HG	2.19	0.43
1:CB:7:ILE:HD12	1:CB:111:ILE:HG13	2.01	0.43
1:CH:7:ILE:HD12	1:CH:111:ILE:HG13	2.01	0.43
1:CX:108:LEU:HA	1:CX:111:ILE:HG22	2.00	0.43
1:DG:111:ILE:O	1:DG:115:LEU:HG	2.19	0.43
1:DJ:111:ILE:O	1:DJ:115:LEU:HG	2.19	0.43
1:AD:111:ILE:O	1:AD:115:LEU:HG	2.19	0.43
1:AH:67:LYS:HD3	1:AH:95:GLU:HB3	2.00	0.43
1:AH:68:GLY:O	1:AH:95:GLU:HA	2.18	0.43
1:AJ:111:ILE:O	1:AJ:115:LEU:HG	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AQ:67:LYS:HD3	1:AQ:95:GLU:HB3	2.00	0.43
1:AT:58:PRO:HB3	1:AT:66:SER:O	2.18	0.43
1:AV:35:LEU:HD22	1:AW:131:ALA:H	1.83	0.43
1:AW:51:ALA:HA	1:AW:73:VAL:O	2.19	0.43
1:BA:114:GLN:NE2	1:CW:16:ILE:HG21	2.34	0.43
1:BO:67:LYS:HD3	1:BO:95:GLU:HB3	2.00	0.43
1:BT:53:PHE:CD1	1:BT:72:VAL:HG22	2.54	0.43
1:BV:7:ILE:HD12	1:BV:111:ILE:HG13	2.01	0.43
1:BZ:53:PHE:CD1	1:BZ:72:VAL:HG22	2.54	0.43
1:CV:58:PRO:HB3	1:CV:66:SER:O	2.18	0.43
1:CW:7:ILE:HD12	1:CW:111:ILE:HG13	2.01	0.43
1:CX:53:PHE:CD1	1:CX:72:VAL:HG22	2.54	0.43
1:DK:68:GLY:O	1:DK:95:GLU:HA	2.18	0.43
1:AA:111:ILE:O	1:AA:115:LEU:HG	2.19	0.43
1:AB:58:PRO:HB3	1:AB:66:SER:O	2.18	0.43
1:AB:67:LYS:HD3	1:AB:95:GLU:HB3	2.00	0.43
1:AE:51:ALA:HA	1:AE:73:VAL:O	2.19	0.43
1:AE:58:PRO:HB3	1:AE:66:SER:O	2.18	0.43
1:AN:67:LYS:HD3	1:AN:95:GLU:HB3	2.00	0.43
1:AP:35:LEU:HD22	1:AQ:131:ALA:H	1.83	0.43
1:AS:53:PHE:CD1	1:AS:72:VAL:HG22	2.54	0.43
1:AU:19:GLY:H	1:CN:11:ALA:HA	1.84	0.43
1:AY:108:LEU:HA	1:AY:111:ILE:HG22	2.00	0.43
1:AZ:58:PRO:HB3	1:AZ:66:SER:O	2.18	0.43
1:BA:19:GLY:O	1:CW:8:LYS:HD2	2.19	0.43
1:BB:26:LEU:HD12	1:BB:36:ASN:O	2.18	0.43
1:BD:14:ALA:HB2	1:CZ:16:ILE:HD12	2.01	0.43
1:BE:35:LEU:HD22	1:BF:131:ALA:H	1.83	0.43
1:BG:7:ILE:HD12	1:BG:111:ILE:HG13	2.01	0.43
1:BO:68:GLY:O	1:BO:95:GLU:HA	2.18	0.43
1:BQ:53:PHE:CD1	1:BQ:72:VAL:HG22	2.54	0.43
1:CD:68:GLY:O	1:CD:95:GLU:HA	2.18	0.43
1:CL:53:PHE:CD1	1:CL:72:VAL:HG22	2.54	0.43
1:CQ:7:ILE:HD12	1:CQ:111:ILE:HG13	2.01	0.43
1:CR:111:ILE:O	1:CR:115:LEU:HG	2.19	0.43
1:CU:53:PHE:CD1	1:CU:72:VAL:HG22	2.54	0.43
1:DA:35:LEU:HD22	1:DB:131:ALA:H	1.83	0.43
1:DA:111:ILE:O	1:DA:115:LEU:HG	2.19	0.43
1:DE:67:LYS:HD3	1:DE:95:GLU:HB3	2.00	0.43
1:AA:53:PHE:CD1	1:AA:72:VAL:HG22	2.54	0.43
1:AD:53:PHE:CD1	1:AD:72:VAL:HG22	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AH:45:TYR:HD2	1:CC:98:PRO:HB3	1.84	0.43
1:AI:100:LEU:HD13	1:BY:88:ASN:ND2	2.34	0.43
1:AM:26:LEU:O	1:CE:129:ALA:HB2	2.19	0.43
1:AM:28:GLY:HA3	1:CE:128:GLY:O	2.19	0.43
1:AR:53:PHE:CD1	1:AR:72:VAL:HG22	2.54	0.43
1:BF:48:ARG:HH22	1:BR:127:LEU:HD23	1.83	0.43
1:BH:53:PHE:CD1	1:BH:72:VAL:HG22	2.54	0.43
1:BK:26:LEU:HB3	1:DI:129:ALA:HB1	2.01	0.43
1:BL:51:ALA:HA	1:BL:73:VAL:O	2.19	0.43
1:BN:35:LEU:HD22	1:BO:131:ALA:H	1.83	0.43
1:BP:7:ILE:HD12	1:BP:111:ILE:HG13	2.01	0.43
1:BX:51:ALA:HA	1:BX:73:VAL:O	2.19	0.43
1:CA:51:ALA:HA	1:CA:73:VAL:O	2.19	0.43
1:CC:35:LEU:HD22	1:CD:131:ALA:H	1.83	0.43
1:CE:39:VAL:HG12	1:CE:42:ASP:H	1.82	0.43
1:CG:51:ALA:HA	1:CG:73:VAL:O	2.19	0.43
1:CJ:67:LYS:HD3	1:CJ:95:GLU:HB3	2.00	0.43
1:CK:53:PHE:CD1	1:CK:72:VAL:HG22	2.54	0.43
1:CY:67:LYS:HD3	1:CY:95:GLU:HB3	2.00	0.43
1:DG:35:LEU:HD22	1:DH:131:ALA:H	1.83	0.43
1:AE:67:LYS:HD3	1:AE:95:GLU:HB3	2.00	0.42
1:AF:7:ILE:O	1:BV:22:ILE:HG22	2.18	0.42
1:AG:53:PHE:CD1	1:AG:72:VAL:HG22	2.54	0.42
1:AI:7:ILE:HD12	1:AI:111:ILE:HG13	2.01	0.42
1:AI:53:PHE:CD1	1:AI:72:VAL:HG22	2.54	0.42
1:AJ:53:PHE:CD1	1:AJ:72:VAL:HG22	2.54	0.42
1:AL:8:LYS:HD2	1:CB:19:GLY:O	2.19	0.42
1:AP:111:ILE:O	1:AP:115:LEU:HG	2.19	0.42
1:BB:111:ILE:O	1:BB:115:LEU:HG	2.19	0.42
1:CD:67:LYS:HD3	1:CD:95:GLU:HB3	2.00	0.42
1:CH:53:PHE:CD1	1:CH:72:VAL:HG22	2.54	0.42
1:CI:35:LEU:HD22	1:CJ:131:ALA:H	1.83	0.42
1:CI:111:ILE:O	1:CI:115:LEU:HG	2.19	0.42
1:CU:108:LEU:HA	1:CU:111:ILE:HG22	2.00	0.42
1:CV:51:ALA:HA	1:CV:73:VAL:O	2.19	0.42
1:CZ:53:PHE:CD1	1:CZ:72:VAL:HG22	2.54	0.42
1:DJ:53:PHE:CD1	1:DJ:72:VAL:HG22	2.54	0.42
1:AC:7:ILE:HD12	1:AC:111:ILE:HG13	2.01	0.42
1:AF:53:PHE:CD1	1:AF:72:VAL:HG22	2.54	0.42
1:AK:48:ARG:HH22	1:BX:127:LEU:HD23	1.84	0.42
1:AK:51:ALA:HA	1:AK:73:VAL:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AN:51:ALA:HA	1:AN:73:VAL:O	2.19	0.42
1:AT:51:ALA:HA	1:AT:73:VAL:O	2.19	0.42
1:AV:111:ILE:O	1:AV:115:LEU:HG	2.19	0.42
1:BG:130:LEU:HD21	1:DC:55:SER:HB3	2.01	0.42
1:BJ:111:ILE:HD13	1:DF:74:PHE:CZ	2.54	0.42
1:BQ:111:ILE:O	1:BQ:115:LEU:HG	2.19	0.42
1:CC:53:PHE:CD1	1:CC:72:VAL:HG22	2.54	0.42
1:CM:67:LYS:HD3	1:CM:95:GLU:HB3	2.00	0.42
1:CT:46:ARG:HA	1:DB:65:TYR:HE2	1.84	0.42
1:DD:108:LEU:HA	1:DD:111:ILE:HG22	2.00	0.42
1:DF:7:ILE:HD12	1:DF:111:ILE:HG13	2.01	0.42
1:DF:39:VAL:HG12	1:DF:42:ASP:H	1.82	0.42
1:DK:67:LYS:HD3	1:DK:95:GLU:HB3	2.00	0.42
1:AC:53:PHE:CD1	1:AC:72:VAL:HG22	2.54	0.42
1:AH:51:ALA:HA	1:AH:73:VAL:O	2.19	0.42
1:AI:114:GLN:NE2	1:BY:16:ILE:HG21	2.34	0.42
1:AK:65:TYR:CE2	1:BY:46:ARG:HG2	2.54	0.42
1:AZ:51:ALA:HA	1:AZ:73:VAL:O	2.19	0.42
1:BA:7:ILE:HD12	1:BA:111:ILE:HG13	2.01	0.42
1:BD:53:PHE:CD1	1:BD:72:VAL:HG22	2.54	0.42
1:BG:53:PHE:CD1	1:BG:72:VAL:HG22	2.54	0.42
1:BP:53:PHE:CD1	1:BP:72:VAL:HG22	2.54	0.42
1:BW:53:PHE:CD1	1:BW:72:VAL:HG22	2.54	0.42
1:CM:51:ALA:HA	1:CM:73:VAL:O	2.19	0.42
1:CO:111:ILE:O	1:CO:115:LEU:HG	2.19	0.42
1:DD:53:PHE:CD1	1:DD:72:VAL:HG22	2.54	0.42
1:DI:53:PHE:CD1	1:DI:72:VAL:HG22	2.55	0.42
1:AC:125:TRP:HZ3	1:BP:70:ASN:O	2.03	0.42
1:BH:35:LEU:HD22	1:BI:131:ALA:H	1.83	0.42
1:CI:53:PHE:CD1	1:CI:72:VAL:HG22	2.54	0.42
1:CN:53:PHE:CD1	1:CN:72:VAL:HG22	2.54	0.42
1:DL:53:PHE:CD1	1:DL:72:VAL:HG22	2.54	0.42
1:AC:46:ARG:HG2	1:DH:65:TYR:CD2	2.54	0.42
1:AM:108:LEU:HA	1:AM:111:ILE:HG22	2.00	0.42
1:BN:111:ILE:O	1:BN:115:LEU:HG	2.19	0.42
1:CF:53:PHE:CD1	1:CF:72:VAL:HG22	2.54	0.42
1:CJ:51:ALA:HA	1:CJ:73:VAL:O	2.19	0.42
1:CL:111:ILE:O	1:CL:115:LEU:HG	2.19	0.42
1:CS:74:PHE:HB3	1:CS:90:VAL:HG13	2.02	0.42
1:CW:53:PHE:CD1	1:CW:72:VAL:HG22	2.55	0.42
1:DK:51:ALA:HA	1:DK:73:VAL:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AB:74:PHE:HB3	1:AB:90:VAL:HG13	2.02	0.42
1:AR:7:ILE:HD12	1:AR:111:ILE:HG13	2.01	0.42
1:AU:16:ILE:HD12	1:CN:14:ALA:HB2	2.01	0.42
1:AY:35:LEU:HD22	1:AZ:131:ALA:H	1.83	0.42
1:BI:51:ALA:HA	1:BI:73:VAL:O	2.19	0.42
1:BI:74:PHE:HB3	1:BI:90:VAL:HG13	2.02	0.42
1:BR:51:ALA:HA	1:BR:73:VAL:O	2.19	0.42
1:BU:33:ASN:OD1	1:DI:30:THR:OG1	2.36	0.42
1:BW:111:ILE:O	1:BW:115:LEU:HG	2.19	0.42
1:CR:53:PHE:CD1	1:CR:72:VAL:HG22	2.54	0.42
1:DF:53:PHE:CD1	1:DF:72:VAL:HG22	2.54	0.42
1:DH:51:ALA:HA	1:DH:73:VAL:O	2.19	0.42
1:AF:7:ILE:HD12	1:AF:111:ILE:HG13	2.01	0.42
1:AI:16:ILE:HA	1:BY:14:ALA:HA	2.02	0.42
1:AQ:51:ALA:HA	1:AQ:73:VAL:O	2.19	0.42
1:AT:74:PHE:HB3	1:AT:90:VAL:HG13	2.02	0.42
1:AU:114:GLN:NE2	1:CN:16:ILE:HD13	2.23	0.42
1:AX:87:PHE:O	1:CT:67:LYS:NZ	2.29	0.42
1:BA:53:PHE:CD1	1:BA:72:VAL:HG22	2.54	0.42
1:BE:130:LEU:O	1:BS:26:LEU:HB3	2.19	0.42
1:BF:51:ALA:HA	1:BF:73:VAL:O	2.19	0.42
1:BS:53:PHE:CD1	1:BS:72:VAL:HG22	2.54	0.42
1:BU:51:ALA:HA	1:BU:73:VAL:O	2.19	0.42
1:CD:51:ALA:HA	1:CD:73:VAL:O	2.19	0.42
1:CT:53:PHE:CD1	1:CT:72:VAL:HG22	2.54	0.42
1:CX:111:ILE:O	1:CX:115:LEU:HG	2.19	0.42
1:AA:98:PRO:HB3	1:DH:45:TYR:HD2	1.83	0.42
1:AB:46:ARG:HA	1:AS:65:TYR:CE2	2.54	0.42
1:AI:16:ILE:HG21	1:BY:114:GLN:NE2	2.35	0.42
1:AL:7:ILE:HD12	1:AL:111:ILE:HG13	2.01	0.42
1:AL:109:ARG:NH2	1:CB:122:ASP:OD1	2.49	0.42
1:AO:53:PHE:CD1	1:AO:72:VAL:HG22	2.55	0.42
1:AS:111:ILE:O	1:AS:115:LEU:HG	2.19	0.42
1:AX:7:ILE:HD12	1:AX:111:ILE:HG13	2.01	0.42
1:AX:16:ILE:HG21	1:CT:114:GLN:CD	2.44	0.42
1:BM:7:ILE:HD12	1:BM:111:ILE:HG13	2.01	0.42
1:BM:53:PHE:CD1	1:BM:72:VAL:HG22	2.54	0.42
1:BO:51:ALA:HA	1:BO:73:VAL:O	2.19	0.42
1:BX:74:PHE:HB3	1:BX:90:VAL:HG13	2.02	0.42
1:CF:111:ILE:O	1:CF:115:LEU:HG	2.19	0.42
1:CY:74:PHE:HB3	1:CY:90:VAL:HG13	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:DB:74:PHE:HB3	1:DB:90:VAL:HG13	2.02	0.42
1:AL:53:PHE:CD1	1:AL:72:VAL:HG22	2.54	0.42
1:AU:53:PHE:CD1	1:AU:72:VAL:HG22	2.54	0.42
1:BJ:53:PHE:CD1	1:BJ:72:VAL:HG22	2.54	0.42
1:CO:53:PHE:CD1	1:CO:72:VAL:HG22	2.54	0.42
1:CP:51:ALA:HA	1:CP:73:VAL:O	2.19	0.42
1:CU:87:PHE:O	1:CV:67:LYS:NZ	2.41	0.42
1:CZ:7:ILE:HD12	1:CZ:111:ILE:HG13	2.01	0.42
1:DB:51:ALA:HA	1:DB:73:VAL:O	2.19	0.42
1:AE:74:PHE:HB3	1:AE:90:VAL:HG13	2.02	0.42
1:AF:20:SER:O	1:BV:8:LYS:HA	2.20	0.42
1:AG:111:ILE:O	1:AG:115:LEU:HG	2.19	0.42
1:BO:74:PHE:HB3	1:BO:90:VAL:HG13	2.02	0.42
1:CE:53:PHE:CD1	1:CE:72:VAL:HG22	2.54	0.42
1:CQ:53:PHE:CD1	1:CQ:72:VAL:HG22	2.54	0.42
1:DA:87:PHE:O	1:DB:67:LYS:NZ	2.41	0.42
1:DE:51:ALA:HA	1:DE:73:VAL:O	2.19	0.42
1:AR:130:LEU:O	1:CK:35:LEU:HD22	2.20	0.41
1:AW:74:PHE:HB3	1:AW:90:VAL:HG13	2.02	0.41
1:BC:74:PHE:HB3	1:BC:90:VAL:HG13	2.02	0.41
1:CS:51:ALA:HA	1:CS:73:VAL:O	2.19	0.41
1:CY:51:ALA:HA	1:CY:73:VAL:O	2.19	0.41
1:DE:74:PHE:HB3	1:DE:90:VAL:HG13	2.02	0.41
1:AO:7:ILE:HD12	1:AO:111:ILE:HG13	2.01	0.41
1:BY:7:ILE:HD12	1:BY:111:ILE:HG13	2.01	0.41
1:CP:74:PHE:HB3	1:CP:90:VAL:HG13	2.02	0.41
1:DH:74:PHE:HB3	1:DH:90:VAL:HG13	2.02	0.41
1:DK:74:PHE:HB3	1:DK:90:VAL:HG13	2.02	0.41
1:AB:51:ALA:HA	1:AB:73:VAL:O	2.19	0.41
1:AC:19:GLY:O	1:BP:8:LYS:HD2	2.20	0.41
1:AI:74:PHE:CZ	1:BY:111:ILE:HD13	2.55	0.41
1:AU:111:ILE:HD13	1:CN:74:PHE:CZ	2.55	0.41
1:BY:53:PHE:CD1	1:BY:72:VAL:HG22	2.54	0.41
1:DC:53:PHE:CD1	1:DC:72:VAL:HG22	2.54	0.41
1:DL:7:ILE:HD12	1:DL:111:ILE:HG13	2.01	0.41
1:AI:109:ARG:NE	1:BY:116:LEU:O	2.50	0.41
1:AJ:18:GLY:HA2	1:AK:12:THR:H	1.86	0.41
1:AM:18:GLY:HA2	1:AN:12:THR:H	1.86	0.41
1:AU:7:ILE:HD12	1:AU:111:ILE:HG13	2.01	0.41
1:AY:18:GLY:HA2	1:AZ:12:THR:H	1.86	0.41
1:BM:109:ARG:NE	1:DI:116:LEU:O	2.52	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BV:53:PHE:CD1	1:BV:72:VAL:HG22	2.54	0.41
1:CD:74:PHE:HB3	1:CD:90:VAL:HG13	2.02	0.41
1:CI:18:GLY:HA2	1:CJ:12:THR:H	1.86	0.41
1:DC:7:ILE:HD12	1:DC:111:ILE:HG13	2.01	0.41
1:AL:46:ARG:HG3	1:CV:98:PRO:HB2	2.03	0.41
1:AU:16:ILE:HA	1:CN:14:ALA:HA	2.02	0.41
1:AX:14:ALA:HB3	1:CT:117:THR:HB	2.01	0.41
1:AX:53:PHE:CD1	1:AX:72:VAL:HG22	2.54	0.41
1:BH:18:GLY:HA2	1:BI:12:THR:H	1.86	0.41
1:BK:18:GLY:HA2	1:BL:12:THR:H	1.86	0.41
1:BL:74:PHE:HB3	1:BL:90:VAL:HG13	2.02	0.41
1:CX:18:GLY:HA2	1:CY:12:THR:H	1.86	0.41
1:AF:95:GLU:HB2	1:BV:89:SER:HB2	2.02	0.41
1:AG:57:VAL:HG21	1:CS:29:GLN:HG3	2.03	0.41
1:AH:74:PHE:HB3	1:AH:90:VAL:HG13	2.02	0.41
1:AJ:65:TYR:CD2	1:CV:46:ARG:HG3	2.55	0.41
1:AK:74:PHE:HB3	1:AK:90:VAL:HG13	2.02	0.41
1:CF:18:GLY:HA2	1:CG:12:THR:H	1.86	0.41
1:AC:46:ARG:HA	1:DH:65:TYR:CE2	2.49	0.41
1:AF:111:ILE:HD13	1:BV:74:PHE:CZ	2.56	0.41
1:AI:46:ARG:HG2	1:CS:65:TYR:HD2	1.84	0.41
1:AR:122:ASP:OD1	1:CK:109:ARG:NH2	2.52	0.41
1:AX:122:ASP:OD1	1:CT:109:ARG:NH2	2.49	0.41
1:BC:51:ALA:HA	1:BC:73:VAL:O	2.19	0.41
1:CB:53:PHE:CD1	1:CB:72:VAL:HG22	2.54	0.41
1:CU:18:GLY:HA2	1:CV:12:THR:H	1.86	0.41
1:DA:18:GLY:HA2	1:DB:12:THR:H	1.86	0.41
1:DG:18:GLY:HA2	1:DH:12:THR:H	1.86	0.41
1:AC:12:THR:OG1	1:BP:17:THR:O	2.36	0.41
1:AE:127:LEU:HD23	1:DE:48:ARG:HH22	1.85	0.41
1:AM:17:THR:O	1:AN:12:THR:OG1	2.37	0.41
1:AM:87:PHE:O	1:AN:67:LYS:NZ	2.41	0.41
1:AP:18:GLY:HA2	1:AQ:12:THR:H	1.86	0.41
1:BQ:16:ILE:CD1	1:BR:114:GLN:HE21	2.26	0.41
1:CM:74:PHE:HB3	1:CM:90:VAL:HG13	2.02	0.41
1:CO:17:THR:O	1:CP:12:THR:OG1	2.37	0.41
1:CO:18:GLY:HA2	1:CP:12:THR:H	1.86	0.41
1:DJ:18:GLY:HA2	1:DK:12:THR:H	1.86	0.41
1:AG:18:GLY:HA2	1:AH:12:THR:H	1.86	0.41
1:AO:122:ASP:OD1	1:CE:109:ARG:NH1	2.45	0.41
1:AQ:74:PHE:HB3	1:AQ:90:VAL:HG13	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BD:19:GLY:O	1:CZ:8:LYS:HD2	2.21	0.41
1:BG:16:ILE:HD13	1:DC:114:GLN:NE2	2.22	0.41
1:BM:12:THR:H	1:DI:18:GLY:HA2	1.86	0.41
1:BQ:18:GLY:HA2	1:BR:12:THR:H	1.86	0.41
1:BR:74:PHE:HB3	1:BR:90:VAL:HG13	2.02	0.41
1:BY:22:ILE:HA	1:BY:41:GLU:CD	2.46	0.41
1:CA:74:PHE:HB3	1:CA:90:VAL:HG13	2.02	0.41
1:CC:18:GLY:HA2	1:CD:12:THR:H	1.86	0.41
1:CG:74:PHE:HB3	1:CG:90:VAL:HG13	2.02	0.41
1:CL:130:LEU:O	1:CZ:26:LEU:HB3	2.20	0.41
1:CV:74:PHE:HB3	1:CV:90:VAL:HG13	2.02	0.41
1:DD:18:GLY:HA2	1:DE:12:THR:H	1.86	0.41
1:AC:6:SER:HA	1:BP:22:ILE:O	2.21	0.41
1:AD:18:GLY:HA2	1:AE:12:THR:H	1.86	0.41
1:AI:46:ARG:HA	1:CS:65:TYR:HE2	1.84	0.41
1:AS:18:GLY:HA2	1:AT:12:THR:H	1.86	0.41
1:AX:109:ARG:HD2	1:CT:125:TRP:HB3	2.03	0.41
1:BF:74:PHE:HB3	1:BF:90:VAL:HG13	2.02	0.41
1:BS:22:ILE:HA	1:BS:41:GLU:CD	2.46	0.41
1:CA:65:TYR:HE2	1:CH:46:ARG:CA	2.33	0.41
1:DC:22:ILE:HA	1:DC:41:GLU:CD	2.47	0.41
1:DF:22:ILE:HA	1:DF:41:GLU:CD	2.46	0.41
1:AC:22:ILE:HA	1:AC:41:GLU:CD	2.46	0.40
1:AC:116:LEU:HG	1:BP:92:ILE:HG21	2.03	0.40
1:AL:22:ILE:HA	1:AL:41:GLU:CD	2.46	0.40
1:AL:26:LEU:HB3	1:CU:130:LEU:O	2.22	0.40
1:AL:46:ARG:HG2	1:CV:65:TYR:HD2	1.84	0.40
1:AR:55:SER:HB3	1:CK:130:LEU:HD21	2.01	0.40
1:AR:87:PHE:O	1:CK:67:LYS:NZ	2.40	0.40
1:AZ:74:PHE:HB3	1:AZ:90:VAL:HG13	2.02	0.40
1:BM:22:ILE:HA	1:BM:41:GLU:CD	2.46	0.40
1:CH:22:ILE:HA	1:CH:41:GLU:CD	2.46	0.40
1:CS:110:LEU:O	1:CS:114:GLN:HG3	2.22	0.40
1:CZ:22:ILE:HA	1:CZ:41:GLU:CD	2.46	0.40
1:AN:74:PHE:HB3	1:AN:90:VAL:HG13	2.02	0.40
1:AR:122:ASP:OD1	1:CK:109:ARG:NH1	2.52	0.40
1:AU:74:PHE:CZ	1:CN:111:ILE:HD13	2.56	0.40
1:AX:22:ILE:HA	1:AX:41:GLU:CD	2.46	0.40
1:BO:110:LEU:O	1:BO:114:GLN:HG3	2.22	0.40
1:CJ:74:PHE:HB3	1:CJ:90:VAL:HG13	2.02	0.40
1:CM:110:LEU:O	1:CM:114:GLN:HG3	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CT:22:ILE:HA	1:CT:41:GLU:CD	2.47	0.40
1:AA:18:GLY:HA2	1:AB:12:THR:H	1.86	0.40
1:AK:26:LEU:O	1:BX:129:ALA:HB2	2.21	0.40
1:AR:22:ILE:HA	1:AR:41:GLU:CD	2.46	0.40
1:AT:48:ARG:HH22	1:AW:127:LEU:HD23	1.85	0.40
1:BA:22:ILE:HA	1:BA:41:GLU:CD	2.46	0.40
1:BB:18:GLY:HA2	1:BC:12:THR:H	1.86	0.40
1:BN:18:GLY:HA2	1:BO:12:THR:H	1.86	0.40
1:BX:110:LEU:O	1:BX:114:GLN:HG3	2.22	0.40
1:CI:11:ALA:HA	1:CJ:18:GLY:HA2	2.04	0.40
1:CU:11:ALA:HA	1:CV:18:GLY:HA2	2.04	0.40
1:CY:110:LEU:O	1:CY:114:GLN:HG3	2.22	0.40
1:DG:11:ALA:HA	1:DH:18:GLY:HA2	2.04	0.40
1:AF:92:ILE:HA	1:BV:91:ARG:O	2.22	0.40
1:AO:129:ALA:HB2	1:CC:26:LEU:O	2.21	0.40
1:AT:85:THR:N	1:AV:63:GLY:O	2.47	0.40
1:AV:18:GLY:HA2	1:AW:12:THR:H	1.86	0.40
1:BA:26:LEU:HD13	1:BA:35:LEU:HD11	2.04	0.40
1:BC:65:TYR:HE2	1:BP:46:ARG:CA	2.32	0.40
1:BD:22:ILE:HA	1:BD:41:GLU:CD	2.47	0.40
1:BG:22:ILE:HA	1:BG:41:GLU:CD	2.46	0.40
1:BJ:22:ILE:HA	1:BJ:41:GLU:CD	2.46	0.40
1:BL:110:LEU:O	1:BL:114:GLN:HG3	2.22	0.40
1:BR:42:ASP:CB	1:BR:48:ARG:HG2	2.50	0.40
1:BU:110:LEU:O	1:BU:114:GLN:HG3	2.22	0.40
1:BW:11:ALA:HA	1:BX:18:GLY:HA2	2.04	0.40
1:CB:22:ILE:HA	1:CB:41:GLU:CD	2.46	0.40
1:CK:22:ILE:HA	1:CK:41:GLU:CD	2.46	0.40
1:CM:8:LYS:HB2	1:CM:114:GLN:NE2	2.32	0.40
1:CN:22:ILE:HA	1:CN:41:GLU:CD	2.46	0.40
1:CR:18:GLY:HA2	1:CS:12:THR:H	1.86	0.40
1:CX:31:VAL:HG13	1:CX:54:LYS:HE2	2.04	0.40
1:DA:31:VAL:HG13	1:DA:54:LYS:HE2	2.04	0.40
1:DE:110:LEU:O	1:DE:114:GLN:HG3	2.22	0.40
1:DG:31:VAL:HG13	1:DG:54:LYS:HE2	2.04	0.40
1:DL:22:ILE:HA	1:DL:41:GLU:CD	2.46	0.40
1:AJ:31:VAL:HG13	1:AJ:54:LYS:HE2	2.04	0.40
1:AU:22:ILE:HA	1:AU:41:GLU:CD	2.47	0.40
1:AY:71:GLU:HG2	1:AY:93:ALA:HB2	2.04	0.40
1:BA:122:ASP:OD1	1:CW:109:ARG:NH2	2.53	0.40
1:BB:71:GLU:HG2	1:BB:93:ALA:HB2	2.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BC:42:ASP:CB	1:BC:48:ARG:HG2	2.50	0.40
1:BE:18:GLY:HA2	1:BF:12:THR:H	1.86	0.40
1:BK:17:THR:O	1:BL:12:THR:OG1	2.37	0.40
1:BR:110:LEU:O	1:BR:114:GLN:HG3	2.22	0.40
1:BT:18:GLY:HA2	1:BU:12:THR:H	1.86	0.40
1:BZ:18:GLY:HA2	1:CA:12:THR:H	1.86	0.40
1:CC:11:ALA:HA	1:CD:18:GLY:HA2	2.04	0.40
1:CC:31:VAL:HG13	1:CC:54:LYS:HE2	2.04	0.40
1:CF:87:PHE:O	1:CG:67:LYS:NZ	2.41	0.40
1:CI:31:VAL:HG13	1:CI:54:LYS:HE2	2.04	0.40
1:CL:11:ALA:HA	1:CM:18:GLY:HA2	2.04	0.40
1:CL:18:GLY:HA2	1:CM:12:THR:H	1.86	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BS:17:THR:O	1:CH:12:THR:OG1[2_555]	2.03	0.17
1:AR:30:THR:OG1	1:CG:33:ASN:OD1[2_555]	2.04	0.16
1:AP:33:ASN:OD1	1:CG:30:THR:OG1[2_555]	2.13	0.07

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	AA	129/131 (98%)	127 (98%)	2 (2%)	0	100	100
1	AB	129/131 (98%)	128 (99%)	1 (1%)	0	100	100
1	AC	129/131 (98%)	129 (100%)	0	0	100	100
1	AD	129/131 (98%)	127 (98%)	2 (2%)	0	100	100
1	AE	129/131 (98%)	128 (99%)	1 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	AF	129/131 (98%)	129 (100%)	0	0	100	100
1	AG	129/131 (98%)	127 (98%)	2 (2%)	0	100	100
1	AH	129/131 (98%)	128 (99%)	1 (1%)	0	100	100
1	AI	129/131 (98%)	129 (100%)	0	0	100	100
1	AJ	129/131 (98%)	127 (98%)	2 (2%)	0	100	100
1	AK	129/131 (98%)	128 (99%)	1 (1%)	0	100	100
1	AL	129/131 (98%)	129 (100%)	0	0	100	100
1	AM	129/131 (98%)	127 (98%)	2 (2%)	0	100	100
1	AN	129/131 (98%)	128 (99%)	1 (1%)	0	100	100
1	AO	129/131 (98%)	129 (100%)	0	0	100	100
1	AP	129/131 (98%)	127 (98%)	2 (2%)	0	100	100
1	AQ	129/131 (98%)	128 (99%)	1 (1%)	0	100	100
1	AR	129/131 (98%)	129 (100%)	0	0	100	100
1	AS	129/131 (98%)	127 (98%)	2 (2%)	0	100	100
1	AT	129/131 (98%)	128 (99%)	1 (1%)	0	100	100
1	AU	129/131 (98%)	129 (100%)	0	0	100	100
1	AV	129/131 (98%)	127 (98%)	2 (2%)	0	100	100
1	AW	129/131 (98%)	128 (99%)	1 (1%)	0	100	100
1	AX	129/131 (98%)	129 (100%)	0	0	100	100
1	AY	129/131 (98%)	127 (98%)	2 (2%)	0	100	100
1	AZ	129/131 (98%)	128 (99%)	1 (1%)	0	100	100
1	BA	129/131 (98%)	129 (100%)	0	0	100	100
1	BB	129/131 (98%)	127 (98%)	2 (2%)	0	100	100
1	BC	129/131 (98%)	128 (99%)	1 (1%)	0	100	100
1	BD	129/131 (98%)	129 (100%)	0	0	100	100
1	BE	129/131 (98%)	127 (98%)	2 (2%)	0	100	100
1	BF	129/131 (98%)	128 (99%)	1 (1%)	0	100	100
1	BG	129/131 (98%)	129 (100%)	0	0	100	100
1	BH	129/131 (98%)	127 (98%)	2 (2%)	0	100	100
1	BI	129/131 (98%)	128 (99%)	1 (1%)	0	100	100
1	BJ	129/131 (98%)	129 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	BK	129/131 (98%)	127 (98%)	2 (2%)	0	100	100
1	BL	129/131 (98%)	128 (99%)	1 (1%)	0	100	100
1	BM	129/131 (98%)	129 (100%)	0	0	100	100
1	BN	129/131 (98%)	127 (98%)	2 (2%)	0	100	100
1	BO	129/131 (98%)	128 (99%)	1 (1%)	0	100	100
1	BP	129/131 (98%)	129 (100%)	0	0	100	100
1	BQ	129/131 (98%)	127 (98%)	2 (2%)	0	100	100
1	BR	129/131 (98%)	128 (99%)	1 (1%)	0	100	100
1	BS	129/131 (98%)	129 (100%)	0	0	100	100
1	BT	129/131 (98%)	127 (98%)	2 (2%)	0	100	100
1	BU	129/131 (98%)	128 (99%)	1 (1%)	0	100	100
1	BV	129/131 (98%)	129 (100%)	0	0	100	100
1	BW	129/131 (98%)	127 (98%)	2 (2%)	0	100	100
1	BX	129/131 (98%)	128 (99%)	1 (1%)	0	100	100
1	BY	129/131 (98%)	129 (100%)	0	0	100	100
1	BZ	129/131 (98%)	127 (98%)	2 (2%)	0	100	100
1	CA	129/131 (98%)	128 (99%)	1 (1%)	0	100	100
1	CB	129/131 (98%)	129 (100%)	0	0	100	100
1	CC	129/131 (98%)	127 (98%)	2 (2%)	0	100	100
1	CD	129/131 (98%)	128 (99%)	1 (1%)	0	100	100
1	CE	129/131 (98%)	129 (100%)	0	0	100	100
1	CF	129/131 (98%)	127 (98%)	2 (2%)	0	100	100
1	CG	129/131 (98%)	128 (99%)	1 (1%)	0	100	100
1	CH	129/131 (98%)	129 (100%)	0	0	100	100
1	CI	129/131 (98%)	127 (98%)	2 (2%)	0	100	100
1	CJ	129/131 (98%)	128 (99%)	1 (1%)	0	100	100
1	CK	129/131 (98%)	129 (100%)	0	0	100	100
1	CL	129/131 (98%)	127 (98%)	2 (2%)	0	100	100
1	CM	129/131 (98%)	128 (99%)	1 (1%)	0	100	100
1	CN	129/131 (98%)	129 (100%)	0	0	100	100
1	CO	129/131 (98%)	127 (98%)	2 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	CP	129/131 (98%)	128 (99%)	1 (1%)	0	100	100
1	CQ	129/131 (98%)	129 (100%)	0	0	100	100
1	CR	129/131 (98%)	127 (98%)	2 (2%)	0	100	100
1	CS	129/131 (98%)	128 (99%)	1 (1%)	0	100	100
1	CT	129/131 (98%)	129 (100%)	0	0	100	100
1	CU	129/131 (98%)	127 (98%)	2 (2%)	0	100	100
1	CV	129/131 (98%)	128 (99%)	1 (1%)	0	100	100
1	CW	129/131 (98%)	129 (100%)	0	0	100	100
1	CX	129/131 (98%)	127 (98%)	2 (2%)	0	100	100
1	CY	129/131 (98%)	128 (99%)	1 (1%)	0	100	100
1	CZ	129/131 (98%)	129 (100%)	0	0	100	100
1	DA	129/131 (98%)	127 (98%)	2 (2%)	0	100	100
1	DB	129/131 (98%)	128 (99%)	1 (1%)	0	100	100
1	DC	129/131 (98%)	129 (100%)	0	0	100	100
1	DD	129/131 (98%)	127 (98%)	2 (2%)	0	100	100
1	DE	129/131 (98%)	128 (99%)	1 (1%)	0	100	100
1	DF	129/131 (98%)	129 (100%)	0	0	100	100
1	DG	129/131 (98%)	127 (98%)	2 (2%)	0	100	100
1	DH	129/131 (98%)	128 (99%)	1 (1%)	0	100	100
1	DI	129/131 (98%)	129 (100%)	0	0	100	100
1	DJ	129/131 (98%)	127 (98%)	2 (2%)	0	100	100
1	DK	129/131 (98%)	128 (99%)	1 (1%)	0	100	100
1	DL	129/131 (98%)	129 (100%)	0	0	100	100
All	All	11610/11790 (98%)	11520 (99%)	90 (1%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	AA	108/108 (100%)	108 (100%)	0	100	100
1	AB	108/108 (100%)	108 (100%)	0	100	100
1	AC	108/108 (100%)	108 (100%)	0	100	100
1	AD	108/108 (100%)	108 (100%)	0	100	100
1	AE	108/108 (100%)	108 (100%)	0	100	100
1	AF	108/108 (100%)	108 (100%)	0	100	100
1	AG	108/108 (100%)	108 (100%)	0	100	100
1	AH	108/108 (100%)	108 (100%)	0	100	100
1	AI	108/108 (100%)	108 (100%)	0	100	100
1	AJ	108/108 (100%)	108 (100%)	0	100	100
1	AK	108/108 (100%)	108 (100%)	0	100	100
1	AL	108/108 (100%)	108 (100%)	0	100	100
1	AM	108/108 (100%)	108 (100%)	0	100	100
1	AN	108/108 (100%)	108 (100%)	0	100	100
1	AO	108/108 (100%)	108 (100%)	0	100	100
1	AP	108/108 (100%)	108 (100%)	0	100	100
1	AQ	108/108 (100%)	108 (100%)	0	100	100
1	AR	108/108 (100%)	108 (100%)	0	100	100
1	AS	108/108 (100%)	108 (100%)	0	100	100
1	AT	108/108 (100%)	108 (100%)	0	100	100
1	AU	108/108 (100%)	108 (100%)	0	100	100
1	AV	108/108 (100%)	108 (100%)	0	100	100
1	AW	108/108 (100%)	108 (100%)	0	100	100
1	AX	108/108 (100%)	108 (100%)	0	100	100
1	AY	108/108 (100%)	108 (100%)	0	100	100
1	AZ	108/108 (100%)	108 (100%)	0	100	100
1	BA	108/108 (100%)	108 (100%)	0	100	100
1	BB	108/108 (100%)	108 (100%)	0	100	100
1	BC	108/108 (100%)	108 (100%)	0	100	100
1	BD	108/108 (100%)	108 (100%)	0	100	100
1	BE	108/108 (100%)	108 (100%)	0	100	100
1	BF	108/108 (100%)	108 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	BG	108/108 (100%)	108 (100%)	0	100	100
1	BH	108/108 (100%)	108 (100%)	0	100	100
1	BI	108/108 (100%)	108 (100%)	0	100	100
1	BJ	108/108 (100%)	108 (100%)	0	100	100
1	BK	108/108 (100%)	108 (100%)	0	100	100
1	BL	108/108 (100%)	108 (100%)	0	100	100
1	BM	108/108 (100%)	108 (100%)	0	100	100
1	BN	108/108 (100%)	108 (100%)	0	100	100
1	BO	108/108 (100%)	108 (100%)	0	100	100
1	BP	108/108 (100%)	108 (100%)	0	100	100
1	BQ	108/108 (100%)	108 (100%)	0	100	100
1	BR	108/108 (100%)	108 (100%)	0	100	100
1	BS	108/108 (100%)	108 (100%)	0	100	100
1	BT	108/108 (100%)	108 (100%)	0	100	100
1	BU	108/108 (100%)	108 (100%)	0	100	100
1	BV	108/108 (100%)	108 (100%)	0	100	100
1	BW	108/108 (100%)	108 (100%)	0	100	100
1	BX	108/108 (100%)	108 (100%)	0	100	100
1	BY	108/108 (100%)	108 (100%)	0	100	100
1	BZ	108/108 (100%)	108 (100%)	0	100	100
1	CA	108/108 (100%)	108 (100%)	0	100	100
1	CB	108/108 (100%)	108 (100%)	0	100	100
1	CC	108/108 (100%)	108 (100%)	0	100	100
1	CD	108/108 (100%)	108 (100%)	0	100	100
1	CE	108/108 (100%)	108 (100%)	0	100	100
1	CF	108/108 (100%)	108 (100%)	0	100	100
1	CG	108/108 (100%)	108 (100%)	0	100	100
1	CH	108/108 (100%)	108 (100%)	0	100	100
1	CI	108/108 (100%)	108 (100%)	0	100	100
1	CJ	108/108 (100%)	108 (100%)	0	100	100
1	CK	108/108 (100%)	108 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	CL	108/108 (100%)	108 (100%)	0	100	100
1	CM	108/108 (100%)	108 (100%)	0	100	100
1	CN	108/108 (100%)	108 (100%)	0	100	100
1	CO	108/108 (100%)	108 (100%)	0	100	100
1	CP	108/108 (100%)	108 (100%)	0	100	100
1	CQ	108/108 (100%)	108 (100%)	0	100	100
1	CR	108/108 (100%)	108 (100%)	0	100	100
1	CS	108/108 (100%)	108 (100%)	0	100	100
1	CT	108/108 (100%)	108 (100%)	0	100	100
1	CU	108/108 (100%)	108 (100%)	0	100	100
1	CV	108/108 (100%)	108 (100%)	0	100	100
1	CW	108/108 (100%)	108 (100%)	0	100	100
1	CX	108/108 (100%)	108 (100%)	0	100	100
1	CY	108/108 (100%)	108 (100%)	0	100	100
1	CZ	108/108 (100%)	108 (100%)	0	100	100
1	DA	108/108 (100%)	108 (100%)	0	100	100
1	DB	108/108 (100%)	108 (100%)	0	100	100
1	DC	108/108 (100%)	108 (100%)	0	100	100
1	DD	108/108 (100%)	108 (100%)	0	100	100
1	DE	108/108 (100%)	108 (100%)	0	100	100
1	DF	108/108 (100%)	108 (100%)	0	100	100
1	DG	108/108 (100%)	108 (100%)	0	100	100
1	DH	108/108 (100%)	108 (100%)	0	100	100
1	DI	108/108 (100%)	108 (100%)	0	100	100
1	DJ	108/108 (100%)	108 (100%)	0	100	100
1	DK	108/108 (100%)	108 (100%)	0	100	100
1	DL	108/108 (100%)	108 (100%)	0	100	100
All	All	9720/9720 (100%)	9720 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (72) such sidechains are listed below:

Mol	Chain	Res	Type
1	AA	70	ASN
1	AB	29	GLN
1	AB	36	ASN
1	AE	29	GLN
1	AE	36	ASN
1	AG	70	ASN
1	AH	29	GLN
1	AH	36	ASN
1	AJ	70	ASN
1	AK	29	GLN
1	AK	36	ASN
1	AN	29	GLN
1	AN	36	ASN
1	AQ	29	GLN
1	AQ	36	ASN
1	AR	70	ASN
1	AS	70	ASN
1	AT	29	GLN
1	AT	36	ASN
1	AW	29	GLN
1	AW	36	ASN
1	AY	70	ASN
1	AZ	29	GLN
1	AZ	36	ASN
1	BC	29	GLN
1	BC	36	ASN
1	BE	70	ASN
1	BF	29	GLN
1	BF	36	ASN
1	BI	29	GLN
1	BI	36	ASN
1	BL	29	GLN
1	BL	36	ASN
1	BN	70	ASN
1	BO	29	GLN
1	BO	36	ASN
1	BR	29	GLN
1	BR	36	ASN
1	BU	29	GLN
1	BU	36	ASN
1	BV	70	ASN
1	BX	29	GLN
1	BX	36	ASN

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Mol	Chain	Res	Type
1	CA	29	GLN
1	CA	36	ASN
1	CD	29	GLN
1	CD	36	ASN
1	CG	29	GLN
1	CG	36	ASN
1	CJ	29	GLN
1	CJ	36	ASN
1	CM	29	GLN
1	CM	36	ASN
1	CP	29	GLN
1	CP	36	ASN
1	CS	29	GLN
1	CS	36	ASN
1	CT	70	ASN
1	CV	29	GLN
1	CV	36	ASN
1	CY	29	GLN
1	CY	36	ASN
1	DA	70	ASN
1	DB	29	GLN
1	DB	36	ASN
1	DE	29	GLN
1	DE	36	ASN
1	DF	70	ASN
1	DH	29	GLN
1	DH	36	ASN
1	DK	29	GLN
1	DK	36	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ > 2	OWAB(Å ²)	Q < 0.9
1	AA	131/131 (100%)	0.38	6 (4%) 37 21	64, 111, 139, 160	0
1	AB	131/131 (100%)	0.19	2 (1%) 72 48	64, 109, 138, 177	0
1	AC	131/131 (100%)	0.34	6 (4%) 37 21	64, 110, 142, 191	0
1	AD	131/131 (100%)	0.23	6 (4%) 37 21	64, 111, 139, 160	0
1	AE	131/131 (100%)	0.22	3 (2%) 61 37	64, 109, 138, 177	0
1	AF	131/131 (100%)	0.21	6 (4%) 37 21	64, 110, 142, 191	0
1	AG	131/131 (100%)	0.25	5 (3%) 44 25	64, 111, 139, 160	0
1	AH	131/131 (100%)	0.27	7 (5%) 32 19	64, 109, 138, 177	0
1	AI	131/131 (100%)	0.40	10 (7%) 20 13	64, 110, 142, 191	0
1	AJ	131/131 (100%)	0.13	6 (4%) 37 21	64, 111, 139, 160	0
1	AK	131/131 (100%)	0.00	2 (1%) 72 48	64, 109, 138, 177	0
1	AL	131/131 (100%)	0.03	3 (2%) 61 37	64, 110, 142, 191	0
1	AM	131/131 (100%)	0.24	4 (3%) 51 30	64, 111, 139, 160	0
1	AN	131/131 (100%)	0.09	2 (1%) 72 48	64, 109, 138, 177	0
1	AO	131/131 (100%)	0.41	6 (4%) 37 21	64, 110, 142, 191	0
1	AP	131/131 (100%)	0.04	6 (4%) 37 21	64, 111, 139, 160	0
1	AQ	131/131 (100%)	0.21	7 (5%) 32 19	64, 109, 138, 177	0
1	AR	131/131 (100%)	0.18	4 (3%) 51 30	64, 110, 142, 191	0
1	AS	131/131 (100%)	0.07	4 (3%) 51 30	64, 111, 139, 160	0
1	AT	131/131 (100%)	0.06	3 (2%) 61 37	64, 109, 138, 177	0
1	AU	131/131 (100%)	0.33	14 (10%) 11 8	64, 110, 142, 191	0
1	AV	131/131 (100%)	0.15	3 (2%) 61 37	64, 111, 139, 160	0
1	AW	131/131 (100%)	0.02	2 (1%) 72 48	64, 109, 138, 177	0
1	AX	131/131 (100%)	0.34	11 (8%) 17 12	64, 110, 142, 191	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å ²)	Q<0.9
1	AY	131/131 (100%)	0.10	4 (3%)	51	30	64, 111, 139, 160	0
1	AZ	131/131 (100%)	0.09	3 (2%)	61	37	64, 109, 138, 177	0
1	BA	131/131 (100%)	0.22	3 (2%)	61	37	64, 110, 142, 191	0
1	BB	131/131 (100%)	0.20	2 (1%)	72	48	64, 111, 139, 160	0
1	BC	131/131 (100%)	0.26	3 (2%)	61	37	64, 109, 138, 177	0
1	BD	131/131 (100%)	0.09	2 (1%)	72	48	64, 110, 142, 191	0
1	BE	131/131 (100%)	0.25	5 (3%)	44	25	64, 111, 139, 160	0
1	BF	131/131 (100%)	0.32	5 (3%)	44	25	64, 109, 138, 177	0
1	BG	131/131 (100%)	0.16	5 (3%)	44	25	64, 110, 142, 191	0
1	BH	131/131 (100%)	0.26	13 (9%)	13	9	64, 111, 139, 160	0
1	BI	131/131 (100%)	0.18	7 (5%)	32	19	64, 109, 138, 177	0
1	BJ	131/131 (100%)	0.14	3 (2%)	61	37	64, 110, 142, 191	0
1	BK	131/131 (100%)	0.18	8 (6%)	27	17	64, 111, 139, 160	0
1	BL	131/131 (100%)	0.05	0	100	100	64, 109, 138, 177	0
1	BM	131/131 (100%)	0.50	12 (9%)	14	10	64, 110, 142, 191	0
1	BN	131/131 (100%)	0.14	7 (5%)	32	19	64, 111, 139, 160	0
1	BO	131/131 (100%)	0.20	7 (5%)	32	19	64, 109, 138, 177	0
1	BP	131/131 (100%)	0.28	6 (4%)	37	21	64, 110, 142, 191	0
1	BQ	131/131 (100%)	0.25	7 (5%)	32	19	64, 111, 139, 160	0
1	BR	131/131 (100%)	0.15	5 (3%)	44	25	64, 109, 138, 177	0
1	BS	131/131 (100%)	0.21	3 (2%)	61	37	64, 110, 142, 191	0
1	BT	131/131 (100%)	0.11	1 (0%)	82	61	64, 111, 139, 160	0
1	BU	131/131 (100%)	0.14	6 (4%)	37	21	64, 109, 138, 177	0
1	BV	131/131 (100%)	0.19	6 (4%)	37	21	64, 110, 142, 191	0
1	BW	131/131 (100%)	0.22	4 (3%)	51	30	64, 111, 139, 160	0
1	BX	131/131 (100%)	0.05	2 (1%)	72	48	64, 109, 138, 177	0
1	BY	131/131 (100%)	0.34	6 (4%)	37	21	64, 110, 142, 191	0
1	BZ	131/131 (100%)	0.19	7 (5%)	32	19	64, 111, 139, 160	0
1	CA	131/131 (100%)	0.08	0	100	100	64, 109, 138, 177	0
1	CB	131/131 (100%)	0.06	3 (2%)	61	37	64, 110, 142, 191	0
1	CC	131/131 (100%)	0.14	3 (2%)	61	37	64, 111, 139, 160	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å ²)	Q<0.9
1	CD	131/131 (100%)	0.18	6 (4%)	37	21	64, 109, 138, 177	0
1	CE	131/131 (100%)	0.27	6 (4%)	37	21	64, 110, 142, 191	0
1	CF	131/131 (100%)	0.34	8 (6%)	27	17	64, 111, 139, 160	0
1	CG	131/131 (100%)	0.15	2 (1%)	72	48	64, 109, 138, 177	0
1	CH	131/131 (100%)	0.45	10 (7%)	20	13	64, 110, 142, 191	0
1	CI	131/131 (100%)	0.24	3 (2%)	61	37	64, 111, 139, 160	0
1	CJ	131/131 (100%)	0.13	5 (3%)	44	25	64, 109, 138, 177	0
1	CK	131/131 (100%)	0.24	9 (6%)	23	15	64, 110, 142, 191	0
1	CL	131/131 (100%)	0.09	3 (2%)	61	37	64, 111, 139, 160	0
1	CM	131/131 (100%)	-0.06	2 (1%)	72	48	64, 109, 138, 177	0
1	CN	131/131 (100%)	0.06	3 (2%)	61	37	64, 110, 142, 191	0
1	CO	131/131 (100%)	0.13	3 (2%)	61	37	64, 111, 139, 160	0
1	CP	131/131 (100%)	-0.04	0	100	100	64, 109, 138, 177	0
1	CQ	131/131 (100%)	0.26	2 (1%)	72	48	64, 110, 142, 191	0
1	CR	131/131 (100%)	0.19	7 (5%)	32	19	64, 111, 139, 160	0
1	CS	131/131 (100%)	0.05	5 (3%)	44	25	64, 109, 138, 177	0
1	CT	131/131 (100%)	0.46	11 (8%)	17	12	64, 110, 142, 191	0
1	CU	131/131 (100%)	-0.15	3 (2%)	61	37	64, 111, 139, 160	0
1	CV	131/131 (100%)	-0.00	3 (2%)	61	37	64, 109, 138, 177	0
1	CW	131/131 (100%)	0.26	3 (2%)	61	37	64, 110, 142, 191	0
1	CX	131/131 (100%)	0.04	5 (3%)	44	25	64, 111, 139, 160	0
1	CY	131/131 (100%)	-0.11	3 (2%)	61	37	64, 109, 138, 177	0
1	CZ	131/131 (100%)	-0.06	1 (0%)	82	61	64, 110, 142, 191	0
1	DA	131/131 (100%)	0.16	3 (2%)	61	37	64, 111, 139, 160	0
1	DB	131/131 (100%)	0.04	2 (1%)	72	48	64, 109, 138, 177	0
1	DC	131/131 (100%)	0.29	2 (1%)	72	48	64, 110, 142, 191	0
1	DD	131/131 (100%)	0.46	9 (6%)	23	15	64, 111, 139, 160	0
1	DE	131/131 (100%)	0.39	8 (6%)	27	17	64, 109, 138, 177	0
1	DF	131/131 (100%)	0.33	6 (4%)	37	21	64, 110, 142, 191	0
1	DG	131/131 (100%)	0.13	3 (2%)	61	37	64, 111, 139, 160	0
1	DH	131/131 (100%)	0.02	4 (3%)	51	30	64, 109, 138, 177	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	DI	131/131 (100%)	0.46	10 (7%) 20 13	64, 110, 142, 191	0
1	DJ	131/131 (100%)	0.02	0 100 100	64, 111, 139, 160	0
1	DK	131/131 (100%)	0.14	5 (3%) 44 25	64, 109, 138, 177	0
1	DL	131/131 (100%)	0.02	1 (0%) 82 61	64, 110, 142, 191	0
All	All	11790/11790 (100%)	0.18	432 (3%) 45 26	64, 110, 141, 191	0

All (432) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	BH	61	VAL	7.1
1	AU	54	LYS	6.5
1	DE	131	ALA	5.9
1	BM	131	ALA	5.7
1	BH	62	ASN	5.2
1	BI	1	SER	5.1
1	AQ	60	VAL	5.0
1	AP	62	ASN	4.9
1	CT	40	SER	4.9
1	CF	30	THR	4.9
1	BJ	130	LEU	4.8
1	BZ	32	THR	4.6
1	AH	131	ALA	4.5
1	AA	131	ALA	4.5
1	AX	92	ILE	4.5
1	BR	131	ALA	4.4
1	BQ	62	ASN	4.4
1	CE	131	ALA	4.4
1	BY	40	SER	4.4
1	AF	71	GLU	4.3
1	BN	62	ASN	4.3
1	BO	29	GLN	4.3
1	CH	130	LEU	4.3
1	BW	71	GLU	4.3
1	BH	65	TYR	4.2
1	CF	6	SER	4.2
1	AU	55	SER	4.1
1	CI	131	ALA	4.1
1	AY	32	THR	4.1
1	DE	126	THR	4.1
1	AA	60	VAL	4.0
1	BO	34	GLY	4.0

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Mol	Chain	Res	Type	RSRZ
1	CV	131	ALA	4.0
1	DD	65	TYR	4.0
1	DG	4	GLY	3.9
1	CY	93	ALA	3.9
1	AI	35	LEU	3.9
1	CE	27	THR	3.9
1	AS	113	ALA	3.9
1	BI	131	ALA	3.9
1	CR	71	GLU	3.8
1	AF	53	PHE	3.8
1	AO	33	ASN	3.8
1	AU	40	SER	3.8
1	CE	53	PHE	3.8
1	CD	131	ALA	3.7
1	AJ	61	VAL	3.7
1	AQ	131	ALA	3.7
1	DF	35	LEU	3.7
1	BA	113	ALA	3.6
1	BP	129	ALA	3.6
1	CF	54	LYS	3.6
1	AA	62	ASN	3.6
1	CC	62	ASN	3.5
1	CT	92	ILE	3.5
1	AO	53	PHE	3.5
1	BG	72	VAL	3.5
1	CH	105	VAL	3.5
1	BH	55	SER	3.5
1	DI	40	SER	3.5
1	BQ	3	ILE	3.5
1	AC	54	LYS	3.4
1	AU	33	ASN	3.4
1	CS	88	ASN	3.4
1	AH	1	SER	3.4
1	AK	131	ALA	3.4
1	AU	71	GLU	3.3
1	BR	1	SER	3.3
1	CM	131	ALA	3.3
1	BI	130	LEU	3.3
1	AU	53	PHE	3.3
1	AM	40	SER	3.3
1	BH	60	VAL	3.3
1	AZ	131	ALA	3.2

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Mol	Chain	Res	Type	RSRZ
1	AL	106	LYS	3.2
1	BM	105	VAL	3.2
1	BU	126	THR	3.2
1	AB	131	ALA	3.2
1	DK	93	ALA	3.2
1	CF	65	TYR	3.2
1	CE	61	VAL	3.2
1	DE	40	SER	3.2
1	AO	70	ASN	3.2
1	BN	64	ASN	3.2
1	DK	33	ASN	3.2
1	AY	131	ALA	3.2
1	AJ	34	GLY	3.1
1	AP	71	GLU	3.1
1	BY	131	ALA	3.1
1	DE	125	TRP	3.1
1	BN	61	VAL	3.1
1	AX	31	VAL	3.1
1	AC	89	SER	3.1
1	CF	40	SER	3.1
1	CT	90	VAL	3.1
1	CL	64	ASN	3.1
1	AM	131	ALA	3.1
1	AU	52	THR	3.1
1	BI	55	SER	3.1
1	CL	61	VAL	3.1
1	CO	55	SER	3.1
1	AD	88	ASN	3.0
1	BT	62	ASN	3.0
1	CK	88	ASN	3.0
1	AX	71	GLU	3.0
1	AD	77	PRO	3.0
1	CZ	93	ALA	3.0
1	AN	60	VAL	3.0
1	CN	30	THR	3.0
1	BH	53	PHE	3.0
1	BE	53	PHE	3.0
1	BZ	30	THR	3.0
1	DH	98	PRO	3.0
1	AI	40	SER	3.0
1	DB	93	ALA	3.0
1	AA	65	TYR	3.0

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Mol	Chain	Res	Type	RSRZ
1	DF	53	PHE	3.0
1	BZ	33	ASN	3.0
1	CR	70	ASN	3.0
1	AD	27	THR	3.0
1	BQ	85	THR	3.0
1	CH	106	LYS	2.9
1	AV	89	SER	2.9
1	DI	71	GLU	2.9
1	BP	18	GLY	2.9
1	AU	30	THR	2.9
1	BB	54	LYS	2.9
1	AO	105	VAL	2.9
1	AQ	1	SER	2.9
1	CH	74	PHE	2.9
1	BD	30	THR	2.9
1	AO	71	GLU	2.9
1	DD	131	ALA	2.9
1	BF	53	PHE	2.9
1	DH	24	PHE	2.9
1	CG	40	SER	2.9
1	DD	106	LYS	2.9
1	AT	93	ALA	2.8
1	CN	106	LYS	2.8
1	CR	53	PHE	2.8
1	BM	1	SER	2.8
1	DK	129	ALA	2.8
1	AP	61	VAL	2.8
1	AI	33	ASN	2.8
1	AY	4	GLY	2.8
1	CJ	35	LEU	2.8
1	CX	62	ASN	2.8
1	DD	70	ASN	2.8
1	CX	91	ARG	2.8
1	CT	72	VAL	2.8
1	AQ	124	PHE	2.8
1	AQ	4	GLY	2.8
1	CF	29	GLN	2.8
1	DF	106	LYS	2.8
1	DI	54	LYS	2.8
1	DF	130	LEU	2.8
1	BM	30	THR	2.8
1	BE	63	GLY	2.7

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Mol	Chain	Res	Type	RSRZ
1	AF	88	ASN	2.7
1	BH	64	ASN	2.7
1	BK	65	TYR	2.7
1	DL	106	LYS	2.7
1	AX	93	ALA	2.7
1	BK	62	ASN	2.7
1	CB	129	ALA	2.7
1	AQ	35	LEU	2.7
1	AC	53	PHE	2.7
1	AD	87	PHE	2.7
1	BO	87	PHE	2.7
1	CK	53	PHE	2.7
1	AP	131	ALA	2.7
1	AX	11	ALA	2.7
1	DF	131	ALA	2.7
1	CG	60	VAL	2.7
1	DB	55	SER	2.7
1	CF	131	ALA	2.7
1	AI	37	VAL	2.7
1	BE	62	ASN	2.7
1	CK	62	ASN	2.7
1	CD	63	GLY	2.6
1	AX	72	VAL	2.6
1	BW	130	LEU	2.6
1	CS	60	VAL	2.6
1	CV	93	ALA	2.6
1	DC	94	LEU	2.6
1	BF	29	GLN	2.6
1	BI	74	PHE	2.6
1	BM	29	GLN	2.6
1	AU	34	GLY	2.6
1	AU	35	LEU	2.6
1	BY	61	VAL	2.6
1	DF	37	VAL	2.6
1	CT	89	SER	2.6
1	DD	69	LYS	2.6
1	CF	53	PHE	2.6
1	AJ	127	LEU	2.6
1	BY	93	ALA	2.6
1	CT	105	VAL	2.6
1	CH	123	SER	2.6
1	BM	33	ASN	2.6

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Mol	Chain	Res	Type	RSRZ
1	AI	41	GLU	2.6
1	AD	61	VAL	2.6
1	BJ	131	ALA	2.6
1	BQ	131	ALA	2.6
1	CB	131	ALA	2.6
1	AS	106	LYS	2.6
1	BV	92	ILE	2.6
1	CN	3	ILE	2.6
1	CC	89	SER	2.6
1	AU	130	LEU	2.6
1	BG	53	PHE	2.5
1	DI	53	PHE	2.5
1	AP	70	ASN	2.5
1	CH	8	LYS	2.5
1	BS	131	ALA	2.5
1	CT	3	ILE	2.5
1	AI	89	SER	2.5
1	AK	35	LEU	2.5
1	AZ	127	LEU	2.5
1	BR	6	SER	2.5
1	BU	38	SER	2.5
1	CD	1	SER	2.5
1	CK	130	LEU	2.5
1	BK	61	VAL	2.5
1	BZ	31	VAL	2.5
1	CY	60	VAL	2.5
1	DG	119	ALA	2.5
1	BZ	45	TYR	2.5
1	AG	130	LEU	2.5
1	AT	35	LEU	2.5
1	CT	93	ALA	2.5
1	DI	131	ALA	2.5
1	AI	130	LEU	2.5
1	BO	91	ARG	2.5
1	CK	35	LEU	2.5
1	BE	57	VAL	2.5
1	CE	60	VAL	2.5
1	CJ	131	ALA	2.5
1	CU	131	ALA	2.5
1	CY	131	ALA	2.5
1	AY	3	ILE	2.5
1	CR	131	ALA	2.4

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Mol	Chain	Res	Type	RSRZ
1	AF	92	ILE	2.4
1	AC	124	PHE	2.4
1	DH	74	PHE	2.4
1	DD	77	PRO	2.4
1	CD	45	TYR	2.4
1	DK	45	TYR	2.4
1	BK	14	ALA	2.4
1	BU	131	ALA	2.4
1	AH	82	SER	2.4
1	AD	62	ASN	2.4
1	BP	30	THR	2.4
1	CO	32	THR	2.4
1	CQ	107	ASP	2.4
1	BY	73	VAL	2.4
1	AI	53	PHE	2.4
1	CX	6	SER	2.4
1	DI	20	SER	2.4
1	BM	127	LEU	2.4
1	CW	130	LEU	2.4
1	BM	129	ALA	2.4
1	CD	3	ILE	2.4
1	AX	53	PHE	2.4
1	BH	71	GLU	2.4
1	AZ	1	SER	2.4
1	BH	42	ASP	2.4
1	BD	106	LYS	2.4
1	BR	24	PHE	2.4
1	AH	127	LEU	2.4
1	BU	127	LEU	2.4
1	BV	1	SER	2.4
1	AM	60	VAL	2.4
1	DG	32	THR	2.3
1	AR	131	ALA	2.3
1	BB	96	ILE	2.3
1	BS	70	ASN	2.3
1	CR	33	ASN	2.3
1	CX	131	ALA	2.3
1	AC	130	LEU	2.3
1	AP	89	SER	2.3
1	CM	55	SER	2.3
1	BM	10	GLY	2.3
1	DI	111	ILE	2.3

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Mol	Chain	Res	Type	RSRZ
1	BA	53	PHE	2.3
1	DI	41	GLU	2.3
1	AH	55	SER	2.3
1	BG	92	ILE	2.3
1	BK	63	GLY	2.3
1	DA	76	ILE	2.3
1	AJ	126	THR	2.3
1	BJ	103	ALA	2.3
1	AH	130	LEU	2.3
1	CJ	26	LEU	2.3
1	CK	110	LEU	2.3
1	DI	130	LEU	2.3
1	AE	131	ALA	2.3
1	AL	131	ALA	2.3
1	AI	1	SER	2.3
1	BG	1	SER	2.3
1	CH	1	SER	2.3
1	CT	6	SER	2.3
1	DE	45	TYR	2.3
1	AX	29	GLN	2.3
1	BK	29	GLN	2.3
1	CE	81	ASP	2.3
1	BF	131	ALA	2.3
1	BZ	131	ALA	2.3
1	AG	40	SER	2.3
1	CD	38	SER	2.3
1	BV	33	ASN	2.3
1	BX	95	GLU	2.2
1	AX	35	LEU	2.2
1	CI	117	THR	2.2
1	BE	40	SER	2.2
1	CR	89	SER	2.2
1	BC	60	VAL	2.2
1	BH	70	ASN	2.2
1	AR	71	GLU	2.2
1	BG	131	ALA	2.2
1	BN	58	PRO	2.2
1	BX	51	ALA	2.2
1	CH	131	ALA	2.2
1	AR	124	PHE	2.2
1	AW	24	PHE	2.2
1	AX	30	THR	2.2

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Mol	Chain	Res	Type	RSRZ
1	AQ	121	TYR	2.2
1	BR	123	SER	2.2
1	DD	89	SER	2.2
1	DE	1	SER	2.2
1	BC	37	VAL	2.2
1	BC	50	ASN	2.2
1	DH	84	GLU	2.2
1	AF	130	LEU	2.2
1	AR	130	LEU	2.2
1	BF	93	ALA	2.2
1	BI	127	LEU	2.2
1	CL	130	LEU	2.2
1	CQ	130	LEU	2.2
1	AH	53	PHE	2.2
1	BW	85	THR	2.2
1	BU	45	TYR	2.2
1	BQ	61	VAL	2.2
1	BZ	29	GLN	2.2
1	CH	16	ILE	2.2
1	BF	33	ASN	2.2
1	BP	71	GLU	2.2
1	CS	126	THR	2.2
1	BH	37	VAL	2.2
1	AS	89	SER	2.2
1	BM	102	SER	2.2
1	CK	89	SER	2.2
1	AV	131	ALA	2.2
1	DA	131	ALA	2.2
1	AC	58	PRO	2.2
1	AF	54	LYS	2.2
1	BV	53	PHE	2.2
1	AJ	59	THR	2.1
1	BP	106	LYS	2.1
1	BW	54	LYS	2.1
1	CC	67	LYS	2.1
1	BP	70	ASN	2.1
1	CH	70	ASN	2.1
1	BV	72	VAL	2.1
1	DD	60	VAL	2.1
1	AG	126	THR	2.1
1	BO	30	THR	2.1
1	CR	85	THR	2.1

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Mol	Chain	Res	Type	RSRZ
1	DC	92	ILE	2.1
1	CI	115	LEU	2.1
1	AU	93	ALA	2.1
1	BK	131	ALA	2.1
1	CW	113	ALA	2.1
1	CX	93	ALA	2.1
1	BQ	55	SER	2.1
1	CV	24	PHE	2.1
1	CW	24	PHE	2.1
1	CK	72	VAL	2.1
1	AA	26	LEU	2.1
1	AN	30	THR	2.1
1	CS	129	ALA	2.1
1	AG	41	GLU	2.1
1	BA	112	GLY	2.1
1	AS	87	PHE	2.1
1	CU	89	SER	2.1
1	AB	60	VAL	2.1
1	DE	60	VAL	2.1
1	AJ	106	LYS	2.1
1	AV	106	LYS	2.1
1	AI	2	ILE	2.1
1	BM	130	LEU	2.1
1	BH	59	THR	2.1
1	BM	27	THR	2.1
1	AO	51	ALA	2.1
1	AW	93	ALA	2.1
1	BH	89	SER	2.1
1	BY	1	SER	2.1
1	CU	82	SER	2.1
1	AX	90	VAL	2.1
1	BO	31	VAL	2.1
1	CK	70	ASN	2.1
1	AA	3	ILE	2.1
1	BK	130	LEU	2.1
1	CJ	92	ILE	2.1
1	AU	131	ALA	2.1
1	DK	131	ALA	2.1
1	CT	53	PHE	2.1
1	AG	106	LYS	2.0
1	AU	69	LYS	2.0
1	AM	89	SER	2.0

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Mol	Chain	Res	Type	RSRZ
1	BV	40	SER	2.0
1	DA	89	SER	2.0
1	CB	130	LEU	2.0
1	BN	59	THR	2.0
1	BS	30	THR	2.0
1	CO	131	ALA	2.0
1	AL	77	PRO	2.0
1	AE	35	LEU	2.0
1	BI	3	ILE	2.0
1	DD	3	ILE	2.0
1	BQ	78	MET	2.0
1	CT	64	ASN	2.0
1	BU	125	TRP	2.0
1	BO	131	ALA	2.0
1	CJ	14	ALA	2.0
1	BN	45	TYR	2.0
1	AE	74	PHE	2.0
1	DI	56	ARG	2.0
1	BN	130	LEU	2.0
1	CS	130	LEU	2.0
1	AT	3	ILE	2.0
1	DE	2	ILE	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.