



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 10:55 AM GMT

PDB ID : 3NH7
Title : Crystal structure of the neutralizing Fab fragment AbD1556 bound to the BMP type I receptor IA
Authors : Mueller, T.D.; Harth, S.; Sebald, W.
Deposited on : 2010-06-14
Resolution : 2.70 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

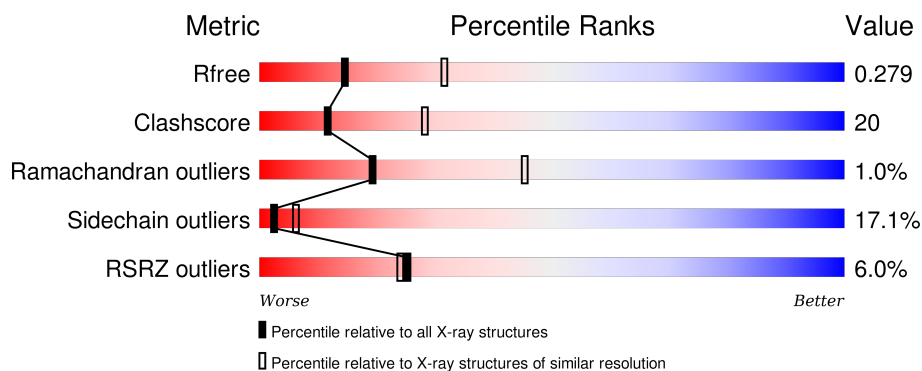
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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}	91344	2103 (2.70-2.70)
Clashscore	102246	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)
RSRZ outliers	91569	2107 (2.70-2.70)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	H	234	<div> <div>3%</div> <div>47% 34% 8% 12%</div> </div>
1	I	234	<div> <div>5%</div> <div>58% 26% 7% 8%</div> </div>
1	J	234	<div> <div>6%</div> <div>59% 29% 6% 6%</div> </div>
1	K	234	<div> <div>4%</div> <div>61% 24% 6% 8%</div> </div>
2	L	213	<div> <div>3%</div> <div>62% 30% 7%</div> </div>

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Mol	Chain	Length	Quality of chain
2	M	213	<div><div>3%</div><div><div></div><div></div><div></div></div><div>64%</div><div>27%</div><div>7%</div><div></div></div>
2	N	213	<div><div>3%</div><div><div></div><div></div><div></div></div><div>64%</div><div>26%</div><div>8%</div><div></div></div>
2	O	213	<div><div>2%</div><div><div></div><div></div><div></div></div><div>62%</div><div>28%</div><div>8%</div><div></div></div>
3	A	129	<div><div>9%</div><div><div></div><div></div><div></div></div><div>41%</div><div>18%</div><div>7%</div><div>34%</div></div>
3	B	129	<div><div>9%</div><div><div></div><div></div><div></div></div><div>39%</div><div>20%</div><div>7%</div><div>34%</div></div>
3	C	129	<div><div>10%</div><div><div></div><div></div><div></div></div><div>40%</div><div>18%</div><div>8%</div><div>34%</div></div>
3	D	129	<div><div>16%</div><div><div></div><div></div><div></div></div><div>38%</div><div>21%</div><div>7%</div><div>34%</div></div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 15350 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 Antibody fragment Fab AbD1556, heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	H	207	Total	C	N	O	S	0	0	0
			1569	992	269	303	5			
1	I	215	Total	C	N	O	S	0	0	0
			1620	1021	278	316	5			
1	J	221	Total	C	N	O	S	0	0	0
			1661	1045	286	325	5			
1	K	215	Total	C	N	O	S	0	0	0
			1625	1025	279	316	5			

- Molecule 2 is a protein called Antibody fragment Fab AbD1556, light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	210	Total	C	N	O	S	0	0	0
			1558	973	258	322	5			
2	M	210	Total	C	N	O	S	0	0	0
			1558	973	258	322	5			
2	N	210	Total	C	N	O	S	0	0	0
			1558	973	258	322	5			
2	O	210	Total	C	N	O	S	0	0	0
			1558	973	258	322	5			

- Molecule 3 is a protein called Bone morphogenetic protein receptor type-1A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	85	Total	C	N	O	S	0	0	0
			654	400	112	131	11			
3	B	85	Total	C	N	O	S	0	0	0
			654	400	112	131	11			
3	C	85	Total	C	N	O	S	0	0	0
			654	400	112	131	11			
3	D	85	Total	C	N	O	S	0	0	0
			654	400	112	131	11			

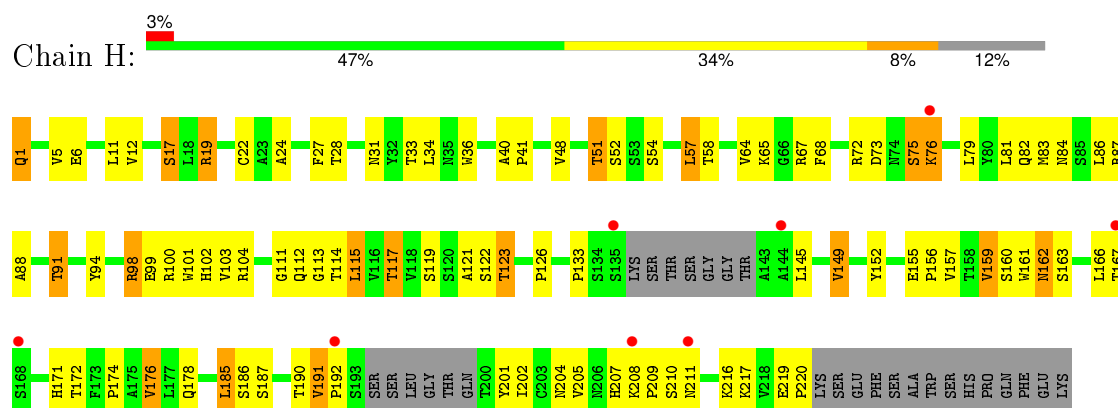
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	H	2	Total 2	O 2	0	0
4	L	3	Total 3	O 3	0	0
4	A	2	Total 2	O 2	0	0
4	I	5	Total 5	O 5	0	0
4	M	4	Total 4	O 4	0	0
4	B	1	Total 1	O 1	0	0
4	J	2	Total 2	O 2	0	0
4	N	1	Total 1	O 1	0	0
4	K	2	Total 2	O 2	0	0
4	O	4	Total 4	O 4	0	0
4	D	1	Total 1	O 1	0	0

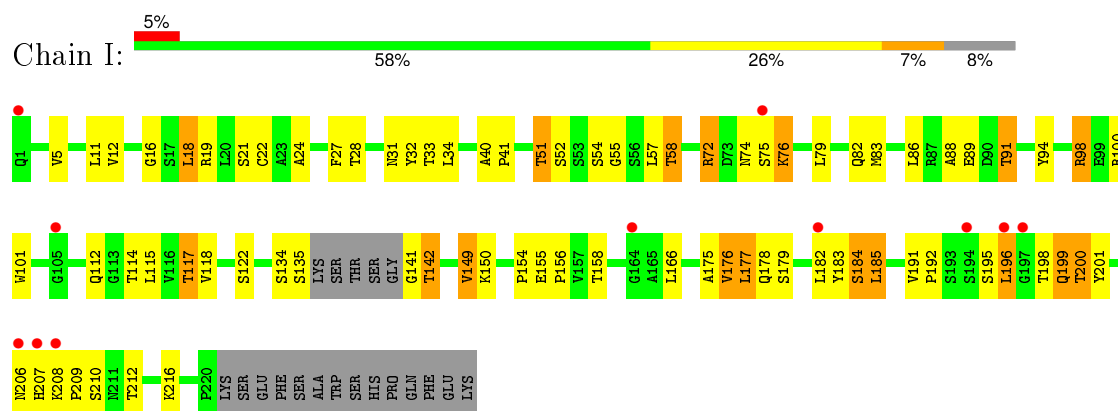
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

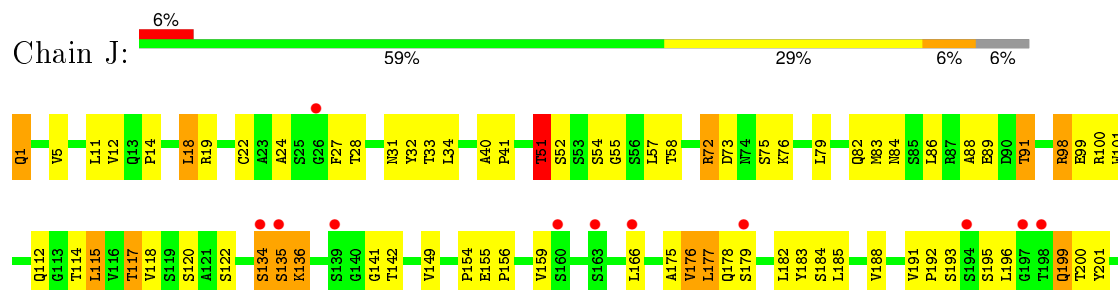
- Molecule 1: Antibody fragment Fab AbD1556, heavy chain



- Molecule 1: Antibody fragment Fab AbD1556, heavy chain

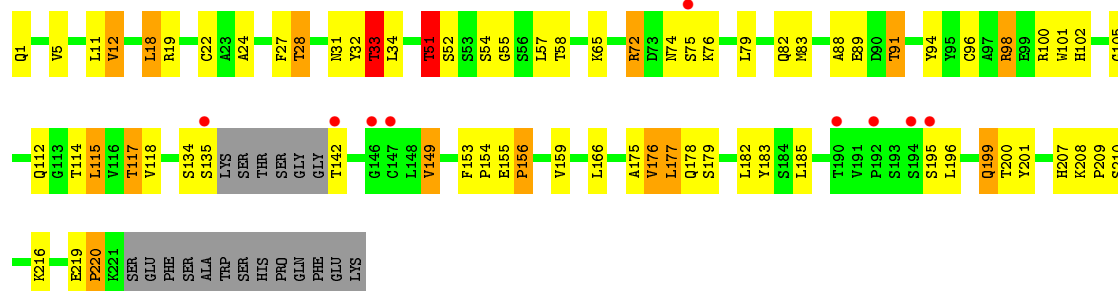


- Molecule 1: Antibody fragment Fab AbD1556, heavy chain

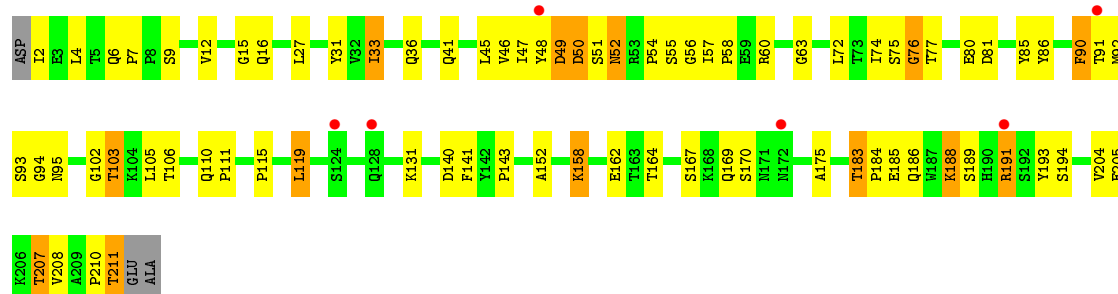




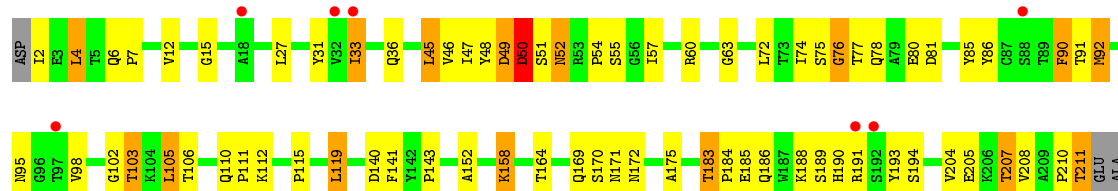
- Molecule 1: Antibody fragment Fab AbD1556, heavy chain



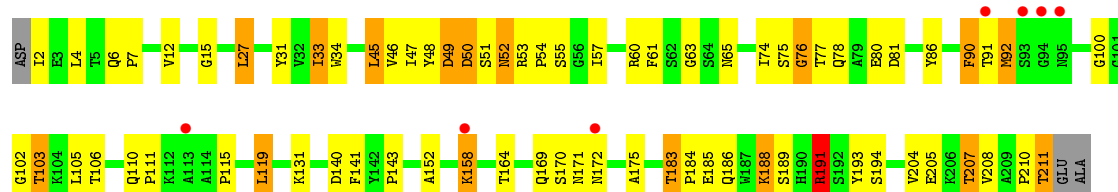
- Molecule 2: Antibody fragment Fab AbD1556, light chain



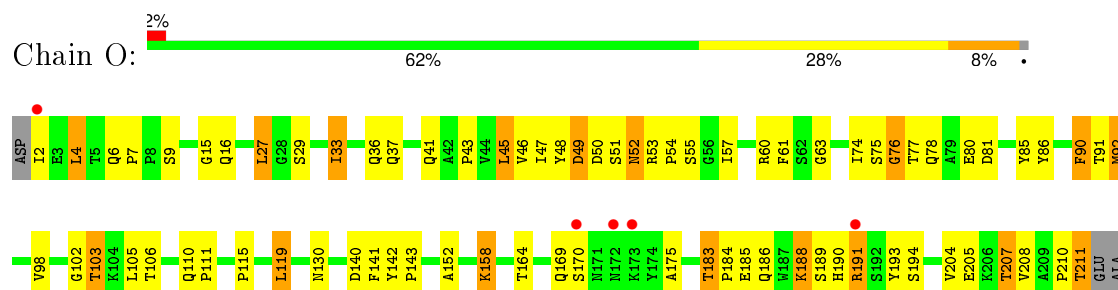
- Molecule 2: Antibody fragment Fab AbD1556, light chain



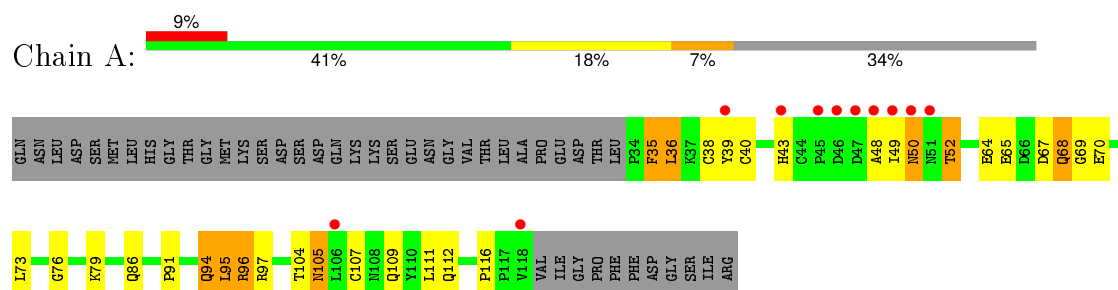
- Molecule 2: Antibody fragment Fab AbD1556, light chain



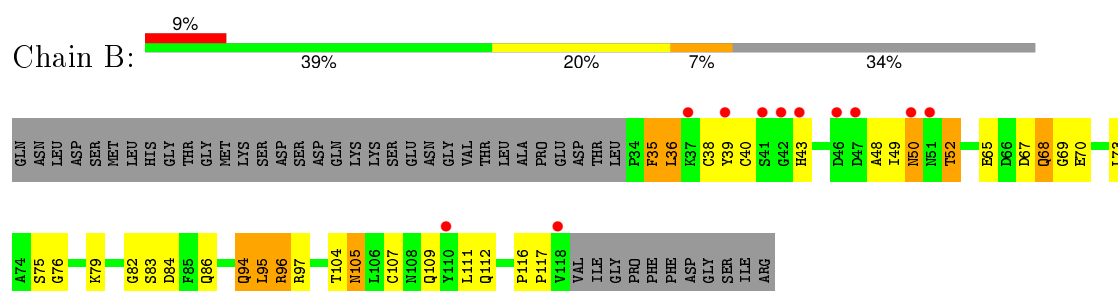
- Molecule 2: Antibody fragment Fab AbD1556, light chain



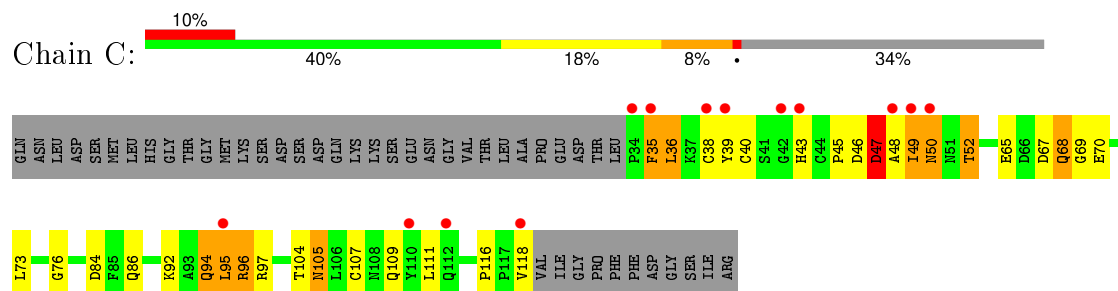
- Molecule 3: Bone morphogenetic protein receptor type-1A



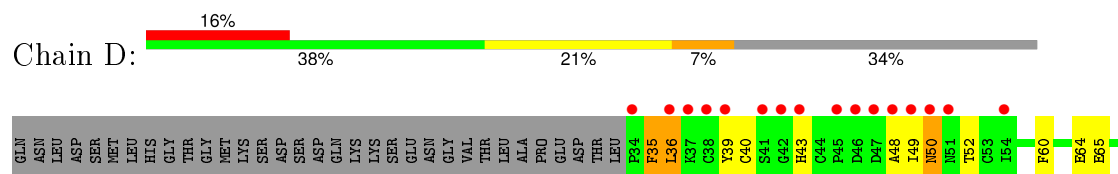
- Molecule 3: Bone morphogenetic protein receptor type-1A

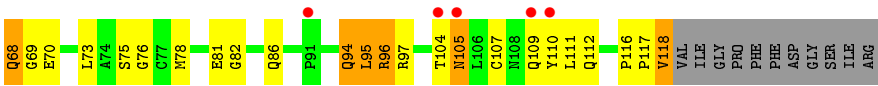


- Molecule 3: Bone morphogenetic protein receptor type-1A



- Molecule 3: Bone morphogenetic protein receptor type-1A





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	89.32Å 129.25Å 100.24Å 90.00° 92.27° 90.00°	Depositor
Resolution (Å)	30.40 – 2.70 30.38 – 2.70	Depositor EDS
% Data completeness (in resolution range)	95.3 (30.40-2.70) 95.3 (30.38-2.70)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.65 (at 2.68Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.234 , 0.280 0.235 , 0.279	Depositor DCC
R_{free} test set	3020 reflections (5.34%)	DCC
Wilson B-factor (Å ²)	50.5	Xtriage
Anisotropy	0.040	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 61.3	EDS
Estimated twinning fraction	0.025 for h,-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 59533 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	15350	wwPDB-VP
Average B, all atoms (Å ²)	64.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.33% 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, 0.375 respectively for untwinned datasets, and 0.333, 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	H	0.65	0/1608	0.80	0/2190
1	I	0.65	0/1660	0.76	0/2262
1	J	0.60	0/1702	0.76	1/2318 (0.0%)
1	K	0.73	2/1665 (0.1%)	0.82	3/2268 (0.1%)
2	L	0.69	1/1595 (0.1%)	0.99	3/2180 (0.1%)
2	M	0.64	0/1595	0.80	3/2180 (0.1%)
2	N	0.69	0/1595	1.08	5/2180 (0.2%)
2	O	0.69	0/1595	0.80	3/2180 (0.1%)
3	A	0.59	0/667	0.72	0/903
3	B	0.54	0/667	0.71	0/903
3	C	0.59	1/667 (0.1%)	0.86	3/903 (0.3%)
3	D	0.53	0/667	0.72	0/903
All	All	0.65	4/15683 (0.0%)	0.84	21/21370 (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
2	L	0	1
3	A	0	1
3	B	0	1
3	C	0	1
3	D	0	1
All	All	0	5

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	L	191	ARG	CZ-NH2	6.19	1.41	1.33
3	C	47	ASP	CB-CG	-5.60	1.40	1.51
1	K	33	THR	CB-CG2	-5.28	1.34	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	K	96	CYS	CB-SG	-5.14	1.73	1.81

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	N	191	ARG	NE-CZ-NH1	-28.45	106.08	120.30
2	L	191	ARG	NE-CZ-NH2	23.52	132.06	120.30
2	N	191	ARG	NE-CZ-NH2	18.50	129.55	120.30
2	L	191	ARG	NE-CZ-NH1	-15.99	112.31	120.30
2	M	191	ARG	NE-CZ-NH1	-11.30	114.65	120.30
3	C	47	ASP	CB-CG-OD2	-10.14	109.18	118.30
2	N	50	ASP	CB-CG-OD2	9.77	127.09	118.30
1	K	98	ARG	NE-CZ-NH1	-7.91	116.34	120.30
2	O	49	ASP	N-CA-C	-7.61	90.45	111.00
2	N	49	ASP	N-CA-C	-7.21	91.53	111.00
2	O	191	ARG	NE-CZ-NH1	-6.98	116.81	120.30
1	K	98	ARG	NE-CZ-NH2	6.75	123.67	120.30
3	C	47	ASP	CB-CA-C	-6.39	97.61	110.40
2	L	49	ASP	N-CA-C	-6.01	94.78	111.00
2	M	49	ASP	N-CA-C	-5.58	95.92	111.00
2	O	191	ARG	CG-CD-NE	5.51	123.36	111.80
2	N	50	ASP	CB-CG-OD1	-5.32	113.51	118.30
1	J	51	THR	CB-CA-C	-5.24	97.45	111.60
3	C	47	ASP	N-CA-CB	-5.23	101.18	110.60
2	M	50	ASP	N-CA-CB	5.08	119.74	110.60
1	K	51	THR	CB-CA-C	-5.06	97.93	111.60

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	A	48	ALA	Peptide
3	B	48	ALA	Peptide
3	C	48	ALA	Peptide
3	D	48	ALA	Peptide
2	L	93	SER	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	H	1569	0	1526	89	0
1	I	1620	0	1576	64	0
1	J	1661	0	1623	58	0
1	K	1625	0	1586	51	0
2	L	1558	0	1507	67	0
2	M	1558	0	1507	75	0
2	N	1558	0	1507	74	0
2	O	1558	0	1507	69	0
3	A	654	0	608	38	0
3	B	654	0	608	37	0
3	C	654	0	608	37	0
3	D	654	0	608	36	0
4	A	2	0	0	0	0
4	B	1	0	0	1	0
4	D	1	0	0	0	0
4	H	2	0	0	1	0
4	I	5	0	0	1	0
4	J	2	0	0	0	0
4	K	2	0	0	0	0
4	L	3	0	0	0	0
4	M	4	0	0	2	0
4	N	1	0	0	0	0
4	O	4	0	0	0	0
All	All	15350	0	14771	610	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 20.

All (610) 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)
2:O:52:ASN:ND2	3:D:43:HIS:CD2	1.99	1.31
2:L:52:ASN:ND2	3:A:43:HIS:CD2	2.07	1.22
1:I:33:THR:CG2	1:I:52:SER:HA	1.70	1.20
2:O:52:ASN:ND2	3:D:43:HIS:HD2	1.37	1.18
2:M:52:ASN:ND2	3:B:43:HIS:CD2	2.12	1.17
2:L:52:ASN:ND2	3:A:43:HIS:HD2	1.42	1.16
1:I:33:THR:HG22	1:I:52:SER:CA	1.78	1.12
1:J:1:GLN:HA	1:J:1:GLN:OE1	1.46	1.12

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:33:THR:CG2	1:J:52:SER:HA	1.86	1.06
1:K:33:THR:HG23	1:K:52:SER:HA	1.34	1.05
2:M:60:ARG:NH1	2:M:81:ASP:OD2	1.88	1.05
1:H:207:HIS:HD2	1:H:210:SER:OG	1.35	1.05
2:M:52:ASN:ND2	3:B:43:HIS:HD2	1.49	1.04
1:H:207:HIS:CD2	1:H:210:SER:OG	2.10	1.04
1:J:33:THR:HG22	1:J:52:SER:CA	1.91	1.01
1:J:134:SER:HB2	1:J:136:LYS:HD2	1.43	0.99
2:N:52:ASN:ND2	3:C:43:HIS:CD2	2.32	0.98
2:L:52:ASN:CG	3:A:43:HIS:CD2	2.38	0.98
1:J:33:THR:HG22	1:J:52:SER:HA	1.00	0.97
3:A:94:GLN:NE2	3:A:97:ARG:HB2	1.79	0.95
2:O:52:ASN:HD21	3:D:43:HIS:HD2	0.99	0.94
2:M:169:GLN:HE21	2:M:175:ALA:HB2	1.31	0.94
2:M:183:THR:HG22	2:M:186:GLN:H	1.33	0.94
3:C:94:GLN:NE2	3:C:97:ARG:HB2	1.83	0.94
2:N:191:ARG:CG	2:N:191:ARG:HH11	1.70	0.93
1:H:17:SER:HB3	1:H:84:ASN:HD22	1.34	0.92
1:I:33:THR:HG22	1:I:52:SER:HA	0.92	0.92
1:K:207:HIS:HD2	1:K:210:SER:OG	1.52	0.91
2:N:169:GLN:HE21	2:N:175:ALA:HB2	1.36	0.91
2:N:183:THR:HG21	2:N:185:GLU:HG2	1.53	0.91
2:O:183:THR:HG22	2:O:186:GLN:H	1.33	0.91
2:O:60:ARG:NH1	2:O:81:ASP:OD2	2.02	0.90
2:N:183:THR:HG22	2:N:186:GLN:H	1.35	0.90
2:L:60:ARG:NH1	2:L:81:ASP:OD2	2.04	0.90
3:A:94:GLN:HE22	3:A:97:ARG:HB2	1.32	0.90
2:M:52:ASN:HD21	3:B:43:HIS:HD2	1.13	0.90
2:L:183:THR:HG21	2:L:185:GLU:HG2	1.51	0.90
2:O:169:GLN:HE21	2:O:175:ALA:HB2	1.34	0.89
1:I:195:SER:O	1:I:199:GLN:HB2	1.71	0.89
2:N:90:PHE:C	2:N:90:PHE:CD1	2.47	0.89
2:M:90:PHE:C	2:M:90:PHE:CD1	2.45	0.89
2:N:52:ASN:ND2	3:C:43:HIS:HD2	1.70	0.88
2:N:60:ARG:NH1	2:N:81:ASP:OD2	2.05	0.88
2:L:52:ASN:HD21	3:A:43:HIS:HD2	1.19	0.88
2:L:6:GLN:HE21	2:L:103:THR:HG23	1.37	0.88
1:H:1:GLN:O	1:H:1:GLN:OE1	1.91	0.88
2:M:90:PHE:C	2:M:90:PHE:HD1	1.77	0.88
2:L:6:GLN:HE21	2:L:103:THR:CG2	1.86	0.87
3:C:94:GLN:HE22	3:C:97:ARG:HB2	1.37	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:88:ALA:O	1:I:91:THR:HG23	1.75	0.87
1:J:134:SER:CB	1:J:136:LYS:HD2	2.05	0.86
3:D:94:GLN:HE22	3:D:97:ARG:HB2	1.39	0.85
1:K:88:ALA:O	1:K:91:THR:HG23	1.76	0.85
2:M:6:GLN:HB3	2:M:103:THR:HG22	1.59	0.85
1:H:185:LEU:C	1:H:185:LEU:HD12	1.96	0.85
2:N:90:PHE:C	2:N:90:PHE:HD1	1.78	0.84
2:M:183:THR:HG21	2:M:185:GLU:HG2	1.57	0.84
1:H:126:PRO:HB3	1:H:152:TYR:HB3	1.60	0.83
3:B:94:GLN:NE2	3:B:97:ARG:HB2	1.93	0.83
2:L:41:GLN:OE1	1:I:122:SER:HB2	1.77	0.82
2:L:90:PHE:CD1	2:L:90:PHE:C	2.53	0.82
2:L:169:GLN:HE21	2:L:175:ALA:HB2	1.45	0.82
2:O:90:PHE:CD1	2:O:90:PHE:C	2.52	0.81
1:J:88:ALA:O	1:J:91:THR:CG2	2.29	0.81
3:D:94:GLN:NE2	3:D:97:ARG:CB	2.43	0.81
3:B:94:GLN:HE22	3:B:97:ARG:HB2	1.46	0.81
2:L:90:PHE:C	2:L:90:PHE:HD1	1.85	0.80
2:L:7:PRO:O	2:L:103:THR:HB	1.82	0.80
3:D:94:GLN:NE2	3:D:97:ARG:HB2	1.94	0.80
2:L:6:GLN:HB3	2:L:103:THR:HG22	1.64	0.80
2:O:90:PHE:C	2:O:90:PHE:HD1	1.85	0.80
2:L:183:THR:CG2	2:L:185:GLU:HG2	2.11	0.79
2:L:183:THR:HG22	2:L:186:GLN:H	1.46	0.79
1:I:88:ALA:O	1:I:91:THR:CG2	2.30	0.79
2:O:183:THR:HG21	2:O:185:GLU:HG2	1.62	0.79
1:K:88:ALA:O	1:K:91:THR:CG2	2.30	0.79
1:H:101:TRP:CZ3	3:A:86:GLN:NE2	2.50	0.79
3:A:94:GLN:NE2	3:A:97:ARG:CB	2.44	0.79
1:I:207:HIS:HD2	1:I:210:SER:OG	1.65	0.79
1:J:1:GLN:CA	1:J:1:GLN:OE1	2.31	0.78
2:O:74:ILE:HG22	2:O:77:THR:HG22	1.64	0.78
1:H:174:PRO:HG2	2:L:167:SER:OG	1.83	0.77
3:C:65:GLU:OE2	3:C:96:ARG:HD3	1.84	0.77
1:I:101:TRP:HH2	3:B:86:GLN:HE22	1.33	0.77
2:L:6:GLN:NE2	2:L:103:THR:HG23	2.00	0.77
2:O:7:PRO:O	2:O:103:THR:HB	1.85	0.77
2:N:52:ASN:CG	3:C:43:HIS:CD2	2.58	0.76
2:N:6:GLN:HB3	2:N:103:THR:HG22	1.66	0.76
2:L:52:ASN:HA	3:A:43:HIS:NE2	2.01	0.76
3:C:94:GLN:NE2	3:C:97:ARG:CB	2.48	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:100:ARG:HG3	3:D:94:GLN:OE1	1.85	0.76
2:L:194:SER:OG	2:L:207:THR:HB	1.86	0.75
2:N:191:ARG:NH1	2:N:191:ARG:CG	2.42	0.75
1:J:207:HIS:HD2	1:J:210:SER:OG	1.68	0.75
2:N:183:THR:CG2	2:N:185:GLU:HG2	2.15	0.75
1:J:88:ALA:O	1:J:91:THR:HG23	1.87	0.74
1:H:17:SER:HB3	1:H:84:ASN:ND2	2.02	0.74
2:M:6:GLN:HE21	2:M:103:THR:HG23	1.50	0.74
2:N:7:PRO:O	2:N:103:THR:HB	1.88	0.74
1:I:31:ASN:HB3	4:B:130:HOH:O	1.87	0.74
1:H:31:ASN:HA	3:A:95:LEU:HD22	1.70	0.74
1:H:176:VAL:HG13	2:L:164:THR:CG2	2.18	0.74
2:O:183:THR:CG2	2:O:185:GLU:HG2	2.17	0.74
2:N:50:ASP:OD1	2:N:65:ASN:HB3	1.87	0.73
2:O:119:LEU:HD13	2:O:208:VAL:HG13	1.69	0.73
2:M:52:ASN:CG	3:B:43:HIS:CD2	2.60	0.73
1:H:178:GLN:HG2	2:L:162:GLU:HG2	1.71	0.73
1:H:17:SER:CB	1:H:84:ASN:HD22	2.02	0.73
2:M:51:SER:HB3	2:M:63:GLY:O	1.89	0.73
3:D:94:GLN:HE21	3:D:97:ARG:HB3	1.54	0.71
1:H:101:TRP:HZ3	3:A:86:GLN:NE2	1.88	0.71
2:N:106:THR:HG21	2:N:143:PRO:HB3	1.71	0.71
2:O:52:ASN:CG	3:D:43:HIS:CD2	2.63	0.71
3:B:65:GLU:OE2	3:B:96:ARG:HD3	1.91	0.71
2:M:106:THR:HG21	2:M:143:PRO:HB3	1.73	0.71
1:K:33:THR:CG2	1:K:52:SER:HA	2.18	0.71
1:H:185:LEU:HD12	1:H:186:SER:N	2.05	0.71
2:M:183:THR:CG2	2:M:185:GLU:HG2	2.20	0.71
2:N:191:ARG:NH1	2:N:191:ARG:HG3	2.06	0.70
2:O:6:GLN:HB3	2:O:103:THR:HG22	1.72	0.70
2:N:74:ILE:HG22	2:N:77:THR:HG22	1.74	0.70
2:M:6:GLN:HE21	2:M:103:THR:CG2	2.05	0.70
1:H:88:ALA:O	1:H:91:THR:HG23	1.91	0.70
1:H:73:ASP:OD1	1:H:75:SER:HB2	1.91	0.70
3:B:94:GLN:NE2	3:B:97:ARG:CB	2.56	0.69
3:A:96:ARG:NH2	3:A:116:PRO:O	2.25	0.69
3:C:116:PRO:HB2	3:C:118:VAL:HG22	1.75	0.69
2:M:119:LEU:HD13	2:M:208:VAL:HG13	1.74	0.69
1:H:52:SER:HB3	1:H:57:LEU:HB2	1.75	0.69
1:J:155:GLU:HG3	1:J:156:PRO:HA	1.75	0.69
2:N:110:GLN:HB2	2:N:111:PRO:HD2	1.74	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:N:52:ASN:HD21	3:C:43:HIS:HD2	1.40	0.68
2:L:49:ASP:O	2:L:51:SER:N	2.26	0.68
1:H:33:THR:C	1:H:34:LEU:HD22	2.13	0.68
1:H:34:LEU:HD23	1:H:72:ARG:NH2	2.09	0.68
1:H:176:VAL:HG13	2:L:164:THR:HG23	1.73	0.68
2:M:6:GLN:NE2	2:M:103:THR:HG23	2.08	0.68
2:O:45:LEU:HD13	2:O:54:PRO:HG3	1.75	0.67
1:J:134:SER:HB2	1:J:136:LYS:CD	2.23	0.67
2:M:60:ARG:NH1	2:M:81:ASP:CG	2.47	0.67
1:H:17:SER:CB	1:H:84:ASN:ND2	2.58	0.67
1:H:17:SER:HB3	1:H:84:ASN:HA	1.76	0.67
2:L:6:GLN:NE2	2:L:103:THR:CG2	2.57	0.67
2:L:119:LEU:HD13	2:L:208:VAL:HG13	1.76	0.66
2:M:74:ILE:HG22	2:M:77:THR:HG22	1.76	0.66
1:I:101:TRP:CZ3	3:B:86:GLN:NE2	2.64	0.66
2:L:106:THR:HG21	2:L:143:PRO:HB3	1.76	0.66
2:N:194:SER:OG	2:N:207:THR:HB	1.95	0.66
1:I:58:THR:HG21	4:I:239:HOH:O	1.95	0.66
2:M:60:ARG:HH12	2:M:81:ASP:CG	1.99	0.66
1:J:134:SER:C	1:J:136:LYS:N	2.49	0.66
2:N:6:GLN:HE21	2:N:103:THR:CG2	2.09	0.66
2:N:191:ARG:HG3	2:N:191:ARG:HH11	1.59	0.65
2:O:46:VAL:O	2:O:47:ILE:HG13	1.96	0.65
2:M:7:PRO:O	2:M:103:THR:HB	1.97	0.65
2:N:60:ARG:HH12	2:N:81:ASP:CG	2.01	0.65
2:O:110:GLN:HB2	2:O:111:PRO:HD2	1.78	0.64
2:O:184:PRO:O	2:O:188:LYS:HG2	1.97	0.64
2:M:90:PHE:HD1	2:M:90:PHE:O	1.78	0.64
2:M:194:SER:OG	2:M:207:THR:HB	1.97	0.64
2:O:15:GLY:O	2:O:76:GLY:HA2	1.98	0.64
1:K:101:TRP:CZ3	3:D:86:GLN:NE2	2.66	0.64
1:I:98:ARG:NH2	3:B:67:ASP:OD2	2.30	0.64
3:A:65:GLU:OE2	3:A:96:ARG:HD3	1.98	0.64
3:D:65:GLU:OE2	3:D:96:ARG:HD3	1.98	0.63
1:J:88:ALA:O	1:J:91:THR:HG22	1.97	0.63
3:C:96:ARG:NH2	3:C:116:PRO:O	2.31	0.63
2:L:31:TYR:CZ	3:A:79:LYS:HG2	2.33	0.63
2:N:6:GLN:HE21	2:N:103:THR:HG23	1.63	0.63
1:K:176:VAL:HG13	2:O:164:THR:CG2	2.28	0.63
1:J:89:GLU:OE1	1:J:89:GLU:N	2.29	0.63
2:M:152:ALA:HB2	2:M:193:TYR:CE1	2.33	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:101:TRP:CH2	3:B:86:GLN:NE2	2.66	0.63
2:O:90:PHE:O	2:O:90:PHE:HD1	1.82	0.63
1:K:102:HIS:ND1	3:D:64:GLU:OE1	2.31	0.62
2:O:194:SER:OG	2:O:207:THR:HB	1.99	0.62
2:N:90:PHE:HD1	2:N:90:PHE:O	1.82	0.62
1:K:195:SER:O	1:K:199:GLN:HB2	1.99	0.62
3:C:45:PRO:C	3:C:47:ASP:H	2.03	0.62
1:K:33:THR:HG22	1:K:51:THR:O	1.99	0.62
2:O:106:THR:HG21	2:O:143:PRO:HB3	1.82	0.62
2:M:183:THR:HG22	2:M:186:GLN:N	2.11	0.62
2:O:6:GLN:HE21	2:O:103:THR:CG2	2.13	0.62
1:J:122:SER:HB2	2:O:41:GLN:OE1	2.00	0.62
1:I:100:ARG:HG3	3:B:94:GLN:OE1	2.00	0.62
1:J:177:LEU:HG	1:J:183:TYR:CE2	2.35	0.62
1:K:51:THR:HG22	1:K:55:GLY:HA2	1.81	0.61
1:H:155:GLU:HG3	1:H:156:PRO:HA	1.81	0.61
3:A:94:GLN:HE21	3:A:97:ARG:CB	2.13	0.61
2:M:183:THR:HB	2:M:186:GLN:OE1	1.99	0.61
2:L:41:GLN:CD	1:I:122:SER:HB2	2.19	0.61
1:K:176:VAL:HG13	2:O:164:THR:HG23	1.82	0.61
2:O:33:ILE:HD13	2:O:48:TYR:O	2.00	0.61
2:M:90:PHE:CE1	2:M:91:THR:O	2.53	0.61
2:O:183:THR:HG22	2:O:186:GLN:N	2.09	0.61
2:L:183:THR:HB	2:L:186:GLN:OE1	2.00	0.61
2:L:110:GLN:HB2	2:L:111:PRO:CD	2.31	0.61
1:K:207:HIS:CD2	1:K:210:SER:OG	2.44	0.61
2:L:110:GLN:HB2	2:L:111:PRO:HD2	1.83	0.60
2:L:46:VAL:O	2:L:47:ILE:HG13	2.00	0.60
3:C:104:THR:HB	3:C:107:CYS:HB3	1.83	0.60
1:K:89:GLU:OE1	1:K:89:GLU:N	2.34	0.60
1:J:178:GLN:HG3	1:J:182:LEU:O	2.01	0.60
1:H:219:GLU:HB3	1:H:220:PRO:CD	2.32	0.60
2:N:52:ASN:CG	3:C:43:HIS:HD2	2.03	0.60
1:H:91:THR:HB	1:H:117:THR:HA	1.82	0.60
1:K:31:ASN:HA	3:D:95:LEU:HD22	1.83	0.60
1:I:19:ARG:HD3	1:I:82:GLN:NE2	2.16	0.60
3:C:94:GLN:HE21	3:C:97:ARG:CB	2.13	0.60
1:H:159:VAL:HG21	1:H:187:SER:HB2	1.84	0.60
1:I:185:LEU:C	1:I:185:LEU:HD12	2.21	0.60
1:H:100:ARG:HG3	3:A:94:GLN:OE1	2.02	0.60
3:B:65:GLU:HG2	3:B:96:ARG:HB3	1.84	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:184:PRO:O	2:L:188:LYS:HG2	2.01	0.59
1:H:208:LYS:HB2	1:H:209:PRO:HD3	1.83	0.59
2:N:140:ASP:H	2:N:169:GLN:HE22	1.50	0.59
1:I:178:GLN:HG3	1:I:182:LEU:O	2.01	0.59
3:D:35:PHE:N	3:D:35:PHE:CD2	2.69	0.59
1:I:51:THR:HG22	1:I:55:GLY:HA2	1.84	0.59
1:I:176:VAL:HG13	2:M:164:THR:HG23	1.85	0.59
2:L:49:ASP:O	2:L:50:ASP:C	2.35	0.59
1:H:34:LEU:HB3	1:H:79:LEU:HD22	1.85	0.59
2:M:52:ASN:HA	3:B:43:HIS:NE2	2.17	0.58
2:M:110:GLN:HB2	2:M:111:PRO:HD2	1.84	0.58
2:L:60:ARG:NH1	2:L:81:ASP:CG	2.56	0.58
1:H:217:LYS:NZ	1:H:219:GLU:OE2	2.34	0.58
1:H:219:GLU:HB3	1:H:220:PRO:HD2	1.84	0.58
1:H:19:ARG:HD3	1:H:82:GLN:NE2	2.19	0.58
1:J:101:TRP:CZ3	3:C:86:GLN:NE2	2.71	0.58
1:H:94:TYR:O	1:H:113:GLY:HA2	2.03	0.58
2:O:60:ARG:NH1	2:O:81:ASP:CG	2.55	0.58
2:L:90:PHE:O	2:L:90:PHE:HD1	1.85	0.58
2:M:110:GLN:HB2	2:M:111:PRO:CD	2.33	0.58
2:N:52:ASN:HD21	3:C:43:HIS:CD2	2.17	0.58
2:N:90:PHE:CE1	2:N:91:THR:O	2.57	0.58
1:I:91:THR:HB	1:I:117:THR:HA	1.85	0.58
1:J:83:MET:C	1:J:84:ASN:HD22	2.06	0.58
2:N:110:GLN:HB2	2:N:111:PRO:CD	2.33	0.58
3:A:35:PHE:CD2	3:A:35:PHE:N	2.70	0.58
2:O:183:THR:HB	2:O:186:GLN:OE1	2.04	0.57
1:H:102:HIS:ND1	3:A:64:GLU:OE1	2.38	0.57
2:N:54:PRO:HG2	2:N:57:ILE:HG12	1.85	0.57
2:O:74:ILE:CG2	2:O:77:THR:HG22	2.33	0.57
1:H:101:TRP:HH2	3:A:86:GLN:HE22	1.48	0.57
2:N:51:SER:HB3	2:N:63:GLY:O	2.03	0.57
3:A:65:GLU:HG2	3:A:96:ARG:HB3	1.87	0.57
2:N:53:ARG:HD3	2:N:61:PHE:O	2.05	0.57
3:C:68:GLN:O	3:C:70:GLU:N	2.36	0.57
2:M:6:GLN:HB3	2:M:103:THR:CG2	2.31	0.57
2:L:52:ASN:CG	3:A:43:HIS:HD2	1.94	0.56
2:L:6:GLN:HB3	2:L:103:THR:CG2	2.34	0.56
2:L:140:ASP:H	2:L:169:GLN:HE22	1.52	0.56
2:L:56:GLY:HA2	1:I:12:VAL:HG13	1.86	0.56
1:I:155:GLU:HG3	1:I:156:PRO:HA	1.86	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:36:GLN:HB2	2:L:85:TYR:CE2	2.40	0.56
3:A:94:GLN:HE21	3:A:97:ARG:HB3	1.70	0.56
2:O:60:ARG:HH12	2:O:81:ASP:CG	2.08	0.56
1:I:175:ALA:HB2	1:I:185:LEU:HB3	1.86	0.56
2:M:210:PRO:O	2:M:211:THR:C	2.44	0.56
3:C:35:PHE:H	3:C:35:PHE:HD2	1.54	0.56
2:M:90:PHE:CZ	2:M:92:MET:SD	2.98	0.56
2:M:6:GLN:NE2	2:M:103:THR:CG2	2.67	0.56
1:K:101:TRP:HZ3	3:D:86:GLN:NE2	2.02	0.56
1:H:155:GLU:HG3	1:H:156:PRO:CA	2.36	0.56
1:J:134:SER:C	1:J:136:LYS:H	2.09	0.56
2:M:90:PHE:CE1	2:M:92:MET:SD	2.99	0.56
2:N:15:GLY:O	2:N:76:GLY:HA2	2.06	0.56
1:H:34:LEU:HD22	1:H:34:LEU:N	2.21	0.56
3:D:96:ARG:NH2	3:D:116:PRO:O	2.39	0.56
1:J:101:TRP:HH2	3:C:86:GLN:HE22	1.54	0.56
1:H:185:LEU:C	1:H:185:LEU:CD1	2.68	0.55
1:I:149:VAL:CG2	1:I:149:VAL:O	2.54	0.55
3:C:94:GLN:HE21	3:C:97:ARG:HB3	1.71	0.55
2:L:60:ARG:HH12	2:L:81:ASP:CG	2.10	0.55
3:A:35:PHE:H	3:A:35:PHE:HD2	1.54	0.55
3:D:35:PHE:N	3:D:35:PHE:HD2	2.02	0.55
2:L:115:PRO:HB3	2:L:141:PHE:HB3	1.88	0.55
2:M:31:TYR:CZ	3:B:79:LYS:HG2	2.41	0.55
2:L:210:PRO:O	2:L:211:THR:C	2.44	0.55
1:H:122:SER:O	1:H:123:THR:O	2.24	0.55
2:N:60:ARG:NH1	2:N:81:ASP:CG	2.58	0.55
2:O:185:GLU:O	2:O:189:SER:HB3	2.06	0.55
2:L:185:GLU:O	2:L:189:SER:HB3	2.06	0.55
2:N:6:GLN:NE2	2:N:103:THR:HG23	2.21	0.55
1:H:34:LEU:HD23	1:H:72:ARG:HH22	1.69	0.55
1:K:178:GLN:HG3	1:K:182:LEU:O	2.07	0.55
3:C:35:PHE:CD2	3:C:35:PHE:N	2.71	0.55
1:H:24:ALA:HB1	1:H:27:PHE:CE1	2.42	0.54
1:I:177:LEU:HG	1:I:183:TYR:CE2	2.42	0.54
1:H:87:ARG:HD3	4:H:235:HOH:O	2.06	0.54
3:C:65:GLU:HG2	3:C:96:ARG:HB3	1.89	0.54
2:L:52:ASN:ND2	3:A:43:HIS:NE2	2.55	0.54
1:H:19:ARG:HB2	1:H:82:GLN:HE22	1.72	0.54
3:C:35:PHE:HD2	3:C:35:PHE:N	2.05	0.54
1:J:19:ARG:HD3	1:J:82:GLN:NE2	2.22	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:74:ILE:HG22	2:L:77:THR:HG22	1.90	0.54
1:H:98:ARG:NH2	3:A:67:ASP:OD2	2.39	0.54
2:N:119:LEU:HD13	2:N:208:VAL:HG13	1.90	0.54
3:B:104:THR:HB	3:B:107:CYS:HB3	1.90	0.54
1:J:185:LEU:HD12	1:J:185:LEU:C	2.27	0.54
2:N:210:PRO:O	2:N:211:THR:C	2.44	0.54
1:K:185:LEU:C	1:K:185:LEU:HD12	2.28	0.54
1:I:83:MET:HE1	1:I:94:TYR:CZ	2.43	0.54
3:B:35:PHE:N	3:B:35:PHE:CD2	2.73	0.54
2:N:49:ASP:O	2:N:51:SER:N	2.41	0.54
1:H:101:TRP:HZ3	3:A:86:GLN:HE21	1.56	0.53
2:N:6:GLN:HB3	2:N:103:THR:CG2	2.37	0.53
2:O:140:ASP:H	2:O:169:GLN:HE22	1.56	0.53
1:I:154:PRO:O	1:I:207:HIS:HE1	1.89	0.53
2:M:95:ASN:HB2	4:M:217:HOH:O	2.08	0.53
2:M:86:TYR:CE2	2:M:102:GLY:HA3	2.43	0.53
1:H:119:SER:OG	1:H:121:ALA:HB3	2.09	0.53
1:H:172:THR:HA	1:H:187:SER:HA	1.89	0.53
1:I:89:GLU:OE1	1:I:89:GLU:N	2.39	0.53
1:K:24:ALA:HB1	1:K:27:PHE:CE1	2.44	0.53
3:D:35:PHE:HD2	3:D:35:PHE:H	1.54	0.53
3:A:35:PHE:HD2	3:A:35:PHE:N	2.07	0.53
2:M:54:PRO:HG2	2:M:57:ILE:HG12	1.90	0.53
1:H:185:LEU:CD1	1:H:186:SER:N	2.71	0.53
1:H:57:LEU:HD23	3:A:91:PRO:HB2	1.90	0.53
3:B:40:CYS:HB3	3:B:76:GLY:HA2	1.90	0.53
3:B:35:PHE:HD2	3:B:35:PHE:N	2.07	0.52
2:L:58:PRO:HA	1:I:16:GLY:HA2	1.91	0.52
2:N:90:PHE:CZ	2:N:92:MET:SD	3.01	0.52
1:K:19:ARG:HD3	1:K:82:GLN:NE2	2.24	0.52
2:L:15:GLY:O	2:L:76:GLY:HA2	2.08	0.52
2:N:183:THR:HG22	2:N:186:GLN:N	2.15	0.52
1:K:149:VAL:O	1:K:149:VAL:CG2	2.58	0.52
1:H:162:ASN:HD21	1:H:201:TYR:HA	1.74	0.52
1:H:176:VAL:HG13	2:L:164:THR:HG22	1.90	0.52
1:I:83:MET:CE	1:I:94:TYR:CZ	2.93	0.52
1:H:162:ASN:ND2	1:H:202:ILE:H	2.08	0.52
3:A:104:THR:HB	3:A:107:CYS:HB3	1.89	0.52
1:K:75:SER:C	1:K:76:LYS:O	2.45	0.52
2:N:184:PRO:O	2:N:188:LYS:HG2	2.09	0.52
3:C:36:LEU:HD23	3:C:105:ASN:HD22	1.74	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:N:6:GLN:NE2	2:N:103:THR:CG2	2.73	0.52
1:J:100:ARG:HG3	3:C:94:GLN:OE1	2.09	0.52
1:H:176:VAL:CG1	2:L:164:THR:HG22	2.39	0.52
2:L:54:PRO:HG2	2:L:57:ILE:HG12	1.92	0.52
2:M:33:ILE:HD13	2:M:48:TYR:O	2.09	0.52
1:H:83:MET:HE2	1:H:86:LEU:HD21	1.91	0.52
2:L:183:THR:HG22	2:L:185:GLU:N	2.25	0.51
3:B:94:GLN:HE21	3:B:97:ARG:CB	2.23	0.51
3:B:94:GLN:HE21	3:B:97:ARG:HB3	1.75	0.51
1:H:161:TRP:O	1:H:162:ASN:C	2.47	0.51
2:N:152:ALA:HB2	2:N:193:TYR:CE1	2.45	0.51
1:I:200:THR:HG22	1:I:200:THR:O	2.10	0.51
2:O:90:PHE:CE1	2:O:92:MET:SD	3.03	0.51
1:H:88:ALA:O	1:H:91:THR:CG2	2.58	0.51
1:H:205:VAL:O	1:H:205:VAL:HG12	2.09	0.51
1:K:177:LEU:HG	1:K:183:TYR:CE2	2.46	0.51
2:O:51:SER:HB3	2:O:63:GLY:O	2.10	0.51
1:I:34:LEU:HB3	1:I:79:LEU:HD22	1.91	0.51
2:M:140:ASP:H	2:M:169:GLN:HE22	1.58	0.51
1:H:145:LEU:C	1:H:145:LEU:HD12	2.32	0.51
2:N:33:ILE:HD13	2:N:48:TYR:O	2.11	0.51
1:J:192:PRO:O	1:J:195:SER:OG	2.20	0.51
2:M:45:LEU:HD13	2:M:54:PRO:HG3	1.92	0.51
1:K:91:THR:HB	1:K:117:THR:HA	1.93	0.50
1:H:48:VAL:HG13	1:H:64:VAL:HG21	1.93	0.50
2:N:90:PHE:CE1	2:N:92:MET:SD	3.04	0.50
2:L:33:ILE:HD13	2:L:48:TYR:O	2.12	0.50
3:A:36:LEU:HD23	3:A:105:ASN:HD22	1.77	0.50
1:J:31:ASN:HA	3:C:95:LEU:HD22	1.94	0.50
1:K:101:TRP:HH2	3:D:86:GLN:HE22	1.54	0.50
2:M:115:PRO:HB3	2:M:141:PHE:HB3	1.93	0.50
2:N:49:ASP:O	2:N:50:ASP:C	2.48	0.50
1:H:75:SER:C	1:H:76:LYS:O	2.47	0.50
3:B:35:PHE:HD2	3:B:35:PHE:H	1.59	0.50
1:J:51:THR:HG22	1:J:55:GLY:HA2	1.94	0.50
2:N:183:THR:HB	2:N:186:GLN:OE1	2.11	0.50
1:H:176:VAL:CG1	2:L:164:THR:CG2	2.88	0.50
2:L:75:SER:O	2:L:76:GLY:C	2.49	0.50
2:M:190:HIS:HE1	3:D:82:GLY:O	1.95	0.50
3:D:65:GLU:HG2	3:D:96:ARG:HB3	1.93	0.50
3:D:50:ASN:ND2	3:D:50:ASN:O	2.39	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:40:CYS:HB3	3:D:76:GLY:HA2	1.94	0.50
2:L:90:PHE:CE1	2:L:91:THR:O	2.65	0.49
1:J:91:THR:HB	1:J:117:THR:HA	1.93	0.49
1:J:175:ALA:HB2	1:J:185:LEU:HB3	1.94	0.49
3:C:50:ASN:HD22	3:C:50:ASN:C	2.15	0.49
1:J:34:LEU:HB3	1:J:79:LEU:HD22	1.93	0.49
2:M:90:PHE:CD1	2:M:91:THR:N	2.79	0.49
2:O:6:GLN:NE2	2:O:103:THR:CG2	2.74	0.49
1:J:154:PRO:O	1:J:207:HIS:HE1	1.95	0.49
2:O:110:GLN:HB2	2:O:111:PRO:CD	2.42	0.49
1:I:176:VAL:HG13	2:M:164:THR:CG2	2.41	0.49
2:N:46:VAL:O	2:N:47:ILE:HG13	2.12	0.49
1:K:34:LEU:HB3	1:K:79:LEU:HD22	1.94	0.49
2:O:6:GLN:HE21	2:O:103:THR:HG23	1.76	0.49
2:M:75:SER:O	2:M:76:GLY:C	2.51	0.49
2:M:184:PRO:O	2:M:188:LYS:HG2	2.13	0.49
2:O:53:ARG:HD3	2:O:61:PHE:O	2.12	0.49
2:M:15:GLY:O	2:M:76:GLY:HA2	2.12	0.49
1:K:72:ARG:HD3	1:K:74:ASN:OD1	2.13	0.49
2:M:52:ASN:N	2:M:52:ASN:HD22	2.11	0.48
1:I:83:MET:HB3	1:I:86:LEU:HD21	1.93	0.48
1:I:185:LEU:HD12	1:I:185:LEU:O	2.14	0.48
2:O:115:PRO:HB3	2:O:141:PHE:HB3	1.95	0.48
1:J:24:ALA:HB1	1:J:27:PHE:CE1	2.47	0.48
2:O:210:PRO:O	2:O:211:THR:C	2.52	0.48
3:D:104:THR:HB	3:D:107:CYS:HB3	1.96	0.48
2:M:90:PHE:CD1	2:M:90:PHE:O	2.61	0.48
2:O:54:PRO:HG2	2:O:57:ILE:HG12	1.96	0.48
1:J:98:ARG:NH2	3:C:67:ASP:OD2	2.40	0.48
1:J:191:VAL:HG21	1:J:201:TYR:CE1	2.49	0.48
1:H:40:ALA:O	1:H:41:PRO:C	2.52	0.48
1:K:154:PRO:O	1:K:207:HIS:HE1	1.97	0.48
3:C:68:GLN:C	3:C:70:GLU:H	2.17	0.48
2:N:90:PHE:CD1	2:N:91:THR:N	2.81	0.48
1:H:36:TRP:NE1	1:H:81:LEU:HB2	2.29	0.48
2:M:52:ASN:ND2	3:B:43:HIS:NE2	2.59	0.48
3:B:36:LEU:HD23	3:B:105:ASN:HD22	1.79	0.48
1:I:75:SER:C	1:I:76:LYS:O	2.52	0.48
1:H:191:VAL:HG11	1:H:201:TYR:CE2	2.48	0.48
1:H:208:LYS:HB2	1:H:209:PRO:CD	2.44	0.47
2:N:45:LEU:HD13	2:N:54:PRO:HG3	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:40:ALA:O	1:I:41:PRO:C	2.51	0.47
1:H:160:SER:OG	1:H:204:ASN:HB2	2.14	0.47
1:I:208:LYS:N	1:I:209:PRO:HD2	2.28	0.47
1:I:33:THR:HG22	1:I:52:SER:C	2.35	0.47
1:H:208:LYS:N	1:H:209:PRO:HD2	2.29	0.47
1:J:101:TRP:HZ3	3:C:86:GLN:NE2	2.10	0.47
1:I:24:ALA:HB1	1:I:27:PHE:CE1	2.50	0.47
2:N:185:GLU:O	2:N:189:SER:HB3	2.13	0.47
2:N:34:TRP:O	2:N:46:VAL:HG22	2.14	0.47
2:N:75:SER:O	2:N:76:GLY:C	2.52	0.47
3:B:111:LEU:HD12	3:B:111:LEU:HA	1.70	0.47
3:B:50:ASN:O	3:B:50:ASN:ND2	2.33	0.47
3:D:36:LEU:HD23	3:D:105:ASN:HD22	1.78	0.47
3:C:111:LEU:HA	3:C:111:LEU:HD12	1.62	0.47
1:K:33:THR:C	1:K:34:LEU:HD22	2.35	0.47
2:N:74:ILE:CG2	2:N:77:THR:HG22	2.43	0.47
1:K:32:TYR:O	1:K:72:ARG:NH2	2.46	0.47
2:O:52:ASN:HA	3:D:43:HIS:NE2	2.30	0.47
1:K:175:ALA:HB2	1:K:185:LEU:HB3	1.95	0.47
2:O:152:ALA:HB2	2:O:193:TYR:CE1	2.50	0.47
1:K:149:VAL:O	1:K:149:VAL:HG23	2.15	0.47
1:K:34:LEU:HD13	1:K:98:ARG:HA	1.97	0.47
2:M:60:ARG:NH1	2:M:81:ASP:OD1	2.46	0.47
2:O:6:GLN:NE2	2:O:103:THR:HG23	2.29	0.47
3:A:40:CYS:HB3	3:A:76:GLY:HA2	1.96	0.47
2:N:140:ASP:H	2:N:169:GLN:NE2	2.13	0.46
1:H:33:THR:HB	1:H:99:GLU:HB3	1.96	0.46
2:N:115:PRO:HB3	2:N:141:PHE:HB3	1.96	0.46
1:I:88:ALA:O	1:I:91:THR:HG22	2.11	0.46
1:K:219:GLU:HA	1:K:220:PRO:HD2	1.69	0.46
2:N:46:VAL:HA	2:N:57:ILE:HG13	1.98	0.46
2:L:56:GLY:CA	1:I:12:VAL:HG13	2.46	0.46
2:M:4:LEU:HD22	2:M:98:VAL:HG13	1.98	0.46
1:K:83:MET:CE	1:K:94:TYR:CZ	2.98	0.46
2:L:60:ARG:NH1	2:L:81:ASP:OD1	2.48	0.46
2:L:51:SER:HB3	2:L:63:GLY:O	2.16	0.46
1:K:195:SER:O	1:K:199:GLN:N	2.47	0.46
1:I:149:VAL:HG23	1:I:149:VAL:O	2.16	0.46
3:B:82:GLY:O	2:O:190:HIS:HE1	1.99	0.46
1:H:101:TRP:CH2	3:A:86:GLN:NE2	2.73	0.46
1:J:134:SER:O	1:J:136:LYS:N	2.48	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:155:GLU:HG3	1:K:156:PRO:HA	1.98	0.46
3:B:68:GLN:O	3:B:70:GLU:N	2.48	0.46
3:C:40:CYS:HB3	3:C:76:GLY:HA2	1.98	0.46
2:O:37:GLN:HE21	2:O:43:PRO:HG3	1.81	0.46
2:M:52:ASN:HD21	3:B:43:HIS:CD2	2.00	0.45
1:J:208:LYS:N	1:J:209:PRO:HD2	2.31	0.45
2:M:171:ASN:O	2:M:172:ASN:HB2	2.15	0.45
2:M:95:ASN:CB	4:M:217:HOH:O	2.64	0.45
1:J:73:ASP:C	1:J:73:ASP:OD1	2.55	0.45
1:H:155:GLU:CG	1:H:156:PRO:HA	2.46	0.45
2:O:52:ASN:ND2	3:D:43:HIS:NE2	2.56	0.45
2:N:52:ASN:ND2	3:C:43:HIS:NE2	2.64	0.45
1:K:88:ALA:O	1:K:91:THR:HG22	2.11	0.45
2:O:90:PHE:CE1	2:O:91:THR:O	2.69	0.45
1:H:6:GLU:OE2	1:H:111:GLY:HA3	2.16	0.45
2:M:46:VAL:HA	2:M:57:ILE:HG13	1.98	0.45
2:M:183:THR:HG22	2:M:185:GLU:N	2.30	0.45
2:N:60:ARG:NH1	2:N:81:ASP:OD1	2.49	0.45
1:J:34:LEU:HD13	1:J:98:ARG:HA	1.99	0.45
1:I:175:ALA:HB1	1:I:184:SER:O	2.17	0.45
1:I:91:THR:HB	1:I:118:VAL:H	1.82	0.45
1:I:33:THR:C	1:I:34:LEU:HD22	2.37	0.45
1:H:185:LEU:HD12	1:H:186:SER:CA	2.47	0.45
1:H:162:ASN:HD21	1:H:202:ILE:H	1.65	0.45
3:D:68:GLN:OE1	3:D:68:GLN:N	2.50	0.45
2:L:86:TYR:CE2	2:L:102:GLY:HA3	2.52	0.45
3:B:50:ASN:HD22	3:B:50:ASN:C	2.07	0.44
3:A:38:CYS:O	3:A:52:THR:HA	2.17	0.44
1:J:18:LEU:HD23	1:J:18:LEU:HA	1.78	0.44
3:D:94:GLN:NE2	3:D:97:ARG:HB3	2.15	0.44
1:H:103:VAL:HG12	1:H:104:ARG:HG2	2.00	0.44
1:H:6:GLU:CD	1:H:113:GLY:H	2.21	0.44
3:D:117:PRO:O	3:D:118:VAL:C	2.56	0.44
2:N:86:TYR:CE2	2:N:102:GLY:HA3	2.53	0.44
1:K:91:THR:HB	1:K:118:VAL:H	1.82	0.44
2:M:49:ASP:O	2:M:50:ASP:C	2.54	0.44
2:L:158:LYS:HA	2:L:158:LYS:HD2	1.77	0.44
2:O:60:ARG:NH1	2:O:81:ASP:OD1	2.50	0.44
2:O:90:PHE:CZ	2:O:92:MET:SD	3.11	0.44
2:M:74:ILE:CG2	2:M:77:THR:HG22	2.43	0.44
1:K:12:VAL:HG21	1:K:18:LEU:HG	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:94:GLN:HG3	3:D:94:GLN:H	1.27	0.44
2:N:90:PHE:O	2:N:90:PHE:CD1	2.66	0.43
1:J:99:GLU:OE1	3:C:92:LYS:HE2	2.18	0.43
2:L:140:ASP:H	2:L:169:GLN:NE2	2.14	0.43
2:O:37:GLN:NE2	2:O:43:PRO:HG3	2.32	0.43
1:I:191:VAL:HG21	1:I:201:TYR:CE1	2.53	0.43
2:M:6:GLN:CB	2:M:103:THR:HG22	2.41	0.43
3:C:38:CYS:O	3:C:52:THR:HA	2.18	0.43
1:H:157:VAL:HG22	1:H:207:HIS:HB2	2.00	0.43
3:A:94:GLN:HG3	3:A:94:GLN:H	1.23	0.43
1:J:155:GLU:HG3	1:J:156:PRO:CA	2.46	0.43
1:I:31:ASN:HA	3:B:95:LEU:HD22	1.99	0.43
3:B:38:CYS:O	3:B:52:THR:HA	2.18	0.43
3:D:68:GLN:O	3:D:70:GLU:N	2.52	0.43
2:O:36:GLN:HB2	2:O:85:TYR:CE2	2.54	0.43
1:I:101:TRP:HZ3	3:B:86:GLN:NE2	2.12	0.43
2:N:31:TYR:CD2	2:N:50:ASP:OD2	2.72	0.43
1:H:33:THR:HG23	1:H:51:THR:O	2.18	0.43
1:I:150:LYS:HA	1:I:184:SER:OG	2.18	0.43
1:J:84:ASN:HD22	1:J:84:ASN:N	2.13	0.43
1:H:68:PHE:CD1	1:H:68:PHE:N	2.87	0.43
2:O:52:ASN:HD22	2:O:52:ASN:N	2.16	0.42
1:K:105:GLY:O	2:O:33:ILE:HD12	2.19	0.42
1:K:28:THR:O	1:K:31:ASN:HB2	2.20	0.42
1:H:19:ARG:HD3	1:H:82:GLN:HE22	1.83	0.42
1:J:176:VAL:HG13	2:N:164:THR:HG23	2.01	0.42
1:K:115:LEU:HD13	1:K:117:THR:HG22	2.01	0.42
1:H:149:VAL:HG13	1:H:185:LEU:CD1	2.49	0.42
2:N:46:VAL:O	2:N:54:PRO:HD2	2.19	0.42
1:H:34:LEU:CD2	1:H:34:LEU:N	2.83	0.42
1:I:196:LEU:C	1:I:198:THR:H	2.22	0.42
2:L:152:ALA:HB2	2:L:193:TYR:CE1	2.55	0.42
2:N:183:THR:HG22	2:N:185:GLU:N	2.35	0.42
1:I:207:HIS:HB3	1:I:212:THR:HB	2.01	0.42
2:O:142:TYR:HA	2:O:143:PRO:C	2.40	0.42
1:J:84:ASN:ND2	1:J:84:ASN:N	2.67	0.42
2:O:86:TYR:CE2	2:O:102:GLY:HA3	2.55	0.42
2:N:90:PHE:CE1	2:N:91:THR:C	2.93	0.42
1:H:115:LEU:HD22	1:H:117:THR:HG22	2.02	0.42
1:H:67:ARG:C	1:H:68:PHE:HD1	2.23	0.42
1:I:72:ARG:HD3	1:I:74:ASN:OD1	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:32:TYR:O	1:J:72:ARG:NH2	2.48	0.42
2:N:6:GLN:OE1	2:N:100:GLY:HA3	2.18	0.42
2:O:75:SER:O	2:O:76:GLY:C	2.58	0.42
2:O:48:TYR:C	2:O:49:ASP:O	2.51	0.42
1:J:75:SER:C	1:J:76:LYS:O	2.58	0.42
3:D:111:LEU:HA	3:D:111:LEU:HD12	1.65	0.42
2:M:105:LEU:HD23	2:M:105:LEU:HA	1.86	0.42
1:J:115:LEU:HA	1:J:115:LEU:HD23	1.84	0.42
1:K:176:VAL:CG1	2:O:164:THR:HG22	2.50	0.42
2:M:158:LYS:HA	2:M:158:LYS:HD2	1.80	0.42
1:H:133:PRO:HB3	1:H:145:LEU:HB3	2.01	0.41
2:M:48:TYR:C	2:M:49:ASP:O	2.57	0.41
1:I:185:LEU:C	1:I:185:LEU:CD1	2.87	0.41
1:J:195:SER:O	1:J:199:GLN:CB	2.69	0.41
1:I:196:LEU:C	1:I:198:THR:N	2.72	0.41
1:J:40:ALA:O	1:J:41:PRO:C	2.58	0.41
1:I:191:VAL:HA	1:I:192:PRO:HD2	1.90	0.41
1:J:91:THR:HB	1:J:118:VAL:H	1.86	0.41
1:H:115:LEU:HD13	1:H:117:THR:HG22	2.02	0.41
3:A:111:LEU:HD12	3:A:111:LEU:HA	1.74	0.41
1:K:153:PHE:HA	1:K:154:PRO:HA	1.74	0.41
2:L:47:ILE:HD12	2:L:72:LEU:CD1	2.50	0.41
2:O:158:LYS:HD2	2:O:158:LYS:HA	1.77	0.41
1:J:33:THR:C	1:J:34:LEU:HD22	2.41	0.41
3:C:49:ILE:HG12	3:C:50:ASN:N	2.36	0.41
3:A:50:ASN:HD22	3:A:50:ASN:C	2.16	0.41
2:M:140:ASP:H	2:M:169:GLN:NE2	2.18	0.41
2:M:86:TYR:CD2	2:M:102:GLY:HA3	2.56	0.41
1:K:83:MET:HE1	1:K:94:TYR:CZ	2.55	0.41
3:B:68:GLN:C	3:B:70:GLU:H	2.24	0.41
1:J:176:VAL:HG13	2:N:164:THR:CG2	2.51	0.41
3:D:60:PHE:HB3	3:D:78:MET:HE3	2.03	0.41
2:M:36:GLN:HB2	2:M:85:TYR:CE2	2.56	0.41
3:A:68:GLN:O	3:A:70:GLU:N	2.53	0.41
2:N:158:LYS:HD2	2:N:158:LYS:HA	1.86	0.41
1:I:199:GLN:HB3	1:I:201:TYR:CE2	2.55	0.41
2:M:49:ASP:O	2:M:51:SER:N	2.54	0.41
2:M:185:GLU:O	2:M:189:SER:HB3	2.20	0.41
2:L:183:THR:HG22	2:L:186:GLN:N	2.23	0.41
1:I:195:SER:O	1:I:199:GLN:CB	2.57	0.41
1:J:207:HIS:HB3	1:J:212:THR:HB	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:12:VAL:HG21	1:I:18:LEU:HG	2.03	0.41
3:B:68:GLN:N	3:B:68:GLN:OE1	2.54	0.41
1:K:208:LYS:N	1:K:209:PRO:HD2	2.36	0.41
2:O:29:SER:O	3:D:81:GLU:HG2	2.21	0.41
2:N:27:LEU:HA	2:N:27:LEU:HD23	1.79	0.41
3:C:94:GLN:H	3:C:94:GLN:HG3	1.27	0.41
1:J:83:MET:HB3	1:J:86:LEU:HD21	2.03	0.41
1:H:119:SER:OG	1:H:121:ALA:CB	2.69	0.41
3:A:68:GLN:C	3:A:70:GLU:H	2.25	0.41
2:N:171:ASN:O	2:N:172:ASN:HB2	2.20	0.41
2:M:47:ILE:HD12	2:M:72:LEU:CD1	2.51	0.41
2:M:90:PHE:CE1	2:M:91:THR:C	2.94	0.40
3:B:96:ARG:NH2	3:B:116:PRO:O	2.54	0.40
1:K:199:GLN:HB3	1:K:201:TYR:CE2	2.57	0.40
1:H:133:PRO:HD2	1:H:219:GLU:O	2.21	0.40
1:K:176:VAL:CG1	2:O:164:THR:CG2	2.99	0.40
1:H:159:VAL:O	1:H:159:VAL:HG23	2.21	0.40
1:I:32:TYR:O	1:I:72:ARG:NH2	2.48	0.40
1:I:158:THR:OG1	1:I:206:ASN:HB3	2.22	0.40
2:O:27:LEU:HD23	2:O:27:LEU:HA	1.84	0.40
3:D:94:GLN:NE2	3:D:97:ARG:CD	2.84	0.40
2:O:90:PHE:CD1	2:O:91:THR:N	2.89	0.40
2:M:188:LYS:HE3	2:M:188:LYS:HB3	1.97	0.40
1:I:141:GLY:C	1:I:142:THR:HG22	2.42	0.40
2:O:4:LEU:HD22	2:O:98:VAL:HG13	2.04	0.40
1:J:141:GLY:C	1:J:193:SER:OG	2.60	0.40
1:K:105:GLY:O	2:O:33:ILE:CD1	2.69	0.40
1:J:14:PRO:HD2	1:J:120:SER:HB3	2.04	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	H	201/234 (86%)	171 (85%)	24 (12%)	6 (3%)	5	13
1	I	211/234 (90%)	202 (96%)	8 (4%)	1 (0%)	34	63
1	J	219/234 (94%)	205 (94%)	13 (6%)	1 (0%)	34	63
1	K	211/234 (90%)	201 (95%)	10 (5%)	0	100	100
2	L	208/213 (98%)	199 (96%)	7 (3%)	2 (1%)	19	45
2	M	208/213 (98%)	200 (96%)	7 (3%)	1 (0%)	34	63
2	N	208/213 (98%)	202 (97%)	5 (2%)	1 (0%)	34	63
2	O	208/213 (98%)	202 (97%)	5 (2%)	1 (0%)	34	63
3	A	83/129 (64%)	74 (89%)	8 (10%)	1 (1%)	16	39
3	B	83/129 (64%)	72 (87%)	9 (11%)	2 (2%)	7	19
3	C	83/129 (64%)	74 (89%)	7 (8%)	2 (2%)	7	19
3	D	83/129 (64%)	74 (89%)	7 (8%)	2 (2%)	7	19
All	All	2006/2304 (87%)	1876 (94%)	110 (6%)	20 (1%)	19	45

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	H	123	THR
1	H	192	PRO
2	L	76	GLY
2	L	94	GLY
2	M	76	GLY
2	N	76	GLY
3	C	69	GLY
2	O	76	GLY
1	H	163	SER
1	H	211	ASN
1	J	135	SER
1	H	76	LYS
1	H	162	ASN
3	B	117	PRO
3	C	46	ASP
3	D	110	TYR
3	A	69	GLY
1	I	76	LYS
3	B	69	GLY
3	D	69	GLY

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	H	175/198 (88%)	145 (83%)	30 (17%)	2	6
1	I	181/198 (91%)	150 (83%)	31 (17%)	2	6
1	J	186/198 (94%)	151 (81%)	35 (19%)	2	5
1	K	182/198 (92%)	148 (81%)	34 (19%)	2	5
2	L	176/178 (99%)	148 (84%)	28 (16%)	3	8
2	M	176/178 (99%)	152 (86%)	24 (14%)	5	11
2	N	176/178 (99%)	151 (86%)	25 (14%)	4	10
2	O	176/178 (99%)	149 (85%)	27 (15%)	3	8
3	A	76/114 (67%)	62 (82%)	14 (18%)	2	5
3	B	76/114 (67%)	59 (78%)	17 (22%)	1	3
3	C	76/114 (67%)	61 (80%)	15 (20%)	1	4
3	D	76/114 (67%)	60 (79%)	16 (21%)	1	3
All	All	1732/1960 (88%)	1436 (83%)	296 (17%)	2	6

All (296) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	H	1	GLN
1	H	5	VAL
1	H	11	LEU
1	H	12	VAL
1	H	17	SER
1	H	19	ARG
1	H	22	CYS
1	H	28	THR
1	H	51	THR
1	H	54	SER
1	H	57	LEU
1	H	58	THR
1	H	65	LYS
1	H	75	SER

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Mol	Chain	Res	Type
1	H	91	THR
1	H	98	ARG
1	H	112	GLN
1	H	114	THR
1	H	115	LEU
1	H	117	THR
1	H	149	VAL
1	H	159	VAL
1	H	166	LEU
1	H	167	THR
1	H	171	HIS
1	H	176	VAL
1	H	185	LEU
1	H	190	THR
1	H	191	VAL
1	H	216	LYS
2	L	2	ILE
2	L	4	LEU
2	L	9	SER
2	L	12	VAL
2	L	16	GLN
2	L	27	LEU
2	L	33	ILE
2	L	45	LEU
2	L	50	ASP
2	L	52	ASN
2	L	55	SER
2	L	80	GLU
2	L	90	PHE
2	L	92	MET
2	L	95	ASN
2	L	103	THR
2	L	105	LEU
2	L	119	LEU
2	L	131	LYS
2	L	158	LYS
2	L	170	SER
2	L	183	THR
2	L	188	LYS
2	L	191	ARG
2	L	204	VAL
2	L	205	GLU

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Mol	Chain	Res	Type
2	L	207	THR
2	L	211	THR
3	A	35	PHE
3	A	36	LEU
3	A	39	TYR
3	A	49	ILE
3	A	50	ASN
3	A	52	THR
3	A	68	GLN
3	A	73	LEU
3	A	94	GLN
3	A	95	LEU
3	A	96	ARG
3	A	105	ASN
3	A	109	GLN
3	A	112	GLN
1	I	5	VAL
1	I	11	LEU
1	I	18	LEU
1	I	21	SER
1	I	22	CYS
1	I	28	THR
1	I	51	THR
1	I	54	SER
1	I	57	LEU
1	I	58	THR
1	I	72	ARG
1	I	91	THR
1	I	98	ARG
1	I	112	GLN
1	I	114	THR
1	I	115	LEU
1	I	117	THR
1	I	134	SER
1	I	135	SER
1	I	142	THR
1	I	149	VAL
1	I	166	LEU
1	I	176	VAL
1	I	177	LEU
1	I	179	SER
1	I	184	SER

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Mol	Chain	Res	Type
1	I	185	LEU
1	I	196	LEU
1	I	199	GLN
1	I	200	THR
1	I	216	LYS
2	M	2	ILE
2	M	4	LEU
2	M	12	VAL
2	M	27	LEU
2	M	33	ILE
2	M	45	LEU
2	M	50	ASP
2	M	52	ASN
2	M	55	SER
2	M	78	GLN
2	M	80	GLU
2	M	90	PHE
2	M	92	MET
2	M	103	THR
2	M	105	LEU
2	M	112	LYS
2	M	119	LEU
2	M	158	LYS
2	M	170	SER
2	M	183	THR
2	M	204	VAL
2	M	205	GLU
2	M	207	THR
2	M	211	THR
3	B	35	PHE
3	B	36	LEU
3	B	39	TYR
3	B	49	ILE
3	B	50	ASN
3	B	52	THR
3	B	68	GLN
3	B	73	LEU
3	B	75	SER
3	B	83	SER
3	B	84	ASP
3	B	94	GLN
3	B	95	LEU

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Mol	Chain	Res	Type
3	B	96	ARG
3	B	105	ASN
3	B	109	GLN
3	B	112	GLN
1	J	1	GLN
1	J	5	VAL
1	J	11	LEU
1	J	12	VAL
1	J	18	LEU
1	J	22	CYS
1	J	28	THR
1	J	51	THR
1	J	54	SER
1	J	57	LEU
1	J	58	THR
1	J	72	ARG
1	J	91	THR
1	J	98	ARG
1	J	112	GLN
1	J	114	THR
1	J	115	LEU
1	J	117	THR
1	J	134	SER
1	J	135	SER
1	J	136	LYS
1	J	142	THR
1	J	149	VAL
1	J	159	VAL
1	J	166	LEU
1	J	176	VAL
1	J	177	LEU
1	J	179	SER
1	J	184	SER
1	J	188	VAL
1	J	196	LEU
1	J	199	GLN
1	J	200	THR
1	J	216	LYS
1	J	220	PRO
2	N	2	ILE
2	N	4	LEU
2	N	12	VAL

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Mol	Chain	Res	Type
2	N	27	LEU
2	N	33	ILE
2	N	45	LEU
2	N	52	ASN
2	N	55	SER
2	N	78	GLN
2	N	80	GLU
2	N	90	PHE
2	N	92	MET
2	N	103	THR
2	N	105	LEU
2	N	119	LEU
2	N	131	LYS
2	N	158	LYS
2	N	170	SER
2	N	183	THR
2	N	188	LYS
2	N	191	ARG
2	N	204	VAL
2	N	205	GLU
2	N	207	THR
2	N	211	THR
3	C	35	PHE
3	C	36	LEU
3	C	39	TYR
3	C	47	ASP
3	C	49	ILE
3	C	50	ASN
3	C	52	THR
3	C	68	GLN
3	C	73	LEU
3	C	84	ASP
3	C	94	GLN
3	C	95	LEU
3	C	96	ARG
3	C	105	ASN
3	C	109	GLN
1	K	1	GLN
1	K	5	VAL
1	K	11	LEU
1	K	12	VAL
1	K	18	LEU

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Mol	Chain	Res	Type
1	K	22	CYS
1	K	28	THR
1	K	33	THR
1	K	51	THR
1	K	54	SER
1	K	57	LEU
1	K	58	THR
1	K	65	LYS
1	K	72	ARG
1	K	91	THR
1	K	112	GLN
1	K	114	THR
1	K	115	LEU
1	K	117	THR
1	K	134	SER
1	K	135	SER
1	K	142	THR
1	K	149	VAL
1	K	156	PRO
1	K	159	VAL
1	K	166	LEU
1	K	176	VAL
1	K	177	LEU
1	K	179	SER
1	K	196	LEU
1	K	199	GLN
1	K	200	THR
1	K	216	LYS
1	K	220	PRO
2	O	2	ILE
2	O	4	LEU
2	O	9	SER
2	O	16	GLN
2	O	27	LEU
2	O	33	ILE
2	O	45	LEU
2	O	50	ASP
2	O	52	ASN
2	O	55	SER
2	O	78	GLN
2	O	80	GLU
2	O	90	PHE

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Mol	Chain	Res	Type
2	O	92	MET
2	O	103	THR
2	O	105	LEU
2	O	119	LEU
2	O	130	ASN
2	O	158	LYS
2	O	170	SER
2	O	183	THR
2	O	188	LYS
2	O	191	ARG
2	O	204	VAL
2	O	205	GLU
2	O	207	THR
2	O	211	THR
3	D	35	PHE
3	D	36	LEU
3	D	39	TYR
3	D	49	ILE
3	D	50	ASN
3	D	52	THR
3	D	68	GLN
3	D	73	LEU
3	D	75	SER
3	D	94	GLN
3	D	95	LEU
3	D	96	ARG
3	D	105	ASN
3	D	109	GLN
3	D	112	GLN
3	D	118	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (68) such sidechains are listed below:

Mol	Chain	Res	Type
1	H	1	GLN
1	H	31	ASN
1	H	39	GLN
1	H	82	GLN
1	H	84	ASN
1	H	162	ASN
1	H	207	HIS
2	L	16	GLN

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Mol	Chain	Res	Type
2	L	36	GLN
2	L	52	ASN
2	L	65	ASN
2	L	169	GLN
2	L	190	HIS
3	A	43	HIS
3	A	50	ASN
3	A	51	ASN
3	A	86	GLN
3	A	105	ASN
3	A	109	GLN
1	I	13	GLN
1	I	31	ASN
1	I	39	GLN
1	I	82	GLN
1	I	84	ASN
1	I	162	ASN
1	I	207	HIS
2	M	52	ASN
2	M	65	ASN
2	M	169	GLN
2	M	190	HIS
3	B	43	HIS
3	B	50	ASN
3	B	51	ASN
3	B	86	GLN
3	B	105	ASN
3	B	109	GLN
1	J	13	GLN
1	J	31	ASN
1	J	82	GLN
1	J	84	ASN
1	J	207	HIS
2	N	37	GLN
2	N	52	ASN
2	N	65	ASN
2	N	169	GLN
2	N	190	HIS
3	C	43	HIS
3	C	51	ASN
3	C	86	GLN
3	C	105	ASN

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Mol	Chain	Res	Type
3	C	109	GLN
1	K	13	GLN
1	K	31	ASN
1	K	82	GLN
1	K	84	ASN
1	K	207	HIS
2	O	36	GLN
2	O	37	GLN
2	O	52	ASN
2	O	65	ASN
2	O	169	GLN
2	O	190	HIS
3	D	43	HIS
3	D	50	ASN
3	D	51	ASN
3	D	86	GLN
3	D	105	ASN
3	D	109	GLN

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 carbohydrates 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 ⓘ

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	H	207/234 (88%)	0.04	8 (3%) 43 43	37, 54, 91, 122	0
1	I	215/234 (91%)	0.29	11 (5%) 32 30	38, 56, 89, 114	0
1	J	221/234 (94%)	0.44	13 (5%) 26 24	44, 64, 102, 125	0
1	K	215/234 (91%)	0.22	9 (4%) 40 39	37, 54, 95, 123	0
2	L	210/213 (98%)	0.33	6 (2%) 55 55	44, 61, 86, 112	0
2	M	210/213 (98%)	0.35	7 (3%) 50 50	41, 59, 77, 92	0
2	N	210/213 (98%)	0.23	7 (3%) 50 50	48, 59, 78, 93	0
2	O	210/213 (98%)	0.10	5 (2%) 62 62	44, 57, 72, 86	0
3	A	85/129 (65%)	0.75	11 (12%) 5 4	51, 73, 107, 127	0
3	B	85/129 (65%)	0.82	11 (12%) 5 4	52, 79, 112, 135	0
3	C	85/129 (65%)	0.87	13 (15%) 3 2	58, 78, 108, 127	0
3	D	85/129 (65%)	1.11	21 (24%) 1 1	50, 80, 131, 157	0
All	All	2038/2304 (88%)	0.36	122 (5%) 25 24	37, 61, 99, 157	0

All (122) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	A	46	ASP	7.5
3	B	118	VAL	7.2
3	B	42	GLY	6.7
3	D	48	ALA	6.1
3	C	34	PRO	5.6
3	A	43	HIS	5.3
3	D	39	TYR	5.0
3	D	45	PRO	4.8
3	D	110	TYR	4.8
3	C	110	TYR	4.6
3	C	39	TYR	4.6

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Mol	Chain	Res	Type	RSRZ
3	D	34	PRO	4.5
3	B	110	TYR	4.5
1	J	197	GLY	4.5
2	M	191	ARG	4.3
3	C	43	HIS	4.3
1	H	135	SER	4.2
3	D	54	ILE	4.1
3	B	43	HIS	4.0
1	I	197	GLY	3.9
3	A	50	ASN	3.8
3	D	46	ASP	3.7
3	D	47	ASP	3.6
2	N	93	SER	3.6
1	I	196	LEU	3.6
1	K	195	SER	3.6
3	A	47	ASP	3.5
3	D	51	ASN	3.5
1	K	192	PRO	3.4
3	B	46	ASP	3.4
3	B	37	LYS	3.4
3	C	50	ASN	3.4
3	A	48	ALA	3.3
2	N	172	ASN	3.3
1	J	198	THR	3.3
3	B	39	TYR	3.2
1	I	1	GLN	3.2
3	D	109	GLN	3.2
3	B	50	ASN	3.2
1	J	163	SER	3.1
3	C	35	PHE	3.0
3	A	51	ASN	3.0
1	H	168	SER	3.0
2	O	2	ILE	3.0
2	O	172	ASN	3.0
1	J	214	VAL	2.9
1	K	135	SER	2.9
1	K	142	THR	2.9
1	J	139	SER	2.9
2	L	124	SER	2.9
1	H	192	PRO	2.8
3	D	105	ASN	2.8
3	A	45	PRO	2.8

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Mol	Chain	Res	Type	RSRZ
3	C	95	LEU	2.8
3	C	48	ALA	2.8
3	D	37	LYS	2.7
2	O	191	ARG	2.7
3	C	49	ILE	2.7
3	A	39	TYR	2.7
2	L	191	ARG	2.7
1	I	182	LEU	2.7
3	A	49	ILE	2.7
1	I	75	SER	2.7
2	L	128	GLN	2.6
1	I	194	SER	2.6
2	L	172	ASN	2.6
3	D	50	ASN	2.6
1	J	135	SER	2.5
1	I	105	GLY	2.5
3	B	41	SER	2.5
3	D	42	GLY	2.5
1	J	166	LEU	2.5
2	N	91	THR	2.5
2	N	158	LYS	2.4
3	D	36	LEU	2.4
1	K	75	SER	2.4
1	K	190	THR	2.4
2	M	192	SER	2.4
2	M	18	ALA	2.4
1	J	26	GLY	2.4
1	K	146	GLY	2.4
2	N	94	GLY	2.4
3	B	47	ASP	2.4
3	C	38	CYS	2.4
3	D	43	HIS	2.4
3	C	112	GLN	2.4
3	C	118	VAL	2.4
1	K	194	SER	2.3
1	H	211	ASN	2.3
1	J	160	SER	2.3
2	M	88	SER	2.3
2	N	95	ASN	2.3
1	H	76	LYS	2.3
3	A	106	LEU	2.2
1	J	179	SER	2.2

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Mol	Chain	Res	Type	RSRZ
1	I	206	ASN	2.2
1	H	208	LYS	2.2
3	D	49	ILE	2.2
1	J	194	SER	2.2
2	M	32	VAL	2.2
2	L	91	THR	2.2
2	M	97	THR	2.2
3	B	51	ASN	2.2
1	H	167	THR	2.2
2	O	170	SER	2.2
2	L	48	TYR	2.2
1	H	144	ALA	2.1
2	N	113	ALA	2.1
3	D	104	THR	2.1
1	J	134	SER	2.1
1	I	207	HIS	2.1
1	I	208	LYS	2.1
1	J	212	THR	2.1
3	D	91	PRO	2.1
3	D	38	CYS	2.1
1	K	147	CYS	2.1
3	D	41	SER	2.1
2	M	33	ILE	2.0
3	C	42	GLY	2.0
2	O	173	LYS	2.0
3	A	118	VAL	2.0
1	I	164	GLY	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 carbohydrates 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.