



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:01 AM UTC

PDB ID : 5K98 / pdb_00005k98
Title : Structure of HipA-HipB-O2-O3 complex
Authors : Schumacher, M.
Deposited on : 2016-05-31
Resolution : 3.99 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:

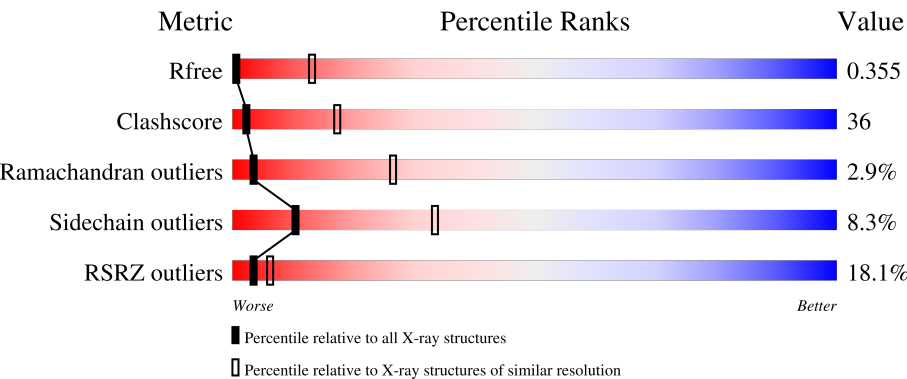
MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)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

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.99 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	1082 (4.20-3.80)
Clashscore	190562	1129 (4.20-3.80)
Ramachandran outliers	187476	1064 (4.20-3.80)
Sidechain outliers	187428	1055 (4.20-3.80)
RSRZ outliers	180081	1082 (4.20-3.80)

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	A	448	19% 46% 40% 5% • 8%
1	D	448	14% 48% 38% 5% • 8%
2	B	91	13% 35% 34% 9% 22%
2	P	91	19% 34% 34% 9% • 22%
3	T	23	9% 17% 78% •

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Mol	Chain	Length	Quality of chain
4	E	23	17% 22% 70% 9%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8605 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 Serine/threonine-protein kinase HipA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	414	Total	C	N	O	S	0	0	0
			3270	2098	570	590	12			
1	D	414	Total	C	N	O	S	0	0	0
			3270	2098	570	590	12			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	MET	-	initiating methionine	UNP P23874
A	-6	HIS	-	expression tag	UNP P23874
A	-5	HIS	-	expression tag	UNP P23874
A	-4	HIS	-	expression tag	UNP P23874
A	-3	HIS	-	expression tag	UNP P23874
A	-2	HIS	-	expression tag	UNP P23874
A	-1	HIS	-	expression tag	UNP P23874
A	0	SER	-	expression tag	UNP P23874
A	1	ARG	-	expression tag	UNP P23874
A	309	GLN	ASP	engineered mutation	UNP P23874
D	-7	MET	-	initiating methionine	UNP P23874
D	-6	HIS	-	expression tag	UNP P23874
D	-5	HIS	-	expression tag	UNP P23874
D	-4	HIS	-	expression tag	UNP P23874
D	-3	HIS	-	expression tag	UNP P23874
D	-2	HIS	-	expression tag	UNP P23874
D	-1	HIS	-	expression tag	UNP P23874
D	0	SER	-	expression tag	UNP P23874
D	1	ARG	-	expression tag	UNP P23874
D	309	GLN	ASP	engineered mutation	UNP P23874

- Molecule 2 is a protein called Antitoxin HipB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	71	Total	C	N	O	S	0	0	0
			564	358	95	108	3			
2	P	71	Total	C	N	O	S	0	0	0
			564	358	95	108	3			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-2	GLY	-	expression tag	UNP M9IJX7
B	-1	SER	-	expression tag	UNP M9IJX7
B	0	HIS	-	expression tag	UNP M9IJX7
P	-2	GLY	-	expression tag	UNP M9IJX7
P	-1	SER	-	expression tag	UNP M9IJX7
P	0	HIS	-	expression tag	UNP M9IJX7

- Molecule 3 is a DNA chain called DNA (5'-D(*TP*CP*CP*CP*TP*AP*TP*CP*CP*CP*CP*TP*TP*AP*AP*GP*GP*GP*GP*AP*TP*AP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	T	23	Total	C	N	O	P	0	0	0
			465	223	83	137	22			

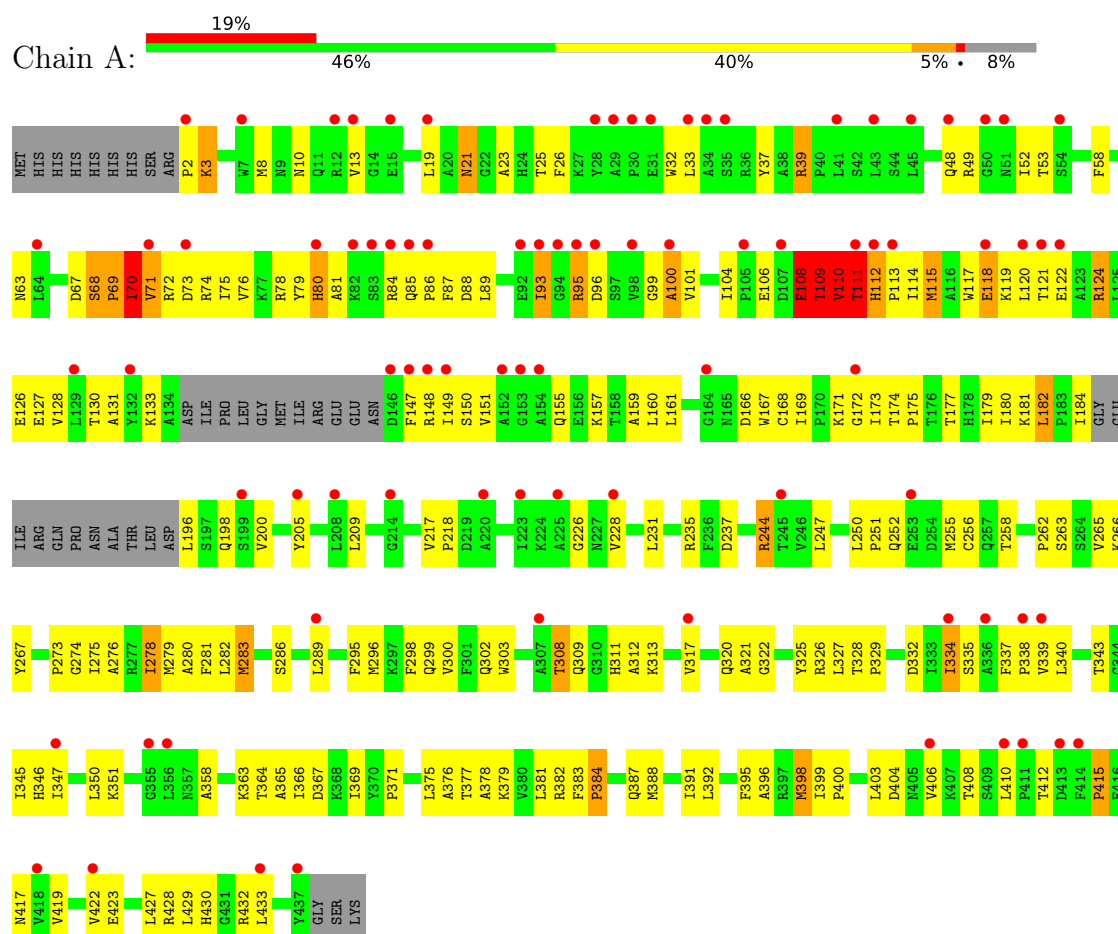
- Molecule 4 is a DNA chain called DNA (5'-D(*CP*TP*AP*TP*CP*CP*CP*CP*TP*TP*AP*AP*GP*GP*GP*GP*AP*TP*AP*GP*GP*GP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	E	23	Total	C	N	O	P	0	0	0
			472	225	90	135	22			

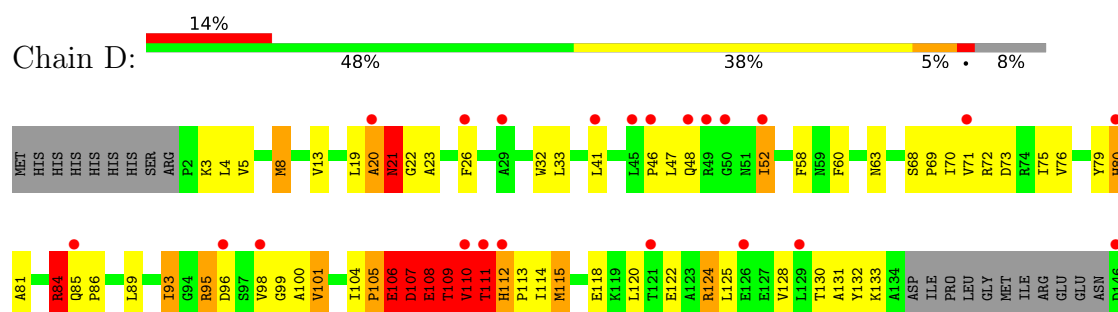
3 Residue-property plots [i](#)

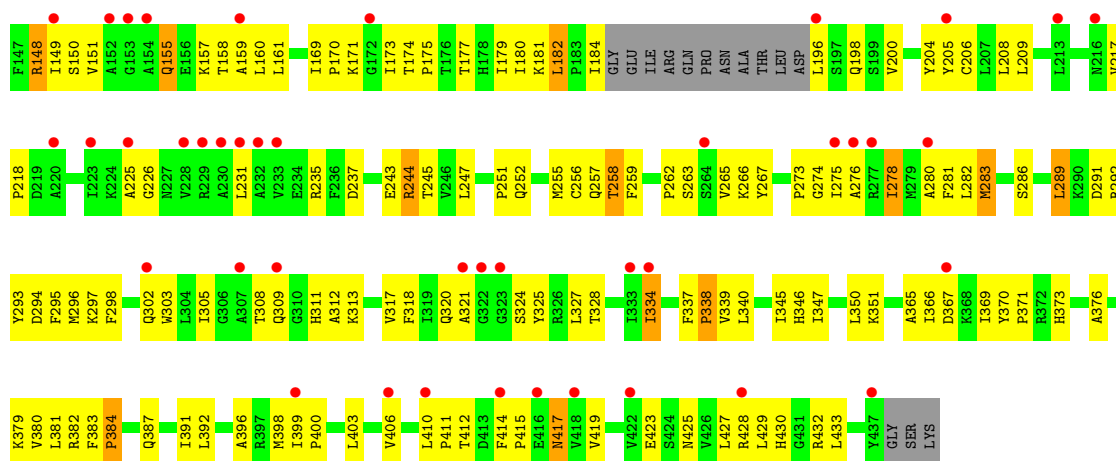
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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.

• Molecule 1: Serine/threonine-protein kinase HipA



• Molecule 1: Serine/threonine-protein kinase HipA

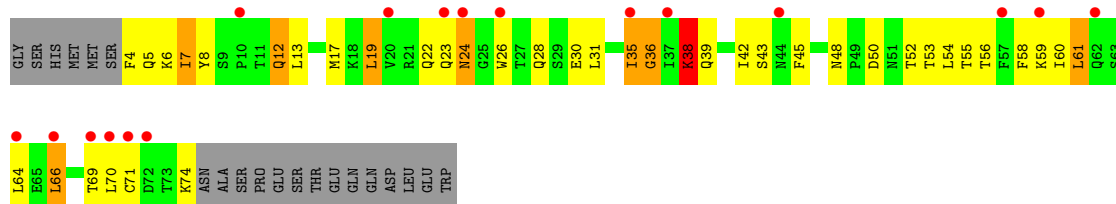




• Molecule 2: Antitoxin HipB



• Molecule 2: Antitoxin HipB

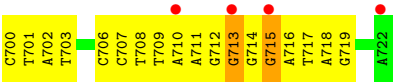


• Molecule 3: DNA (5'-D(*TP*CP*CP*CP*TP*AP*TP*CP*CP*CP*CP*TP*TP*AP*AP*GP*GP*GP*GP*AP*TP*AP*G)-3')



• Molecule 4: DNA (5'-D(*CP*TP*AP*TP*CP*CP*CP*CP*TP*TP*AP*AP*GP*GP*GP*GP*AP*TP*AP*GP*GP*GP*A)-3')





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	214.04Å 146.83Å 53.75Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	121.08 – 3.99 121.08 – 3.99	Depositor EDS
% Data completeness (in resolution range)	92.2 (121.08-3.99) 92.4 (121.08-3.99)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.97 (at 4.01Å)	Xtriage
Refinement program	PHENIX 1.6.4 _486	Depositor
R, R_{free}	0.350 , 0.375 0.329 , 0.355	Depositor DCC
R_{free} test set	1405 reflections (10.00%)	wwPDB-VP
Wilson B-factor (Å ²)	120.1	Xtriage
Anisotropy	0.930	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 128.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.36$, $\langle L^2 \rangle = 0.19$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.83	EDS
Total number of atoms	8605	wwPDB-VP
Average B, all atoms (Å ²)	179.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 24.73 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 3.6243e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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	A	1.17	16/3343 (0.5%)	1.10	19/4530 (0.4%)
1	D	0.92	13/3343 (0.4%)	1.05	23/4530 (0.5%)
2	B	0.71	0/572	0.97	0/771
2	P	0.60	0/572	0.97	3/771 (0.4%)
3	T	0.46	0/520	0.96	1/800 (0.1%)
4	E	0.46	0/530	1.20	4/817 (0.5%)
All	All	0.96	29/8880 (0.3%)	1.07	50/12219 (0.4%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (29) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	110	VAL	CA-CB	46.73	2.18	1.54
1	D	109	THR	CA-CB	23.43	1.93	1.53
1	D	109	THR	CA-C	19.72	1.79	1.52
1	A	110	VAL	CB-CG1	18.30	2.12	1.52
1	A	110	VAL	CA-C	15.62	1.72	1.52
1	A	111	THR	N-CA	15.17	1.66	1.46
1	D	110	VAL	CA-C	12.57	1.68	1.52
1	D	107	ASP	CA-C	9.97	1.66	1.52
1	D	107	ASP	CB-CG	9.94	1.76	1.52
1	D	107	ASP	CA-CB	8.30	1.67	1.53
1	A	111	THR	CA-CB	7.98	1.66	1.53
1	D	283	MET	CG-SD	7.36	1.99	1.80
1	A	109	THR	CA-C	7.17	1.62	1.52
1	D	110	VAL	N-CA	7.09	1.55	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	398	MET	CG-SD	7.06	1.98	1.80
1	A	109	THR	CA-CB	6.86	1.65	1.53
1	D	398	MET	SD-CE	6.83	1.96	1.79
1	A	112	HIS	N-CA	6.81	1.56	1.46
1	A	283	MET	CG-SD	6.78	1.97	1.80
1	D	283	MET	SD-CE	6.71	1.96	1.79
1	D	106	GLU	CA-C	-6.67	1.44	1.52
1	A	283	MET	SD-CE	6.51	1.95	1.79
1	A	398	MET	SD-CE	6.45	1.95	1.79
1	A	110	VAL	C-N	6.43	1.42	1.33
1	A	398	MET	CG-SD	6.30	1.96	1.80
1	D	110	VAL	CA-CB	-5.97	1.46	1.54
1	A	109	THR	C-N	5.51	1.41	1.33
1	A	109	THR	C-O	5.50	1.30	1.24
1	A	228	VAL	CA-CB	5.15	1.59	1.54

All (50) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	111	THR	N-CA-CB	18.82	140.61	109.56
1	A	68	SER	CA-C-N	-13.51	105.82	119.87
1	A	68	SER	C-N-CA	-13.51	105.82	119.87
1	A	111	THR	N-CA-C	-11.40	94.39	110.35
1	D	109	THR	O-C-N	-11.19	107.70	122.59
4	E	713	DG	O4'-C1'-N9	10.56	124.24	108.40
1	A	110	VAL	CB-CA-C	10.53	128.56	111.29
1	D	110	VAL	N-CA-C	9.82	129.77	109.34
1	A	70	ILE	CB-CA-C	-9.72	99.52	111.97
1	D	109	THR	CA-CB-CG2	9.42	126.51	110.50
1	D	106	GLU	N-CA-C	-8.76	101.68	111.14
1	A	110	VAL	CA-CB-CG1	8.47	124.80	110.40
1	D	109	THR	CB-CA-C	8.09	126.52	110.42
1	D	107	ASP	CB-CA-C	8.07	126.48	110.42
4	E	715	DG	O4'-C1'-N9	7.51	119.66	108.40
1	A	111	THR	CA-CB-OG1	7.43	120.74	109.60
1	A	112	HIS	CA-C-N	-7.20	110.84	119.84
1	A	112	HIS	C-N-CA	-7.20	110.84	119.84
1	D	108	GLU	CA-C-N	-6.99	108.19	121.54
1	D	108	GLU	C-N-CA	-6.99	108.19	121.54
1	A	109	THR	CA-C-O	-6.73	110.89	120.51
1	D	109	THR	CA-C-N	-6.72	109.86	121.97
1	D	109	THR	C-N-CA	-6.72	109.86	121.97

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	E	715	DG	O5'-C5'-C4'	-6.56	100.97	110.80
1	D	107	ASP	CA-C-N	6.43	133.82	121.54
1	D	107	ASP	C-N-CA	6.43	133.82	121.54
1	D	106	GLU	CA-C-N	-6.32	109.46	121.54
1	D	106	GLU	C-N-CA	-6.32	109.46	121.54
2	P	48	ASN	CA-C-N	6.32	125.94	119.56
2	P	48	ASN	C-N-CA	6.32	125.94	119.56
1	D	70	ILE	CB-CA-C	-6.12	103.88	112.14
1	A	111	THR	CA-C-O	-6.00	114.59	121.07
1	D	84	ARG	NE-CZ-NH1	-5.91	115.59	121.50
1	D	110	VAL	CB-CA-C	-5.90	101.61	111.29
1	D	68	SER	CA-C-N	-5.84	114.23	121.00
1	D	68	SER	C-N-CA	-5.84	114.23	121.00
1	D	111	THR	N-CA-CB	5.55	119.88	110.49
1	A	108	GLU	CA-C-N	-5.47	111.08	121.54
1	A	108	GLU	C-N-CA	-5.47	111.08	121.54
2	P	38	LYS	CD-CE-NZ	5.46	129.38	111.90
1	A	71	VAL	N-CA-C	-5.44	104.42	110.62
1	D	112	HIS	N-CA-C	-5.44	103.18	109.93
1	A	39	ARG	CA-C-N	-5.43	114.37	119.85
1	A	39	ARG	C-N-CA	-5.43	114.37	119.85
1	A	278	ILE	CB-CA-C	-5.31	104.97	112.14
4	E	713	DG	C4'-C3'-O3'	5.30	117.96	110.00
1	D	106	GLU	N-CA-CB	5.28	117.72	110.07
1	A	111	THR	CB-CA-C	-5.26	101.30	109.61
3	T	708	DT	O4'-C1'-N1	5.23	116.25	108.40
1	D	70	ILE	N-CA-C	5.14	115.86	110.62

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	108	GLU	Peptide

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	A	3270	0	3313	250	10
1	D	3270	0	3313	223	10
2	B	564	0	574	54	0
2	P	564	0	574	49	0
3	T	465	0	261	37	0
4	E	472	0	260	55	0
All	All	8605	0	8295	607	10

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 36.

All (607) 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:D:107:ASP:CG	1:D:107:ASP:CB	1.76	1.53
1:D:109:THR:C	1:D:109:THR:CA	1.79	1.52
1:D:109:THR:CA	1:D:109:THR:CB	1.93	1.42
1:A:110:VAL:CB	1:A:110:VAL:CG1	2.12	1.27
1:A:68:SER:OG	1:A:70:ILE:HD12	1.31	1.25
1:A:110:VAL:CB	1:A:110:VAL:CA	2.18	1.22
1:A:112:HIS:CG	1:A:173:ILE:HG23	1.71	1.22
2:B:38:LYS:HD2	4:E:715:DG:C6	1.76	1.19
1:A:118:GLU:HB2	1:A:171:LYS:HE3	1.24	1.17
1:D:110:VAL:HG12	1:D:110:VAL:O	1.35	1.13
1:D:104:ILE:HG22	1:D:108:GLU:HB3	1.31	1.11
1:D:110:VAL:HG11	1:D:112:HIS:CE1	1.87	1.08
1:D:124:ARG:HH11	1:D:124:ARG:HB3	0.95	1.07
1:D:148:ARG:HG3	1:D:148:ARG:HH11	1.18	1.04
1:D:206:CYS:SG	1:D:305:ILE:HD12	1.97	1.03
1:A:95:ARG:NE	1:A:112:HIS:HE1	1.58	1.01
1:D:106:GLU:O	1:D:107:ASP:C	2.02	1.01
1:A:95:ARG:CD	1:A:112:HIS:CE1	2.45	1.00
1:A:110:VAL:CB	1:A:110:VAL:HA	1.92	0.99
1:D:124:ARG:HB3	1:D:124:ARG:NH1	1.80	0.95
2:P:39:GLN:HE22	4:E:702:DA:H62	1.06	0.95
1:A:112:HIS:HB2	1:A:173:ILE:HG12	1.48	0.94
1:A:345:ILE:HD12	1:A:345:ILE:H	1.33	0.94
1:A:95:ARG:NE	1:A:112:HIS:CE1	2.36	0.93
1:A:95:ARG:HD3	1:A:112:HIS:ND1	1.84	0.93
1:D:345:ILE:HD12	1:D:345:ILE:H	1.33	0.92
1:D:20:ALA:O	1:D:22:GLY:N	2.03	0.91
1:A:112:HIS:CD2	1:A:173:ILE:HG23	2.05	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:429:LEU:HD23	1:A:432:ARG:HH12	1.34	0.90
1:A:112:HIS:CB	1:A:173:ILE:HG23	2.02	0.90
1:A:118:GLU:HB2	1:A:171:LYS:CE	2.01	0.89
1:D:110:VAL:HG11	1:D:112:HIS:NE2	1.86	0.89
1:D:105:PRO:HB2	1:D:108:GLU:HB2	1.54	0.89
1:D:110:VAL:O	1:D:112:HIS:N	2.06	0.89
3:T:713:DG:H2''	2:P:38:LYS:HD3	1.54	0.88
1:D:399:ILE:HB	1:D:430:HIS:HD2	1.38	0.87
1:D:106:GLU:O	1:D:107:ASP:O	1.93	0.87
4:E:707:DC:H2''	4:E:708:DT:H5'	1.54	0.87
3:T:700:DC:H42	4:E:719:DG:H1	1.22	0.87
1:D:399:ILE:HB	1:D:430:HIS:CD2	2.09	0.87
1:D:110:VAL:O	1:D:111:THR:C	2.19	0.86
1:A:399:ILE:HB	1:A:430:HIS:HD2	1.40	0.85
1:D:429:LEU:HD23	1:D:432:ARG:HH12	1.41	0.85
1:A:99:GLY:HA2	1:A:252:GLN:HG2	1.55	0.85
2:B:38:LYS:HD2	4:E:715:DG:C5	2.11	0.85
1:D:124:ARG:HH11	1:D:124:ARG:CB	1.86	0.85
1:A:63:ASN:OD1	1:A:263:SER:HB3	1.76	0.85
2:B:27:THR:OG1	2:B:30:GLU:HG2	1.78	0.83
1:D:104:ILE:HG22	1:D:108:GLU:CB	2.07	0.83
4:E:709:DT:H1'	4:E:710:DA:H5'	1.59	0.83
2:B:38:LYS:HE2	4:E:713:DG:H2''	1.60	0.83
1:A:399:ILE:HB	1:A:430:HIS:CD2	2.13	0.83
1:D:96:ASP:HB3	1:D:149:ILE:HG23	1.59	0.83
1:D:428:ARG:HH11	1:D:428:ARG:HG2	1.42	0.83
1:A:112:HIS:CB	1:A:173:ILE:N	2.42	0.82
1:A:95:ARG:HD3	1:A:112:HIS:HD1	1.43	0.82
1:A:112:HIS:HB3	1:A:173:ILE:N	1.95	0.82
2:B:38:LYS:HE2	4:E:713:DG:C2'	2.10	0.81
1:A:429:LEU:HD23	1:A:432:ARG:NH1	1.94	0.81
1:A:403:LEU:HD13	1:A:423:GLU:HG3	1.63	0.81
4:E:706:DC:H1'	4:E:707:DC:H5'	1.60	0.81
1:A:112:HIS:HB2	1:A:173:ILE:HG23	1.62	0.80
1:A:68:SER:HG	1:A:70:ILE:HD12	1.40	0.80
1:D:99:GLY:HA2	1:D:252:GLN:HG2	1.64	0.80
2:B:38:LYS:CE	4:E:713:DG:H2''	2.11	0.80
3:T:704:DC:O2	4:E:716:DA:H2	1.65	0.79
1:A:119:LYS:HA	1:A:168:CYS:SG	2.22	0.79
1:A:200:VAL:HG13	1:A:231:LEU:HB2	1.65	0.79
1:A:124:ARG:HH11	1:A:124:ARG:HB3	1.48	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:T:702:DA:N1	4:E:718:DA:H2	1.82	0.78
1:D:20:ALA:C	1:D:22:GLY:H	1.91	0.78
3:T:705:DC:C4	4:E:714:DG:O6	2.36	0.78
1:A:76:VAL:HG13	1:A:81:ALA:HB3	1.64	0.77
1:A:95:ARG:HD3	1:A:112:HIS:CE1	2.16	0.77
1:A:112:HIS:HB3	1:A:172:GLY:CA	2.15	0.77
3:T:702:DA:H2''	3:T:703:DT:H5''	1.66	0.77
1:D:110:VAL:O	1:D:110:VAL:CG1	2.21	0.77
1:D:109:THR:CA	1:D:109:THR:HB	2.14	0.77
3:T:698:DC:H2''	3:T:699:DC:C5	2.20	0.77
3:T:712:DG:H4'	3:T:713:DG:H5'	1.66	0.76
1:D:169:ILE:HG23	1:D:170:PRO:HD2	1.69	0.75
1:D:109:THR:O	1:D:110:VAL:HB	1.86	0.75
1:A:112:HIS:HB2	1:A:173:ILE:CG1	2.17	0.74
1:D:105:PRO:CB	1:D:108:GLU:HB2	2.17	0.74
1:D:383:PHE:CD1	1:D:384:PRO:HD2	2.23	0.74
1:A:49:ARG:HD2	2:B:23:GLN:HG2	1.70	0.74
1:A:151:VAL:HG12	1:A:157:LYS:HE3	1.69	0.74
2:B:7:ILE:HG23	2:B:12:GLN:HG2	1.70	0.74
2:P:39:GLN:NE2	4:E:702:DA:H62	1.85	0.74
1:D:69:PRO:HA	1:D:72:ARG:HG3	1.70	0.73
1:A:95:ARG:HD2	1:A:148:ARG:HB2	1.71	0.73
1:A:112:HIS:CB	1:A:173:ILE:H	2.02	0.72
1:D:72:ARG:O	1:D:76:VAL:HG23	1.90	0.72
2:B:10:PRO:HG2	2:B:50:ASP:OD1	1.89	0.72
1:D:76:VAL:HG22	1:D:89:LEU:HD21	1.71	0.72
1:D:392:LEU:HD22	1:D:433:LEU:HD21	1.72	0.71
1:A:21:ASN:HD21	1:A:23:ALA:CB	2.03	0.71
1:D:204:TYR:CZ	1:D:208:LEU:HD11	2.24	0.71
1:D:109:THR:O	1:D:110:VAL:CB	2.36	0.71
1:A:69:PRO:O	1:A:72:ARG:HB2	1.90	0.71
1:A:339:VAL:HG12	1:A:339:VAL:O	1.91	0.70
1:D:76:VAL:HG13	1:D:81:ALA:HB3	1.72	0.70
1:A:111:THR:HG21	1:A:114:ILE:HD11	1.74	0.70
2:B:38:LYS:CD	4:E:713:DG:H2''	2.22	0.69
1:A:399:ILE:CB	1:A:430:HIS:HD2	2.06	0.69
1:A:95:ARG:HG2	1:A:112:HIS:CE1	2.28	0.69
1:A:85:GLN:HE21	1:A:86:PRO:HD2	1.59	0.69
1:A:21:ASN:HD21	1:A:23:ALA:HB3	1.58	0.68
1:D:109:THR:C	1:D:109:THR:HA	2.08	0.68
1:D:120:LEU:HD22	1:D:124:ARG:NH1	2.09	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:384:PRO:HG2	1:D:387:GLN:NE2	2.08	0.68
1:A:71:VAL:O	1:A:75:ILE:HD13	1.94	0.68
1:A:298:PHE:O	1:A:302:GLN:HG3	1.93	0.68
1:D:148:ARG:HH11	1:D:148:ARG:CG	2.03	0.68
1:A:111:THR:HG23	1:A:112:HIS:N	2.08	0.68
3:T:701:DT:O2	3:T:702:DA:C5	2.46	0.68
1:D:76:VAL:HG21	1:D:84:ARG:HG3	1.76	0.68
1:A:280:ALA:O	1:A:283:MET:HB2	1.94	0.68
1:A:71:VAL:O	1:A:74:ARG:HB2	1.94	0.68
1:D:148:ARG:HG3	1:D:148:ARG:NH1	1.98	0.67
1:D:71:VAL:O	1:D:75:ILE:HD13	1.95	0.67
1:D:5:VAL:HG23	1:D:106:GLU:HA	1.76	0.67
1:D:286:SER:HA	2:P:8:TYR:CE2	2.29	0.67
1:A:428:ARG:HH11	1:A:428:ARG:HG2	1.60	0.67
2:B:59:LYS:HA	1:D:283:MET:HE1	1.76	0.66
1:D:76:VAL:HG21	1:D:84:ARG:CG	2.26	0.66
1:D:58:PHE:CE1	1:D:86:PRO:HG2	2.30	0.66
1:A:112:HIS:HB2	1:A:173:ILE:CG2	2.24	0.66
1:A:96:ASP:HB3	1:A:149:ILE:HG23	1.78	0.66
2:P:6:LYS:NZ	2:P:74:LYS:HE3	2.10	0.66
1:A:95:ARG:CG	1:A:112:HIS:CE1	2.78	0.66
1:A:161:LEU:HD12	1:A:177:THR:HG23	1.77	0.66
2:B:56:THR:OG1	4:E:712:DG:H5"	1.96	0.66
2:B:55:THR:O	2:B:59:LYS:HG3	1.96	0.66
1:D:179:ILE:HD11	1:D:235:ARG:CZ	2.26	0.65
1:A:112:HIS:CD2	1:A:173:ILE:CG2	2.79	0.65
3:T:704:DC:O2	4:E:716:DA:C2	2.49	0.65
1:D:26:PHE:O	1:D:52:ILE:HG12	1.97	0.65
1:A:115:MET:HE2	1:A:115:MET:HA	1.79	0.65
2:B:17:MET:SD	2:B:45:PHE:CZ	2.89	0.64
1:A:112:HIS:HB2	1:A:173:ILE:N	2.12	0.64
1:A:13:VAL:HG13	1:A:32:TRP:CE2	2.33	0.64
1:A:383:PHE:CD1	1:A:384:PRO:HD2	2.32	0.64
2:B:37:ILE:HB	4:E:713:DG:OP2	1.97	0.64
1:A:244:ARG:HH11	1:A:244:ARG:HG2	1.61	0.64
1:D:21:ASN:C	1:D:21:ASN:HD22	2.05	0.63
1:D:399:ILE:CB	1:D:430:HIS:HD2	2.08	0.63
2:P:19:LEU:CD2	2:P:23:GLN:HG3	2.29	0.63
1:A:111:THR:CG2	1:A:112:HIS:N	2.59	0.63
3:T:716:DA:H61	4:E:703:DT:H3	1.46	0.63
2:P:35:ILE:HG13	2:P:36:GLY:H	1.63	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:17:MET:SD	2:B:45:PHE:HZ	2.21	0.63
1:D:429:LEU:HD23	1:D:432:ARG:NH1	2.12	0.63
1:A:112:HIS:O	1:A:113:PRO:C	2.38	0.63
1:A:175:PRO:HB2	1:A:247:LEU:HD23	1.81	0.63
1:A:112:HIS:HB2	1:A:173:ILE:CB	2.30	0.62
1:D:3:LYS:O	1:D:106:GLU:N	2.32	0.62
1:A:112:HIS:HB3	1:A:173:ILE:H	1.60	0.62
1:D:182:LEU:HD22	1:D:182:LEU:H	1.64	0.62
2:B:38:LYS:HG3	4:E:713:DG:C2'	2.30	0.62
1:D:345:ILE:HD12	1:D:345:ILE:N	2.12	0.62
1:A:112:HIS:CG	1:A:173:ILE:CG2	2.67	0.62
1:D:345:ILE:H	1:D:345:ILE:CD1	2.09	0.62
2:P:55:THR:O	2:P:59:LYS:HG3	2.00	0.62
1:A:119:LYS:HG3	1:A:166:ASP:CB	2.30	0.62
1:D:95:ARG:HB2	1:D:148:ARG:O	1.99	0.62
1:A:112:HIS:HB3	1:A:172:GLY:C	2.25	0.61
1:A:115:MET:CE	1:A:174:THR:HG23	2.30	0.61
2:P:6:LYS:HZ1	2:P:74:LYS:HE3	1.65	0.61
1:D:179:ILE:HG12	1:D:235:ARG:HG2	1.83	0.61
1:A:68:SER:HB3	1:A:71:VAL:HG23	1.82	0.61
1:A:110:VAL:CA	1:A:110:VAL:HB	2.27	0.61
2:B:59:LYS:HA	1:D:283:MET:CE	2.29	0.61
1:D:151:VAL:HG12	1:D:157:LYS:HE3	1.82	0.61
1:D:244:ARG:HG2	1:D:244:ARG:HH11	1.63	0.61
2:B:64:LEU:O	2:B:66:LEU:HD13	2.00	0.61
1:A:75:ILE:HG22	1:A:89:LEU:HD22	1.83	0.61
1:A:205:TYR:HD1	1:A:410:LEU:HD21	1.65	0.61
1:D:175:PRO:HB2	1:D:247:LEU:HD23	1.81	0.61
4:E:712:DG:C6	4:E:713:DG:C2	2.89	0.61
1:A:159:ALA:O	1:A:160:LEU:HD23	2.00	0.60
1:A:274:GLY:O	1:A:278:ILE:HG12	2.01	0.60
1:A:256:CYS:SG	1:A:313:LYS:HG3	2.42	0.60
2:P:64:LEU:O	2:P:66:LEU:HD13	2.00	0.60
1:A:13:VAL:HG13	1:A:32:TRP:NE1	2.16	0.60
1:A:382:ARG:HD3	3:T:714:DG:OP1	2.01	0.60
2:B:52:THR:HB	2:P:54:LEU:HD12	1.83	0.60
1:D:21:ASN:HD22	1:D:22:GLY:N	1.98	0.60
1:A:108:GLU:O	1:A:109:THR:C	2.42	0.60
1:A:120:LEU:HD22	1:A:124:ARG:HH11	1.66	0.60
1:D:337:PHE:C	1:D:339:VAL:H	2.08	0.60
1:A:110:VAL:O	1:A:111:THR:O	2.20	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:38:LYS:CD	4:E:715:DG:C6	2.70	0.60
1:D:281:PHE:HE2	1:D:317:VAL:HG11	1.66	0.59
1:A:320:GLN:OE1	1:A:326:ARG:HG2	2.02	0.59
1:A:184:ILE:HD12	1:A:196:LEU:HB2	1.84	0.59
2:P:7:ILE:HG23	2:P:12:GLN:HG2	1.83	0.59
1:D:115:MET:HA	1:D:115:MET:HE2	1.83	0.59
1:D:303:TRP:HE1	1:D:369:ILE:HB	1.67	0.59
1:A:58:PHE:CE1	1:A:86:PRO:HG2	2.38	0.59
1:A:21:ASN:ND2	1:A:23:ALA:CB	2.65	0.59
2:B:26:TRP:HA	2:B:30:GLU:OE1	2.03	0.58
1:D:108:GLU:HG3	1:D:110:VAL:CG2	2.33	0.58
3:T:715:DG:O6	2:P:38:LYS:NZ	2.31	0.58
3:T:718:DA:H2	4:E:702:DA:C2	2.21	0.58
1:A:72:ARG:O	1:A:76:VAL:HG23	2.03	0.58
3:T:712:DG:H2"	3:T:713:DG:C8	2.38	0.58
1:D:110:VAL:CG1	1:D:112:HIS:CE1	2.77	0.58
1:D:63:ASN:OD1	1:D:263:SER:HB3	2.03	0.58
1:A:89:LEU:O	1:A:93:ILE:HD13	2.03	0.58
2:B:38:LYS:HE2	4:E:713:DG:C8	2.39	0.58
1:A:339:VAL:O	1:A:345:ILE:HB	2.04	0.58
1:D:159:ALA:O	1:D:160:LEU:HD23	2.03	0.58
1:D:347:ILE:HG13	1:D:367:ASP:HB2	1.86	0.58
3:T:711:DA:N1	4:E:710:DA:H5"	2.19	0.58
1:D:19:LEU:HD12	1:D:23:ALA:HB3	1.85	0.58
1:A:10:ASN:HD21	1:A:173:ILE:HD12	1.69	0.58
1:A:345:ILE:H	1:A:345:ILE:CD1	2.09	0.58
1:A:217:VAL:HB	1:A:218:PRO:HD2	1.85	0.58
1:D:48:GLN:HB2	2:P:22:GLN:HE22	1.68	0.58
1:D:370:TYR:CD2	1:D:432:ARG:HD3	2.39	0.58
1:A:48:GLN:HB2	2:B:22:GLN:HE22	1.68	0.57
1:D:157:LYS:HD3	1:D:181:LYS:HE2	1.86	0.57
1:A:112:HIS:HB2	1:A:173:ILE:H	1.68	0.57
1:D:21:ASN:HD21	1:D:23:ALA:HB3	1.69	0.57
1:D:303:TRP:NE1	1:D:369:ILE:HB	2.20	0.57
1:A:19:LEU:HB2	1:A:23:ALA:HB3	1.87	0.57
1:A:387:GLN:O	1:A:391:ILE:HG13	2.04	0.57
2:P:17:MET:SD	2:P:45:PHE:CZ	2.98	0.57
1:A:85:GLN:HB2	1:A:88:ASP:OD2	2.04	0.57
1:D:291:ASP:O	1:D:292:ARG:C	2.48	0.57
1:A:322:GLY:HA2	2:B:5:GLN:NE2	2.20	0.56
1:A:104:ILE:CD1	1:A:112:HIS:NE2	2.69	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:108:GLU:HG3	1:A:110:VAL:HG23	1.86	0.56
3:T:712:DG:H2''	3:T:713:DG:N7	2.19	0.56
2:B:69:THR:OG1	2:P:4:PHE:HB3	2.05	0.56
1:D:124:ARG:O	1:D:128:VAL:HG23	2.05	0.56
4:E:702:DA:H2''	4:E:703:DT:H71	1.88	0.56
1:A:179:ILE:HD11	1:A:235:ARG:CZ	2.35	0.56
4:E:707:DC:H2''	4:E:708:DT:C5'	2.33	0.56
1:D:298:PHE:O	1:D:302:GLN:HG3	2.06	0.56
1:D:308:THR:HG22	1:D:351:LYS:O	2.06	0.56
1:A:75:ILE:CG2	1:A:89:LEU:HD22	2.36	0.56
1:D:339:VAL:O	1:D:345:ILE:HB	2.06	0.55
1:A:161:LEU:CD1	1:A:177:THR:HG23	2.36	0.55
1:D:20:ALA:C	1:D:22:GLY:N	2.55	0.55
2:P:17:MET:SD	2:P:45:PHE:HZ	2.30	0.55
3:T:718:DA:C2	4:E:702:DA:C2	2.94	0.55
2:P:39:GLN:HE22	4:E:702:DA:N6	1.89	0.55
4:E:714:DG:H2''	4:E:715:DG:O4'	2.06	0.55
1:A:21:ASN:ND2	1:A:23:ALA:HB2	2.21	0.55
1:D:79:TYR:O	1:D:80:HIS:C	2.49	0.55
1:D:115:MET:HE2	1:D:174:THR:HG23	1.88	0.55
1:A:399:ILE:HG22	1:A:400:PRO:N	2.19	0.55
1:D:200:VAL:HG13	1:D:231:LEU:HB2	1.89	0.55
1:A:26:PHE:O	1:A:52:ILE:HG12	2.07	0.55
1:A:281:PHE:CZ	1:A:325:TYR:CE2	2.95	0.55
1:D:295:PHE:O	1:D:296:MET:C	2.49	0.55
1:A:130:THR:HA	1:A:133:LYS:NZ	2.22	0.55
1:D:320:GLN:O	1:D:321:ALA:C	2.50	0.55
1:D:347:ILE:HD12	1:D:350:LEU:HD12	1.89	0.55
1:A:85:GLN:NE2	1:A:86:PRO:HD2	2.21	0.55
1:D:282:LEU:HD21	1:D:327:LEU:HD13	1.88	0.55
1:A:308:THR:HG22	1:A:351:LYS:O	2.07	0.54
4:E:717:DT:O2	4:E:718:DA:C2	2.60	0.54
1:A:19:LEU:HB2	1:A:21:ASN:HD21	1.72	0.54
1:A:384:PRO:HG2	1:A:387:GLN:NE2	2.22	0.54
1:D:130:THR:HA	1:D:133:LYS:NZ	2.22	0.54
1:A:347:ILE:HD12	1:A:350:LEU:HD12	1.88	0.54
1:D:370:TYR:H	1:D:373:HIS:HD1	1.55	0.54
1:A:337:PHE:C	1:A:339:VAL:H	2.15	0.54
4:E:712:DG:C5	4:E:713:DG:C4	2.95	0.54
1:D:339:VAL:O	1:D:339:VAL:HG12	2.07	0.54
1:D:428:ARG:HG2	1:D:428:ARG:NH1	2.16	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:35:ILE:HG23	2:B:63:SER:HB3	1.87	0.54
1:D:85:GLN:HE21	1:D:86:PRO:HD2	1.71	0.54
1:D:104:ILE:HG22	1:D:108:GLU:CG	2.38	0.54
1:D:427:LEU:HA	1:D:430:HIS:HB3	1.90	0.54
1:A:112:HIS:HB3	1:A:172:GLY:HA3	1.87	0.54
4:E:702:DA:C2'	4:E:703:DT:H71	2.38	0.54
1:A:93:ILE:HD13	1:A:93:ILE:H	1.72	0.54
1:A:303:TRP:NE1	1:A:369:ILE:HB	2.23	0.54
3:T:713:DG:C4	3:T:714:DG:N7	2.77	0.53
1:D:52:ILE:HD13	1:D:52:ILE:N	2.22	0.53
1:D:198:GLN:HE22	1:D:415:PRO:HB3	1.74	0.53
2:B:54:LEU:O	2:B:58:PHE:HD1	1.90	0.53
3:T:702:DA:C2'	3:T:703:DT:H5''	2.35	0.53
1:A:104:ILE:HD12	1:A:112:HIS:NE2	2.22	0.53
1:A:278:ILE:HG21	1:A:295:PHE:CE1	2.42	0.53
2:B:35:ILE:HG21	2:B:59:LYS:O	2.09	0.53
1:D:71:VAL:O	1:D:75:ILE:CD1	2.57	0.53
1:A:119:LYS:HG3	1:A:166:ASP:HB3	1.89	0.53
1:D:104:ILE:CG2	1:D:108:GLU:HG2	2.38	0.53
4:E:700:DC:H2''	4:E:701:DT:O5'	2.08	0.53
1:A:19:LEU:HD12	1:A:23:ALA:HB3	1.91	0.53
1:A:120:LEU:HD22	1:A:124:ARG:HB3	1.92	0.52
1:A:182:LEU:H	1:A:182:LEU:HD22	1.73	0.52
1:A:345:ILE:HD12	1:A:345:ILE:N	2.14	0.52
1:D:85:GLN:NE2	1:D:86:PRO:HD2	2.24	0.52
2:P:58:PHE:HA	2:P:61:LEU:HB2	1.91	0.52
1:D:410:LEU:HD13	1:D:419:VAL:HG22	1.90	0.52
1:D:108:GLU:HG3	1:D:110:VAL:HG23	1.90	0.52
1:D:46:PRO:O	1:D:47:LEU:C	2.53	0.52
2:P:19:LEU:HD22	2:P:23:GLN:HG3	1.91	0.52
1:A:95:ARG:HE	1:A:112:HIS:HE1	1.52	0.52
1:D:243:GLU:CB	1:D:245:THR:HG23	2.40	0.52
1:A:19:LEU:HB2	1:A:21:ASN:ND2	2.25	0.52
1:D:21:ASN:ND2	1:D:23:ALA:H	2.07	0.52
1:D:237:ASP:HB3	1:D:252:GLN:OE1	2.10	0.52
2:P:53:THR:HG22	2:P:55:THR:H	1.75	0.52
1:D:370:TYR:CE2	1:D:432:ARG:HD3	2.45	0.51
2:B:35:ILE:CG2	2:B:63:SER:HB3	2.40	0.51
2:B:69:THR:HG22	2:P:69:THR:HG23	1.92	0.51
1:D:96:ASP:CB	1:D:149:ILE:HG23	2.37	0.51
1:A:209:LEU:HA	1:A:406:VAL:HG21	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:E:716:DA:H2''	4:E:717:DT:H5'	1.93	0.51
4:E:717:DT:C2	4:E:718:DA:C6	2.98	0.51
1:A:399:ILE:CB	1:A:430:HIS:CD2	2.87	0.51
1:A:148:ARG:HG3	1:A:148:ARG:HH11	1.76	0.51
1:D:205:TYR:HD1	1:D:410:LEU:HD21	1.75	0.51
2:B:22:GLN:C	2:B:24:ASN:H	2.19	0.51
2:B:38:LYS:CG	4:E:713:DG:H2''	2.41	0.51
2:P:7:ILE:HG23	2:P:12:GLN:CG	2.40	0.51
1:A:110:VAL:O	1:A:111:THR:C	2.54	0.50
1:D:255:MET:HB2	1:D:312:ALA:O	2.10	0.50
1:A:275:ILE:HG22	1:A:381:LEU:HD11	1.93	0.50
1:D:19:LEU:HB2	1:D:21:ASN:HD21	1.76	0.50
1:D:110:VAL:CG1	1:D:112:HIS:NE2	2.66	0.50
1:D:428:ARG:HH11	1:D:428:ARG:CG	2.19	0.50
1:A:95:ARG:CZ	1:A:148:ARG:HG3	2.41	0.50
2:B:31:LEU:O	2:B:35:ILE:HG12	2.12	0.50
2:B:61:LEU:HD21	2:P:70:LEU:HD22	1.92	0.50
1:A:33:LEU:HD11	1:A:48:GLN:O	2.10	0.50
1:A:255:MET:HB2	1:A:312:ALA:O	2.12	0.50
1:A:347:ILE:HG13	1:A:367:ASP:HB2	1.93	0.50
1:D:376:ALA:O	1:D:379:LYS:HB2	2.12	0.50
2:P:53:THR:HG22	2:P:54:LEU:N	2.25	0.50
1:A:175:PRO:HG2	1:A:247:LEU:HB3	1.94	0.50
1:A:428:ARG:HG2	1:A:428:ARG:NH1	2.26	0.50
1:D:380:VAL:HG23	1:D:381:LEU:N	2.27	0.50
1:D:387:GLN:O	1:D:391:ILE:HG13	2.11	0.50
4:E:718:DA:H2'	4:E:718:DA:OP2	2.12	0.50
2:B:7:ILE:HG23	2:B:12:GLN:CG	2.42	0.49
1:D:293:TYR:O	1:D:294:ASP:C	2.55	0.49
1:D:33:LEU:HD11	1:D:48:GLN:O	2.12	0.49
1:D:104:ILE:O	1:D:105:PRO:C	2.55	0.49
1:D:104:ILE:O	1:D:105:PRO:O	2.29	0.49
1:D:161:LEU:CD1	1:D:177:THR:HG23	2.42	0.49
1:D:169:ILE:CG2	1:D:170:PRO:HD2	2.42	0.49
1:D:371:PRO:HB3	1:D:392:LEU:HD13	1.95	0.49
1:A:100:ALA:HB3	1:A:251:PRO:HA	1.95	0.49
1:D:262:PRO:O	1:D:265:VAL:HG22	2.12	0.49
2:P:55:THR:O	2:P:56:THR:C	2.52	0.49
1:A:198:GLN:HE22	1:A:415:PRO:HB3	1.78	0.49
3:T:703:DT:H1'	3:T:704:DC:H5'	1.94	0.49
1:D:21:ASN:ND2	1:D:23:ALA:CB	2.76	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:P:6:LYS:NZ	2:P:74:LYS:HG3	2.28	0.49
1:A:375:LEU:HD23	1:A:388:MET:HG2	1.94	0.49
1:A:376:ALA:O	1:A:379:LYS:HB2	2.13	0.49
1:D:125:LEU:HD23	1:D:225:ALA:HB2	1.93	0.49
1:A:115:MET:HE2	1:A:174:THR:HG23	1.95	0.49
1:D:149:ILE:HG22	1:D:150:SER:H	1.78	0.49
1:D:266:LYS:HD2	1:D:311:HIS:CD2	2.47	0.49
1:A:179:ILE:HG12	1:A:235:ARG:HG2	1.94	0.48
1:A:430:HIS:ND1	1:A:430:HIS:C	2.71	0.48
1:A:151:VAL:CG1	1:A:157:LYS:HE3	2.41	0.48
1:A:237:ASP:HB3	1:A:252:GLN:OE1	2.13	0.48
1:A:117:TRP:CD1	1:A:168:CYS:HG	2.31	0.48
1:A:182:LEU:N	1:A:182:LEU:HD13	2.28	0.48
1:A:275:ILE:CG2	1:A:381:LEU:HD11	2.43	0.48
1:A:320:GLN:O	1:A:321:ALA:C	2.57	0.48
1:A:404:ASP:O	1:A:408:THR:HG23	2.14	0.48
1:A:110:VAL:C	1:A:111:THR:O	2.53	0.48
1:A:384:PRO:HB2	1:A:387:GLN:HB2	1.95	0.48
1:A:340:LEU:HD23	1:A:345:ILE:HG22	1.96	0.48
1:D:115:MET:CE	1:D:174:THR:HG23	2.43	0.48
1:D:161:LEU:HD12	1:D:177:THR:HG23	1.94	0.48
2:P:42:ILE:HD13	2:P:60:ILE:CD1	2.43	0.48
1:A:429:LEU:CD2	1:A:432:ARG:HH12	2.17	0.48
1:D:280:ALA:O	1:D:283:MET:HB2	2.14	0.48
1:D:318:PHE:CD1	1:D:328:THR:HG22	2.49	0.48
2:P:53:THR:CG2	2:P:54:LEU:N	2.76	0.48
4:E:712:DG:O6	4:E:713:DG:C2	2.67	0.48
2:B:17:MET:HG3	2:P:70:LEU:HD11	1.94	0.48
1:D:309:GLN:HG3	1:D:334:ILE:CG2	2.44	0.48
2:B:6:LYS:HZ1	2:B:74:LYS:HG3	1.78	0.47
1:D:93:ILE:O	1:D:93:ILE:HG12	2.13	0.47
1:D:148:ARG:CG	1:D:148:ARG:NH1	2.69	0.47
2:P:31:LEU:O	2:P:35:ILE:HG12	2.14	0.47
1:A:303:TRP:HE1	1:A:369:ILE:HB	1.78	0.47
1:D:104:ILE:HG22	1:D:108:GLU:HG2	1.96	0.47
1:D:209:LEU:HA	1:D:406:VAL:HG21	1.97	0.47
1:A:358:ALA:HB2	1:A:363:LYS:H	1.78	0.47
2:B:6:LYS:HZ2	2:B:74:LYS:HE3	1.79	0.47
1:D:399:ILE:CB	1:D:430:HIS:CD2	2.88	0.47
1:D:414:PHE:CD2	1:D:415:PRO:HD2	2.50	0.47
1:D:414:PHE:CG	1:D:415:PRO:HD2	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:T:705:DC:N4	4:E:714:DG:O6	2.47	0.47
1:A:3:LYS:HD2	1:A:106:GLU:OE1	2.15	0.47
2:B:19:LEU:CD2	2:B:23:GLN:HG3	2.45	0.47
1:D:8:MET:HE3	1:D:101:VAL:HG12	1.97	0.47
1:A:429:LEU:CD2	1:A:432:ARG:NH1	2.72	0.47
2:B:38:LYS:HG3	4:E:713:DG:H2''	1.95	0.47
3:T:703:DT:H2''	3:T:704:DC:OP2	2.14	0.47
3:T:708:DT:H3	4:E:712:DG:H22	1.62	0.47
3:T:710:DA:H5''	4:E:711:DA:C2	2.49	0.47
2:P:56:THR:O	2:P:59:LYS:HB2	2.15	0.47
1:A:351:LYS:HD2	1:A:365:ALA:HA	1.97	0.47
1:A:68:SER:HB3	1:A:71:VAL:CG2	2.45	0.47
1:A:300:VAL:O	1:A:303:TRP:HB3	2.15	0.47
1:A:392:LEU:HD22	1:A:433:LEU:HD21	1.97	0.47
1:D:108:GLU:HG3	1:D:110:VAL:HG21	1.97	0.47
1:D:309:GLN:HG3	1:D:334:ILE:HG23	1.97	0.47
1:A:398:MET:O	1:A:399:ILE:C	2.56	0.46
1:D:60:PHE:CE1	1:D:257:GLN:NE2	2.83	0.46
1:D:256:CYS:SG	1:D:313:LYS:HG3	2.55	0.46
2:P:26:TRP:HA	2:P:30:GLU:OE1	2.15	0.46
1:A:371:PRO:HB3	1:A:392:LEU:HD13	1.97	0.46
2:B:54:LEU:HD12	2:P:52:THR:HB	1.97	0.46
1:D:337:PHE:O	1:D:339:VAL:N	2.49	0.46
1:D:399:ILE:HG22	1:D:400:PRO:N	2.31	0.46
4:E:702:DA:H2''	4:E:703:DT:C6	2.51	0.46
2:B:53:THR:HG22	2:B:55:THR:H	1.80	0.46
1:D:311:HIS:CE1	1:D:313:LYS:HB2	2.51	0.46
1:D:396:ALA:O	1:D:430:HIS:NE2	2.47	0.46
2:P:58:PHE:O	2:P:59:LYS:C	2.56	0.46
2:B:6:LYS:NZ	2:B:74:LYS:HG3	2.30	0.46
4:E:717:DT:O2	4:E:718:DA:N1	2.49	0.46
1:A:375:LEU:HD23	1:A:388:MET:CG	2.45	0.46
1:D:182:LEU:HD22	1:D:182:LEU:N	2.30	0.46
1:D:184:ILE:HD12	1:D:196:LEU:HB2	1.98	0.46
1:D:204:TYR:CE2	1:D:208:LEU:HD11	2.51	0.46
1:A:382:ARG:CD	3:T:714:DG:OP1	2.63	0.46
1:A:399:ILE:HG22	1:A:400:PRO:CD	2.46	0.46
1:A:273:PRO:HD2	1:A:312:ALA:HB2	1.98	0.46
1:A:76:VAL:HG22	1:A:89:LEU:HD21	1.97	0.46
1:D:109:THR:CB	1:D:109:THR:HA	2.27	0.45
1:D:114:ILE:HD13	1:D:173:ILE:HD11	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:25:THR:HG22	1:A:53:THR:HG22	1.99	0.45
1:A:255:MET:HE3	1:A:278:ILE:HG23	1.97	0.45
1:D:158:THR:OG1	1:D:159:ALA:N	2.49	0.45
1:A:396:ALA:O	1:A:430:HIS:NE2	2.48	0.45
1:D:370:TYR:CZ	1:D:432:ARG:NE	2.83	0.45
1:A:2:PRO:O	1:A:87:PHE:HZ	2.00	0.45
1:A:99:GLY:O	1:A:250:LEU:O	2.34	0.45
1:A:296:MET:O	1:A:299:GLN:N	2.50	0.45
3:T:714:DG:OP2	3:T:714:DG:H3'	2.17	0.45
1:D:13:VAL:HG13	1:D:32:TRP:CE2	2.52	0.45
1:D:21:ASN:HD21	1:D:23:ALA:CB	2.29	0.45
1:D:410:LEU:HD13	1:D:419:VAL:CG2	2.46	0.45
1:A:377:THR:O	1:A:378:ALA:C	2.58	0.45
1:D:274:GLY:O	1:D:278:ILE:HG12	2.17	0.45
1:D:297:LYS:HG3	1:D:391:ILE:HG23	1.99	0.45
1:A:126:GLU:O	1:A:130:THR:HG23	2.17	0.45
1:A:37:TYR:HB3	1:A:39:ARG:HH21	1.82	0.45
1:A:126:GLU:OE2	1:A:226:GLY:HA3	2.17	0.45
1:D:100:ALA:HB3	1:D:251:PRO:HA	1.98	0.45
4:E:707:DC:C2'	4:E:708:DT:H5'	2.37	0.45
1:A:115:MET:HA	1:A:115:MET:CE	2.47	0.44
1:A:147:PHE:O	1:A:147:PHE:CD1	2.70	0.44
1:A:351:LYS:CD	1:A:365:ALA:HA	2.47	0.44
3:T:709:DT:H2''	3:T:710:DA:O5'	2.17	0.44
1:D:282:LEU:HD23	1:D:282:LEU:HA	1.89	0.44
4:E:712:DG:N7	4:E:713:DG:C5	2.84	0.44
1:A:427:LEU:HA	1:A:430:HIS:HB3	1.98	0.44
3:T:712:DG:OP1	2:P:53:THR:N	2.34	0.44
1:D:157:LYS:CD	1:D:181:LYS:HE2	2.48	0.44
1:A:104:ILE:HD11	1:A:112:HIS:NE2	2.32	0.44
1:D:410:LEU:CD1	1:D:419:VAL:HG22	2.47	0.44
2:P:54:LEU:O	2:P:58:PHE:HD1	1.99	0.44
1:A:130:THR:HA	1:A:133:LYS:CE	2.48	0.44
1:A:295:PHE:O	1:A:296:MET:C	2.57	0.44
1:A:328:THR:O	1:A:329:PRO:C	2.60	0.44
1:A:429:LEU:O	1:A:432:ARG:HB2	2.17	0.44
1:A:262:PRO:O	1:A:265:VAL:HG22	2.18	0.44
4:E:708:DT:H1'	4:E:709:DT:C4	2.52	0.44
1:A:410:LEU:HD13	1:A:419:VAL:HG22	1.99	0.44
1:D:320:GLN:HB2	1:D:324:SER:HB2	1.99	0.44
1:D:430:HIS:ND1	1:D:430:HIS:C	2.75	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:P:6:LYS:HZ1	2:P:74:LYS:HG3	1.83	0.44
1:A:118:GLU:HB2	1:A:171:LYS:NZ	2.31	0.44
1:A:181:LYS:C	1:A:182:LEU:HD13	2.42	0.44
1:D:273:PRO:HD2	1:D:312:ALA:HB2	1.99	0.44
1:D:340:LEU:HD23	1:D:345:ILE:HG22	1.99	0.44
1:D:425:ASN:HD22	1:D:425:ASN:H	1.66	0.44
1:D:110:VAL:CG1	1:D:112:HIS:CD2	3.01	0.44
1:A:100:ALA:CB	1:A:251:PRO:HA	2.48	0.43
1:A:308:THR:HG22	1:A:351:LYS:C	2.43	0.43
1:D:243:GLU:O	1:D:245:THR:HG23	2.17	0.43
4:E:700:DC:C5	4:E:701:DT:H73	2.53	0.43
1:D:98:VAL:HG12	1:D:99:GLY:N	2.33	0.43
2:B:55:THR:O	2:B:56:THR:C	2.60	0.43
1:D:160:LEU:HG	1:D:180:ILE:HD12	2.00	0.43
1:D:258:THR:HG22	1:D:259:PHE:CD2	2.53	0.43
1:A:244:ARG:HG2	1:A:244:ARG:NH1	2.29	0.43
1:A:266:LYS:HD2	1:A:311:HIS:CD2	2.53	0.43
1:A:279:MET:O	1:A:280:ALA:C	2.61	0.43
1:A:337:PHE:O	1:A:339:VAL:N	2.51	0.43
1:D:120:LEU:HD22	1:D:124:ARG:HH11	1.81	0.43
1:A:95:ARG:NH2	1:A:148:ARG:NH1	2.67	0.43
1:D:98:VAL:HG12	1:D:99:GLY:H	1.84	0.43
1:D:266:LYS:HD2	1:D:311:HIS:CE1	2.53	0.43
4:E:717:DT:C2	4:E:718:DA:N1	2.86	0.43
1:A:124:ARG:O	1:A:128:VAL:HG23	2.19	0.43
2:B:38:LYS:HD2	4:E:715:DG:N1	2.27	0.43
2:P:22:GLN:C	2:P:24:ASN:H	2.27	0.43
1:A:52:ILE:N	1:A:52:ILE:HD13	2.33	0.43
1:D:255:MET:HE3	1:D:278:ILE:HG23	2.00	0.43
2:B:59:LYS:HG2	1:D:283:MET:HE2	2.00	0.43
1:D:337:PHE:C	1:D:339:VAL:N	2.72	0.43
1:A:182:LEU:HD22	1:A:182:LEU:N	2.34	0.43
1:A:279:MET:HG3	1:A:381:LEU:HD13	2.00	0.43
1:A:337:PHE:C	1:A:339:VAL:N	2.74	0.43
1:D:21:ASN:ND2	1:D:23:ALA:N	2.67	0.43
1:D:337:PHE:N	1:D:338:PRO:CD	2.82	0.42
1:A:120:LEU:HD11	1:A:169:ILE:HG13	2.01	0.42
1:A:108:GLU:C	1:A:110:VAL:N	2.74	0.42
1:A:120:LEU:HB2	1:A:167:TRP:O	2.20	0.42
1:A:149:ILE:HG22	1:A:150:SER:H	1.83	0.42
1:A:283:MET:HE1	2:P:59:LYS:HA	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:38:LYS:CG	4:E:713:DG:C2'	2.96	0.42
2:B:65:GLU:O	2:B:66:LEU:HD12	2.18	0.42
1:D:217:VAL:HB	1:D:218:PRO:HD2	2.01	0.42
2:P:38:LYS:HB2	2:P:38:LYS:HE2	1.35	0.42
1:A:296:MET:O	1:A:299:GLN:HB2	2.19	0.42
2:B:62:GLN:NE2	2:P:8:TYR:HB3	2.35	0.42
3:T:701:DT:C2	3:T:702:DA:C5	3.07	0.42
1:A:84:ARG:HE	1:A:84:ARG:HB2	1.69	0.42
1:D:120:LEU:HD11	1:D:169:ILE:HG13	2.02	0.42
1:D:170:PRO:C	1:D:171:LYS:HG3	2.43	0.42
1:A:127:GLU:O	1:A:131:ALA:HB2	2.20	0.42
1:A:275:ILE:O	1:A:276:ALA:C	2.63	0.42
1:A:343:THR:OG1	1:A:345:ILE:HD13	2.19	0.42
1:D:155:GLN:HG2	1:D:184:ILE:HG12	2.02	0.42
3:T:707:DC:H42	4:E:713:DG:H21	1.68	0.42
1:D:428:ARG:NH1	1:D:428:ARG:CG	2.79	0.42
1:A:96:ASP:OD2	1:A:175:PRO:HA	2.19	0.42
1:A:309:GLN:HG3	1:A:334:ILE:HG23	2.01	0.42
1:A:334:ILE:HG12	1:A:335:SER:N	2.34	0.42
3:T:712:DG:P	2:P:53:THR:H	2.42	0.42
1:A:75:ILE:HG22	1:A:89:LEU:CD2	2.50	0.42
1:D:275:ILE:CG2	1:D:381:LEU:HD11	2.50	0.42
1:D:403:LEU:HD13	1:D:423:GLU:HG3	2.01	0.42
1:A:282:LEU:HD21	1:A:327:LEU:HD13	2.01	0.41
3:T:705:DC:N3	4:E:714:DG:O6	2.52	0.41
1:D:13:VAL:HG13	1:D:32:TRP:NE1	2.34	0.41
2:P:39:GLN:HE21	2:P:39:GLN:HB3	1.64	0.41
1:A:89:LEU:O	1:A:93:ILE:CD1	2.68	0.41
1:A:382:ARG:NE	3:T:714:DG:OP1	2.53	0.41
3:T:712:DG:C4'	3:T:713:DG:H5'	2.43	0.41
1:D:21:ASN:HD22	1:D:23:ALA:H	1.66	0.41
1:D:410:LEU:HA	1:D:411:PRO:HD2	1.89	0.41
1:A:205:TYR:HE2	1:A:422:VAL:HG11	1.84	0.41
1:A:365:ALA:O	1:A:366:ILE:C	2.64	0.41
1:D:105:PRO:HG2	1:D:108:GLU:CD	2.45	0.41
1:D:160:LEU:HB2	1:D:180:ILE:HD11	2.02	0.41
1:D:275:ILE:O	1:D:276:ALA:C	2.62	0.41
1:D:351:LYS:CD	1:D:365:ALA:HA	2.49	0.41
1:D:383:PHE:O	1:D:384:PRO:C	2.63	0.41
2:P:28:GLN:OE1	2:P:43:SER:HB2	2.20	0.41
1:D:266:LYS:HE2	1:D:267:TYR:HE1	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:76:VAL:HG21	1:D:84:ARG:HG2	2.00	0.41
1:D:128:VAL:O	1:D:131:ALA:HB3	2.20	0.41
1:D:281:PHE:CZ	1:D:325:TYR:CE2	3.08	0.41
1:A:157:LYS:HD3	1:A:181:LYS:HE2	2.02	0.41
1:D:266:LYS:HD2	1:D:311:HIS:CG	2.56	0.41
1:D:380:VAL:C	1:D:382:ARG:H	2.28	0.41
1:D:180:ILE:HA	1:D:231:LEU:O	2.20	0.41
1:A:21:ASN:C	1:A:21:ASN:HD22	2.28	0.41
1:A:120:LEU:HD22	1:A:124:ARG:NH1	2.35	0.41
1:A:205:TYR:CD1	1:A:410:LEU:HD21	2.52	0.41
1:A:282:LEU:HA	1:A:282:LEU:HD23	1.73	0.41
1:A:286:SER:HA	2:B:8:TYR:CE2	2.56	0.41
2:B:34:LYS:HE2	1:D:289:LEU:HD23	2.02	0.41
3:T:714:DG:N7	2:P:38:LYS:NZ	2.63	0.41
1:D:120:LEU:HD21	1:D:169:ILE:HD12	2.02	0.41
1:D:149:ILE:HG22	1:D:150:SER:N	2.36	0.41
4:E:712:DG:C5	4:E:713:DG:C5	3.09	0.41
1:A:13:VAL:HG13	1:A:32:TRP:CD1	2.55	0.41
1:A:68:SER:O	1:A:69:PRO:C	2.63	0.41
1:A:112:HIS:O	1:A:114:ILE:N	2.53	0.41
1:A:70:ILE:H	1:A:70:ILE:HG13	1.57	0.40
1:A:79:TYR:O	1:A:80:HIS:C	2.64	0.40
1:A:205:TYR:CE2	1:A:422:VAL:HG11	2.55	0.40
2:B:67:SER:N	2:P:71:CYS:O	2.51	0.40
1:D:182:LEU:N	1:D:182:LEU:HD13	2.36	0.40
2:P:60:ILE:O	2:P:64:LEU:HG	2.20	0.40
1:A:85:GLN:HE21	1:A:86:PRO:CD	2.32	0.40
1:A:174:THR:HA	1:A:175:PRO:HD3	1.86	0.40
1:A:196:LEU:N	1:A:196:LEU:HD22	2.36	0.40
1:A:281:PHE:HE2	1:A:317:VAL:HG11	1.86	0.40
3:T:712:DG:OP2	2:P:52:THR:HA	2.22	0.40
1:D:8:MET:HE1	1:D:41:LEU:HD12	2.03	0.40
1:A:180:ILE:HA	1:A:231:LEU:O	2.21	0.40
1:A:181:LYS:NZ	1:A:332:ASP:OD1	2.44	0.40
1:A:266:LYS:HE2	1:A:267:TYR:HE1	1.86	0.40
2:B:34:LYS:HE2	1:D:289:LEU:CD2	2.51	0.40
1:D:4:LEU:HD23	1:D:105:PRO:HA	2.02	0.40
1:D:366:ILE:HG22	1:D:367:ASP:N	2.34	0.40
1:A:21:ASN:ND2	1:A:21:ASN:N	2.67	0.40
1:A:399:ILE:CB	1:A:400:PRO:HD3	2.51	0.40
1:A:100:ALA:HB2	1:A:250:LEU:C	2.46	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:112:HIS:CB	1:A:173:ILE:HG12	2.35	0.40
1:A:121:THR:H	1:A:124:ARG:HB2	1.87	0.40
1:A:395:PHE:O	1:A:396:ALA:C	2.64	0.40
1:D:21:ASN:HD22	1:D:23:ALA:N	2.19	0.40
1:D:337:PHE:HB2	1:D:338:PRO:HD3	2.03	0.40

All (10) 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:A:110:VAL:CB	1:D:109:THR:C[3_456]	1.68	0.52
1:A:110:VAL:CA	1:D:110:VAL:N[3_456]	1.94	0.26
1:A:110:VAL:CG1	1:D:109:THR:O[3_456]	1.95	0.25
1:A:110:VAL:O	1:D:110:VAL:N[3_456]	2.00	0.20
1:A:110:VAL:CB	1:D:110:VAL:N[3_456]	2.09	0.11
1:A:110:VAL:C	1:D:110:VAL:N[3_456]	2.10	0.10
1:A:67:ASP:O	1:D:84:ARG:NH1[3_455]	2.15	0.05
1:A:110:VAL:CB	1:D:109:THR:CA[3_456]	2.15	0.05
1:A:340:LEU:O	1:D:417:ASN:OD1[4_465]	2.15	0.05
1:A:110:VAL:CG1	1:D:109:THR:C[3_456]	2.18	0.02

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	A	408/448 (91%)	353 (86%)	45 (11%)	10 (2%)	4	29
1	D	408/448 (91%)	353 (86%)	41 (10%)	14 (3%)	3	24
2	B	69/91 (76%)	63 (91%)	5 (7%)	1 (1%)	9	39
2	P	69/91 (76%)	63 (91%)	3 (4%)	3 (4%)	2	20
All	All	954/1078 (88%)	832 (87%)	94 (10%)	28 (3%)	3	26

All (28) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	95	ARG
1	A	108	GLU
1	A	110	VAL
1	D	20	ALA
1	D	21	ASN
1	D	95	ARG
1	D	108	GLU
1	A	78	ARG
1	A	80	HIS
1	D	80	HIS
1	D	110	VAL
1	D	111	THR
1	A	100	ALA
2	B	23	GLN
1	D	105	PRO
1	A	384	PRO
1	D	113	PRO
1	A	415	PRO
1	D	109	THR
1	D	132	TYR
1	D	384	PRO
2	P	36	GLY
2	P	50	ASP
1	A	109	THR
1	D	338	PRO
1	A	338	PRO
1	D	226	GLY
2	P	35	ILE

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	A	351/382 (92%)	325 (93%)	26 (7%)	13	36
1	D	351/382 (92%)	326 (93%)	25 (7%)	13	38

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	B	64/83 (77%)	55 (86%)	9 (14%)	3	17
2	P	64/83 (77%)	55 (86%)	9 (14%)	3	17
All	All	830/930 (89%)	761 (92%)	69 (8%)	10	33

All (69) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	LYS
1	A	8	MET
1	A	21	ASN
1	A	69	PRO
1	A	70	ILE
1	A	73	ASP
1	A	93	ILE
1	A	101	VAL
1	A	108	GLU
1	A	110	VAL
1	A	111	THR
1	A	115	MET
1	A	118	GLU
1	A	122	GLU
1	A	124	ARG
1	A	155	GLN
1	A	182	LEU
1	A	244	ARG
1	A	258	THR
1	A	289	LEU
1	A	308	THR
1	A	334	ILE
1	A	346	HIS
1	A	364	THR
1	A	412	THR
1	A	417	ASN
2	B	5	GLN
2	B	7	ILE
2	B	12	GLN
2	B	13	LEU
2	B	19	LEU
2	B	24	ASN
2	B	42	ILE
2	B	61	LEU

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Mol	Chain	Res	Type
2	B	66	LEU
1	D	8	MET
1	D	21	ASN
1	D	52	ILE
1	D	73	ASP
1	D	84	ARG
1	D	93	ILE
1	D	101	VAL
1	D	106	GLU
1	D	107	ASP
1	D	108	GLU
1	D	115	MET
1	D	118	GLU
1	D	122	GLU
1	D	124	ARG
1	D	148	ARG
1	D	155	GLN
1	D	182	LEU
1	D	244	ARG
1	D	258	THR
1	D	278	ILE
1	D	289	LEU
1	D	334	ILE
1	D	346	HIS
1	D	412	THR
1	D	417	ASN
2	P	5	GLN
2	P	7	ILE
2	P	12	GLN
2	P	13	LEU
2	P	19	LEU
2	P	24	ASN
2	P	38	LYS
2	P	61	LEU
2	P	66	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	10	ASN
1	A	21	ASN
1	A	24	HIS

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Mol	Chain	Res	Type
1	A	48	GLN
1	A	85	GLN
1	A	198	GLN
1	A	227	ASN
1	A	299	GLN
1	A	357	ASN
1	A	425	ASN
2	B	5	GLN
2	B	12	GLN
2	B	22	GLN
2	B	39	GLN
2	B	48	ASN
2	B	62	GLN
1	D	10	ASN
1	D	21	ASN
1	D	80	HIS
1	D	85	GLN
1	D	198	GLN
1	D	216	ASN
1	D	302	GLN
1	D	320	GLN
1	D	357	ASN
1	D	387	GLN
1	D	425	ASN
2	P	5	GLN
2	P	12	GLN
2	P	22	GLN
2	P	24	ASN
2	P	39	GLN

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	A	414/448 (92%)	1.16	85 (20%) 2 5	155, 179, 191, 199	0
1	D	414/448 (92%)	1.08	64 (15%) 5 9	157, 179, 191, 206	0
2	B	71/91 (78%)	1.34	12 (16%) 4 7	162, 176, 183, 185	0
2	P	71/91 (78%)	1.49	17 (23%) 2 4	158, 175, 187, 190	0
3	T	23/23 (100%)	1.16	2 (8%) 16 17	168, 186, 195, 198	0
4	E	23/23 (100%)	1.35	4 (17%) 4 7	166, 191, 196, 204	0
All	All	1016/1124 (90%)	1.17	184 (18%) 3 6	155, 179, 191, 206	0

All (184) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	230	ALA	8.7
2	P	72	ASP	8.6
2	P	20	VAL	5.8
1	D	205	TYR	5.7
1	A	118	GLU	5.3
2	B	37	ILE	5.2
1	A	84	ARG	5.2
1	D	220	ALA	5.2
1	D	307	ALA	4.7
2	B	70	LEU	4.6
1	A	113	PRO	4.6
1	D	406	VAL	4.5
1	A	41	LEU	4.5
1	D	334	ILE	4.5
1	D	228	VAL	4.2
1	A	83	SER	4.1
1	D	50	GLY	4.0
1	D	223	ILE	3.9
1	A	29	ALA	3.9

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Mol	Chain	Res	Type	RSRZ
1	A	111	THR	3.9
2	P	23	GLN	3.8
2	P	70	LEU	3.8
1	D	129	LEU	3.8
1	A	172	GLY	3.8
1	A	214	GLY	3.8
1	A	45	LEU	3.7
1	A	96	ASP	3.6
1	D	422	VAL	3.6
1	A	414	PHE	3.6
1	D	399	ILE	3.5
1	A	86	PRO	3.5
2	P	66	LEU	3.4
1	A	205	TYR	3.4
1	D	52	ILE	3.4
1	A	220	ALA	3.4
1	A	422	VAL	3.4
1	A	146	ASP	3.4
1	A	406	VAL	3.4
2	P	37	ILE	3.4
1	D	96	ASP	3.3
1	D	154	ALA	3.3
1	D	172	GLY	3.3
1	A	223	ILE	3.3
1	D	41	LEU	3.3
1	D	233	VAL	3.3
2	B	68	MET	3.3
1	D	126	GLU	3.3
1	A	112	HIS	3.2
2	B	57	PHE	3.2
1	A	154	ALA	3.2
1	A	245	THR	3.1
1	D	48	GLN	3.1
1	A	28	TYR	3.1
2	B	72	ASP	3.1
1	A	418	VAL	3.1
1	A	149	ILE	3.1
1	A	334	ILE	3.1
1	A	120	LEU	3.1
1	D	110	VAL	3.0
1	A	31	GLU	3.0
1	A	73	ASP	3.0

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Mol	Chain	Res	Type	RSRZ
2	P	57	PHE	3.0
1	D	45	LEU	3.0
1	D	367	ASP	3.0
1	D	231	LEU	2.9
1	D	85	GLN	2.9
1	D	437	TYR	2.9
2	B	48	ASN	2.9
1	A	339	VAL	2.9
1	D	410	LEU	2.9
2	B	44	ASN	2.9
1	A	12	ARG	2.9
1	A	122	GLU	2.9
1	D	196	LEU	2.8
4	E	713	DG	2.8
1	A	2	PRO	2.8
1	A	437	TYR	2.8
1	D	280	ALA	2.8
2	P	69	THR	2.8
1	A	153	GLY	2.7
1	A	93	ILE	2.7
1	D	112	HIS	2.7
1	A	411	PRO	2.7
1	A	33	LEU	2.7
1	D	152	ALA	2.7
1	D	121	THR	2.7
2	P	62	GLN	2.6
3	T	719	DG	2.6
1	D	153	GLY	2.6
1	A	289	LEU	2.6
1	D	229	ARG	2.6
4	E	722	DA	2.6
1	D	322	GLY	2.6
1	D	49	ARG	2.6
1	D	80	HIS	2.6
1	D	276	ALA	2.6
1	D	333	ILE	2.6
2	P	64	LEU	2.6
1	A	43	LEU	2.6
1	A	85	GLN	2.6
1	A	164	GLY	2.6
1	A	71	VAL	2.5
1	A	410	LEU	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	48	GLN	2.5
1	D	20	ALA	2.5
1	A	98	VAL	2.5
1	D	321	ALA	2.5
3	T	718	DA	2.5
1	A	19	LEU	2.5
1	A	433	LEU	2.5
1	D	26	PHE	2.5
2	B	69	THR	2.5
1	A	148	ARG	2.5
1	A	92	GLU	2.5
1	A	34	ALA	2.5
1	A	100	ALA	2.4
1	A	307	ALA	2.4
1	D	277	ARG	2.4
1	A	336	ALA	2.4
2	B	23	GLN	2.4
2	B	39	GLN	2.4
1	A	94	GLY	2.4
1	D	264	SER	2.4
1	A	147	PHE	2.4
1	D	98	VAL	2.4
1	D	213	LEU	2.4
1	D	232	ALA	2.4
1	A	30	PRO	2.3
2	B	66	LEU	2.3
1	A	132	TYR	2.3
1	A	51	ASN	2.3
2	P	24	ASN	2.3
1	A	356	LEU	2.3
1	D	302	GLN	2.3
4	E	715	DG	2.3
1	A	317	VAL	2.3
1	D	149	ILE	2.3
1	A	228	VAL	2.2
1	D	71	VAL	2.2
1	A	105	PRO	2.2
4	E	710	DA	2.2
2	P	26	TRP	2.2
1	A	129	LEU	2.2
1	D	111	THR	2.2
1	A	199	SER	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	338	PRO	2.2
2	P	35	ILE	2.2
2	P	44	ASN	2.2
2	P	71	CYS	2.2
1	A	80	HIS	2.2
1	A	413	ASP	2.2
1	D	159	ALA	2.2
1	A	35	SER	2.2
1	A	54	SER	2.2
1	A	82	LYS	2.2
1	A	225	ALA	2.1
1	D	46	PRO	2.1
1	D	323	GLY	2.1
1	A	15	GLU	2.1
1	D	216	ASN	2.1
2	B	34	LYS	2.1
2	P	10	PRO	2.1
1	A	95	ARG	2.1
1	A	13	VAL	2.1
1	A	208	LEU	2.1
1	A	121	THR	2.1
1	D	416	GLU	2.1
1	D	225	ALA	2.1
1	D	275	ILE	2.1
1	D	309	GLN	2.1
1	D	414	PHE	2.1
1	A	253	GLU	2.1
1	D	29	ALA	2.1
1	D	418	VAL	2.1
1	A	355	GLY	2.0
1	A	107	ASP	2.0
1	D	146	ASP	2.0
1	D	428	ARG	2.0
1	A	50	GLY	2.0
1	A	347	ILE	2.0
1	A	64	LEU	2.0
2	P	59	LYS	2.0
1	A	7	TRP	2.0
1	A	152	ALA	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.