Full wwPDB X-ray Structure Validation Report (i)

Oct 15, 2023 – 02:15 PM EDT

PDB ID	:	2E2M
Title	:	Crystal structure of archaeal peroxiredoxin, thioredoxin peroxidase from Aeropyrum pernix K1 (sulfinic acid form)
Authors	:	Nakamura, T.; Yamamoto, T.; Abe, M.; Matsumura, H.; Hagihara, Y.; Goto, T.; Yamaguchi, T.; Inoue, T.
Deposited on	:	2006-11-14
Resolution	:	2.60 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 (i) 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 (i)) were used in the production of this report:

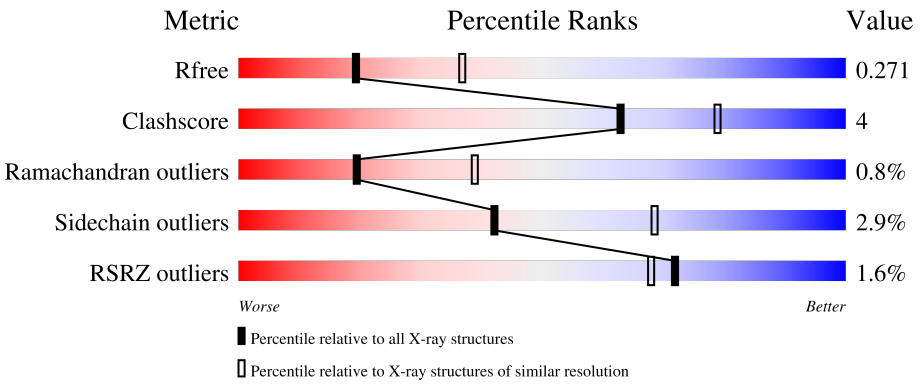
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 2.60 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}	130704	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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 <=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	250	80% 14% 6%
1	B	250	83% 11% • 5%
1	C	250	2% 81% 15% • •
1	D	250	% 80% 14% 7%
1	E	250	4% 78% 16% 5%

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Mol	Chain	Length	Quality of chain
1	F	250	2% 82% 13% 5%
1	G	250	2% 80% 14% 5%
1	H	250	% 79% 17%
1	I	250	2% 82% 12% 5%
1	J	250	% 79% 16% 5%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 19750 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
			Total	C	N	O	S			
1	A	236	1920	1236	337	341	6	0	0	0
1	B	237	1928	1241	338	343	6	0	0	0
1	C	242	1964	1261	345	352	6	0	0	0
1	D	233	1897	1222	334	335	6	0	0	0
1	E	238	1934	1244	339	345	6	0	0	0
1	F	238	1934	1244	339	345	6	0	0	0
1	G	238	1934	1244	339	345	6	0	0	0
1	H	241	1954	1255	342	351	6	0	0	0
1	I	237	1929	1241	338	344	6	0	0	0
1	J	238	1934	1244	339	345	6	0	0	0

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 water.

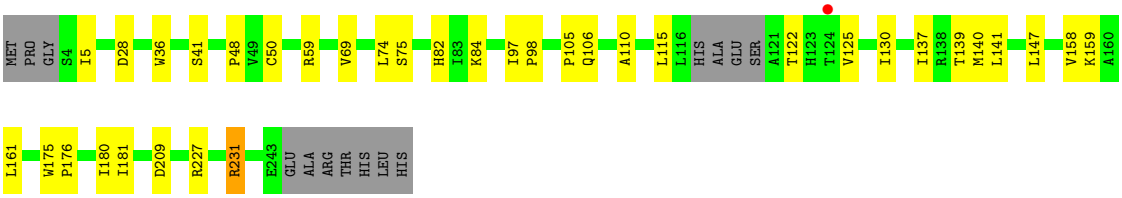
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	43	43	43	0	0
2	B	62	62	62	0	0
2	C	47	47	47	0	0
2	D	57	57	57	0	0
2	E	37	37	37	0	0
2	F	40	40	40	0	0
2	G	47	47	47	0	0
2	H	32	32	32	0	0
2	I	34	34	34	0	0
2	J	23	23	23	0	0

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.

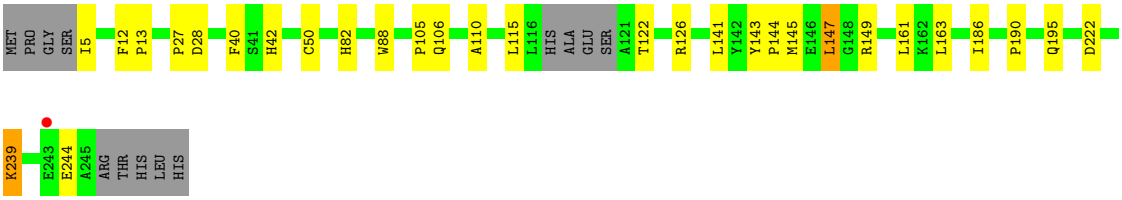
• Molecule 1: Probable peroxiredoxin

Chain A:



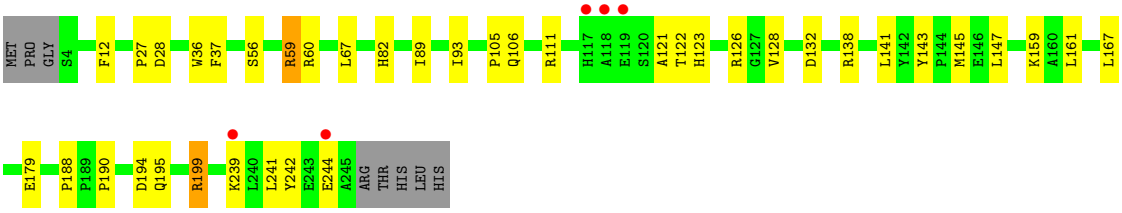
• Molecule 1: Probable peroxiredoxin

Chain B:



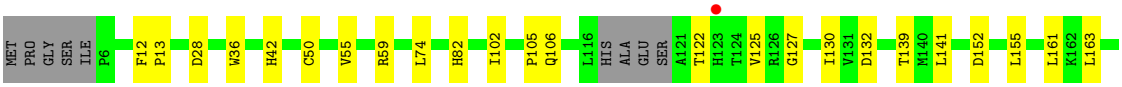
• Molecule 1: Probable peroxiredoxin

Chain C:



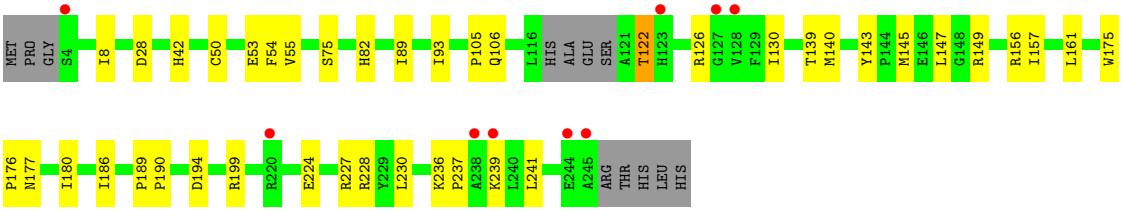
• Molecule 1: Probable peroxiredoxin

Chain D:



• Molecule 1: Probable peroxiredoxin

Chain E:

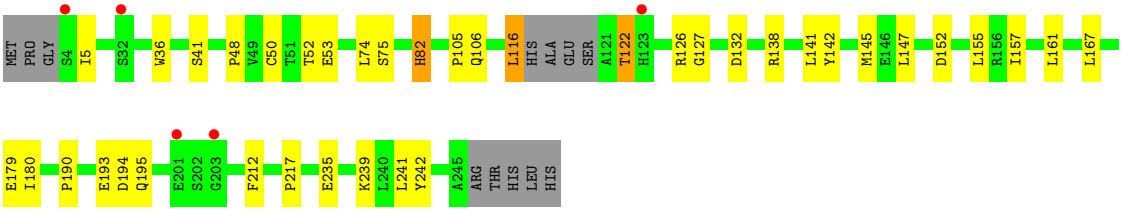


• Molecule 1: Probable peroxiredoxin

Chain F:

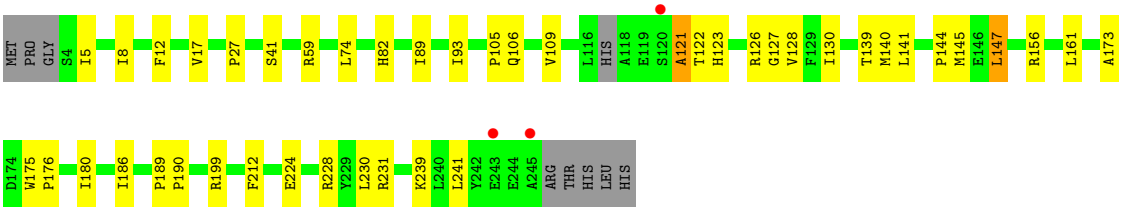
• Molecule 1: Probable peroxiredoxin

Chain G:



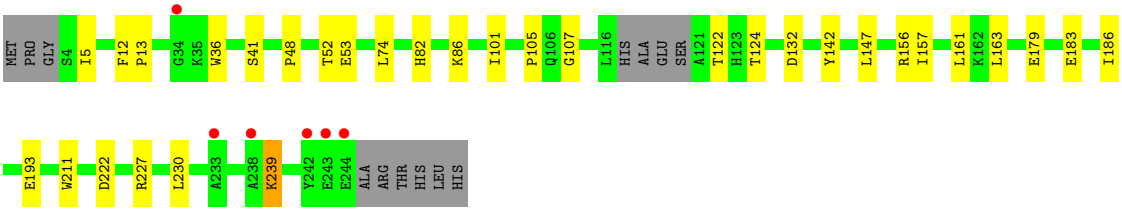
• Molecule 1: Probable peroxiredoxin

Chain H:



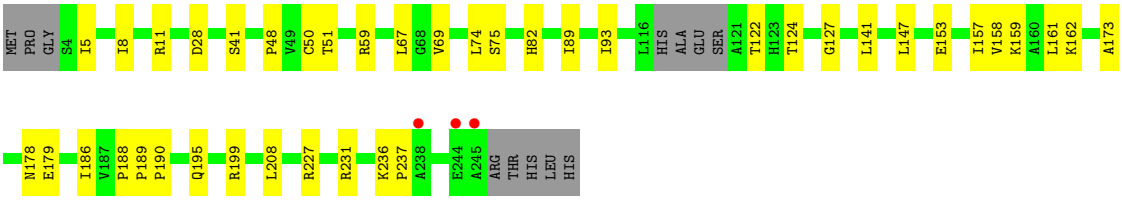
• Molecule 1: Probable peroxiredoxin

Chain I:



- Molecule 1: Probable peroxiredoxin

Chain J:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	76.24Å 103.13Å 104.68Å 105.88° 105.12° 92.63°	Depositor
Resolution (Å)	39.37 – 2.60 39.37 – 2.60	Depositor EDS
% Data completeness (in resolution range)	90.5 (39.37-2.60) 90.5 (39.37-2.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.71 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.231 , 0.280 0.225 , 0.271	Depositor DCC
R_{free} test set	4086 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å^2)	35.2	Xtriage
Anisotropy	0.221	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 44.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.91	EDS
Total number of atoms	19750	wwPDB-VP
Average B, all atoms (Å^2)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 12.79% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: CSD

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	0.32	0/1964	0.49	0/2669
1	B	0.32	0/1972	0.49	0/2680
1	C	0.32	0/2010	0.48	0/2733
1	D	0.32	0/1941	0.48	0/2637
1	E	0.32	0/1978	0.47	0/2688
1	F	0.32	0/1978	0.47	0/2688
1	G	0.32	0/1978	0.48	0/2688
1	H	0.32	0/1998	0.47	0/2715
1	I	0.32	0/1973	0.48	0/2681
1	J	0.32	0/1978	0.49	0/2688
All	All	0.32	0/19770	0.48	0/26867

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (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	1920	0	1913	20	0
1	B	1928	0	1919	18	0
1	C	1964	0	1948	18	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	1897	0	1892	15	0
1	E	1934	0	1924	26	0
1	F	1934	0	1924	20	0
1	G	1934	0	1924	23	0
1	H	1954	0	1940	25	0
1	I	1929	0	1919	18	0
1	J	1934	0	1924	22	0
2	A	43	0	0	0	0
2	B	62	0	0	0	0
2	C	47	0	0	0	0
2	D	57	0	0	0	0
2	E	37	0	0	0	0
2	F	40	0	0	0	0
2	G	47	0	0	0	0
2	H	32	0	0	0	0
2	I	34	0	0	0	0
2	J	23	0	0	0	0
All	All	19750	0	19227	167	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 4.

All (167) 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:G:142:TYR:HB3	1:H:8:ILE:HD11	1.48	0.95
1:A:5:ILE:HG12	1:B:5:ILE:HG12	1.61	0.83
1:E:147:LEU:HD12	1:F:161:LEU:HD13	1.72	0.72
1:E:177:ASN:HD21	1:E:227:ARG:HE	1.40	0.69
1:G:147:LEU:HD12	1:H:161:LEU:HD13	1.75	0.69
1:B:106:GLN:HE21	1:C:111:ARG:HH21	1.43	0.67
1:B:105:PRO:O	1:B:106:GLN:HB2	1.95	0.66
1:J:67:LEU:O	1:J:162:LYS:HE3	1.95	0.66
1:D:105:PRO:O	1:D:106:GLN:HB2	1.96	0.64
1:F:105:PRO:O	1:F:106:GLN:HB2	1.99	0.62
1:G:126:ARG:HD2	1:G:145:MET:HA	1.81	0.62
1:I:147:LEU:HD12	1:J:161:LEU:HD13	1.82	0.61
1:C:105:PRO:O	1:C:106:GLN:HB2	2.01	0.60
1:H:17:VAL:HG12	1:H:109:VAL:HG11	1.84	0.59
1:I:186:ILE:HG21	1:J:48:PRO:HB2	1.85	0.58
1:C:12:PHE:HB3	1:C:27:PRO:HG3	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:126:ARG:HD2	1:E:145:MET:HA	1.87	0.56
1:H:228:ARG:HG2	1:H:231:ARG:HH12	1.71	0.56
1:C:241:LEU:HB2	1:D:180:ILE:HA	1.87	0.56
1:D:74:LEU:HD23	1:D:102:ILE:HB	1.88	0.55
1:C:190:PRO:HB3	1:C:195:GLN:HB3	1.89	0.55
1:F:43:PRO:HD2	1:F:50:CSD:OD2	2.07	0.55
1:B:126:ARG:HH21	1:B:149:ARG:CZ	2.20	0.55
1:I:179:GLU:HG2	1:J:59:ARG:HD3	1.89	0.54
1:A:209:ASP:HB2	1:B:88:TRP:CZ2	2.42	0.54
1:G:179:GLU:HG2	1:H:59:ARG:HD3	1.88	0.54
1:A:137:ILE:HG21	1:A:140:MET:HE3	1.89	0.54
1:G:5:ILE:HG12	1:H:5:ILE:HG12	1.89	0.54
1:D:127:GLY:HA2	1:D:141:LEU:O	2.08	0.54
1:J:179:GLU:H	1:J:179:GLU:CD	2.11	0.54
1:I:5:ILE:HG12	1:J:5:ILE:HG12	1.89	0.54
1:E:105:PRO:O	1:E:106:GLN:HB2	2.08	0.53
1:A:41:SER:HA	1:A:74:LEU:HB3	1.90	0.53
1:B:110:ALA:HA	1:B:115:LEU:HD12	1.91	0.53
1:A:105:PRO:O	1:A:106:GLN:HB2	2.09	0.53
1:I:48:PRO:HB2	1:J:186:ILE:HG12	1.90	0.52
1:A:69:VAL:HG21	1:A:158:VAL:HG11	1.91	0.52
1:C:126:ARG:HB3	1:C:143:TYR:O	2.10	0.51
1:F:157:ILE:O	1:F:161:LEU:HB2	2.10	0.51
1:A:137:ILE:HG21	1:A:140:MET:CE	2.41	0.51
1:C:121:ALA:C	1:C:123:HIS:H	2.13	0.51
1:H:12:PHE:HB3	1:H:27:PRO:HG3	1.93	0.51
1:I:156:ARG:HD3	1:I:230:LEU:HD21	1.93	0.51
1:I:163:LEU:HD11	1:I:222:ASP:HB3	1.93	0.51
1:J:75:SER:HB3	1:J:82:HIS:CE1	2.46	0.51
1:B:190:PRO:HB3	1:B:195:GLN:HB3	1.92	0.51
1:J:190:PRO:HB3	1:J:195:GLN:HB3	1.93	0.51
1:H:127:GLY:HA2	1:H:141:LEU:O	2.11	0.50
1:F:7:LEU:HD23	1:F:8:ILE:H	1.77	0.50
1:C:128:VAL:HB	1:C:141:LEU:HB2	1.92	0.50
1:H:105:PRO:O	1:H:106:GLN:HB2	2.12	0.50
1:H:175:TRP:CG	1:H:176:PRO:HA	2.47	0.50
1:G:116:LEU:HD21	1:G:122:THR:H	1.77	0.49
1:E:186:ILE:HG12	1:F:48:PRO:HB2	1.93	0.49
1:H:121:ALA:C	1:H:123:HIS:H	2.16	0.49
1:G:167:LEU:HD23	1:G:217:PRO:HG2	1.94	0.49
1:E:180:ILE:HA	1:F:241:LEU:HB2	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:48:PRO:HB2	1:H:186:ILE:HD13	1.94	0.49
1:A:209:ASP:HB2	1:B:88:TRP:HZ2	1.78	0.49
1:B:239:LYS:HE2	1:B:244:GLU:HG2	1.94	0.49
1:E:8:ILE:HD11	1:F:142:TYR:HB3	1.94	0.49
1:E:177:ASN:HD21	1:E:227:ARG:NE	2.08	0.49
1:E:75:SER:HB3	1:E:82:HIS:CE1	2.48	0.49
1:G:41:SER:HA	1:G:74:LEU:HB3	1.95	0.48
1:E:89:ILE:HG23	1:E:93:ILE:HD12	1.95	0.48
1:H:156:ARG:HD3	1:H:230:LEU:HD21	1.96	0.48
1:D:152:ASP:HA	1:D:155:LEU:HD12	1.96	0.48
1:F:41:SER:HA	1:F:74:LEU:HB3	1.95	0.48
1:G:157:ILE:O	1:G:161:LEU:HB2	2.13	0.48
1:I:142:TYR:HB3	1:J:8:ILE:HD11	1.95	0.48
1:I:227:ARG:HH21	1:J:236:LYS:HD3	1.78	0.48
1:E:53:GLU:OE2	1:E:149:ARG:HG3	2.13	0.47
1:E:156:ARG:HD3	1:E:230:LEU:HD21	1.97	0.47
1:G:152:ASP:HA	1:G:155:LEU:HD12	1.96	0.47
1:A:36:TRP:HB2	1:A:69:VAL:HG22	1.96	0.47
1:J:127:GLY:HA2	1:J:141:LEU:O	2.14	0.47
1:D:163:LEU:HD11	1:D:222:ASP:HB3	1.96	0.46
1:G:52:THR:HG22	1:H:180:ILE:HD12	1.97	0.46
1:G:190:PRO:HB3	1:G:195:GLN:HB3	1.97	0.46
1:E:130:ILE:HD12	1:E:139:THR:HB	1.98	0.46
1:J:227:ARG:HD3	1:J:231:ARG:HH22	1.81	0.46
1:A:130:ILE:HD12	1:A:139:THR:HB	1.97	0.46
1:G:127:GLY:HA2	1:G:141:LEU:O	2.16	0.46
1:I:86:LYS:HD2	1:I:101:ILE:HD12	1.96	0.46
1:E:236:LYS:HD3	1:F:227:ARG:NH2	2.30	0.46
1:C:56:SER:O	1:C:60:ARG:HD3	2.15	0.46
1:I:41:SER:HA	1:I:74:LEU:HB3	1.98	0.46
1:J:157:ILE:O	1:J:161:LEU:HB2	2.16	0.46
1:B:163:LEU:HD11	1:B:222:ASP:HB3	1.97	0.46
1:G:53:GLU:HG3	1:H:173:ALA:HB2	1.98	0.45
1:I:53:GLU:HG2	1:J:173:ALA:HB2	1.98	0.45
1:D:36:TRP:CD2	1:D:132:ASP:HA	2.51	0.45
1:I:157:ILE:O	1:I:161:LEU:HB2	2.16	0.45
1:E:175:TRP:CG	1:E:176:PRO:HA	2.52	0.45
1:A:175:TRP:CG	1:A:176:PRO:HA	2.51	0.45
1:B:12:PHE:HB3	1:B:27:PRO:HG3	1.98	0.45
1:D:175:TRP:CG	1:D:176:PRO:HA	2.52	0.45
1:J:69:VAL:HG21	1:J:158:VAL:HG11	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:105:PRO:HB2	1:E:122:THR:HG21	1.99	0.45
1:G:105:PRO:O	1:G:106:GLN:HB2	2.17	0.45
1:G:138:ARG:HB3	1:G:161:LEU:HD21	1.99	0.45
1:H:144:PRO:HG2	1:H:147:LEU:HB2	1.99	0.45
1:A:48:PRO:HB2	1:B:186:ILE:HD13	1.99	0.44
1:E:75:SER:HB3	1:E:82:HIS:HE1	1.82	0.44
1:H:41:SER:HA	1:H:74:LEU:HB3	1.99	0.44
1:E:236:LYS:HD2	1:E:237:PRO:HD2	1.99	0.44
1:G:180:ILE:HA	1:H:241:LEU:HB2	1.99	0.44
1:C:36:TRP:CD2	1:C:132:ASP:HA	2.52	0.44
1:F:189:PRO:HA	1:F:190:PRO:HD3	1.90	0.44
1:C:89:ILE:HG23	1:C:93:ILE:HD12	2.00	0.44
1:D:130:ILE:HD12	1:D:139:THR:HB	1.99	0.44
1:J:236:LYS:HD2	1:J:237:PRO:HD2	2.00	0.44
1:F:115:LEU:HB3	1:F:125:VAL:HA	1.99	0.43
1:C:59:ARG:HH21	1:D:179:GLU:HG2	1.81	0.43
1:D:225:GLU:HA	1:D:228:ARG:HD3	2.00	0.43
1:F:128:VAL:O	1:F:140:MET:HA	2.18	0.43
1:F:188:PRO:HA	1:F:189:PRO:HD3	1.85	0.43
1:H:224:GLU:O	1:H:228:ARG:HG3	2.17	0.43
1:E:241:LEU:HB2	1:F:180:ILE:HA	2.00	0.43
1:A:97:ILE:HA	1:A:98:PRO:HD3	1.87	0.43
1:E:55:VAL:HG21	1:F:180:ILE:HD11	2.00	0.43
1:G:75:SER:HB3	1:G:82:HIS:CE1	2.54	0.43
1:G:241:LEU:HB2	1:H:180:ILE:HA	2.01	0.43
1:A:110:ALA:HA	1:A:115:LEU:HD12	2.01	0.43
1:H:89:ILE:HG23	1:H:93:ILE:HD12	2.00	0.43
1:A:75:SER:HB3	1:A:82:HIS:CE1	2.54	0.42
1:H:189:PRO:HA	1:H:190:PRO:HD3	1.94	0.42
1:J:89:ILE:HG23	1:J:93:ILE:HD12	2.01	0.42
1:E:189:PRO:HA	1:E:190:PRO:HD3	1.92	0.42
1:C:138:ARG:HB3	1:C:161:LEU:HD21	2.01	0.42
1:G:36:TRP:CD2	1:G:132:ASP:HA	2.55	0.42
1:I:211:TRP:HZ3	1:J:51:THR:HG1	1.66	0.42
1:H:130:ILE:HD12	1:H:139:THR:HB	2.02	0.42
1:I:52:THR:HB	1:J:178:ASN:HD21	1.85	0.42
1:J:188:PRO:HA	1:J:189:PRO:HD2	1.90	0.42
1:A:227:ARG:O	1:A:231:ARG:HB2	2.19	0.42
1:C:67:LEU:HD21	1:C:159:LYS:HD3	2.02	0.42
1:F:5:ILE:HG22	1:F:114:GLY:HA3	2.02	0.42
1:A:84:LYS:HA	1:A:84:LYS:HD2	1.90	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:42:HIS:NE2	1:E:54:PHE:HE1	2.18	0.42
1:E:126:ARG:HD3	1:E:143:TYR:O	2.20	0.42
1:I:12:PHE:HA	1:I:13:PRO:HD3	1.96	0.41
1:G:126:ARG:HD2	1:G:145:MET:CA	2.48	0.41
1:B:12:PHE:HA	1:B:13:PRO:HD3	1.94	0.41
1:E:180:ILE:HD12	1:F:52:THR:HG22	2.02	0.41
1:I:105:PRO:C	1:I:107:GLY:H	2.23	0.41
1:A:141:LEU:HD22	1:B:141:LEU:HD22	2.03	0.41
1:C:239:LYS:HE2	1:C:244:GLU:HG2	2.03	0.41
1:J:41:SER:HA	1:J:74:LEU:HB3	2.02	0.41
1:C:27:PRO:HB3	1:C:37:PHE:HE2	1.85	0.41
1:E:224:GLU:O	1:E:228:ARG:HG3	2.21	0.41
1:H:128:VAL:O	1:H:140:MET:HA	2.20	0.41
1:I:36:TRP:CD2	1:I:132:ASP:HA	2.55	0.41
1:B:144:PRO:HG2	1:B:147:LEU:HB2	2.02	0.41
1:A:139:THR:HG23	1:B:143:TYR:HA	2.01	0.41
1:C:188:PRO:O	1:C:199:ARG:NH2	2.53	0.41
1:A:180:ILE:HG22	1:A:181:ILE:HG23	2.03	0.41
1:E:157:ILE:O	1:E:161:LEU:HB2	2.21	0.41
1:F:160:ALA:HB1	1:F:171:VAL:HG11	2.03	0.41
1:H:126:ARG:HD2	1:H:145:MET:HA	2.03	0.41
1:C:126:ARG:HH11	1:C:143:TYR:HB2	1.86	0.41
1:D:42:HIS:N	1:D:74:LEU:O	2.44	0.41
1:F:122:THR:HG21	1:G:105:PRO:HB2	2.02	0.40
1:B:40:PHE:HD1	1:B:42:HIS:NE2	2.20	0.40
1:D:12:PHE:HA	1:D:13:PRO:HD3	1.96	0.40
1:D:55:VAL:O	1:D:59:ARG:HG3	2.21	0.40
1:B:126:ARG:HD2	1:B:145:MET:HA	2.03	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	231/250 (92%)	223 (96%)	6 (3%)	2 (1%)	17	35
1	B	232/250 (93%)	224 (97%)	6 (3%)	2 (1%)	17	35
1	C	239/250 (96%)	227 (95%)	11 (5%)	1 (0%)	34	57
1	D	228/250 (91%)	221 (97%)	5 (2%)	2 (1%)	17	35
1	E	233/250 (93%)	227 (97%)	4 (2%)	2 (1%)	17	35
1	F	233/250 (93%)	226 (97%)	5 (2%)	2 (1%)	17	35
1	G	233/250 (93%)	223 (96%)	8 (3%)	2 (1%)	17	35
1	H	236/250 (94%)	227 (96%)	6 (2%)	3 (1%)	12	24
1	I	232/250 (93%)	223 (96%)	7 (3%)	2 (1%)	17	35
1	J	233/250 (93%)	223 (96%)	9 (4%)	1 (0%)	34	57
All	All	2330/2500 (93%)	2244 (96%)	67 (3%)	19 (1%)	19	39

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	122	THR
1	G	239	LYS
1	H	122	THR
1	I	239	LYS
1	J	122	THR
1	E	122	THR
1	F	122	THR
1	G	122	THR
1	B	122	THR
1	C	122	THR
1	D	122	THR
1	E	239	LYS
1	F	239	LYS
1	H	239	LYS
1	I	122	THR
1	B	239	LYS
1	H	121	ALA
1	A	125	VAL
1	D	125	VAL

5.3.2 Protein sidechains (i)

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	204/215 (95%)	198 (97%)	6 (3%)	42	68
1	B	204/215 (95%)	200 (98%)	4 (2%)	55	78
1	C	208/215 (97%)	198 (95%)	10 (5%)	25	49
1	D	201/215 (94%)	195 (97%)	6 (3%)	41	67
1	E	205/215 (95%)	201 (98%)	4 (2%)	55	78
1	F	205/215 (95%)	199 (97%)	6 (3%)	42	68
1	G	205/215 (95%)	198 (97%)	7 (3%)	37	63
1	H	207/215 (96%)	203 (98%)	4 (2%)	57	79
1	I	205/215 (95%)	200 (98%)	5 (2%)	49	74
1	J	205/215 (95%)	197 (96%)	8 (4%)	32	58
All	All	2049/2150 (95%)	1989 (97%)	60 (3%)	42	68

All (60) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	ASP
1	A	59	ARG
1	A	147	LEU
1	A	159	LYS
1	A	161	LEU
1	A	231	ARG
1	B	28	ASP
1	B	82	HIS
1	B	147	LEU
1	B	161	LEU
1	C	28	ASP
1	C	59	ARG
1	C	82	HIS
1	C	145	MET
1	C	147	LEU
1	C	167	LEU
1	C	179	GLU
1	C	194	ASP
1	C	199	ARG
1	C	242	TYR

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Mol	Chain	Res	Type
1	D	28	ASP
1	D	82	HIS
1	D	161	LEU
1	D	167	LEU
1	D	212	PHE
1	D	242	TYR
1	E	28	ASP
1	E	140	MET
1	E	194	ASP
1	E	199	ARG
1	F	7	LEU
1	F	62	GLU
1	F	82	HIS
1	F	147	LEU
1	F	153	GLU
1	F	199	ARG
1	G	82	HIS
1	G	116	LEU
1	G	193	GLU
1	G	194	ASP
1	G	212	PHE
1	G	235	GLU
1	G	242	TYR
1	H	82	HIS
1	H	147	LEU
1	H	199	ARG
1	H	212	PHE
1	I	82	HIS
1	I	124	THR
1	I	183	GLU
1	I	193	GLU
1	I	239	LYS
1	J	11	ARG
1	J	28	ASP
1	J	124	THR
1	J	147	LEU
1	J	153	GLU
1	J	159	LYS
1	J	199	ARG
1	J	208	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	21	HIS
1	A	92	HIS
1	A	106	GLN
1	B	106	GLN
1	B	195	GLN
1	C	65	GLN
1	C	106	GLN
1	C	117	HIS
1	C	195	GLN
1	D	106	GLN
1	E	177	ASN
1	F	106	GLN
1	G	195	GLN
1	I	195	GLN
1	I	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)

10 non-standard protein/DNA/RNA residues 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	$\#\|Z\| > 2$
1	CSD	F	50	1	3,7,8	0.58	0	1,8,10	1.20	0
1	CSD	B	50	1	3,7,8	0.61	0	1,8,10	2.14	1 (100%)
1	CSD	H	50	1	3,7,8	0.72	0	1,8,10	1.50	0
1	CSD	J	50	1	3,7,8	0.59	0	1,8,10	2.64	1 (100%)
1	CSD	D	50	1	3,7,8	0.66	0	1,8,10	3.14	1 (100%)
1	CSD	E	50	1	3,7,8	0.67	0	1,8,10	2.12	1 (100%)
1	CSD	A	50	1	3,7,8	0.66	0	1,8,10	2.05	1 (100%)
1	CSD	G	50	1	3,7,8	0.71	0	1,8,10	3.57	1 (100%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
1	CSD	I	50	1	3,7,8	0.60	0	1,8,10	0.43	0
1	CSD	C	50	1	3,7,8	0.63	0	1,8,10	1.52	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	CSD	F	50	1	-	0/2/6/8	-
1	CSD	B	50	1	-	0/2/6/8	-
1	CSD	H	50	1	-	1/2/6/8	-
1	CSD	J	50	1	-	0/2/6/8	-
1	CSD	D	50	1	-	1/2/6/8	-
1	CSD	E	50	1	-	0/2/6/8	-
1	CSD	A	50	1	-	0/2/6/8	-
1	CSD	G	50	1	-	1/2/6/8	-
1	CSD	I	50	1	-	1/2/6/8	-
1	CSD	C	50	1	-	1/2/6/8	-

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	G	50	CSD	OD1-SG-CB	3.57	112.34	105.54
1	D	50	CSD	OD1-SG-CB	3.14	111.51	105.54
1	J	50	CSD	OD1-SG-CB	2.64	110.57	105.54
1	B	50	CSD	OD1-SG-CB	2.14	109.61	105.54
1	E	50	CSD	OD1-SG-CB	2.12	109.56	105.54
1	A	50	CSD	OD1-SG-CB	2.05	109.44	105.54

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	C	50	CSD	CA-CB-SG-OD1
1	D	50	CSD	CA-CB-SG-OD1
1	G	50	CSD	CA-CB-SG-OD1
1	H	50	CSD	CA-CB-SG-OD1
1	I	50	CSD	CA-CB-SG-OD1

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	F	50	CSD	1	0

5.5 Carbohydrates (i)

There are no monosaccharides 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 (i)

6.1 Protein, DNA and RNA chains (i)

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	235/250 (94%)	-0.24	1 (0%)	92	91	22, 29, 43, 48	0
1	B	236/250 (94%)	-0.24	1 (0%)	92	91	22, 30, 44, 53	0
1	C	241/250 (96%)	-0.13	5 (2%)	63	58	23, 32, 49, 53	0
1	D	232/250 (92%)	-0.15	2 (0%)	84	82	23, 32, 44, 46	0
1	E	237/250 (94%)	0.06	9 (3%)	40	33	29, 37, 54, 58	0
1	F	237/250 (94%)	-0.11	4 (1%)	70	66	26, 34, 52, 60	0
1	G	237/250 (94%)	-0.18	5 (2%)	63	58	25, 32, 46, 53	0
1	H	240/250 (96%)	-0.15	3 (1%)	77	73	25, 33, 49, 59	0
1	I	236/250 (94%)	-0.10	6 (2%)	57	51	23, 32, 49, 55	0
1	J	237/250 (94%)	-0.00	3 (1%)	77	73	24, 33, 47, 53	0
All	All	2368/2500 (94%)	-0.13	39 (1%)	72	68	22, 33, 48, 60	0

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	120	SER	4.8
1	F	238	ALA	4.0
1	E	238	ALA	3.8
1	C	118	ALA	3.4
1	G	123	HIS	3.2
1	D	123	HIS	3.1
1	I	243	GLU	3.1
1	I	242	TYR	3.0
1	E	239	LYS	2.9
1	C	119	GLU	2.8
1	H	245	ALA	2.8
1	G	201	GLU	2.8
1	F	4	SER	2.7

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Mol	Chain	Res	Type	RSRZ
1	F	245	ALA	2.6
1	J	238	ALA	2.6
1	G	4	SER	2.5
1	E	123	HIS	2.5
1	H	243	GLU	2.4
1	I	238	ALA	2.4
1	J	244	GLU	2.4
1	C	117	HIS	2.3
1	E	127	GLY	2.3
1	E	128	VAL	2.3
1	E	220	ARG	2.3
1	F	239	LYS	2.3
1	E	244	GLU	2.3
1	D	238	ALA	2.2
1	C	244	GLU	2.2
1	G	203	GLY	2.2
1	E	245	ALA	2.2
1	I	34	GLY	2.2
1	E	4	SER	2.1
1	I	244	GLU	2.1
1	B	243	GLU	2.1
1	C	239	LYS	2.0
1	J	245	ALA	2.0
1	A	124	THR	2.0
1	G	32	SER	2.0
1	I	233	ALA	2.0

6.2 Non-standard residues in protein, DNA, RNA chains (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(Å^2)	Q<0.9
1	CSD	D	50	8/9	0.94	0.13	28,28,30,31	0
1	CSD	C	50	8/9	0.95	0.14	30,30,31,31	0
1	CSD	I	50	8/9	0.95	0.14	32,32,33,34	0
1	CSD	G	50	8/9	0.96	0.11	28,28,30,30	0
1	CSD	H	50	8/9	0.96	0.12	28,29,30,30	0
1	CSD	A	50	8/9	0.96	0.15	25,25,26,27	0
1	CSD	J	50	8/9	0.96	0.15	26,26,27,27	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
1	CSD	E	50	8/9	0.97	0.13	29,29,30,31	0
1	CSD	F	50	8/9	0.97	0.17	27,27,28,28	0
1	CSD	B	50	8/9	0.97	0.12	24,24,24,24	0

6.3 Carbohydrates (i)

There are no monosaccharides 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.