



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 10:22 PM UTC

PDB ID : 3C60 / pdb_00003c60
Title : Crystal structure of mouse MHC class II I-Ab/3K peptide complexed with mouse TCR YAc62
Authors : Dai, S.
Deposited on : 2008-02-01
Resolution : 3.05 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

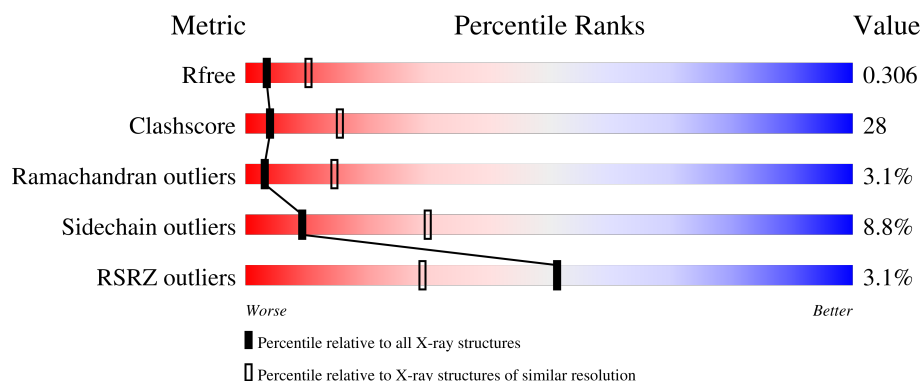
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.05 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	2469 (3.10-3.02)
Clashscore	190562	2569 (3.10-3.02)
Ramachandran outliers	187476	2424 (3.10-3.02)
Sidechain outliers	187428	2423 (3.10-3.02)
RSRZ outliers	180081	2469 (3.10-3.02)

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	199	5% 47% 45% 8%
1	E	199	6% 48% 45% 8%
2	B	236	5% 51% 41% 8%
2	F	236	4% 48% 43% 8%
3	C	182	2% 59% 36% ...

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Mol	Chain	Length	Quality of chain
3	G	182	57%38%5%
4	D	217	% 38%47%7%7%
4	H	217	% 39%46%8%7%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 13102 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 TCR YAc62 alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	199	Total	C	N	O	S	0	0	0
			1554	971	252	323	8			
1	E	199	Total	C	N	O	S	0	0	0
			1554	971	252	323	8			

- Molecule 2 is a protein called TCR YAc62 beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	236	Total	C	N	O	S	0	0	0
			1869	1177	325	361	6			
2	F	236	Total	C	N	O	S	0	0	0
			1869	1177	325	361	6			

- Molecule 3 is a protein called H-2 class II histocompatibility antigen, A-B alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	182	Total	C	N	O	S	0	0	0
			1459	944	230	282	3			
3	G	182	Total	C	N	O	S	0	0	0
			1459	944	230	282	3			

- Molecule 4 is a protein called 3K peptide, Linker, and H-2 class II histocompatibility antigen (A beta chain).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	201	Total	C	N	O	S	0	0	0
			1669	1045	303	314	7			
4	H	201	Total	C	N	O	S	0	0	0
			1669	1045	303	314	7			

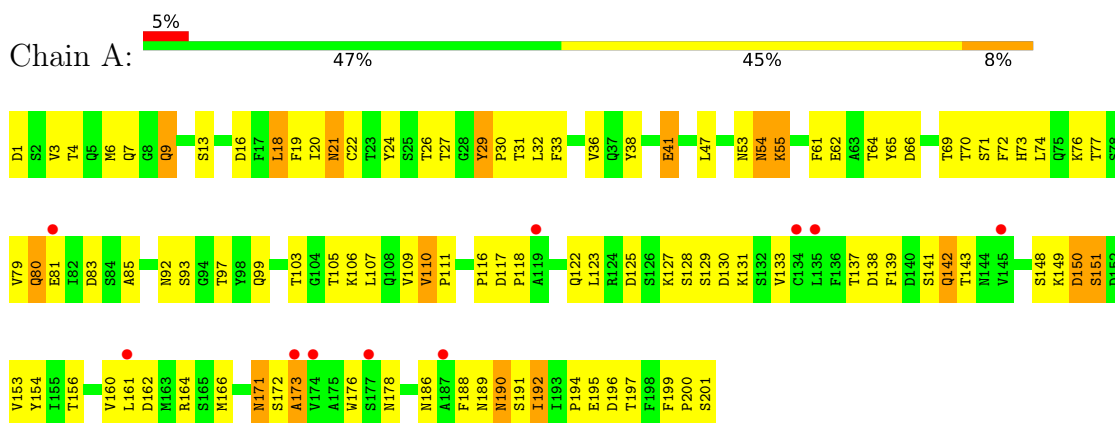
There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	14	GLY	-	linker	UNP P14483
D	15	GLY	-	linker	UNP P14483
D	16	GLY	-	linker	UNP P14483
D	17	GLY	-	linker	UNP P14483
D	18	SER	-	linker	UNP P14483
D	19	LEU	-	linker	UNP P14483
D	20	VAL	-	linker	UNP P14483
D	21	PRO	-	linker	UNP P14483
D	22	ARG	-	linker	UNP P14483
D	23	GLY	-	linker	UNP P14483
D	24	SER	-	linker	UNP P14483
D	25	GLY	-	linker	UNP P14483
D	26	GLY	-	linker	UNP P14483
D	27	GLY	-	linker	UNP P14483
D	28	GLY	-	linker	UNP P14483
D	216	LYS	ARG	engineered mutation	UNP P14483
H	14	GLY	-	linker	UNP P14483
H	15	GLY	-	linker	UNP P14483
H	16	GLY	-	linker	UNP P14483
H	17	GLY	-	linker	UNP P14483
H	18	SER	-	linker	UNP P14483
H	19	LEU	-	linker	UNP P14483
H	20	VAL	-	linker	UNP P14483
H	21	PRO	-	linker	UNP P14483
H	22	ARG	-	linker	UNP P14483
H	23	GLY	-	linker	UNP P14483
H	24	SER	-	linker	UNP P14483
H	25	GLY	-	linker	UNP P14483
H	26	GLY	-	linker	UNP P14483
H	27	GLY	-	linker	UNP P14483
H	28	GLY	-	linker	UNP P14483
H	216	LYS	ARG	engineered mutation	UNP P14483

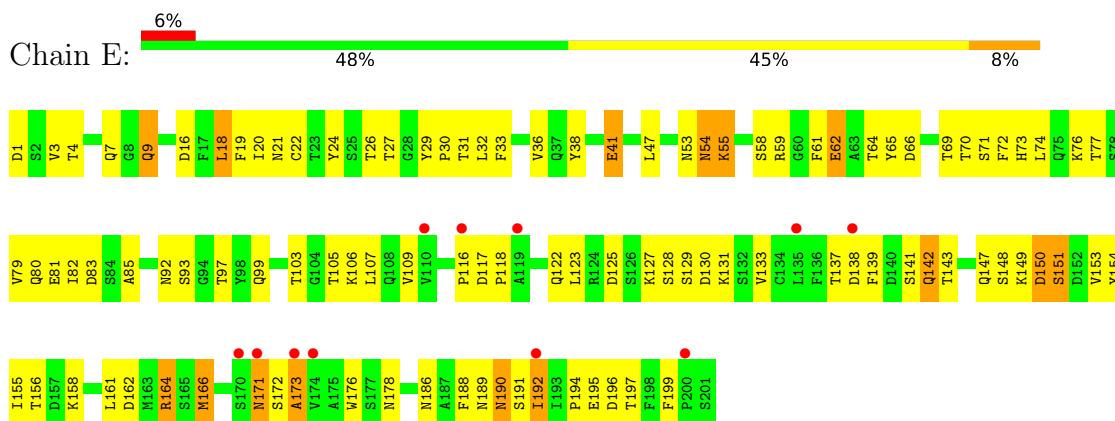
3 Residue-property plots

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.

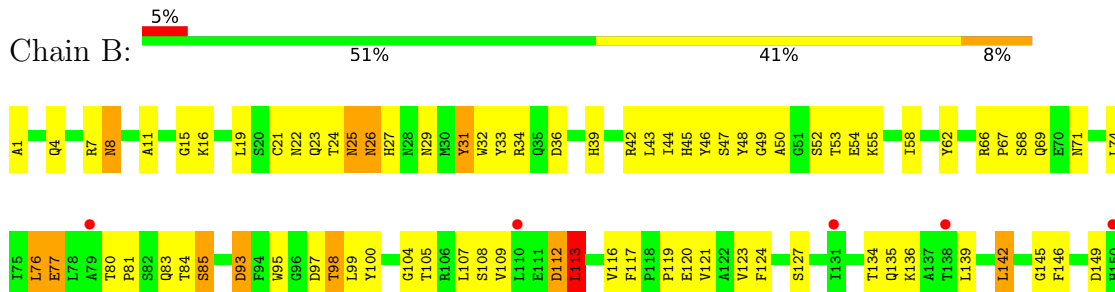
• Molecule 1: TCR YAE62 alpha chain

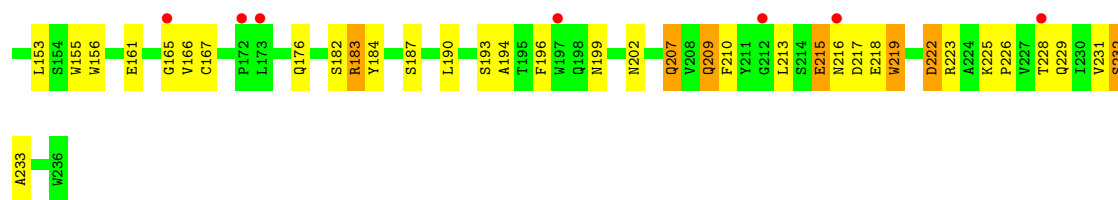


• Molecule 1: TCR YAE62 alpha chain

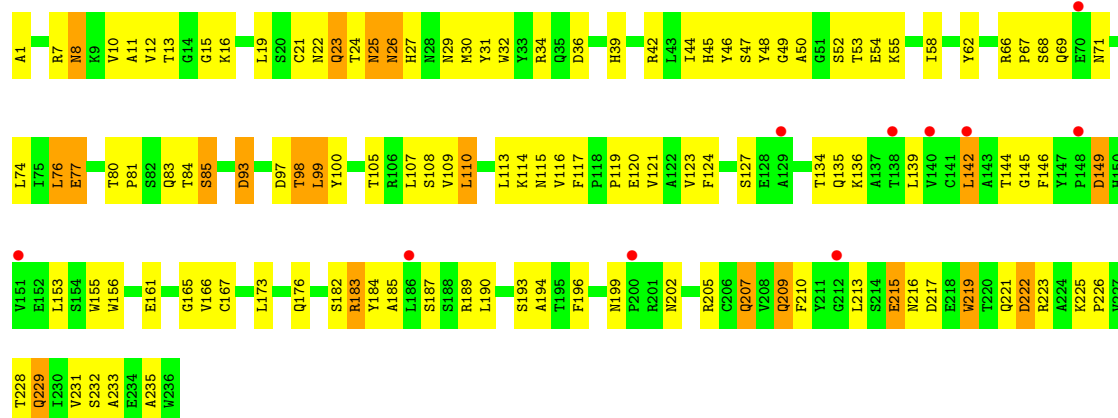


• Molecule 2: TCR YAE62 beta chain

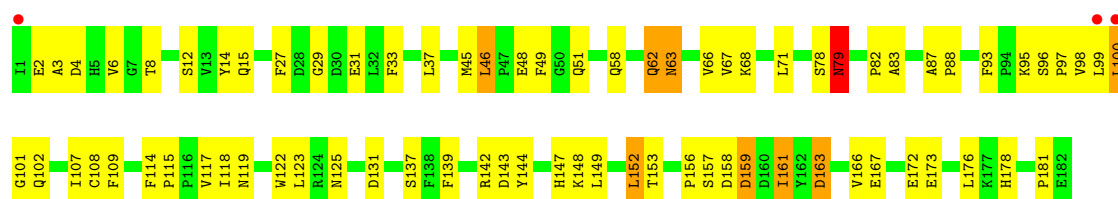




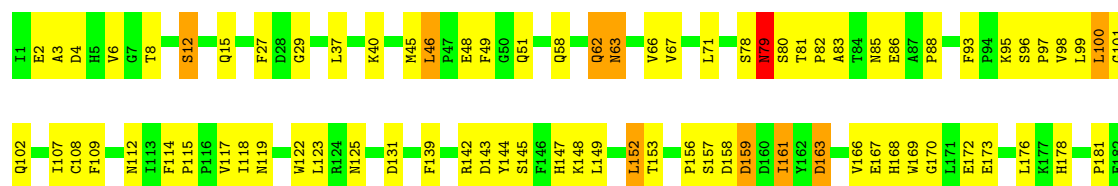
• Molecule 2: TCR YAc62 beta chain



• Molecule 3: H-2 class II histocompatibility antigen, A-B alpha chain

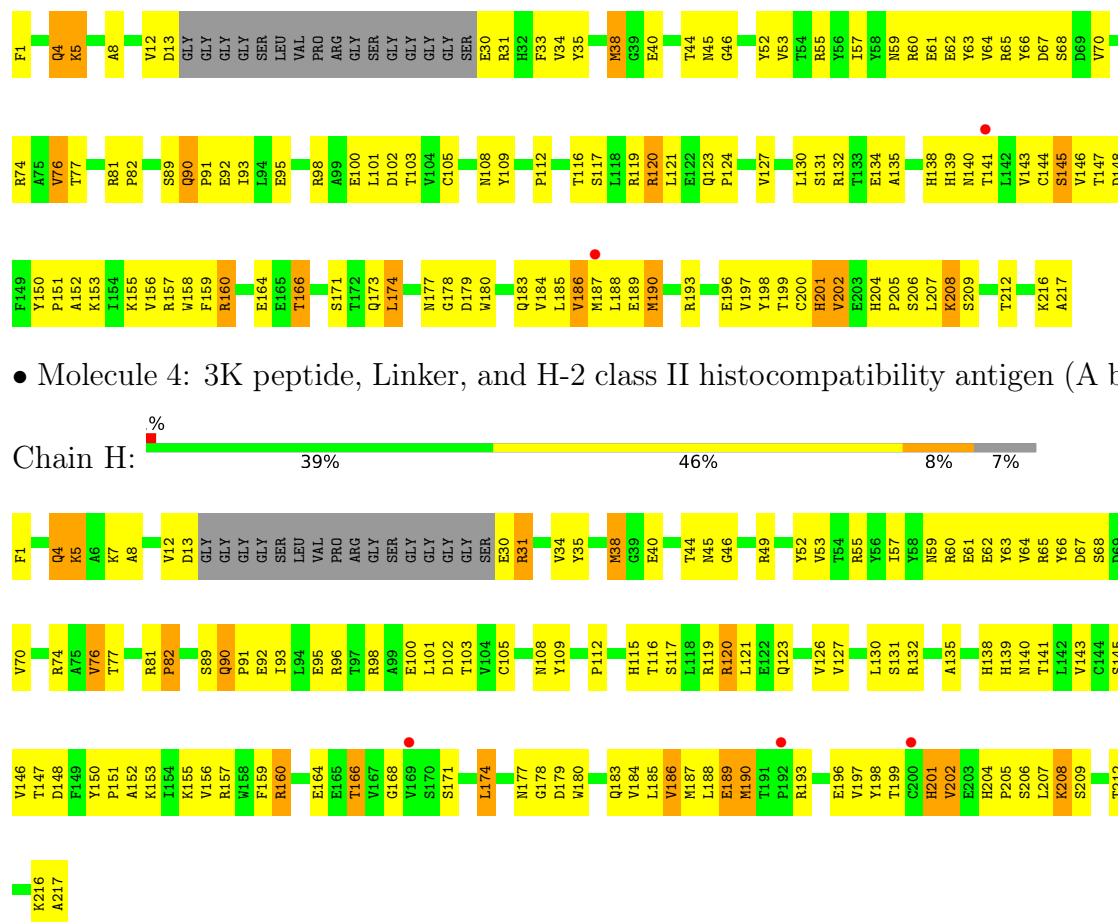


• Molecule 3: H-2 class II histocompatibility antigen, A-B alpha chain



• Molecule 4: 3K peptide, Linker, and H-2 class II histocompatibility antigen (A beta chain)





- Molecule 4: 3K peptide, Linker, and H-2 class II histocompatibility antigen (A beta chain)

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	66.06Å 126.17Å 277.18Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.92 – 3.05 40.92 – 3.05	Depositor EDS
% Data completeness (in resolution range)	92.6 (40.92-3.05) 92.6 (40.92-3.05)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.69 (at 2.90Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.267 , 0.309 0.266 , 0.306	Depositor DCC
R_{free} test set	2372 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	73.1	Xtriage
Anisotropy	0.307	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 61.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	13102	wwPDB-VP
Average B, all atoms (Å ²)	78.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.51% 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 [i](#)

5.1 Standard geometry [i](#)

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	0.54	0/1588	0.92	7/2157 (0.3%)
1	E	0.53	0/1588	0.92	6/2157 (0.3%)
2	B	0.53	0/1922	0.86	1/2620 (0.0%)
2	F	0.50	0/1922	0.84	0/2620
3	C	0.69	0/1504	0.96	0/2054
3	G	0.66	0/1504	0.98	4/2054 (0.2%)
4	D	0.58	0/1709	0.91	5/2316 (0.2%)
4	H	0.57	0/1709	0.91	4/2316 (0.2%)
All	All	0.57	0/13446	0.91	27/18294 (0.1%)

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
3	C	0	1
3	G	0	1
All	All	0	2

There are no bond length outliers.

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	29	TYR	CA-C-N	6.48	126.44	120.03
1	E	29	TYR	C-N-CA	6.48	126.44	120.03
1	A	29	TYR	CA-C-N	6.28	126.25	120.03
1	A	29	TYR	C-N-CA	6.28	126.25	120.03
1	A	55	LYS	N-CA-C	5.56	119.12	110.17
1	E	55	LYS	N-CA-C	5.51	119.05	110.17
4	D	57	ILE	N-CA-C	5.42	115.70	108.11
1	A	27	THR	N-CA-C	-5.42	105.48	113.61

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	232	SER	N-CA-C	5.41	115.38	108.24
3	G	80	SER	N-CA-C	5.36	118.60	111.30
1	E	27	THR	N-CA-C	-5.31	105.64	113.61
4	H	57	ILE	N-CA-C	5.28	115.50	108.11
4	D	145	SER	N-CA-C	5.26	117.20	107.73
4	D	202	VAL	N-CA-C	5.25	115.83	108.17
1	E	31	THR	N-CA-C	-5.18	101.26	109.76
3	G	81	THR	CA-C-N	5.16	125.14	120.03
3	G	81	THR	C-N-CA	5.16	125.14	120.03
1	A	110	VAL	N-CA-C	5.16	113.82	109.02
1	A	128	SER	CB-CA-C	-5.12	109.00	116.53
4	H	202	VAL	N-CA-C	5.12	115.64	108.17
1	A	31	THR	N-CA-C	-5.11	101.38	109.76
4	H	90	GLN	CA-C-N	5.11	125.39	119.47
4	H	90	GLN	C-N-CA	5.11	125.39	119.47
1	E	128	SER	CB-CA-C	-5.09	109.05	116.53
4	D	90	GLN	CA-C-N	5.03	125.31	119.47
4	D	90	GLN	C-N-CA	5.03	125.31	119.47
3	G	145	SER	N-CA-C	-5.02	102.15	109.63

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	C	144	TYR	Sidechain
3	G	144	TYR	Sidechain

5.2 Too-close contacts [i](#)

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	1554	0	1454	81	0
1	E	1554	0	1454	78	0
2	B	1869	0	1763	98	0
2	F	1869	0	1763	121	0
3	C	1459	0	1386	70	0
3	G	1459	0	1386	69	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	1669	0	1600	127	0
4	H	1669	0	1600	126	0
All	All	13102	0	12406	711	0

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 28.

All (711) 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)
4:D:208:LYS:HE3	4:D:208:LYS:H	1.14	1.12
4:H:208:LYS:H	4:H:208:LYS:HE3	1.18	1.09
4:D:174:LEU:HD13	4:D:174:LEU:H	1.34	0.93
2:F:48:TYR:CE1	3:G:62:GLN:HG2	2.05	0.91
2:F:16:LYS:HD2	2:F:77:GLU:HA	1.51	0.90
2:B:16:LYS:HD2	2:B:77:GLU:HA	1.53	0.90
2:B:48:TYR:CE1	3:C:62:GLN:HG2	2.06	0.90
2:F:113:LEU:HD21	2:F:213:LEU:HD21	1.54	0.89
4:H:174:LEU:HD13	4:H:174:LEU:H	1.35	0.88
4:H:31:ARG:H	4:H:31:ARG:HD3	1.37	0.87
4:H:112:PRO:O	4:H:116:THR:HG22	1.75	0.86
4:D:112:PRO:O	4:D:116:THR:HG22	1.77	0.83
4:D:138:HIS:CE1	3:G:142:ARG:HH22	1.97	0.82
3:G:172:GLU:CD	3:G:172:GLU:H	1.89	0.81
1:E:18:LEU:HB3	1:E:76:LYS:HB3	1.62	0.80
1:A:18:LEU:HB3	1:A:76:LYS:HB3	1.63	0.80
2:B:80:THR:HB	2:B:81:PRO:HD2	1.63	0.79
4:D:208:LYS:HE3	4:D:208:LYS:N	1.96	0.79
2:F:80:THR:HB	2:F:81:PRO:HD2	1.62	0.79
1:A:97:THR:HG22	3:C:62:GLN:HG3	1.64	0.79
1:E:30:PRO:HB3	1:E:93:SER:HB3	1.66	0.78
2:B:113:LEU:HD11	2:B:213:LEU:HD21	1.64	0.77
1:A:30:PRO:HB3	1:A:93:SER:HB3	1.67	0.77
2:B:123:VAL:HG23	2:B:233:ALA:HB3	1.67	0.76
4:H:143:VAL:HG22	4:H:187:MET:HG2	1.68	0.76
4:D:143:VAL:HG22	4:D:187:MET:HG2	1.66	0.76
3:C:172:GLU:CD	3:C:172:GLU:H	1.94	0.76
3:G:118:ILE:HD11	3:G:166:VAL:HG13	1.68	0.76
2:B:84:THR:HG23	2:B:108:SER:HA	1.68	0.74
4:H:159:PHE:HB2	4:H:199:THR:HG23	1.70	0.74
4:D:208:LYS:H	4:D:208:LYS:CE	1.97	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:159:PHE:HB2	4:D:199:THR:HG23	1.70	0.74
1:E:83:ASP:O	1:E:107:LEU:HD23	1.88	0.73
2:B:209:GLN:HE21	2:B:209:GLN:C	1.96	0.73
1:E:97:THR:HG22	3:G:62:GLN:HG3	1.71	0.73
4:D:166:THR:HG21	4:H:156:VAL:H	1.53	0.73
1:E:54:ASN:N	1:E:54:ASN:HD22	1.87	0.73
4:H:208:LYS:H	4:H:208:LYS:CE	2.00	0.73
4:H:63:TYR:CD2	4:H:64:VAL:HG23	2.24	0.72
4:H:62:GLU:HG2	4:H:76:VAL:HG11	1.72	0.72
4:D:157:ARG:HB2	4:D:201:HIS:HD2	1.53	0.72
3:G:12:SER:HB2	3:G:67:VAL:HG21	1.72	0.72
4:D:63:TYR:CD2	4:D:64:VAL:HG23	2.24	0.72
3:C:118:ILE:HD11	3:C:166:VAL:HG13	1.70	0.72
2:F:84:THR:HG23	2:F:108:SER:HA	1.72	0.72
4:H:157:ARG:HB2	4:H:201:HIS:HD2	1.55	0.71
1:A:81:GLU:OE2	1:A:111:PRO:HG3	1.91	0.71
4:D:120:ARG:HG3	4:D:120:ARG:HH11	1.55	0.71
4:H:208:LYS:HE3	4:H:208:LYS:N	2.00	0.71
1:A:54:ASN:N	1:A:54:ASN:HD22	1.87	0.70
1:E:161:LEU:HD13	1:E:162:ASP:N	2.08	0.69
4:D:116:THR:HG23	4:D:117:SER:N	2.08	0.69
3:C:161:ILE:HD13	3:C:161:ILE:H	1.58	0.68
1:E:192:ILE:HD12	1:E:192:ILE:H	1.59	0.68
1:A:83:ASP:O	1:A:107:LEU:HD23	1.93	0.68
4:D:120:ARG:O	4:D:121:LEU:HD12	1.94	0.68
2:F:48:TYR:CD1	3:G:62:GLN:NE2	2.61	0.68
3:C:12:SER:HB2	3:C:67:VAL:HG21	1.76	0.68
2:F:209:GLN:HE21	2:F:209:GLN:C	2.02	0.68
1:A:111:PRO:HG2	1:A:160:VAL:HG11	1.76	0.68
3:G:161:ILE:HD13	3:G:161:ILE:H	1.58	0.68
4:D:116:THR:HG23	4:D:117:SER:H	1.59	0.67
4:D:156:VAL:HG22	4:D:202:VAL:HG12	1.77	0.67
4:H:120:ARG:O	4:H:121:LEU:HD12	1.95	0.66
4:D:31:ARG:HD2	4:D:33:PHE:HZ	1.59	0.66
4:D:156:VAL:HG13	4:D:202:VAL:HG12	1.77	0.66
4:H:116:THR:HG23	4:H:117:SER:N	2.11	0.66
4:D:62:GLU:HG2	4:D:76:VAL:HG11	1.78	0.66
1:A:161:LEU:HD13	1:A:162:ASP:N	2.11	0.66
1:A:192:ILE:HD12	1:A:192:ILE:H	1.60	0.66
2:F:123:VAL:HG11	2:F:235:ALA:HB2	1.79	0.65
3:G:148:LYS:C	3:G:149:LEU:HD22	2.22	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:156:VAL:HG22	4:H:202:VAL:HG12	1.79	0.65
2:F:48:TYR:OH	3:G:58:GLN:NE2	2.29	0.65
4:H:156:VAL:HG13	4:H:202:VAL:HG12	1.78	0.64
4:D:38:MET:HE2	4:D:40:GLU:OE1	1.97	0.64
4:H:120:ARG:HG3	4:H:120:ARG:HH11	1.62	0.64
1:E:122:GLN:O	2:F:127:SER:HB3	1.96	0.64
2:B:84:THR:HG23	2:B:107:LEU:O	1.97	0.64
4:D:12:VAL:HG12	4:D:13:ASP:N	2.13	0.64
3:C:125:ASN:OD1	3:C:161:ILE:HD13	1.98	0.64
1:A:38:TYR:O	1:A:41:GLU:HG3	1.98	0.64
1:A:122:GLN:C	1:A:123:LEU:HD12	2.23	0.64
4:D:159:PHE:HE2	4:D:164:GLU:HB2	1.63	0.63
3:C:15:GLN:NE2	3:C:115:PRO:HB2	2.13	0.63
4:D:193:ARG:HB2	4:D:196:GLU:CD	2.24	0.63
3:G:15:GLN:HE22	3:G:117:VAL:HG23	1.62	0.63
4:H:12:VAL:HG12	4:H:13:ASP:N	2.12	0.63
2:B:76:LEU:HD23	2:B:83:GLN:OE1	1.98	0.63
4:D:67:ASP:HB3	4:D:70:VAL:HG23	1.80	0.63
4:H:12:VAL:HG12	4:H:13:ASP:H	1.63	0.63
2:B:176:GLN:HB2	2:B:182:SER:CB	2.28	0.63
2:F:84:THR:HG23	2:F:107:LEU:O	1.99	0.63
4:H:4:GLN:HG2	4:H:103:THR:O	1.98	0.63
1:E:122:GLN:C	1:E:123:LEU:HD12	2.24	0.62
1:A:122:GLN:O	2:B:127:SER:HB3	1.99	0.62
1:E:38:TYR:O	1:E:41:GLU:HG3	1.99	0.62
4:D:108:ASN:HA	4:D:112:PRO:HD2	1.82	0.62
4:H:159:PHE:HE2	4:H:164:GLU:HB2	1.65	0.62
2:F:176:GLN:HB2	2:F:182:SER:CB	2.30	0.62
1:E:81:GLU:HA	1:E:109:VAL:HB	1.83	0.61
2:F:7:ARG:O	2:F:105:THR:HA	2.01	0.61
4:H:38:MET:HE2	4:H:40:GLU:OE1	2.00	0.61
4:H:116:THR:HG23	4:H:117:SER:H	1.65	0.61
2:B:29:ASN:N	2:B:29:ASN:HD22	1.98	0.61
1:E:131:LYS:NZ	1:E:178:ASN:HA	2.16	0.61
2:F:76:LEU:HD23	2:F:83:GLN:OE1	2.01	0.61
3:G:95:LYS:O	4:H:183:GLN:NE2	2.33	0.61
1:E:154:TYR:HD2	1:E:176:TRP:NE1	1.99	0.61
2:F:123:VAL:HG23	2:F:233:ALA:HB3	1.83	0.61
4:D:164:GLU:HB3	4:H:155:LYS:NZ	2.16	0.61
4:D:174:LEU:HD22	4:D:174:LEU:N	2.13	0.61
4:D:12:VAL:HG12	4:D:13:ASP:H	1.66	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:48:GLU:HG3	3:G:49:PHE:N	2.15	0.61
2:F:113:LEU:CD2	2:F:213:LEU:HD21	2.28	0.60
2:B:7:ARG:O	2:B:105:THR:HA	2.02	0.60
2:F:93:ASP:OD2	2:F:93:ASP:N	2.26	0.60
4:H:171:SER:HB3	4:H:186:VAL:HG13	1.84	0.60
4:H:67:ASP:HB3	4:H:70:VAL:HG23	1.83	0.60
1:A:47:LEU:C	1:A:47:LEU:HD12	2.27	0.60
4:D:141:THR:HG21	4:D:187:MET:HE2	1.84	0.60
4:D:160:ARG:HG3	4:D:198:TYR:CE2	2.36	0.60
3:G:125:ASN:OD1	3:G:161:ILE:HD13	2.01	0.60
4:H:160:ARG:HG3	4:H:198:TYR:CE2	2.37	0.59
4:H:108:ASN:HA	4:H:112:PRO:HD2	1.84	0.59
4:H:55:ARG:HG2	4:H:55:ARG:HH11	1.67	0.59
1:A:149:LYS:O	1:A:149:LYS:HD3	2.02	0.59
1:E:47:LEU:HD12	1:E:47:LEU:C	2.28	0.59
4:H:193:ARG:HB2	4:H:196:GLU:CD	2.27	0.59
4:D:155:LYS:NZ	4:H:164:GLU:HB3	2.18	0.59
2:F:54:GLU:OE2	3:G:40:LYS:HD2	2.03	0.59
2:B:15:GLY:O	2:B:77:GLU:O	2.21	0.59
4:H:141:THR:HG21	4:H:187:MET:HE2	1.84	0.59
1:A:154:TYR:HD2	1:A:176:TRP:NE1	2.00	0.59
1:A:131:LYS:NZ	1:A:178:ASN:HA	2.18	0.58
4:D:171:SER:HB3	4:D:186:VAL:HG13	1.85	0.58
2:F:12:VAL:HA	2:F:110:LEU:O	2.03	0.58
4:D:147:THR:HG22	4:D:183:GLN:HG3	1.85	0.58
4:D:197:VAL:HG22	4:D:216:LYS:HA	1.84	0.58
4:H:63:TYR:C	4:H:76:VAL:HG13	2.28	0.58
4:H:174:LEU:N	4:H:174:LEU:HD22	2.18	0.58
4:H:147:THR:HG22	4:H:183:GLN:HG3	1.85	0.58
1:A:81:GLU:HA	1:A:109:VAL:HB	1.86	0.58
4:D:63:TYR:C	4:D:76:VAL:HG13	2.29	0.58
4:D:89:SER:O	4:D:91:PRO:HD3	2.03	0.58
3:G:85:ASN:ND2	3:G:169:TRP:HB2	2.19	0.58
1:A:1:ASP:HA	1:A:26:THR:HG22	1.86	0.57
1:A:186:ASN:HA	1:A:189:ASN:ND2	2.19	0.57
2:B:48:TYR:OH	3:C:58:GLN:NE2	2.34	0.57
3:G:15:GLN:NE2	3:G:115:PRO:HB2	2.19	0.57
1:E:149:LYS:O	1:E:149:LYS:HD3	2.03	0.57
4:D:159:PHE:CB	4:D:199:THR:HG23	2.35	0.57
4:D:156:VAL:H	4:H:166:THR:HG21	1.70	0.57
2:F:16:LYS:CD	2:F:77:GLU:HA	2.29	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:5:LYS:HD2	4:H:5:LYS:O	2.04	0.57
4:D:65:ARG:HG2	4:D:66:TYR:N	2.20	0.57
1:A:9:GLN:HE21	1:A:9:GLN:C	2.12	0.56
3:C:48:GLU:HG3	3:C:49:PHE:N	2.19	0.56
2:B:22:ASN:HD22	2:B:71:ASN:ND2	2.03	0.56
2:B:209:GLN:NE2	2:B:210:PHE:N	2.53	0.56
3:C:82:PRO:HB3	4:D:31:ARG:HB2	1.87	0.56
4:D:119:ARG:HB2	4:D:119:ARG:HH11	1.71	0.56
1:E:186:ASN:HA	1:E:189:ASN:ND2	2.20	0.56
2:F:22:ASN:HD22	2:F:71:ASN:ND2	2.04	0.56
4:H:65:ARG:HG2	4:H:66:TYR:N	2.20	0.56
3:G:83:ALA:HB2	4:H:59:ASN:OD1	2.05	0.56
2:B:99:LEU:O	2:B:99:LEU:HD12	2.06	0.56
2:F:135:GLN:C	2:F:194:ALA:HB2	2.30	0.56
4:H:119:ARG:HH11	4:H:119:ARG:HB2	1.70	0.56
2:B:48:TYR:CD1	3:C:62:GLN:NE2	2.73	0.56
1:E:190:ASN:N	1:E:190:ASN:HD22	2.03	0.56
4:H:44:THR:O	4:H:45:ASN:HB3	2.06	0.56
4:H:52:TYR:CD1	4:H:105:CYS:SG	2.98	0.56
1:E:1:ASP:HA	1:E:26:THR:HG22	1.87	0.56
1:A:200:PRO:O	1:A:201:SER:HB2	2.05	0.56
4:D:130:LEU:HD23	4:D:130:LEU:H	1.71	0.56
3:C:157:SER:HB2	3:C:159:ASP:OD2	2.06	0.56
2:F:29:ASN:N	2:F:29:ASN:HD22	2.01	0.56
3:C:96:SER:HB2	3:C:97:PRO:HD2	1.88	0.55
1:E:9:GLN:HG2	1:E:106:LYS:HB3	1.89	0.55
2:F:76:LEU:N	2:F:76:LEU:HD12	2.20	0.55
3:G:118:ILE:HG12	3:G:119:ASN:N	2.20	0.55
4:H:156:VAL:HG11	4:H:186:VAL:HG21	1.88	0.55
1:A:9:GLN:HG2	1:A:106:LYS:HB3	1.87	0.55
1:A:190:ASN:N	1:A:190:ASN:HD22	2.01	0.55
3:C:63:ASN:HD21	4:D:8:ALA:H	1.54	0.55
2:F:123:VAL:HG23	2:F:233:ALA:CB	2.36	0.55
2:F:156:TRP:NE1	2:F:207:GLN:NE2	2.54	0.55
3:G:114:PHE:HA	3:G:115:PRO:C	2.31	0.55
4:H:159:PHE:CB	4:H:199:THR:HG23	2.36	0.55
2:B:113:LEU:CD1	2:B:213:LEU:HD11	2.36	0.55
3:G:67:VAL:HG13	4:H:35:TYR:CD2	2.42	0.55
2:F:121:VAL:HG23	2:F:231:VAL:CG1	2.37	0.55
4:H:197:VAL:HG22	4:H:216:LYS:HA	1.88	0.55
2:B:49:GLY:O	2:B:52:SER:HB3	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:113:LEU:HD11	2:B:213:LEU:HD11	1.87	0.55
3:G:45:MET:HE3	4:H:178:GLY:O	2.06	0.55
3:G:63:ASN:HD21	4:H:8:ALA:H	1.53	0.55
2:B:16:LYS:CD	2:B:77:GLU:HA	2.29	0.55
2:B:84:THR:O	2:B:85:SER:HB3	2.07	0.54
2:B:199:ASN:HB3	2:B:202:ASN:ND2	2.22	0.54
3:C:62:GLN:O	3:C:66:VAL:HG23	2.08	0.54
4:D:116:THR:CG2	4:D:117:SER:H	2.19	0.54
2:F:99:LEU:HD12	2:F:99:LEU:O	2.08	0.54
4:H:108:ASN:O	4:H:112:PRO:HD2	2.07	0.54
2:B:156:TRP:NE1	2:B:207:GLN:NE2	2.55	0.54
3:C:15:GLN:HE22	3:C:117:VAL:HG23	1.72	0.54
3:C:119:ASN:HB2	3:C:167:GLU:HB3	1.88	0.54
1:E:55:LYS:HB3	1:E:64:THR:OG1	2.07	0.54
1:E:192:ILE:HD12	1:E:192:ILE:N	2.22	0.54
1:A:192:ILE:HD12	1:A:192:ILE:N	2.22	0.54
2:B:135:GLN:C	2:B:194:ALA:HB2	2.32	0.54
3:C:67:VAL:HG13	4:D:35:TYR:CD2	2.42	0.54
3:C:118:ILE:HG12	3:C:119:ASN:N	2.23	0.54
3:C:148:LYS:C	3:C:149:LEU:HD22	2.32	0.54
4:D:159:PHE:CE2	4:D:164:GLU:HB2	2.42	0.54
4:H:130:LEU:HD23	4:H:130:LEU:H	1.72	0.54
4:H:157:ARG:HD3	4:H:164:GLU:CD	2.32	0.54
4:D:120:ARG:HG3	4:D:120:ARG:NH1	2.22	0.54
3:C:83:ALA:HB2	4:D:59:ASN:OD1	2.08	0.54
2:F:199:ASN:HB3	2:F:202:ASN:ND2	2.22	0.54
1:E:7:GLN:HG3	1:E:20:ILE:HG23	1.90	0.54
4:H:52:TYR:H	4:H:68:SER:CB	2.21	0.53
4:H:132:ARG:H	4:H:140:ASN:HD21	1.56	0.53
2:B:134:THR:HB	2:B:136:LYS:HE3	1.88	0.53
1:A:7:GLN:HG3	1:A:20:ILE:HG23	1.90	0.53
3:G:157:SER:HB2	3:G:159:ASP:OD2	2.08	0.53
4:D:164:GLU:O	4:H:155:LYS:HE2	2.09	0.53
1:A:143:THR:HG21	1:A:194:PRO:HD3	1.90	0.53
3:G:98:VAL:C	3:G:99:LEU:HD22	2.33	0.53
2:B:116:VAL:HG12	2:B:226:PRO:HB2	1.91	0.53
3:C:98:VAL:C	3:C:99:LEU:HD22	2.34	0.53
4:D:52:TYR:CD1	4:D:105:CYS:SG	3.02	0.53
4:D:204:HIS:ND1	4:D:205:PRO:HD2	2.23	0.53
4:H:116:THR:CG2	4:H:117:SER:H	2.22	0.53
1:A:47:LEU:HD12	1:A:47:LEU:O	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:44:THR:O	4:D:45:ASN:HB3	2.09	0.53
4:H:76:VAL:HG22	4:H:77:THR:HG23	1.89	0.53
4:H:190:MET:C	4:H:190:MET:HE3	2.34	0.53
1:A:29:TYR:CD2	4:D:103:THR:HG22	2.43	0.53
4:D:132:ARG:H	4:D:140:ASN:HD21	1.57	0.53
1:E:47:LEU:HD12	1:E:47:LEU:O	2.08	0.53
1:E:161:LEU:HD13	1:E:162:ASP:H	1.73	0.53
2:F:15:GLY:O	2:F:77:GLU:O	2.27	0.53
2:F:134:THR:HB	2:F:136:LYS:HE3	1.90	0.53
1:A:36:VAL:HG13	1:A:36:VAL:O	2.09	0.53
2:B:113:LEU:HD11	2:B:213:LEU:CD2	2.35	0.53
2:B:8:ASN:N	2:B:8:ASN:ND2	2.57	0.52
4:D:76:VAL:HG22	4:D:77:THR:HG23	1.91	0.52
4:H:53:VAL:HA	4:H:66:TYR:O	2.09	0.52
2:F:209:GLN:C	2:F:209:GLN:NE2	2.68	0.52
3:G:96:SER:HB2	3:G:97:PRO:HD2	1.90	0.52
2:B:219:TRP:HB2	2:B:225:LYS:HG2	1.92	0.52
4:D:52:TYR:H	4:D:68:SER:CB	2.23	0.52
4:D:150:TYR:CD1	4:D:151:PRO:HA	2.44	0.52
2:F:156:TRP:HE1	2:F:207:GLN:NE2	2.08	0.52
4:H:101:LEU:O	4:H:105:CYS:HB2	2.09	0.52
2:F:156:TRP:HE1	2:F:207:GLN:HE21	1.57	0.52
3:G:163:ASP:OD2	3:G:163:ASP:N	2.43	0.52
2:B:210:PHE:O	2:B:228:THR:HG23	2.09	0.52
4:D:188:LEU:HG	4:D:190:MET:HG3	1.92	0.52
1:E:79:VAL:HG13	1:E:79:VAL:O	2.09	0.52
3:C:98:VAL:O	3:C:99:LEU:HD13	2.09	0.52
2:F:84:THR:O	2:F:85:SER:HB3	2.10	0.52
3:G:86:GLU:O	3:G:170:GLY:HA3	2.10	0.52
3:G:98:VAL:O	3:G:99:LEU:HD13	2.09	0.52
4:H:188:LEU:HG	4:H:190:MET:HG3	1.91	0.52
3:G:100:LEU:HD12	3:G:100:LEU:H	1.74	0.52
2:B:117:PHE:O	2:B:146:PHE:HA	2.09	0.52
3:C:163:ASP:OD2	3:C:163:ASP:N	2.42	0.52
2:B:156:TRP:HE1	2:B:207:GLN:HE21	1.58	0.52
3:C:123:LEU:HB2	3:C:163:ASP:OD2	2.10	0.52
4:D:31:ARG:HD2	4:D:33:PHE:CZ	2.43	0.52
2:F:100:TYR:N	2:F:100:TYR:CD1	2.78	0.52
2:F:121:VAL:HG23	2:F:231:VAL:HG12	1.92	0.52
2:F:209:GLN:NE2	2:F:210:PHE:N	2.58	0.52
3:G:123:LEU:HB2	3:G:163:ASP:OD2	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:76:LEU:HD12	2:B:76:LEU:N	2.25	0.51
3:C:93:PHE:CE2	3:C:107:ILE:HD12	2.44	0.51
4:H:89:SER:O	4:H:91:PRO:HD3	2.09	0.51
2:B:95:TRP:N	3:C:62:GLN:OE1	2.35	0.51
2:B:155:TRP:O	2:B:161:GLU:HA	2.09	0.51
3:C:142:ARG:HH22	4:H:138:HIS:CE1	2.28	0.51
2:F:8:ASN:N	2:F:8:ASN:ND2	2.58	0.51
2:F:123:VAL:HG11	2:F:235:ALA:CB	2.39	0.51
2:F:119:PRO:HG2	2:F:231:VAL:HB	1.91	0.51
4:H:112:PRO:O	4:H:116:THR:CG2	2.55	0.51
4:H:204:HIS:ND1	4:H:205:PRO:HD2	2.23	0.51
4:D:116:THR:CG2	4:D:117:SER:N	2.73	0.51
4:D:157:ARG:HD3	4:D:164:GLU:CD	2.35	0.51
1:E:143:THR:HG21	1:E:194:PRO:HD3	1.92	0.51
4:D:112:PRO:O	4:D:116:THR:CG2	2.54	0.51
4:H:159:PHE:CE2	4:H:164:GLU:HB2	2.44	0.51
2:B:156:TRP:HE1	2:B:207:GLN:NE2	2.08	0.51
3:G:119:ASN:HB2	3:G:167:GLU:HB3	1.92	0.51
2:B:100:TYR:CD1	2:B:100:TYR:N	2.78	0.51
2:B:209:GLN:C	2:B:209:GLN:NE2	2.65	0.51
4:D:5:LYS:HD2	4:D:5:LYS:O	2.10	0.51
2:B:145:GLY:O	2:B:183:ARG:HB3	2.11	0.51
4:D:155:LYS:HE2	4:H:164:GLU:O	2.10	0.51
4:H:46:GLY:HA2	4:H:109:TYR:OH	2.10	0.51
2:B:80:THR:O	2:B:109:VAL:HG11	2.11	0.51
3:C:95:LYS:O	4:D:183:GLN:NE2	2.44	0.51
3:G:100:LEU:HD12	3:G:100:LEU:N	2.26	0.51
4:H:116:THR:CG2	4:H:117:SER:N	2.74	0.51
2:F:155:TRP:O	2:F:161:GLU:HA	2.11	0.50
2:F:219:TRP:HB2	2:F:225:LYS:HG2	1.92	0.50
1:A:79:VAL:O	1:A:79:VAL:HG13	2.11	0.50
1:A:176:TRP:CE3	2:B:142:LEU:HD11	2.46	0.50
2:F:219:TRP:HA	2:F:219:TRP:CE3	2.47	0.50
1:A:161:LEU:HD13	1:A:162:ASP:H	1.75	0.50
3:C:139:PHE:HB2	3:C:147:HIS:NE2	2.27	0.50
1:A:69:THR:O	1:A:69:THR:HG22	2.09	0.50
3:C:93:PHE:HE2	3:C:107:ILE:HD12	1.76	0.50
4:D:53:VAL:HA	4:D:66:TYR:O	2.11	0.50
4:D:159:PHE:HB2	4:D:199:THR:O	2.12	0.50
1:E:30:PRO:CB	1:E:93:SER:HB3	2.39	0.50
1:E:53:ASN:HA	1:E:65:TYR:O	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:176:TRP:CE3	2:F:142:LEU:HD11	2.47	0.50
3:G:142:ARG:HG3	3:G:143:ASP:N	2.26	0.50
4:H:204:HIS:CD2	4:H:206:SER:OG	2.65	0.50
3:C:100:LEU:HD12	3:C:100:LEU:H	1.75	0.50
2:F:119:PRO:O	2:F:231:VAL:HG11	2.12	0.50
4:H:152:ALA:O	4:H:153:LYS:C	2.54	0.50
3:C:15:GLN:HG2	4:D:34:VAL:HG22	1.94	0.49
4:D:196:GLU:O	4:D:217:ALA:HB3	2.12	0.49
1:E:156:THR:HG21	2:F:187:SER:OG	2.11	0.49
1:A:18:LEU:HD22	1:A:19:PHE:N	2.27	0.49
3:G:15:GLN:HG2	4:H:34:VAL:HG22	1.94	0.49
4:H:204:HIS:HD2	4:H:206:SER:OG	1.95	0.49
1:A:53:ASN:HA	1:A:65:TYR:O	2.12	0.49
2:B:219:TRP:HA	2:B:219:TRP:CE3	2.46	0.49
3:C:100:LEU:HD12	3:C:100:LEU:N	2.26	0.49
2:B:11:ALA:O	2:B:109:VAL:HA	2.11	0.49
3:C:45:MET:HE3	4:D:178:GLY:O	2.12	0.49
4:D:1:PHE:HD1	4:D:1:PHE:H3	1.61	0.49
4:D:59:ASN:O	4:D:61:GLU:N	2.46	0.49
4:D:101:LEU:O	4:D:105:CYS:HB2	2.11	0.49
4:D:164:GLU:HB3	4:H:155:LYS:HZ1	1.77	0.49
1:E:69:THR:HG22	1:E:69:THR:O	2.13	0.49
1:E:131:LYS:HZ3	1:E:178:ASN:HA	1.77	0.49
1:A:30:PRO:CB	1:A:93:SER:HB3	2.40	0.49
2:F:116:VAL:HG12	2:F:226:PRO:HB2	1.95	0.49
2:F:205:ARG:HE	2:F:232:SER:HB2	1.77	0.49
4:H:190:MET:HE3	4:H:190:MET:O	2.11	0.49
2:B:156:TRP:HA	2:B:161:GLU:HA	1.95	0.49
2:F:22:ASN:HB2	2:F:71:ASN:OD1	2.12	0.49
4:H:55:ARG:HG2	4:H:55:ARG:NH1	2.27	0.49
4:H:150:TYR:CD1	4:H:151:PRO:HA	2.48	0.49
4:H:179:ASP:O	4:H:180:TRP:HB2	2.13	0.49
3:C:142:ARG:HG3	3:C:143:ASP:N	2.27	0.49
4:D:108:ASN:O	4:D:112:PRO:HD2	2.13	0.49
3:C:67:VAL:HG13	4:D:35:TYR:HD2	1.78	0.49
1:E:149:LYS:HB2	1:E:190:ASN:OD1	2.12	0.49
2:F:50:ALA:HA	2:F:66:ARG:HG3	1.95	0.49
2:B:19:LEU:HD22	2:B:105:THR:HG21	1.95	0.48
2:F:210:PHE:O	2:F:228:THR:HG23	2.13	0.48
3:C:158:ASP:HA	3:C:181:PRO:HG2	1.94	0.48
1:E:9:GLN:HE21	1:E:9:GLN:C	2.20	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:54:GLU:HG2	3:G:58:GLN:NE2	2.27	0.48
3:G:158:ASP:HA	3:G:181:PRO:HG2	1.95	0.48
3:C:139:PHE:HB2	3:C:147:HIS:CE1	2.49	0.48
4:D:55:ARG:HH11	4:D:55:ARG:HG2	1.78	0.48
4:D:127:VAL:HG23	4:D:145:SER:HB2	1.95	0.48
2:F:166:VAL:HG12	2:F:167:CYS:N	2.28	0.48
4:H:120:ARG:HG3	4:H:120:ARG:NH1	2.26	0.48
3:G:93:PHE:CE2	3:G:107:ILE:HD12	2.48	0.48
3:G:172:GLU:CD	3:G:172:GLU:N	2.67	0.48
1:A:197:THR:HB	1:A:199:PHE:HE1	1.78	0.48
2:B:209:GLN:NE2	2:B:209:GLN:CA	2.77	0.48
4:D:152:ALA:O	4:D:153:LYS:C	2.57	0.48
1:E:24:TYR:C	1:E:24:TYR:CD1	2.91	0.48
1:E:61:PHE:HB3	1:E:74:LEU:HD11	1.94	0.48
3:G:71:LEU:HD13	4:H:35:TYR:HB2	1.95	0.48
1:A:156:THR:HG21	2:B:187:SER:OG	2.13	0.48
2:B:146:PHE:CE2	2:B:184:TYR:HB2	2.49	0.48
1:E:36:VAL:HG13	1:E:36:VAL:O	2.13	0.48
1:A:54:ASN:N	1:A:54:ASN:ND2	2.61	0.47
4:D:4:GLN:HG2	4:D:103:THR:O	2.14	0.47
1:E:3:VAL:HG12	1:E:4:THR:N	2.29	0.47
2:F:119:PRO:HG2	2:F:231:VAL:CG2	2.44	0.47
4:D:156:VAL:HG22	4:D:202:VAL:CG1	2.43	0.47
1:E:69:THR:O	1:E:70:THR:C	2.56	0.47
2:F:114:LYS:HE2	2:F:221:GLN:NE2	2.28	0.47
1:A:66:ASP:OD2	1:A:69:THR:OG1	2.26	0.47
2:B:62:TYR:HB3	2:B:74:LEU:HD11	1.96	0.47
2:B:176:GLN:HB2	2:B:182:SER:HB3	1.95	0.47
4:D:190:MET:HE3	4:D:190:MET:C	2.39	0.47
4:D:204:HIS:CD2	4:D:206:SER:OG	2.68	0.47
4:D:207:LEU:HD12	4:D:207:LEU:O	2.14	0.47
3:G:67:VAL:HG13	4:H:35:TYR:HD2	1.78	0.47
4:H:59:ASN:O	4:H:61:GLU:N	2.47	0.47
1:A:24:TYR:CD1	1:A:24:TYR:C	2.93	0.47
2:F:39:HIS:HB3	2:F:42:ARG:HD3	1.96	0.47
4:H:127:VAL:HG23	4:H:145:SER:HB2	1.96	0.47
1:A:133:VAL:HG23	2:B:124:PHE:CE2	2.50	0.47
4:D:123:GLN:HA	4:D:206:SER:OG	2.15	0.47
4:D:190:MET:HE3	4:D:190:MET:O	2.15	0.47
1:E:194:PRO:C	1:E:196:ASP:H	2.22	0.47
2:F:98:THR:O	2:F:98:THR:OG1	2.31	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:117:PHE:O	2:F:146:PHE:HA	2.14	0.47
2:F:215:GLU:O	2:F:217:ASP:N	2.48	0.47
4:H:81:ARG:HG3	4:H:81:ARG:HH11	1.80	0.47
4:H:207:LEU:HD12	4:H:207:LEU:O	2.14	0.47
1:A:3:VAL:HG12	1:A:4:THR:N	2.30	0.47
3:C:114:PHE:HA	3:C:115:PRO:C	2.40	0.47
1:A:149:LYS:HB2	1:A:190:ASN:OD1	2.15	0.47
3:C:139:PHE:HB2	3:C:147:HIS:CD2	2.50	0.47
2:F:10:VAL:HG13	2:F:110:LEU:HD21	1.96	0.47
2:F:19:LEU:HD22	2:F:105:THR:HG21	1.96	0.47
2:F:199:ASN:HB3	2:F:202:ASN:HD22	1.79	0.47
4:H:196:GLU:O	4:H:217:ALA:HB3	2.14	0.47
1:A:131:LYS:HZ3	1:A:178:ASN:HA	1.79	0.47
2:B:139:LEU:HD12	2:B:139:LEU:N	2.30	0.47
4:D:208:LYS:HG2	4:D:209:SER:N	2.30	0.47
4:H:159:PHE:HB2	4:H:199:THR:O	2.15	0.47
4:D:156:VAL:HG11	4:D:186:VAL:HG21	1.96	0.46
1:E:66:ASP:OD2	1:E:69:THR:OG1	2.25	0.46
2:F:29:ASN:HA	2:F:47:SER:O	2.16	0.46
2:F:156:TRP:HA	2:F:161:GLU:HA	1.96	0.46
3:G:78:SER:O	3:G:79:ASN:C	2.57	0.46
4:H:5:LYS:HD2	4:H:5:LYS:C	2.40	0.46
2:B:7:ARG:NH1	2:B:7:ARG:HB3	2.30	0.46
1:E:125:ASP:HA	2:F:124:PHE:HA	1.98	0.46
2:F:46:TYR:CE2	2:F:54:GLU:HB3	2.50	0.46
4:H:52:TYR:H	4:H:68:SER:HB3	1.81	0.46
2:B:123:VAL:HG23	2:B:233:ALA:CB	2.43	0.46
2:B:166:VAL:HG12	2:B:167:CYS:N	2.30	0.46
2:B:199:ASN:HB3	2:B:202:ASN:HD22	1.80	0.46
4:H:208:LYS:HG2	4:H:209:SER:N	2.30	0.46
1:A:6:MET:HB3	1:A:21:ASN:OD1	2.15	0.46
2:B:134:THR:HB	2:B:136:LYS:CE	2.45	0.46
4:D:64:VAL:HG12	4:D:65:ARG:N	2.31	0.46
1:E:16:ASP:O	1:E:77:THR:O	2.33	0.46
1:E:32:LEU:HD13	1:E:72:PHE:HB2	1.98	0.46
1:A:55:LYS:HB3	1:A:64:THR:OG1	2.15	0.46
1:A:125:ASP:HA	2:B:124:PHE:HA	1.98	0.46
2:B:93:ASP:OD2	2:B:93:ASP:N	2.25	0.46
2:F:44:ILE:HG23	2:F:58:ILE:O	2.16	0.46
1:A:125:ASP:OD1	1:A:127:LYS:HB3	2.15	0.46
1:A:192:ILE:H	1:A:192:ILE:CD1	2.29	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:156:VAL:HG13	4:D:202:VAL:CG1	2.45	0.46
4:D:199:THR:OG1	4:D:212:THR:HG22	2.15	0.46
1:E:18:LEU:HD22	1:E:19:PHE:N	2.30	0.46
2:F:110:LEU:N	2:F:110:LEU:HD23	2.30	0.46
3:G:125:ASN:HA	3:G:161:ILE:HD11	1.98	0.46
1:A:194:PRO:C	1:A:196:ASP:H	2.23	0.46
2:B:98:THR:O	2:B:98:THR:OG1	2.33	0.46
3:C:125:ASN:HA	3:C:161:ILE:HD11	1.97	0.46
4:D:46:GLY:HA2	4:D:109:TYR:OH	2.15	0.46
1:E:172:SER:O	1:E:173:ALA:HB2	2.15	0.46
4:H:100:GLU:HA	4:H:103:THR:OG1	2.16	0.46
3:C:46:LEU:O	3:C:49:PHE:HB2	2.15	0.46
2:F:1:ALA:N	2:F:24:THR:HB	2.31	0.46
1:A:29:TYR:CG	4:D:103:THR:HG22	2.51	0.46
4:D:150:TYR:HA	4:D:151:PRO:O	2.16	0.46
1:E:197:THR:HB	1:E:199:PHE:HE1	1.81	0.46
2:F:58:ILE:O	2:F:58:ILE:HG13	2.15	0.46
2:F:223:ARG:HH11	2:F:223:ARG:HG2	1.80	0.46
1:E:139:PHE:CZ	1:E:171:ASN:HB3	2.51	0.46
2:B:215:GLU:O	2:B:217:ASP:N	2.49	0.45
4:D:158:TRP:CZ3	4:D:200:CYS:HB2	2.52	0.45
1:E:3:VAL:HG13	1:E:22:CYS:SG	2.57	0.45
4:H:52:TYR:HD1	4:H:105:CYS:SG	2.38	0.45
1:A:172:SER:O	1:A:173:ALA:HB2	2.16	0.45
2:B:153:LEU:C	2:B:153:LEU:HD23	2.41	0.45
2:F:80:THR:HB	2:F:81:PRO:CD	2.42	0.45
2:F:145:GLY:O	2:F:183:ARG:HB3	2.17	0.45
4:H:135:ALA:O	4:H:138:HIS:HB2	2.16	0.45
2:B:67:PRO:HG2	2:B:68:SER:H	1.81	0.45
3:C:152:LEU:HD23	3:C:153:THR:N	2.31	0.45
2:F:54:GLU:HG2	3:G:58:GLN:HE22	1.81	0.45
3:G:161:ILE:HD13	3:G:161:ILE:N	2.29	0.45
4:H:197:VAL:HG13	4:H:216:LYS:HA	1.97	0.45
2:B:29:ASN:N	2:B:29:ASN:ND2	2.64	0.45
1:E:62:GLU:O	1:E:62:GLU:HG3	2.16	0.45
2:F:11:ALA:O	2:F:109:VAL:HA	2.15	0.45
2:F:153:LEU:C	2:F:153:LEU:HD23	2.42	0.45
3:G:101:GLY:N	3:G:156:PRO:HG2	2.31	0.45
4:H:1:PHE:HD1	4:H:1:PHE:H3	1.63	0.45
2:B:45:HIS:HD2	2:B:55:LYS:HA	1.81	0.45
2:F:31:TYR:N	2:F:31:TYR:CD1	2.84	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:100:GLU:HA	4:D:103:THR:OG1	2.16	0.45
4:D:190:MET:HE3	4:D:190:MET:H	1.82	0.45
4:D:197:VAL:HG13	4:D:216:LYS:HA	1.98	0.45
1:E:20:ILE:HD12	1:E:74:LEU:HD23	1.99	0.45
2:F:68:SER:OG	2:F:69:GLN:N	2.50	0.45
3:G:163:ASP:HB3	3:G:178:HIS:HA	1.98	0.45
2:B:165:GLY:O	2:B:190:LEU:HD12	2.17	0.45
3:C:78:SER:C	3:C:79:ASN:ND2	2.74	0.45
3:G:48:GLU:O	3:G:51:GLN:HB2	2.17	0.45
2:B:29:ASN:HA	2:B:47:SER:O	2.17	0.45
2:B:39:HIS:HB3	2:B:42:ARG:HD3	1.98	0.45
2:B:58:ILE:HG13	2:B:58:ILE:O	2.17	0.45
3:C:71:LEU:HD13	4:D:35:TYR:HB2	1.99	0.45
2:F:62:TYR:HB3	2:F:74:LEU:HD11	1.98	0.45
2:F:134:THR:HB	2:F:136:LYS:CE	2.47	0.45
3:G:93:PHE:HE2	3:G:107:ILE:HD12	1.82	0.45
4:D:140:ASN:CG	4:D:141:THR:H	2.25	0.45
2:B:22:ASN:HB2	2:B:71:ASN:OD1	2.17	0.44
4:D:204:HIS:HD2	4:D:206:SER:OG	2.00	0.44
1:E:148:SER:OG	1:E:153:VAL:HB	2.17	0.44
2:F:176:GLN:HB2	2:F:182:SER:HB3	1.97	0.44
3:G:148:LYS:O	3:G:149:LEU:HD22	2.17	0.44
2:F:22:ASN:HD22	2:F:71:ASN:HD21	1.65	0.44
4:H:64:VAL:HG12	4:H:65:ARG:N	2.33	0.44
2:B:22:ASN:HD22	2:B:71:ASN:HD21	1.65	0.44
2:B:44:ILE:HG23	2:B:58:ILE:O	2.16	0.44
3:C:78:SER:O	3:C:79:ASN:C	2.60	0.44
3:C:163:ASP:HB3	3:C:178:HIS:HA	1.99	0.44
4:D:155:LYS:HZ1	4:H:164:GLU:HB3	1.83	0.44
3:G:82:PRO:HB3	4:H:31:ARG:HB2	1.98	0.44
4:H:199:THR:OG1	4:H:212:THR:HG22	2.17	0.44
1:A:148:SER:OG	1:A:153:VAL:HB	2.18	0.44
2:B:1:ALA:N	2:B:24:THR:HB	2.33	0.44
1:E:133:VAL:HG23	2:F:124:PHE:CE2	2.52	0.44
2:F:52:SER:O	2:F:53:THR:HG23	2.17	0.44
2:B:31:TYR:CD1	2:B:31:TYR:N	2.85	0.44
2:B:46:TYR:CE2	2:B:54:GLU:HB3	2.53	0.44
3:C:161:ILE:HD13	3:C:161:ILE:N	2.29	0.44
4:D:135:ALA:O	4:D:138:HIS:HB2	2.17	0.44
1:E:188:PHE:HB3	1:E:191:SER:HB2	1.99	0.44
2:F:26:ASN:N	2:F:26:ASN:HD22	2.13	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:46:LEU:O	3:G:49:PHE:HB2	2.17	0.44
4:H:156:VAL:HG22	4:H:202:VAL:CG1	2.44	0.44
1:E:7:GLN:HG3	1:E:20:ILE:CG2	2.47	0.44
1:E:125:ASP:OD1	1:E:127:LYS:HB3	2.18	0.44
2:F:45:HIS:HD2	2:F:55:LYS:HA	1.82	0.44
2:F:80:THR:O	2:F:109:VAL:HG11	2.18	0.44
2:F:121:VAL:HA	2:F:142:LEU:O	2.18	0.44
1:E:147:GLN:O	1:E:155:ILE:HD12	2.17	0.44
2:F:209:GLN:NE2	2:F:209:GLN:CA	2.80	0.44
1:A:61:PHE:HB3	1:A:74:LEU:HD11	1.99	0.44
1:A:139:PHE:CZ	1:A:171:ASN:HB3	2.52	0.44
4:H:123:GLN:HA	4:H:206:SER:OG	2.18	0.44
2:B:213:LEU:HD12	2:B:226:PRO:HG2	1.99	0.43
3:C:108:CYS:HB2	3:C:122:TRP:CH2	2.52	0.43
2:F:146:PHE:CE2	2:F:184:TYR:HB2	2.53	0.43
3:G:157:SER:O	3:G:181:PRO:HG3	2.18	0.43
1:A:9:GLN:C	1:A:9:GLN:NE2	2.76	0.43
1:A:32:LEU:HD13	1:A:72:PHE:HB2	1.99	0.43
1:A:188:PHE:HB3	1:A:191:SER:HB2	1.99	0.43
2:B:50:ALA:HA	2:B:66:ARG:HG3	2.00	0.43
2:B:219:TRP:HA	2:B:219:TRP:HE3	1.83	0.43
2:F:21:CYS:HB2	2:F:32:TRP:CZ2	2.53	0.43
2:F:49:GLY:O	2:F:66:ARG:HG2	2.18	0.43
2:F:139:LEU:N	2:F:139:LEU:HD12	2.33	0.43
3:G:88:PRO:HD3	3:G:168:HIS:HD2	1.83	0.43
4:H:174:LEU:H	4:H:174:LEU:CD1	2.11	0.43
2:B:68:SER:OG	2:B:69:GLN:N	2.51	0.43
2:B:121:VAL:HA	2:B:142:LEU:O	2.18	0.43
2:B:223:ARG:HG2	2:B:223:ARG:HH11	1.82	0.43
4:D:81:ARG:HG3	4:D:81:ARG:HH11	1.83	0.43
4:D:124:PRO:HD3	4:D:204:HIS:HD2	1.83	0.43
3:G:139:PHE:HB2	3:G:147:HIS:CD2	2.53	0.43
3:C:14:TYR:CD2	3:C:68:LYS:HA	2.53	0.43
2:F:123:VAL:CG1	2:F:235:ALA:HB2	2.48	0.43
3:C:31:GLU:OE1	3:C:137:SER:HB2	2.18	0.43
3:C:157:SER:O	3:C:181:PRO:HG3	2.19	0.43
4:D:119:ARG:HB2	4:D:119:ARG:NH1	2.33	0.43
1:E:81:GLU:HA	1:E:109:VAL:CG1	2.48	0.43
2:F:219:TRP:HA	2:F:219:TRP:HE3	1.83	0.43
4:H:31:ARG:HD3	4:H:31:ARG:N	2.20	0.43
4:H:55:ARG:NH1	4:H:62:GLU:OE1	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:119:ARG:HB2	4:H:119:ARG:NH1	2.33	0.43
4:H:147:THR:O	4:H:148:ASP:HB2	2.19	0.43
1:A:92:ASN:HB2	1:A:99:GLN:HA	2.01	0.43
2:B:112:ASP:N	2:B:112:ASP:OD1	2.51	0.43
1:E:118:PRO:HB2	1:E:197:THR:OG1	2.18	0.43
3:G:12:SER:CB	3:G:67:VAL:HG21	2.45	0.43
1:A:3:VAL:HG13	1:A:22:CYS:SG	2.59	0.43
1:A:154:TYR:HD2	1:A:176:TRP:HE1	1.67	0.43
3:C:87:ALA:HA	3:C:88:PRO:HD2	1.88	0.43
3:C:101:GLY:N	3:C:156:PRO:HG2	2.32	0.43
4:D:146:VAL:HB	4:D:184:VAL:CG1	2.48	0.43
3:G:27:PHE:C	3:G:29:GLY:N	2.76	0.43
4:H:190:MET:HE3	4:H:190:MET:H	1.84	0.43
1:A:69:THR:O	1:A:70:THR:C	2.62	0.43
1:A:97:THR:CG2	3:C:62:GLN:HG3	2.42	0.43
4:H:44:THR:HB	4:H:49:ARG:HD3	2.00	0.43
1:A:16:ASP:O	1:A:79:VAL:HG12	2.19	0.43
2:B:80:THR:HB	2:B:81:PRO:CD	2.41	0.43
4:D:164:GLU:HB3	4:H:155:LYS:HZ3	1.84	0.43
1:E:164:ARG:O	1:E:166:MET:N	2.47	0.43
1:E:192:ILE:H	1:E:192:ILE:CD1	2.28	0.43
2:F:7:ARG:HB3	2:F:7:ARG:NH1	2.33	0.43
4:H:90:GLN:HB2	4:H:93:ILE:HD12	2.01	0.43
1:A:190:ASN:N	1:A:190:ASN:ND2	2.67	0.42
3:C:118:ILE:CG1	3:C:119:ASN:N	2.82	0.42
2:F:1:ALA:H2	2:F:24:THR:HB	1.84	0.42
2:F:165:GLY:O	2:F:190:LEU:HD12	2.19	0.42
3:G:108:CYS:HB2	3:G:122:TRP:CH2	2.53	0.42
3:C:12:SER:CB	3:C:67:VAL:HG21	2.46	0.42
3:C:27:PHE:C	3:C:29:GLY:N	2.76	0.42
4:D:5:LYS:HD2	4:D:5:LYS:C	2.44	0.42
4:D:177:ASN:N	4:D:177:ASN:HD22	2.16	0.42
1:E:38:TYR:CD2	1:E:85:ALA:HB2	2.54	0.42
4:D:74:ARG:HA	4:D:74:ARG:HD3	1.89	0.42
4:D:124:PRO:HD3	4:D:204:HIS:CD2	2.54	0.42
2:F:167:CYS:HB3	2:F:189:ARG:HD3	1.99	0.42
3:G:2:GLU:O	3:G:3:ALA:HB2	2.18	0.42
1:A:111:PRO:HG2	1:A:160:VAL:CG1	2.48	0.42
4:D:90:GLN:HB2	4:D:93:ILE:HD12	2.02	0.42
1:E:141:SER:O	1:E:142:GLN:HG2	2.19	0.42
2:F:49:GLY:O	2:F:52:SER:HB3	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:121:VAL:HG21	2:F:231:VAL:O	2.19	0.42
4:H:177:ASN:N	4:H:177:ASN:HD22	2.17	0.42
1:A:7:GLN:HB2	1:A:105:THR:OG1	2.18	0.42
2:B:21:CYS:HB2	2:B:32:TRP:CZ2	2.55	0.42
3:C:6:VAL:HG13	3:C:6:VAL:O	2.19	0.42
3:C:15:GLN:HE21	3:C:115:PRO:HB2	1.83	0.42
1:E:7:GLN:HB2	1:E:105:THR:OG1	2.19	0.42
1:E:137:THR:OG1	1:E:138:ASP:N	2.52	0.42
2:F:23:GLN:HE21	2:F:30:MET:HE3	1.84	0.42
2:F:114:LYS:HE2	2:F:221:GLN:HE21	1.84	0.42
2:B:26:ASN:N	2:B:26:ASN:HD22	2.17	0.42
3:C:100:LEU:H	3:C:100:LEU:CD1	2.29	0.42
4:D:81:ARG:N	4:D:82:PRO:CD	2.83	0.42
1:E:190:ASN:HD22	1:E:190:ASN:H	1.66	0.42
2:F:34:ARG:HB2	2:F:44:ILE:HD11	2.01	0.42
3:G:62:GLN:O	3:G:66:VAL:HG23	2.20	0.42
4:H:7:LYS:HD2	4:H:96:ARG:HD3	2.02	0.42
4:H:146:VAL:HB	4:H:184:VAL:HG12	2.02	0.42
4:H:150:TYR:HA	4:H:151:PRO:O	2.19	0.42
1:A:81:GLU:HA	1:A:109:VAL:CG1	2.49	0.42
2:B:52:SER:O	2:B:53:THR:HG23	2.20	0.42
3:C:4:ASP:CG	4:D:45:ASN:H	2.27	0.42
4:D:147:THR:O	4:D:148:ASP:HB2	2.20	0.42
4:D:179:ASP:O	4:D:180:TRP:HB2	2.19	0.42
2:F:26:ASN:N	2:F:26:ASN:ND2	2.68	0.42
2:F:119:PRO:HG2	2:F:231:VAL:CB	2.50	0.42
2:F:119:PRO:HG2	2:F:231:VAL:HG21	2.01	0.42
2:F:193:SER:OG	2:F:196:PHE:HB2	2.20	0.42
3:G:4:ASP:CG	4:H:45:ASN:H	2.27	0.42
4:H:146:VAL:HB	4:H:184:VAL:CG1	2.49	0.42
1:A:13:SER:HA	1:A:110:VAL:CG2	2.50	0.42
1:A:137:THR:OG1	1:A:138:ASP:N	2.53	0.42
4:D:52:TYR:H	4:D:68:SER:HB3	1.85	0.42
1:E:150:ASP:O	1:E:151:SER:HB2	2.20	0.42
1:E:156:THR:HG22	2:F:173:LEU:HD21	2.01	0.42
2:F:25:ASN:O	2:F:27:HIS:N	2.53	0.42
4:H:1:PHE:CD1	4:H:1:PHE:N	2.87	0.42
1:E:71:SER:HB2	1:E:73:HIS:CE1	2.55	0.42
1:E:190:ASN:N	1:E:190:ASN:ND2	2.68	0.42
2:F:229:GLN:HE21	2:F:229:GLN:HB2	1.54	0.42
4:H:156:VAL:HG13	4:H:202:VAL:CG1	2.47	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:38:TYR:CD2	1:A:85:ALA:HB2	2.55	0.41
3:C:33:PHE:CD1	3:C:33:PHE:C	2.98	0.41
3:C:109:PHE:HA	3:C:149:LEU:HD13	2.01	0.41
4:D:146:VAL:HB	4:D:184:VAL:HG12	2.02	0.41
1:E:58:SER:O	1:E:59:ARG:C	2.62	0.41
1:E:118:PRO:HG2	1:E:196:ASP:O	2.20	0.41
2:F:144:THR:HG22	2:F:185:ALA:HB1	2.01	0.41
1:E:156:THR:HG21	2:F:187:SER:HG	1.84	0.41
2:F:29:ASN:N	2:F:29:ASN:ND2	2.67	0.41
2:F:149:ASP:OD1	2:F:149:ASP:N	2.46	0.41
4:H:12:VAL:CG1	4:H:13:ASP:N	2.82	0.41
4:H:140:ASN:CG	4:H:141:THR:H	2.27	0.41
2:B:193:SER:OG	2:B:196:PHE:HB2	2.21	0.41
4:D:12:VAL:CG1	4:D:13:ASP:N	2.82	0.41
4:D:173:GLN:H	4:D:173:GLN:HG2	1.69	0.41
1:E:92:ASN:HB2	1:E:99:GLN:HA	2.02	0.41
3:G:78:SER:C	3:G:79:ASN:ND2	2.78	0.41
4:H:81:ARG:N	4:H:82:PRO:CD	2.83	0.41
1:A:190:ASN:HD22	1:A:190:ASN:H	1.64	0.41
4:D:55:ARG:HG2	4:D:55:ARG:NH1	2.35	0.41
2:F:67:PRO:HG2	2:F:68:SER:H	1.86	0.41
3:G:86:GLU:HB3	3:G:112:ASN:ND2	2.36	0.41
1:A:150:ASP:O	1:A:151:SER:HB2	2.21	0.41
2:B:34:ARG:HB2	2:B:44:ILE:HD11	2.03	0.41
2:B:119:PRO:HG2	2:B:231:VAL:CG2	2.50	0.41
3:C:2:GLU:O	3:C:3:ALA:HB2	2.20	0.41
3:G:6:VAL:O	3:G:6:VAL:HG13	2.21	0.41
3:G:27:PHE:C	3:G:29:GLY:H	2.29	0.41
3:C:27:PHE:C	3:C:29:GLY:H	2.29	0.41
4:D:184:VAL:HG22	4:D:185:LEU:N	2.35	0.41
1:A:16:ASP:O	1:A:77:THR:O	2.38	0.41
2:B:4:GLN:NE2	2:B:104:GLY:H	2.19	0.41
2:B:219:TRP:CB	2:B:225:LYS:HG2	2.50	0.41
4:D:52:TYR:HB2	4:D:101:LEU:HD13	2.01	0.41
4:H:74:ARG:HA	4:H:74:ARG:HD3	1.88	0.41
1:A:7:GLN:HG3	1:A:20:ILE:CG2	2.51	0.41
1:A:71:SER:HB2	1:A:73:HIS:CE1	2.55	0.41
1:A:141:SER:O	1:A:142:GLN:HG2	2.21	0.41
2:F:213:LEU:HD12	2:F:226:PRO:HG2	2.02	0.41
3:G:109:PHE:HA	3:G:149:LEU:HD13	2.03	0.41
4:H:1:PHE:CZ	4:H:115:HIS:ND1	2.89	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:110:VAL:HA	1:A:111:PRO:HD3	1.90	0.40
1:A:118:PRO:HB2	1:A:197:THR:OG1	2.21	0.40
2:B:218:GLU:O	2:B:218:GLU:HG3	2.21	0.40
2:B:231:VAL:HG12	2:B:232:SER:N	2.36	0.40
3:C:48:GLU:O	3:C:51:GLN:HB2	2.21	0.40
3:C:152:LEU:HD23	3:C:153:THR:H	1.86	0.40
1:E:54:ASN:N	1:E:54:ASN:ND2	2.60	0.40
2:F:119:PRO:HD3	2:F:210:PHE:CG	2.56	0.40
3:G:85:ASN:ND2	3:G:169:TRP:CB	2.84	0.40
4:H:126:VAL:O	4:H:127:VAL:HG13	2.21	0.40
2:B:25:ASN:O	2:B:27:HIS:N	2.54	0.40
2:B:33:TYR:CE2	2:B:43:LEU:HB2	2.56	0.40
4:D:55:ARG:NH1	4:D:62:GLU:OE1	2.54	0.40
4:D:144:CYS:HB2	4:D:158:TRP:CZ2	2.57	0.40
1:E:172:SER:HG	2:F:189:ARG:HE	1.69	0.40
2:F:13:THR:N	2:F:110:LEU:O	2.50	0.40
3:G:139:PHE:HB2	3:G:147:HIS:NE2	2.36	0.40
4:D:52:TYR:HD1	4:D:105:CYS:SG	2.44	0.40
1:E:32:LEU:HB3	1:E:72:PHE:CD1	2.56	0.40
3:G:152:LEU:HD23	3:G:153:THR:N	2.37	0.40
4:H:168:GLY:O	4:H:189:GLU:HB2	2.21	0.40
1:A:80:GLN:O	1:A:109:VAL:HG11	2.21	0.40
1:E:82:ILE:H	1:E:82:ILE:HG13	1.75	0.40
4:H:184:VAL:HG22	4:H:185:LEU:N	2.37	0.40

There are no symmetry-related clashes.

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	A	197/199 (99%)	159 (81%)	27 (14%)	11 (6%)	1 7

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	E	197/199 (99%)	160 (81%)	25 (13%)	12 (6%)	1	6
2	B	234/236 (99%)	192 (82%)	33 (14%)	9 (4%)	2	12
2	F	234/236 (99%)	193 (82%)	33 (14%)	8 (3%)	3	13
3	C	180/182 (99%)	162 (90%)	17 (9%)	1 (1%)	21	48
3	G	180/182 (99%)	161 (89%)	18 (10%)	1 (1%)	21	48
4	D	197/217 (91%)	166 (84%)	27 (14%)	4 (2%)	6	21
4	H	197/217 (91%)	166 (84%)	27 (14%)	4 (2%)	6	21
All	All	1616/1668 (97%)	1359 (84%)	207 (13%)	50 (3%)	3	14

All (50) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	142	GLN
1	E	142	GLN
2	F	215	GLU
1	A	150	ASP
1	A	151	SER
1	A	166	MET
1	A	173	ALA
2	B	26	ASN
2	B	113	LEU
2	B	215	GLU
2	B	216	ASN
4	D	60	ARG
4	D	131	SER
1	E	129	SER
1	E	150	ASP
1	E	151	SER
1	E	166	MET
1	E	173	ALA
2	F	26	ASN
2	F	216	ASN
2	F	222	ASP
4	H	60	ARG
4	H	131	SER
1	A	129	SER
1	A	164	ARG
1	A	171	ASN
2	B	85	SER
2	B	222	ASP

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Mol	Chain	Res	Type
3	C	79	ASN
4	D	160	ARG
1	E	164	ARG
1	E	171	ASN
2	F	77	GLU
2	F	85	SER
2	F	115	ASN
3	G	79	ASN
4	H	160	ARG
1	A	116	PRO
1	A	117	ASP
2	B	77	GLU
2	B	112	ASP
2	B	149	ASP
1	E	116	PRO
1	E	117	ASP
2	F	149	ASP
1	A	192	ILE
4	D	134	GLU
1	E	158	LYS
1	E	192	ILE
4	H	82	PRO

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	178/178 (100%)	166 (93%)	12 (7%)	15	39
1	E	178/178 (100%)	166 (93%)	12 (7%)	15	39
2	B	203/203 (100%)	185 (91%)	18 (9%)	9	29
2	F	203/203 (100%)	185 (91%)	18 (9%)	9	29
3	C	163/163 (100%)	148 (91%)	15 (9%)	8	28
3	G	163/163 (100%)	147 (90%)	16 (10%)	7	26
4	D	182/189 (96%)	164 (90%)	18 (10%)	7	26

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
4	H	182/189 (96%)	163 (90%)	19 (10%)	7	23
All	All	1452/1466 (99%)	1324 (91%)	128 (9%)	9	30

All (128) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	GLN
1	A	18	LEU
1	A	21	ASN
1	A	33	PHE
1	A	41	GLU
1	A	54	ASN
1	A	62	GLU
1	A	80	GLN
1	A	103	THR
1	A	130	ASP
1	A	190	ASN
1	A	195	GLU
2	B	8	ASN
2	B	23	GLN
2	B	25	ASN
2	B	31	TYR
2	B	36	ASP
2	B	76	LEU
2	B	93	ASP
2	B	97	ASP
2	B	98	THR
2	B	113	LEU
2	B	120	GLU
2	B	142	LEU
2	B	183	ARG
2	B	207	GLN
2	B	209	GLN
2	B	219	TRP
2	B	222	ASP
2	B	229	GLN
3	C	8	THR
3	C	37	LEU
3	C	46	LEU
3	C	62	GLN
3	C	63	ASN
3	C	79	ASN

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Mol	Chain	Res	Type
3	C	100	LEU
3	C	102	GLN
3	C	131	ASP
3	C	152	LEU
3	C	159	ASP
3	C	161	ILE
3	C	163	ASP
3	C	173	GLU
3	C	176	LEU
4	D	4	GLN
4	D	5	LYS
4	D	30	GLU
4	D	38	MET
4	D	76	VAL
4	D	92	GLU
4	D	95	GLU
4	D	98	ARG
4	D	102	ASP
4	D	120	ARG
4	D	139	HIS
4	D	166	THR
4	D	174	LEU
4	D	186	VAL
4	D	189	GLU
4	D	190	MET
4	D	201	HIS
4	D	208	LYS
1	E	9	GLN
1	E	18	LEU
1	E	21	ASN
1	E	33	PHE
1	E	41	GLU
1	E	54	ASN
1	E	62	GLU
1	E	80	GLN
1	E	103	THR
1	E	130	ASP
1	E	190	ASN
1	E	195	GLU
2	F	8	ASN
2	F	23	GLN
2	F	25	ASN

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Mol	Chain	Res	Type
2	F	36	ASP
2	F	76	LEU
2	F	93	ASP
2	F	97	ASP
2	F	98	THR
2	F	99	LEU
2	F	110	LEU
2	F	120	GLU
2	F	142	LEU
2	F	183	ARG
2	F	207	GLN
2	F	209	GLN
2	F	219	TRP
2	F	222	ASP
2	F	229	GLN
3	G	8	THR
3	G	12	SER
3	G	37	LEU
3	G	46	LEU
3	G	62	GLN
3	G	63	ASN
3	G	79	ASN
3	G	100	LEU
3	G	102	GLN
3	G	131	ASP
3	G	152	LEU
3	G	159	ASP
3	G	161	ILE
3	G	163	ASP
3	G	173	GLU
3	G	176	LEU
4	H	4	GLN
4	H	5	LYS
4	H	30	GLU
4	H	31	ARG
4	H	38	MET
4	H	76	VAL
4	H	92	GLU
4	H	95	GLU
4	H	98	ARG
4	H	102	ASP
4	H	120	ARG

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Mol	Chain	Res	Type
4	H	139	HIS
4	H	166	THR
4	H	174	LEU
4	H	186	VAL
4	H	189	GLU
4	H	190	MET
4	H	201	HIS
4	H	208	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (73) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	7	GLN
1	A	44	GLN
1	A	54	ASN
1	A	112	ASN
1	A	186	ASN
1	A	189	ASN
1	A	190	ASN
2	B	4	GLN
2	B	22	ASN
2	B	23	GLN
2	B	25	ASN
2	B	26	ASN
2	B	29	ASN
2	B	45	HIS
2	B	176	GLN
2	B	202	ASN
2	B	207	GLN
2	B	209	GLN
2	B	221	GLN
3	C	15	GLN
3	C	58	GLN
3	C	63	ASN
3	C	79	ASN
3	C	102	GLN
3	C	178	HIS
4	D	4	GLN
4	D	32	HIS
4	D	36	GLN
4	D	45	ASN
4	D	123	GLN

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Mol	Chain	Res	Type
4	D	137	ASN
4	D	138	HIS
4	D	140	ASN
4	D	163	GLN
4	D	173	GLN
4	D	183	GLN
4	D	201	HIS
4	D	204	HIS
1	E	7	GLN
1	E	44	GLN
1	E	54	ASN
1	E	142	GLN
1	E	189	ASN
1	E	190	ASN
2	F	4	GLN
2	F	22	ASN
2	F	23	GLN
2	F	25	ASN
2	F	26	ASN
2	F	29	ASN
2	F	45	HIS
2	F	202	ASN
2	F	207	GLN
2	F	209	GLN
2	F	221	GLN
3	G	15	GLN
3	G	58	GLN
3	G	63	ASN
3	G	79	ASN
3	G	102	GLN
3	G	168	HIS
3	G	178	HIS
4	H	4	GLN
4	H	32	HIS
4	H	36	GLN
4	H	45	ASN
4	H	123	GLN
4	H	137	ASN
4	H	138	HIS
4	H	140	ASN
4	H	183	GLN
4	H	201	HIS

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Mol	Chain	Res	Type
4	H	204	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

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	199/199 (100%)	0.51	10 (5%) 34 17	29, 84, 130, 140	0
1	E	199/199 (100%)	0.61	11 (5%) 30 16	29, 85, 131, 140	0
2	B	236/236 (100%)	0.58	12 (5%) 33 17	26, 89, 139, 159	0
2	F	236/236 (100%)	0.51	10 (4%) 40 21	29, 90, 139, 158	0
3	C	182/182 (100%)	0.03	3 (1%) 70 47	23, 55, 107, 128	0
3	G	182/182 (100%)	-0.01	0 100 100	26, 56, 106, 130	0
4	D	201/217 (92%)	0.37	2 (0%) 79 59	26, 72, 113, 128	0
4	H	201/217 (92%)	0.32	3 (1%) 72 50	30, 72, 112, 127	0
All	All	1636/1668 (98%)	0.38	51 (3%) 51 29	23, 74, 129, 159	0

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	F	140	VAL	5.1
1	E	135	LEU	3.9
2	B	79	ALA	3.3
2	F	212	GLY	3.3
2	B	212	GLY	3.3
1	E	173	ALA	3.2
1	E	174	VAL	3.0
2	B	110	LEU	2.9
1	E	200	PRO	2.8
1	A	135	LEU	2.8
3	C	100	LEU	2.8
1	E	116	PRO	2.8
2	F	129	ALA	2.7
3	C	1	ILE	2.7
2	B	173	LEU	2.6
4	H	192	PRO	2.6

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Mol	Chain	Res	Type	RSRZ
2	F	151	VAL	2.6
1	E	192	ILE	2.5
1	E	138	ASP	2.5
3	C	99	LEU	2.5
4	D	141	THR	2.5
4	H	200	CYS	2.5
2	F	148	PRO	2.5
1	A	187	ALA	2.5
1	E	119	ALA	2.5
2	B	131	ILE	2.5
1	E	110	VAL	2.5
1	A	174	VAL	2.4
1	A	81	GLU	2.4
1	A	177	SER	2.4
1	A	134	CYS	2.4
2	F	200	PRO	2.4
2	B	228	THR	2.3
2	F	138	THR	2.3
1	A	119	ALA	2.3
1	E	170	SER	2.3
2	B	216	ASN	2.3
2	B	150	HIS	2.3
2	B	172	PRO	2.2
1	A	145	VAL	2.2
2	B	197	TRP	2.2
2	B	138	THR	2.1
2	F	70	GLU	2.1
4	H	169	VAL	2.1
2	B	165	GLY	2.1
1	A	161	LEU	2.0
2	F	186	LEU	2.0
4	D	187	MET	2.0
1	A	173	ALA	2.0
2	F	142	LEU	2.0
1	E	171	ASN	2.0

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

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.