



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:31 PM UTC

PDB ID : 3A2V / pdb_00003a2v
Title : Peroxiredoxin (C207S) from *Aeropyrum pernix* K1 complexed with hydrogen peroxide
Authors : Nakamura, T.; Kado, Y.; Yamaguchi, F.; Ishikawa, K.; Matsumura, H.; Inoue, T.
Deposited on : 2009-06-04
Resolution : 1.65 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
Mogul	:	2022.3.0, CSD as543be (2022)
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

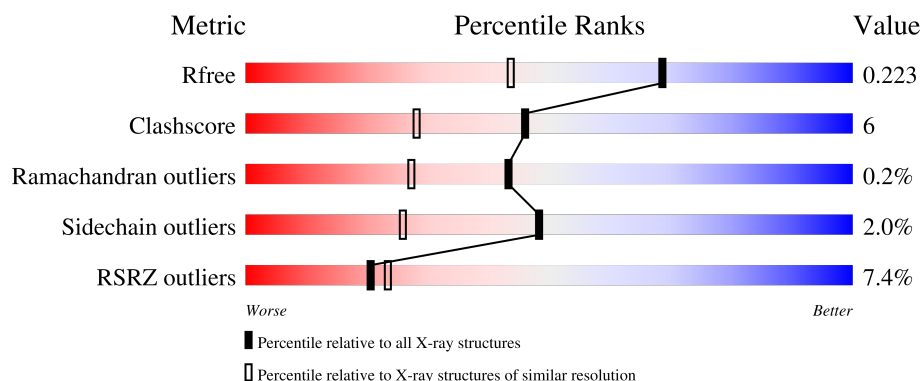
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.65 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.



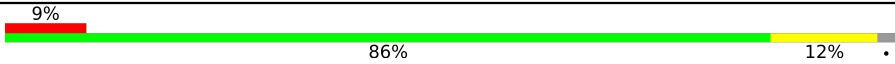
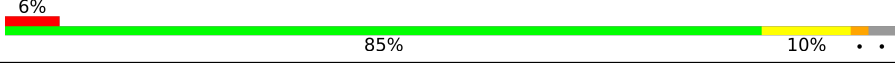
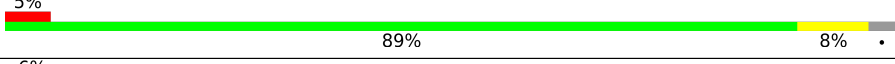
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	2563 (1.66-1.66)
Clashscore	190562	2662 (1.66-1.66)
Ramachandran outliers	187476	2621 (1.66-1.66)
Sidechain outliers	187428	2621 (1.66-1.66)
RSRZ outliers	180081	2564 (1.66-1.66)

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	249	7% 80% 17% ..
1	B	249	8% 87% 10% .
1	C	249	10% 79% 17% ..
1	D	249	5% 86% 10% ..
1	E	249	7% 86% 10% ..

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Mol	Chain	Length	Quality of chain
1	F	249	
1	G	249	
1	H	249	
1	I	249	
1	J	249	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 20999 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 Probable peroxiredoxin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	243	Total	C	N	O	S	0	15	0
			2050	1334	350	359	7			
1	B	242	Total	C	N	O	S	0	6	0
			1993	1287	348	352	6			
1	C	243	Total	C	N	O	S	0	11	0
			2038	1318	358	354	8			
1	D	240	Total	C	N	O	S	0	7	0
			1978	1276	349	346	7			
1	E	242	Total	C	N	O	S	0	5	0
			1990	1281	352	351	6			
1	F	244	Total	C	N	O	S	0	7	0
			2014	1302	348	357	7			
1	G	243	Total	C	N	O	S	0	5	0
			1998	1288	350	353	7			
1	H	242	Total	C	N	O	S	0	5	0
			1996	1292	349	349	6			
1	I	242	Total	C	N	O	S	0	5	0
			1993	1286	352	349	6			
1	J	243	Total	C	N	O	S	0	6	0
			2019	1300	353	358	8			

There are 10 discrepancies between the modelled and reference sequences:

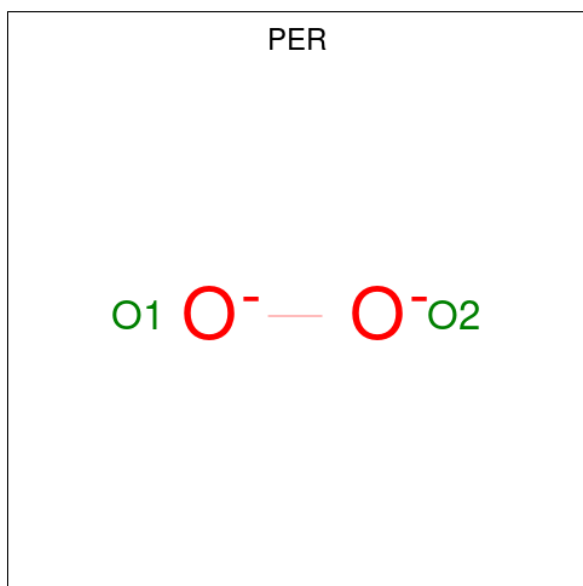
Chain	Residue	Modelled	Actual	Comment	Reference
A	207	SER	CYS	engineered mutation	UNP Q9Y9L0
B	207	SER	CYS	engineered mutation	UNP Q9Y9L0
C	207	SER	CYS	engineered mutation	UNP Q9Y9L0
D	207	SER	CYS	engineered mutation	UNP Q9Y9L0
E	207	SER	CYS	engineered mutation	UNP Q9Y9L0
F	207	SER	CYS	engineered mutation	UNP Q9Y9L0
G	207	SER	CYS	engineered mutation	UNP Q9Y9L0
H	207	SER	CYS	engineered mutation	UNP Q9Y9L0
I	207	SER	CYS	engineered mutation	UNP Q9Y9L0

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Chain	Residue	Modelled	Actual	Comment	Reference
J	207	SER	CYS	engineered mutation	UNP Q9Y9L0

- Molecule 2 is PEROXIDE ION (CCD ID: PER) (formula: O₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total O 2 2	0	0
2	B	1	Total O 2 2	0	0
2	C	1	Total O 2 2	0	0
2	D	1	Total O 2 2	0	0
2	E	1	Total O 2 2	0	0
2	F	1	Total O 2 2	0	0
2	G	1	Total O 2 2	0	0
2	H	1	Total O 2 2	0	0
2	I	1	Total O 2 2	0	0
2	J	1	Total O 2 2	0	0

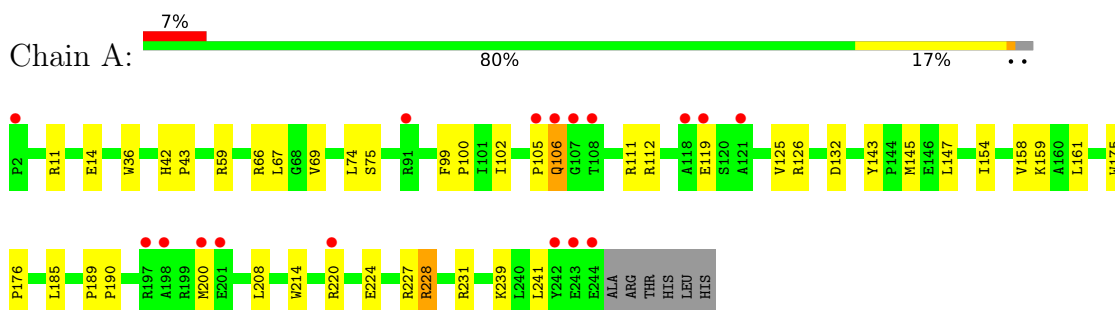
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	106	Total 106	O 106	0	0
3	B	79	Total 79	O 79	0	0
3	C	77	Total 77	O 77	0	0
3	D	88	Total 88	O 88	0	0
3	E	79	Total 79	O 79	0	0
3	F	76	Total 76	O 76	0	0
3	G	93	Total 93	O 93	0	0
3	H	85	Total 85	O 85	0	0
3	I	124	Total 124	O 124	0	0
3	J	103	Total 103	O 103	0	0

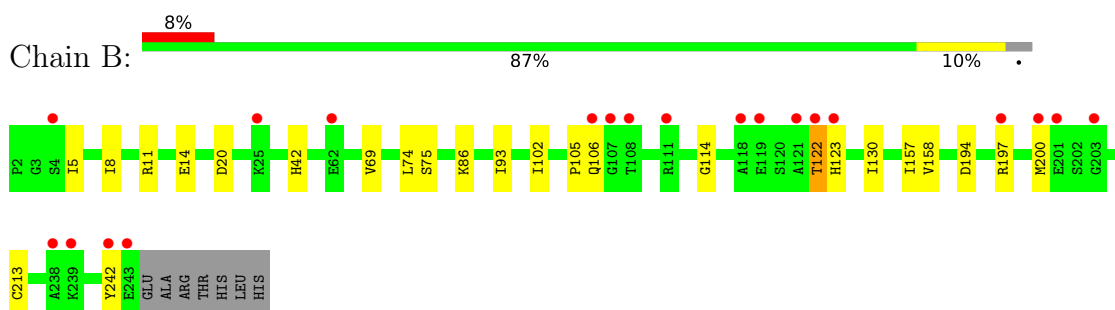
3 Residue-property plots [i](#)

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

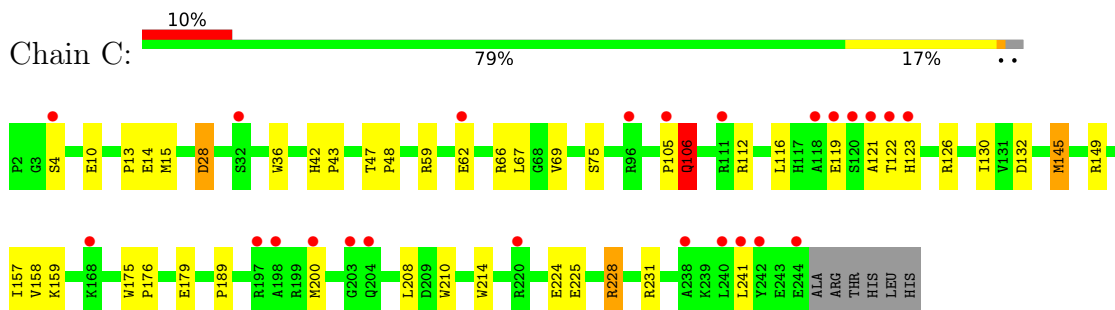
- Molecule 1: Probable peroxiredoxin



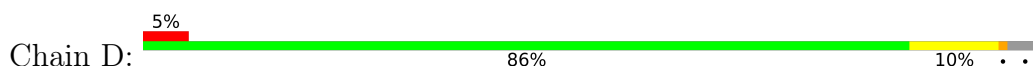
- Molecule 1: Probable peroxiredoxin

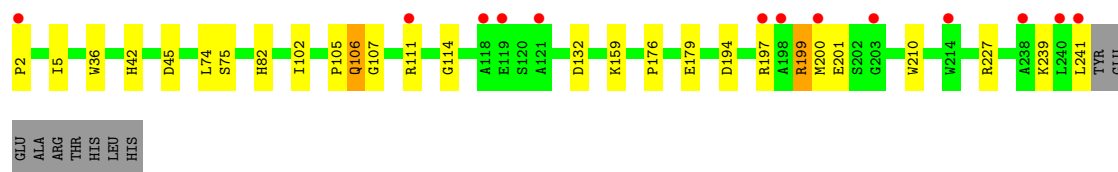


- Molecule 1: Probable peroxiredoxin

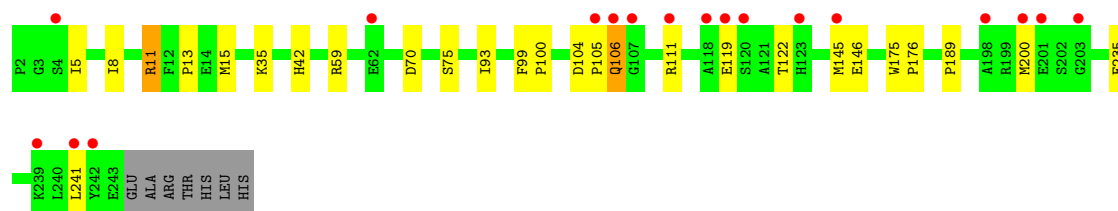
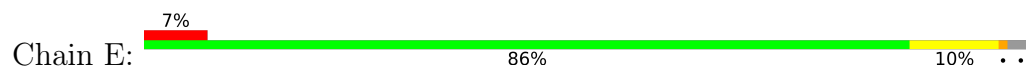


- Molecule 1: Probable peroxiredoxin

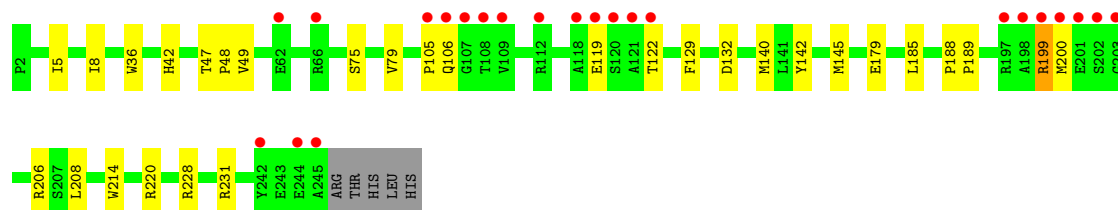
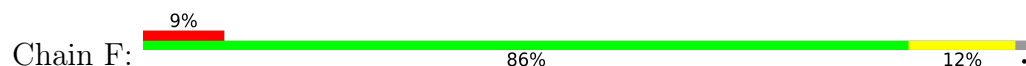




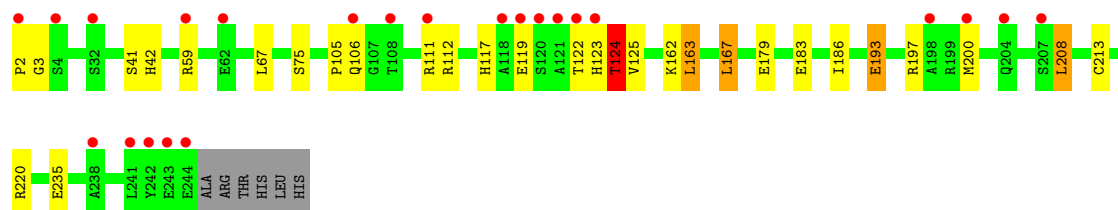
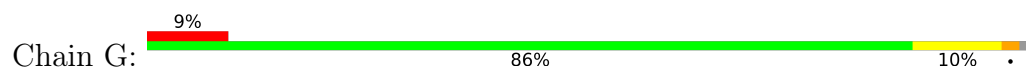
- Molecule 1: Probable peroxiredoxin



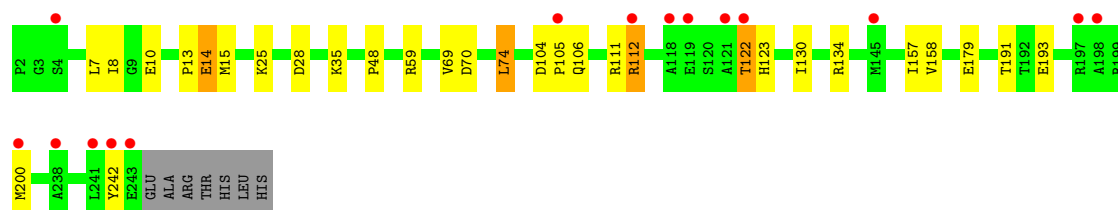
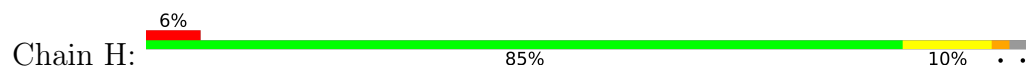
- Molecule 1: Probable peroxiredoxin



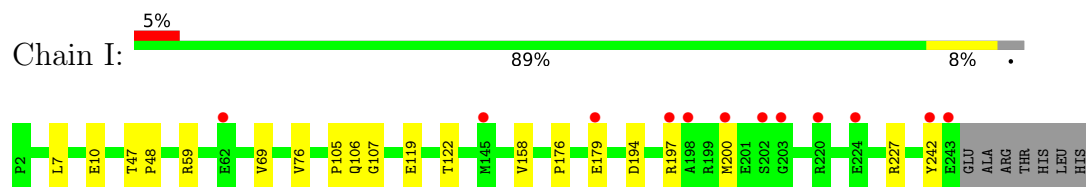
- Molecule 1: Probable peroxiredoxin



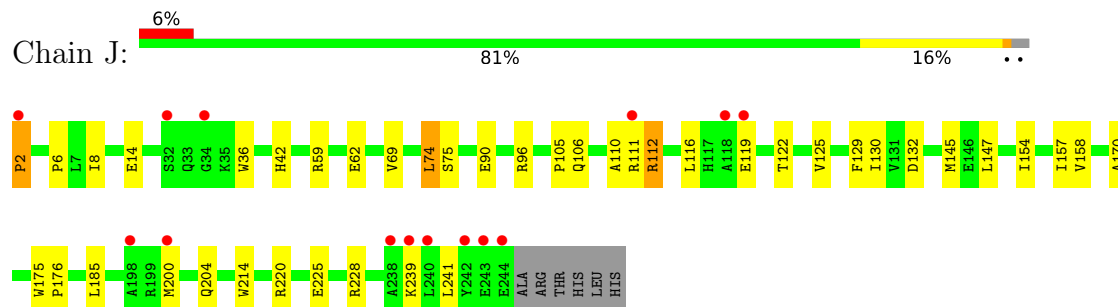
- Molecule 1: Probable peroxiredoxin



● Molecule 1: Probable peroxiredoxin



● Molecule 1: Probable peroxiredoxin



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	76.21Å 103.11Å 104.40Å 106.02° 104.91° 92.97°	Depositor
Resolution (Å)	39.25 – 1.65 39.25 – 1.65	Depositor EDS
% Data completeness (in resolution range)	94.0 (39.25-1.65) 93.9 (39.25-1.65)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.79 (at 1.65Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.194 , 0.223 0.193 , 0.223	Depositor DCC
R_{free} test set	16633 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	23.1	Xtriage
Anisotropy	0.064	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 37.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	20999	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.15% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: PER

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	0.77	0/2153	0.84	0/2927
1	B	0.74	0/2066	0.85	0/2808
1	C	0.71	1/2128 (0.0%)	0.86	0/2889
1	D	0.75	0/2053	0.88	0/2788
1	E	0.75	0/2060	0.87	1/2797 (0.0%)
1	F	0.76	1/2093 (0.0%)	0.86	1/2846 (0.0%)
1	G	0.76	0/2070	0.89	1/2813 (0.0%)
1	H	0.73	0/2069	0.89	2/2814 (0.1%)
1	I	0.81	0/2065	0.92	0/2807
1	J	0.85	0/2076	0.90	2/2822 (0.1%)
All	All	0.76	2/20833 (0.0%)	0.88	7/28311 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	189	PRO	CA-C	7.24	1.55	1.51
1	F	189	PRO	CA-C	5.39	1.54	1.51

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	189	PRO	O-C-N	5.92	123.93	121.15
1	G	124	THR	N-CA-CB	-5.37	102.44	110.17
1	J	170	ALA	CA-C-N	-5.36	118.67	123.33
1	J	170	ALA	C-N-CA	-5.36	118.67	123.33
1	H	104	ASP	CA-C-N	-5.26	114.50	119.76
1	H	104	ASP	C-N-CA	-5.26	114.50	119.76
1	F	49	VAL	N-CA-C	-5.13	105.54	110.72

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2050	0	2084	41	0
1	B	1993	0	2004	27	0
1	C	2038	0	2064	37	0
1	D	1978	0	2001	19	0
1	E	1990	0	1998	26	0
1	F	2014	0	2012	28	0
1	G	1998	0	1996	25	0
1	H	1996	0	2000	30	0
1	I	1993	0	1998	24	0
1	J	2019	0	2001	35	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
2	E	2	0	0	0	0
2	F	2	0	0	0	0
2	G	2	0	0	0	0
2	H	2	0	0	0	0
2	I	2	0	0	0	0
2	J	2	0	0	0	0
3	A	106	0	0	1	0
3	B	79	0	0	0	0
3	C	77	0	0	1	0
3	D	88	0	0	1	0
3	E	79	0	0	1	0
3	F	76	0	0	1	0
3	G	93	0	0	0	0
3	H	85	0	0	1	0
3	I	124	0	0	1	0
3	J	103	0	0	1	0
All	All	20999	0	20158	233	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 6.

All (233) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:242:TYR:CD1	1:J:214[B]:TRP:HH2	1.55	1.23
1:E:59:ARG:HD3	1:F:179[B]:GLU:OE1	1.41	1.20
1:E:59:ARG:CD	1:F:179[B]:GLU:OE1	1.91	1.16
1:E:59:ARG:HD2	1:F:179[B]:GLU:CD	1.71	1.14
1:E:59:ARG:HD2	1:F:179[B]:GLU:OE2	1.49	1.12
1:I:242:TYR:CD1	1:J:214[B]:TRP:CH2	2.40	1.09
1:F:129:PHE:CE2	1:F:140:MET:HE3	1.90	1.07
1:A:214[A]:TRP:HH2	1:B:242:TYR:CD1	1.73	1.05
1:I:242:TYR:CE1	1:J:214[B]:TRP:HH2	1.78	0.99
1:G:105:PRO:O	1:G:106:GLN:HB2	1.66	0.94
1:I:242:TYR:CE1	1:J:214[B]:TRP:CH2	2.54	0.94
1:A:214[A]:TRP:CH2	1:B:242:TYR:CD1	2.56	0.93
1:F:129:PHE:HE2	1:F:140:MET:HE3	1.33	0.92
1:F:105:PRO:O	1:F:106:GLN:HB2	1.69	0.92
1:F:140:MET:HE2	1:F:142:TYR:OH	1.69	0.91
1:C:59:ARG:HD2	1:D:179:GLU:OE2	1.74	0.88
1:B:11[A]:ARG:NH1	1:B:14[A]:GLU:OE2	2.06	0.87
1:B:69:VAL:HG21	1:B:158[A]:VAL:HG11	1.56	0.87
1:C:208[A]:LEU:CD1	1:C:214[A]:TRP:HZ3	1.88	0.85
1:B:194:ASP:OD2	1:B:197:ARG:NH2	2.10	0.85
1:E:105:PRO:O	1:E:106:GLN:HB2	1.73	0.85
1:E:241:LEU:HD11	1:F:179[B]:GLU:OE2	1.78	0.83
1:I:69:VAL:HG21	1:I:158[A]:VAL:HG11	1.61	0.83
1:C:69:VAL:HG21	1:C:158:VAL:HG11	1.59	0.81
1:J:69:VAL:HG21	1:J:158[A]:VAL:HG21	1.64	0.80
1:J:105:PRO:O	1:J:106:GLN:HB2	1.82	0.79
1:E:59:ARG:CD	1:F:179[B]:GLU:CD	2.45	0.77
1:D:45:ASP:OD2	1:D:82:HIS:HD2	1.68	0.77
1:A:125[A]:VAL:HG22	1:A:143:TYR:O	1.84	0.76
1:E:8:ILE:HG22	1:F:119:GLU:HG3	1.67	0.76
1:I:200:MET:HA	1:I:200:MET:HE2	1.67	0.76
1:C:200[A]:MET:HA	1:C:200[A]:MET:HE2	1.69	0.75
1:I:179:GLU:OE1	1:J:241:LEU:HD11	1.86	0.75
1:C:208[A]:LEU:CD1	1:C:214[A]:TRP:CZ3	2.70	0.74
1:G:106:GLN:O	1:G:111:ARG:NH2	2.20	0.74
1:H:106:GLN:O	1:H:111:ARG:NH2	2.20	0.73
1:B:105:PRO:O	1:B:106:GLN:HB2	1.88	0.72
1:H:134:ARG:NH1	3:H:1139:HOH:O	2.22	0.72
1:C:200[A]:MET:HE3	1:C:210:TRP:HA	1.72	0.72
1:D:105:PRO:O	1:D:106:GLN:HB2	1.88	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:179:GLU:OE1	1:D:241:LEU:HD11	1.90	0.71
1:F:200:MET:HE2	1:F:200:MET:HA	1.72	0.71
1:C:119:GLU:HB3	3:C:445:HOH:O	1.92	0.69
1:G:163:LEU:HD22	1:G:167:LEU:HD22	1.74	0.68
1:D:200:MET:HE3	1:D:210:TRP:HA	1.76	0.68
1:C:43:PRO:HG3	1:C:145[A]:MET:HG3	1.74	0.68
1:H:69:VAL:HG21	1:H:158[B]:VAL:HG11	1.76	0.68
1:A:105:PRO:O	1:A:106:GLN:HB2	1.93	0.68
1:A:214[A]:TRP:HH2	1:B:242:TYR:CE1	2.12	0.67
1:J:105:PRO:O	1:J:106:GLN:CB	2.43	0.67
1:H:122:THR:HG23	1:H:123:HIS:CD2	2.29	0.66
1:C:208[A]:LEU:HD12	1:C:214[A]:TRP:HZ3	1.59	0.66
1:H:35:LYS:HD3	1:H:70:ASP:OD2	1.96	0.66
1:E:11:ARG:HD3	3:E:392:HOH:O	1.95	0.66
1:E:200:MET:HE2	1:E:200:MET:HA	1.78	0.65
1:C:224:GLU:OE2	1:C:231:ARG:NH1	2.29	0.65
1:J:225:GLU:CD	1:J:228:ARG:HH21	2.05	0.65
1:H:122:THR:HG21	3:I:627:HOH:O	1.97	0.65
1:A:214[A]:TRP:CH2	1:B:242:TYR:CE1	2.85	0.64
1:A:224[B]:GLU:OE2	1:A:231:ARG:NH2	2.30	0.64
1:J:200[A]:MET:HA	1:J:200[A]:MET:HE2	1.79	0.64
1:F:122:THR:HB	1:G:105:PRO:HG2	1.78	0.63
1:H:122:THR:CG2	1:H:123:HIS:CD2	2.81	0.63
1:J:110:ALA:HB1	1:J:116:LEU:CD2	2.28	0.63
1:C:105:PRO:O	1:C:106[A]:GLN:HB2	1.99	0.63
1:A:43:PRO:HB3	1:A:145[B]:MET:HE1	1.82	0.62
1:F:140:MET:CE	1:F:142:TYR:OH	2.47	0.61
1:D:105:PRO:O	1:D:106:GLN:CB	2.48	0.61
1:A:105:PRO:HG2	1:J:122:THR:HB	1.84	0.60
1:I:176:PRO:HG2	1:I:227:ARG:HG2	1.85	0.59
1:H:105:PRO:O	1:H:106:GLN:HB2	2.03	0.59
1:I:105:PRO:O	1:I:106:GLN:HB2	2.02	0.58
1:A:106:GLN:O	1:A:111:ARG:NH2	2.36	0.58
1:A:69:VAL:HG21	1:A:158[A]:VAL:HG11	1.84	0.58
1:B:106:GLN:NE2	1:C:116:LEU:HD22	2.18	0.57
1:B:42:HIS:CE1	1:B:75:SER:HB3	2.38	0.57
1:E:35[B]:LYS:HD2	1:E:70:ASP:OD2	2.03	0.57
1:D:106:GLN:O	1:D:111:ARG:NH2	2.37	0.57
1:D:176:PRO:HG2	1:D:227:ARG:HG2	1.85	0.57
1:A:185:LEU:O	1:A:214[A]:TRP:HB2	2.05	0.57
1:A:106:GLN:HG2	1:J:122:THR:HA	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:10:GLU:OE2	1:D:2:PRO:HB2	2.05	0.56
1:B:74:LEU:HD23	1:B:102:ILE:HB	1.87	0.56
1:A:200:MET:HE2	1:A:200:MET:HA	1.88	0.56
1:E:119:GLU:OE2	1:E:145:MET:HG2	2.05	0.56
1:G:124:THR:HG23	1:G:125:VAL:O	2.07	0.55
1:I:105:PRO:O	1:I:106:GLN:CB	2.54	0.55
1:J:74:LEU:C	1:J:74:LEU:HD13	2.32	0.55
1:C:200[A]:MET:HA	1:C:200[A]:MET:CE	2.37	0.54
1:A:105:PRO:O	1:A:106:GLN:CB	2.55	0.54
1:B:122:THR:HG23	1:B:123:HIS:CD2	2.42	0.54
1:H:105:PRO:HG2	1:I:122:THR:HB	1.90	0.54
1:B:130:ILE:HD13	1:B:157:ILE:HG21	1.89	0.54
1:A:125[A]:VAL:HG21	1:B:8:ILE:HD12	1.89	0.53
1:G:200:MET:HA	1:G:200:MET:HE2	1.90	0.53
1:A:119:GLU:HG3	1:B:8:ILE:HG22	1.91	0.53
1:I:200:MET:HA	1:I:200:MET:CE	2.39	0.53
1:H:106:GLN:HE22	1:I:107:GLY:HA3	1.73	0.52
1:C:105:PRO:O	1:C:106[B]:GLN:CB	2.57	0.52
1:G:59:ARG:HG3	1:H:179:GLU:OE1	2.08	0.52
1:H:130:ILE:HD13	1:H:157:ILE:HG21	1.91	0.52
1:D:105:PRO:HG2	1:E:122:THR:HB	1.92	0.52
1:B:200:MET:HA	1:B:200:MET:HE2	1.91	0.52
1:J:90:GLU:OE2	1:J:96:ARG:HD3	2.09	0.52
1:A:11:ARG:NH1	1:A:14[A]:GLU:OE2	2.32	0.52
1:G:200:MET:HE1	1:G:213:CYS:SG	2.50	0.52
1:J:14:GLU:O	1:J:112:ARG:NH2	2.40	0.52
1:A:74[A]:LEU:HD23	1:A:102:ILE:HB	1.92	0.51
1:I:179:GLU:OE1	1:J:241:LEU:CD1	2.56	0.51
1:H:200:MET:HE2	1:H:200:MET:HA	1.92	0.51
1:I:194:ASP:OD1	1:I:197:ARG:NH1	2.43	0.51
1:F:106:GLN:HE22	1:G:111:ARG:HE	1.59	0.51
1:F:105:PRO:HG2	1:G:122:THR:HB	1.93	0.51
1:J:42:HIS:CE1	1:J:75:SER:HB3	2.46	0.51
1:C:208[A]:LEU:HD13	1:C:214[A]:TRP:CZ3	2.46	0.50
1:C:200[A]:MET:HG2	1:C:210:TRP:HB3	1.92	0.50
1:J:185:LEU:O	1:J:214[B]:TRP:HB2	2.11	0.50
1:A:36:TRP:HB2	1:A:69:VAL:HG22	1.94	0.50
1:B:74:LEU:HD23	1:B:102:ILE:CG2	2.42	0.50
1:C:241:LEU:HD11	1:D:179:GLU:OE1	2.12	0.50
1:H:74:LEU:C	1:H:74:LEU:HD13	2.36	0.50
1:A:119:GLU:OE2	1:A:145[B]:MET:HG2	2.11	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:220:ARG:O	1:A:224[A]:GLU:HG3	2.12	0.49
1:C:13:PRO:HB2	1:C:15:MET:HE3	1.94	0.49
1:E:119:GLU:HG3	1:F:8:ILE:HG22	1.94	0.49
1:C:59:ARG:CD	1:D:179:GLU:OE2	2.55	0.49
1:B:5:ILE:HG22	1:B:114:GLY:HA3	1.93	0.49
1:B:200:MET:HE1	1:B:213:CYS:SG	2.52	0.49
1:D:42:HIS:CE1	1:D:75:SER:HB3	2.48	0.49
1:E:175:TRP:CG	1:E:176:PRO:HA	2.48	0.49
1:I:242:TYR:CD1	1:J:214[B]:TRP:CZ2	2.98	0.49
1:B:69:VAL:CG2	1:B:158[A]:VAL:HG11	2.36	0.48
1:C:130:ILE:HD13	1:C:157:ILE:HG21	1.95	0.48
1:B:105:PRO:O	1:B:106:GLN:CB	2.57	0.48
1:E:93:ILE:HD11	1:F:208:LEU:HD13	1.96	0.48
1:H:105:PRO:O	1:H:106:GLN:CB	2.62	0.48
1:I:200:MET:HE2	1:I:200:MET:CA	2.41	0.48
1:D:82:HIS:HE1	1:D:102:ILE:O	1.97	0.48
1:D:199:ARG:HD3	3:D:472:HOH:O	2.14	0.47
1:E:13:PRO:HB2	1:E:15:MET:HE3	1.97	0.47
1:J:119:GLU:OE2	1:J:145[B]:MET:HG2	2.14	0.47
1:F:185:LEU:O	1:F:214[A]:TRP:HB2	2.14	0.47
1:J:175:TRP:CG	1:J:176:PRO:HA	2.50	0.47
1:B:106:GLN:HE22	1:C:116:LEU:HD22	1.78	0.47
1:I:179:GLU:OE2	1:J:59:ARG:HG3	2.14	0.47
1:J:200[A]:MET:HE2	1:J:200[A]:MET:CA	2.45	0.47
1:A:125[A]:VAL:HG22	1:A:126:ARG:H	1.79	0.47
1:J:106:GLN:O	1:J:111:ARG:NH2	2.48	0.47
1:C:42:HIS:CE1	1:C:75:SER:HB3	2.50	0.47
1:C:122:THR:HG23	1:C:123:HIS:CD2	2.51	0.47
1:I:119:GLU:HG3	1:J:8:ILE:HG22	1.96	0.46
1:G:208:LEU:CD1	1:H:242[B]:TYR:CD1	2.99	0.46
1:F:42:HIS:CE1	1:F:75:SER:HB3	2.51	0.46
1:J:6:PRO:HG2	1:J:129:PHE:HE2	1.81	0.46
1:E:5:ILE:HG12	1:F:5:ILE:HG12	1.97	0.46
1:G:119:GLU:HG3	1:H:8:ILE:HG22	1.97	0.46
1:G:179:GLU:OE2	1:H:59:ARG:HG3	2.16	0.46
1:G:193:GLU:O	1:G:197:ARG:HG2	2.15	0.46
1:H:14:GLU:O	1:H:112[B]:ARG:NH2	2.49	0.46
1:G:3:GLY:HA3	1:H:7:LEU:HD21	1.97	0.46
1:A:67:LEU:HD21	1:A:159:LYS:HD3	1.98	0.46
1:E:146:GLU:HG3	3:F:558:HOH:O	2.16	0.46
1:J:204:GLN:NE2	3:J:283:HOH:O	2.48	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:125[A]:VAL:CG2	1:A:143:TYR:O	2.60	0.45
1:G:208:LEU:CD1	1:H:242[B]:TYR:HD1	2.29	0.45
1:E:42:HIS:CE1	1:E:75:SER:HB3	2.52	0.45
1:G:186:ILE:HG21	1:H:48:PRO:HG2	1.99	0.45
1:C:200[B]:MET:HG3	1:C:210:TRP:HB3	1.98	0.45
1:A:99:PHE:HB2	1:A:100:PRO:HD2	1.97	0.44
1:E:99:PHE:HB2	1:E:100:PRO:HD2	1.99	0.44
1:F:47:THR:HB	1:F:48:PRO:HD2	1.97	0.44
1:A:176:PRO:HG2	1:A:227:ARG:HG2	2.00	0.44
1:C:175:TRP:CG	1:C:176:PRO:HA	2.53	0.44
1:D:36:TRP:CD2	1:D:132:ASP:HA	2.52	0.44
1:C:67:LEU:HD21	1:C:159:LYS:HD3	1.99	0.44
1:D:194:ASP:OD1	1:D:197:ARG:NH2	2.49	0.44
1:J:154:ILE:O	1:J:158[B]:VAL:HG23	2.17	0.44
1:F:188:PRO:O	1:F:199:ARG:NH2	2.47	0.43
1:C:36:TRP:CD2	1:C:132:ASP:HA	2.53	0.43
1:C:14:GLU:O	1:C:112:ARG:NH2	2.52	0.43
1:H:13:PRO:HB2	1:H:15:MET:HE3	2.00	0.43
1:A:175:TRP:CG	1:A:176:PRO:HA	2.54	0.43
1:A:214[A]:TRP:CZ2	1:B:242:TYR:CD1	3.06	0.43
1:F:79[B]:VAL:HG11	1:H:191:THR:O	2.19	0.43
1:G:122:THR:OG1	1:G:123:HIS:HD2	2.01	0.43
1:C:47:THR:HB	1:C:48:PRO:HD2	2.01	0.43
1:G:41:SER:HB2	1:G:124:THR:HG21	2.00	0.43
1:D:5:ILE:HG22	1:D:114:GLY:HA3	2.01	0.43
1:E:104:ASP:N	1:E:105:PRO:HD3	2.34	0.43
1:H:106:GLN:HG2	1:I:122:THR:HA	2.01	0.43
1:A:125[A]:VAL:HG22	1:A:126:ARG:N	2.34	0.43
1:J:36:TRP:CD2	1:J:132:ASP:HA	2.54	0.42
1:A:59[A]:ARG:HH21	1:A:241:LEU:HD21	1.85	0.42
1:A:224[B]:GLU:CD	1:A:231:ARG:HH22	2.27	0.42
1:G:208:LEU:HD11	1:H:242[B]:TYR:CD1	2.54	0.42
1:C:28:ASP:OD1	1:C:28:ASP:N	2.51	0.42
1:E:200:MET:HA	1:E:200:MET:CE	2.49	0.42
1:A:74[A]:LEU:HD23	1:A:102:ILE:CG2	2.50	0.42
1:A:154:ILE:O	1:A:158[A]:VAL:HG23	2.20	0.42
1:F:228:ARG:HG3	1:F:231:ARG:NH1	2.35	0.42
1:B:106:GLN:NE2	1:C:116:LEU:CD2	2.83	0.42
1:F:200:MET:HE2	1:F:200:MET:CA	2.45	0.42
1:F:119:GLU:OE2	1:F:145[B]:MET:HG2	2.20	0.41
1:A:42:HIS:CE1	1:A:75:SER:HB3	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:189:PRO:HA	1:A:190:PRO:HD3	1.97	0.41
1:A:228:ARG:HD3	3:A:387:HOH:O	2.20	0.41
1:C:126:ARG:HB3	1:C:149:ARG:CZ	2.50	0.41
1:G:2:PRO:HB2	1:H:10:GLU:OE2	2.20	0.41
1:B:106:GLN:NE2	1:C:121:ALA:O	2.53	0.41
1:C:62:GLU:HG2	1:C:66:ARG:NH1	2.35	0.41
1:A:228:ARG:CZ	1:A:228:ARG:HB3	2.51	0.41
1:D:107:GLY:HA3	1:E:106:GLN:NE2	2.36	0.41
1:G:67:LEU:O	1:G:162:LYS:HE3	2.20	0.41
1:I:59:ARG:HH11	1:I:59:ARG:HD3	1.74	0.41
1:E:175:TRP:CD1	1:E:176:PRO:HA	2.56	0.41
1:G:42:HIS:CE1	1:G:75:SER:HB3	2.56	0.41
1:G:183:GLU:OE2	1:G:220[B]:ARG:NH2	2.42	0.41
1:I:10:GLU:OE2	1:J:2:PRO:HB2	2.21	0.41
1:J:69:VAL:CG2	1:J:158[A]:VAL:HG21	2.42	0.41
1:B:20:ASP:OD2	1:B:86:LYS:NZ	2.50	0.41
1:J:130:ILE:HD13	1:J:157:ILE:HG21	2.02	0.41
1:A:208:LEU:HD13	1:B:93[A]:ILE:CD1	2.51	0.40
1:H:106:GLN:OE1	1:I:76:VAL:HG11	2.22	0.40
1:G:117:HIS:HA	1:H:7:LEU:HD13	2.02	0.40
1:J:110:ALA:CB	1:J:116:LEU:CD2	2.97	0.40
1:F:36:TRP:CD2	1:F:132:ASP:HA	2.57	0.40
1:C:225:GLU:O	1:C:228[B]:ARG:HB3	2.22	0.40
1:H:25:LYS:NZ	1:H:28:ASP:OD2	2.54	0.40
1:A:36:TRP:CD2	1:A:132:ASP:HA	2.56	0.40
1:I:47:THR:HB	1:I:48:PRO:HD2	2.04	0.40
1:J:6:PRO:HG2	1:J:129:PHE:CE2	2.57	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	256/249 (103%)	251 (98%)	4 (2%)	1 (0%)	30	15
1	B	246/249 (99%)	243 (99%)	3 (1%)	0	100	100
1	C	252/249 (101%)	244 (97%)	6 (2%)	2 (1%)	16	4
1	D	245/249 (98%)	239 (98%)	4 (2%)	2 (1%)	16	4
1	E	245/249 (98%)	238 (97%)	7 (3%)	0	100	100
1	F	249/249 (100%)	245 (98%)	4 (2%)	0	100	100
1	G	246/249 (99%)	241 (98%)	5 (2%)	0	100	100
1	H	245/249 (98%)	240 (98%)	5 (2%)	0	100	100
1	I	245/249 (98%)	241 (98%)	4 (2%)	0	100	100
1	J	247/249 (99%)	238 (96%)	7 (3%)	2 (1%)	16	4
All	All	2476/2490 (99%)	2420 (98%)	49 (2%)	7 (0%)	43	21

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	106[A]	GLN
1	C	106[B]	GLN
1	D	106	GLN
1	A	106	GLN
1	D	239	LYS
1	J	239	LYS
1	J	125	VAL

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	225/215 (105%)	217 (96%)	8 (4%)	31	9
1	B	215/215 (100%)	214 (100%)	1 (0%)	81	72
1	C	221/215 (103%)	213 (96%)	8 (4%)	31	9
1	D	214/215 (100%)	210 (98%)	4 (2%)	50	28
1	E	214/215 (100%)	211 (99%)	3 (1%)	59	39

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	F	217/215 (101%)	214 (99%)	3 (1%)	59	39
1	G	215/215 (100%)	208 (97%)	7 (3%)	33	11
1	H	214/215 (100%)	208 (97%)	6 (3%)	38	15
1	I	214/215 (100%)	213 (100%)	1 (0%)	81	72
1	J	216/215 (100%)	210 (97%)	6 (3%)	38	15
All	All	2165/2150 (101%)	2118 (98%)	47 (2%)	48	23

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	66	ARG
1	A	112	ARG
1	A	147[A]	LEU
1	A	147[B]	LEU
1	A	161[A]	LEU
1	A	161[B]	LEU
1	A	228	ARG
1	A	239	LYS
1	B	122	THR
1	C	4	SER
1	C	28	ASP
1	C	106[A]	GLN
1	C	106[B]	GLN
1	C	145[A]	MET
1	C	145[B]	MET
1	C	228[A]	ARG
1	C	228[B]	ARG
1	D	74	LEU
1	D	159	LYS
1	D	199	ARG
1	D	201	GLU
1	E	11	ARG
1	E	106	GLN
1	E	111	ARG
1	F	199	ARG
1	F	206	ARG
1	F	220	ARG
1	G	112	ARG
1	G	124	THR
1	G	163	LEU

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Mol	Chain	Res	Type
1	G	167	LEU
1	G	193	GLU
1	G	208	LEU
1	G	235	GLU
1	H	14	GLU
1	H	74	LEU
1	H	112[A]	ARG
1	H	112[B]	ARG
1	H	122	THR
1	H	193	GLU
1	I	7	LEU
1	J	2	PRO
1	J	62	GLU
1	J	74	LEU
1	J	112	ARG
1	J	147	LEU
1	J	220	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	33	GLN
1	A	65	GLN
1	A	106	GLN
1	A	123	HIS
1	A	204	GLN
1	B	106	GLN
1	B	123	HIS
1	B	204	GLN
1	C	123	HIS
1	C	204	GLN
1	D	65	GLN
1	D	82	HIS
1	D	106	GLN
1	D	204	GLN
1	E	123	HIS
1	F	106	GLN
1	F	123	HIS
1	F	204	GLN
1	G	33	GLN
1	G	106	GLN
1	G	123	HIS

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Mol	Chain	Res	Type
1	G	204	GLN
1	H	106	GLN
1	H	123	HIS
1	H	204	GLN
1	I	204	GLN
1	J	204	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

10 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	PER	F	251	-	1,1,1	0.41	0	-		
2	PER	A	1	-	1,1,1	0.33	0	-		
2	PER	H	251	-	1,1,1	0.41	0	-		
2	PER	G	251	-	1,1,1	0.45	0	-		
2	PER	D	251	-	1,1,1	0.33	0	-		
2	PER	B	251	-	1,1,1	0.43	0	-		
2	PER	C	251	-	1,1,1	0.38	0	-		
2	PER	I	251	-	1,1,1	0.34	0	-		

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	PER	J	251	-	1,1,1	0.39	0	-		
2	PER	E	251	-	1,1,1	0.35	0	-		

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	243/249 (97%)	0.27	17 (6%)	22	25	11, 21, 36, 48	15 (6%)
1	B	242/249 (97%)	0.47	20 (8%)	17	19	13, 24, 40, 48	6 (2%)
1	C	243/249 (97%)	0.74	24 (9%)	13	13	13, 26, 44, 62	11 (4%)
1	D	240/249 (96%)	0.45	13 (5%)	31	35	13, 23, 42, 50	7 (2%)
1	E	242/249 (97%)	0.51	18 (7%)	20	23	11, 24, 39, 44	5 (2%)
1	F	244/249 (97%)	0.51	23 (9%)	14	15	13, 24, 41, 54	7 (2%)
1	G	243/249 (97%)	0.47	23 (9%)	14	14	12, 24, 40, 59	5 (2%)
1	H	242/249 (97%)	0.41	15 (6%)	26	30	13, 24, 36, 47	5 (2%)
1	I	242/249 (97%)	0.17	12 (4%)	34	38	11, 20, 35, 45	5 (2%)
1	J	243/249 (97%)	0.32	14 (5%)	29	32	9, 21, 37, 56	6 (2%)
All	All	2424/2490 (97%)	0.43	179 (7%)	20	23	9, 23, 40, 62	72 (2%)

All (179) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	245	ALA	7.2
1	J	238	ALA	5.8
1	C	122	THR	5.3
1	D	238	ALA	5.1
1	E	242	TYR	4.8
1	C	198	ALA	4.7
1	G	242	TYR	4.7
1	F	121	ALA	4.5
1	J	242	TYR	4.5
1	D	200	MET	4.3
1	H	242[A]	TYR	4.3
1	I	242	TYR	4.2
1	B	243	GLU	4.1

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Mol	Chain	Res	Type	RSRZ
1	C	121	ALA	4.0
1	B	122	THR	4.0
1	G	118	ALA	4.0
1	F	200	MET	3.9
1	B	242	TYR	3.9
1	D	121	ALA	3.9
1	H	118	ALA	3.9
1	G	120	SER	3.8
1	A	244	GLU	3.7
1	H	145	MET	3.7
1	A	2	PRO	3.7
1	G	121	ALA	3.7
1	C	242	TYR	3.7
1	F	244	GLU	3.7
1	D	241	LEU	3.6
1	I	203	GLY	3.6
1	E	106	GLN	3.6
1	A	242[A]	TYR	3.6
1	C	4	SER	3.5
1	G	119	GLU	3.4
1	D	118	ALA	3.4
1	F	198	ALA	3.4
1	D	198	ALA	3.3
1	F	201	GLU	3.3
1	C	200[A]	MET	3.3
1	C	32	SER	3.3
1	E	119	GLU	3.3
1	F	118	ALA	3.3
1	C	203	GLY	3.3
1	I	198	ALA	3.3
1	A	118	ALA	3.2
1	H	243	GLU	3.2
1	C	204	GLN	3.2
1	J	200[A]	MET	3.2
1	F	120	SER	3.2
1	J	34	GLY	3.2
1	B	121	ALA	3.2
1	E	198	ALA	3.2
1	F	202	SER	3.1
1	F	108	THR	3.1
1	A	198	ALA	3.1
1	H	197	ARG	3.0

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Mol	Chain	Res	Type	RSRZ
1	H	198	ALA	3.0
1	J	118	ALA	3.0
1	F	242[A]	TYR	3.0
1	C	123	HIS	3.0
1	A	119	GLU	2.9
1	F	107	GLY	2.9
1	H	200	MET	2.9
1	G	244	GLU	2.9
1	B	197	ARG	2.8
1	E	203	GLY	2.8
1	J	2	PRO	2.8
1	C	220	ARG	2.8
1	E	62	GLU	2.8
1	J	119	GLU	2.8
1	B	200	MET	2.8
1	C	118	ALA	2.7
1	C	197	ARG	2.7
1	F	197	ARG	2.7
1	B	123	HIS	2.7
1	A	197	ARG	2.7
1	D	2	PRO	2.7
1	D	240	LEU	2.7
1	B	4	SER	2.7
1	H	122	THR	2.7
1	J	243	GLU	2.7
1	I	200	MET	2.7
1	B	106	GLN	2.6
1	E	120	SER	2.6
1	F	122	THR	2.6
1	E	145	MET	2.6
1	J	240	LEU	2.6
1	B	25	LYS	2.6
1	C	119	GLU	2.6
1	C	241	LEU	2.6
1	F	119	GLU	2.6
1	B	238	ALA	2.6
1	H	238	ALA	2.6
1	B	239	LYS	2.6
1	B	62	GLU	2.6
1	G	122	THR	2.5
1	C	105	PRO	2.5
1	E	4	SER	2.5

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Mol	Chain	Res	Type	RSRZ
1	G	62	GLU	2.5
1	G	106	GLN	2.5
1	G	241	LEU	2.5
1	D	203	GLY	2.5
1	C	244	GLU	2.5
1	H	4	SER	2.4
1	E	123	HIS	2.4
1	E	200	MET	2.4
1	I	202	SER	2.4
1	A	243	GLU	2.4
1	D	197	ARG	2.4
1	A	105	PRO	2.3
1	F	106	GLN	2.3
1	G	32	SER	2.3
1	B	201	GLU	2.3
1	F	105	PRO	2.3
1	H	121	ALA	2.3
1	J	244	GLU	2.3
1	A	106	GLN	2.3
1	J	198	ALA	2.3
1	G	204	GLN	2.2
1	H	105	PRO	2.2
1	B	108	THR	2.2
1	G	198	ALA	2.2
1	B	111	ARG	2.2
1	C	96[A]	ARG	2.2
1	E	111	ARG	2.2
1	A	107	GLY	2.2
1	G	243	GLU	2.2
1	C	120	SER	2.2
1	B	107	GLY	2.2
1	G	123	HIS	2.2
1	A	108	THR	2.2
1	C	168	LYS	2.2
1	E	118	ALA	2.2
1	A	201	GLU	2.2
1	I	224	GLU	2.2
1	H	241	LEU	2.2
1	A	91	ARG	2.2
1	E	105	PRO	2.2
1	G	200	MET	2.1
1	I	179	GLU	2.1

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Mol	Chain	Res	Type	RSRZ
1	I	243	GLU	2.1
1	C	111	ARG	2.1
1	F	66	ARG	2.1
1	J	111	ARG	2.1
1	C	240	LEU	2.1
1	E	239	LYS	2.1
1	G	108	THR	2.1
1	G	59	ARG	2.1
1	F	109	VAL	2.1
1	F	203	GLY	2.1
1	A	200	MET	2.1
1	D	214	TRP	2.1
1	I	62	GLU	2.1
1	J	32	SER	2.1
1	F	112	ARG	2.1
1	F	199	ARG	2.1
1	G	2	PRO	2.1
1	I	220	ARG	2.1
1	B	203	GLY	2.1
1	B	119	GLU	2.1
1	E	201	GLU	2.1
1	E	241	LEU	2.1
1	G	111	ARG	2.1
1	H	112[A]	ARG	2.1
1	E	107	GLY	2.0
1	A	121	ALA	2.0
1	B	118	ALA	2.0
1	C	62	GLU	2.0
1	H	119	GLU	2.0
1	A	220	ARG	2.0
1	D	111	ARG	2.0
1	G	4	SER	2.0
1	G	207	SER	2.0
1	I	197	ARG	2.0
1	D	119	GLU	2.0
1	F	62	GLU	2.0
1	I	145	MET	2.0
1	C	238	ALA	2.0
1	G	238	ALA	2.0
1	J	239	LYS	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	PER	J	251	2/2	0.88	0.13	22,22,22,29	0
2	PER	C	251	2/2	0.89	0.14	27,27,27,33	0
2	PER	D	251	2/2	0.91	0.11	30,30,30,35	0
2	PER	F	251	2/2	0.92	0.12	28,28,28,37	0
2	PER	B	251	2/2	0.92	0.12	28,28,28,35	0
2	PER	H	251	2/2	0.94	0.12	29,29,29,32	0
2	PER	A	1	2/2	0.94	0.10	25,25,25,31	0
2	PER	E	251	2/2	0.95	0.11	32,32,32,34	0
2	PER	G	251	2/2	0.95	0.10	29,29,29,33	0
2	PER	I	251	2/2	0.97	0.12	22,22,22,27	0

6.5 Other polymers [i](#)

There are no such residues in this entry.