



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 10:32 PM GMT

PDB ID : 1U10
Title : MEPA, active form with ZN in P1
Authors : Marcyjaniak, M.; Odintsov, S.G.; Sabala, I.; Bochtler, M.
Deposited on : 2004-07-14
Resolution : 2.40 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

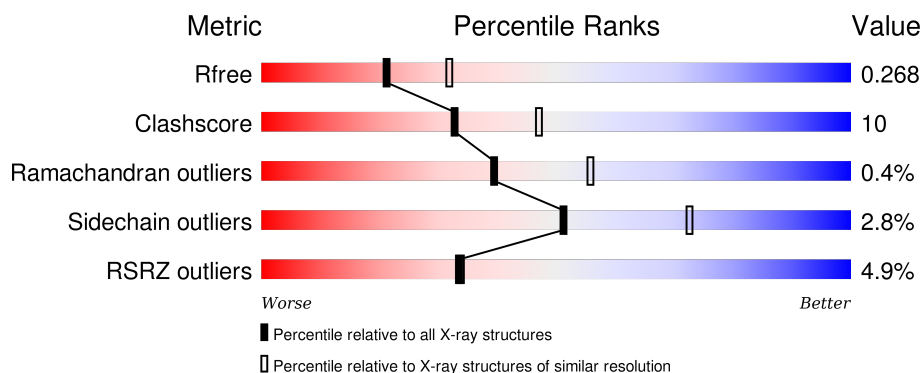
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	2919 (2.40-2.40)
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (2.40-2.40)

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	A	255	
1	B	255	
1	C	255	
1	D	255	
1	E	255	

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Mol	Chain	Length	Quality of chain
1	F	255	<div><div></div><div>4%</div><div>78%</div><div>16%</div><div>• 6%</div></div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11311 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 Penicillin-insensitive murein endopeptidase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	232	Total	C	N	O	S	Se	0	0	0
			1813	1142	329	330	6	6			
1	B	239	Total	C	N	O	S	Se	0	0	0
			1871	1179	339	341	6	6			
1	C	241	Total	C	N	O	S	Se	0	0	0
			1885	1189	341	343	6	6			
1	D	241	Total	C	N	O	S	Se	0	0	0
			1885	1189	341	343	6	6			
1	E	241	Total	C	N	O	S	Se	0	0	0
			1885	1189	341	343	6	6			
1	F	240	Total	C	N	O	S	Se	0	0	0
			1878	1184	340	342	6	6			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	61	MSE	MET	MODIFIED RESIDUE	UNP P14007
A	76	MSE	MET	MODIFIED RESIDUE	UNP P14007
A	90	MSE	MET	MODIFIED RESIDUE	UNP P14007
A	98	MSE	MET	MODIFIED RESIDUE	UNP P14007
A	100	MSE	MET	MODIFIED RESIDUE	UNP P14007
A	210	MSE	MET	MODIFIED RESIDUE	UNP P14007
B	61	MSE	MET	MODIFIED RESIDUE	UNP P14007
B	76	MSE	MET	MODIFIED RESIDUE	UNP P14007
B	90	MSE	MET	MODIFIED RESIDUE	UNP P14007
B	98	MSE	MET	MODIFIED RESIDUE	UNP P14007
B	100	MSE	MET	MODIFIED RESIDUE	UNP P14007
B	210	MSE	MET	MODIFIED RESIDUE	UNP P14007
C	61	MSE	MET	MODIFIED RESIDUE	UNP P14007
C	76	MSE	MET	MODIFIED RESIDUE	UNP P14007
C	90	MSE	MET	MODIFIED RESIDUE	UNP P14007
C	98	MSE	MET	MODIFIED RESIDUE	UNP P14007
C	100	MSE	MET	MODIFIED RESIDUE	UNP P14007

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Chain	Residue	Modelled	Actual	Comment	Reference
C	210	MSE	MET	MODIFIED RESIDUE	UNP P14007
D	61	MSE	MET	MODIFIED RESIDUE	UNP P14007
D	76	MSE	MET	MODIFIED RESIDUE	UNP P14007
D	90	MSE	MET	MODIFIED RESIDUE	UNP P14007
D	98	MSE	MET	MODIFIED RESIDUE	UNP P14007
D	100	MSE	MET	MODIFIED RESIDUE	UNP P14007
D	210	MSE	MET	MODIFIED RESIDUE	UNP P14007
E	61	MSE	MET	MODIFIED RESIDUE	UNP P14007
E	76	MSE	MET	MODIFIED RESIDUE	UNP P14007
E	90	MSE	MET	MODIFIED RESIDUE	UNP P14007
E	98	MSE	MET	MODIFIED RESIDUE	UNP P14007
E	100	MSE	MET	MODIFIED RESIDUE	UNP P14007
E	210	MSE	MET	MODIFIED RESIDUE	UNP P14007
F	61	MSE	MET	MODIFIED RESIDUE	UNP P14007
F	76	MSE	MET	MODIFIED RESIDUE	UNP P14007
F	90	MSE	MET	MODIFIED RESIDUE	UNP P14007
F	98	MSE	MET	MODIFIED RESIDUE	UNP P14007
F	100	MSE	MET	MODIFIED RESIDUE	UNP P14007
F	210	MSE	MET	MODIFIED RESIDUE	UNP P14007

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	D	2	Total Zn 2 2	0	0
2	E	2	Total Zn 2 2	0	0
2	B	1	Total Zn 1 1	0	0
2	C	1	Total Zn 1 1	0	0
2	A	2	Total Zn 2 2	0	0
2	F	1	Total Zn 1 1	0	0

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	E	1	Total	O	S	0	0
			5	4	1		
3	F	1	Total	O	S	0	0
			5	4	1		

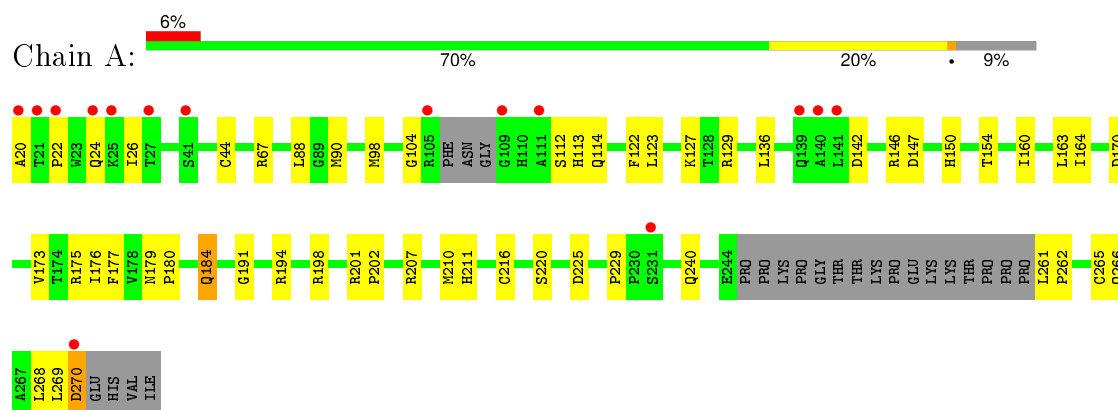
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	6	Total	O	0	0
			6	6		
4	B	12	Total	O	0	0
			12	12		
4	C	9	Total	O	0	0
			9	9		
4	D	8	Total	O	0	0
			8	8		
4	E	12	Total	O	0	0
			12	12		
4	F	8	Total	O	0	0
			8	8		

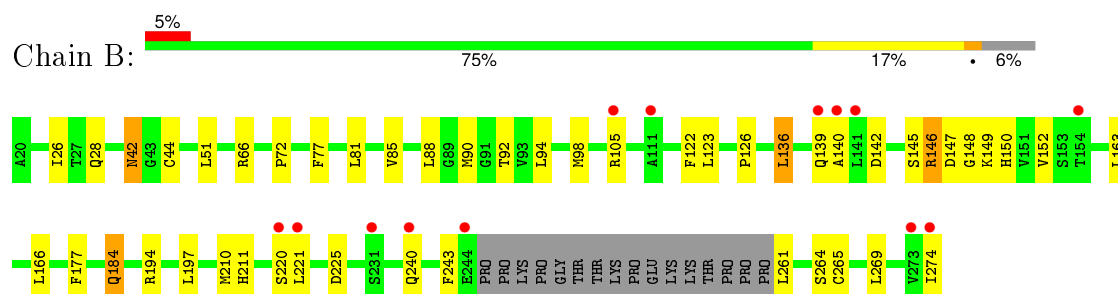
3 Residue-property plots [i](#)

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.

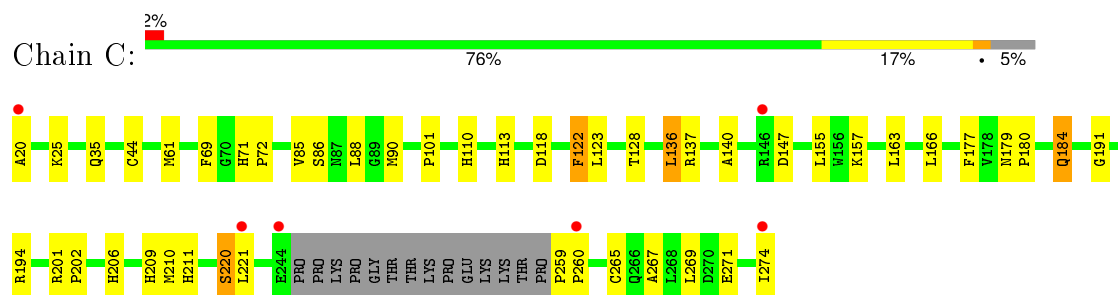
- Molecule 1: Penicillin-insensitive murein endopeptidase



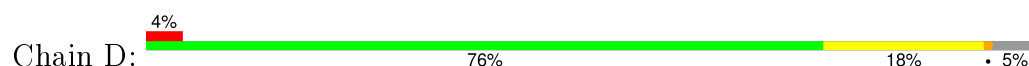
- Molecule 1: Penicillin-insensitive murein endopeptidase

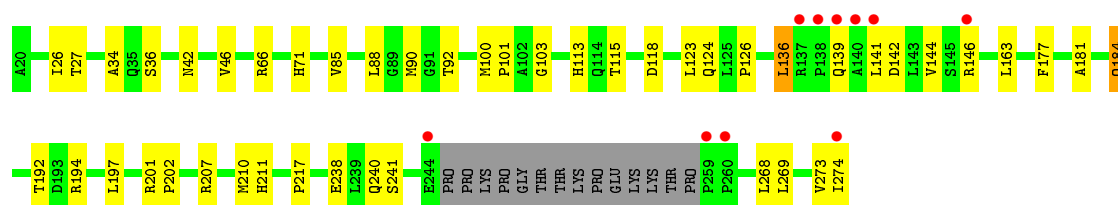


- Molecule 1: Penicillin-insensitive murein endopeptidase

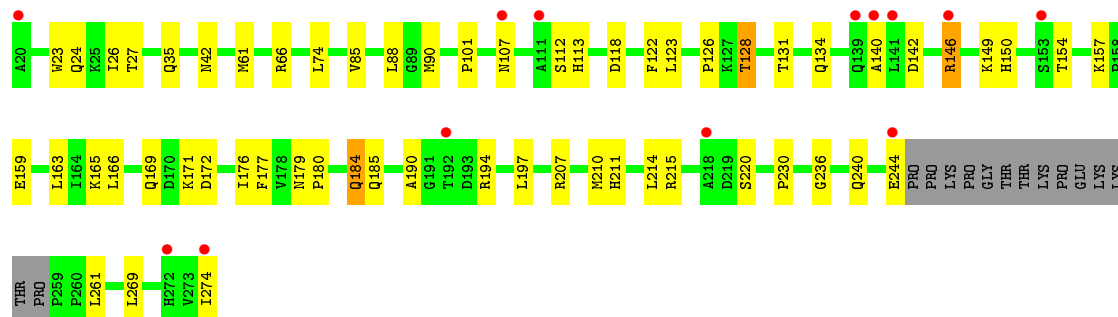


- Molecule 1: Penicillin-insensitive murein endopeptidase

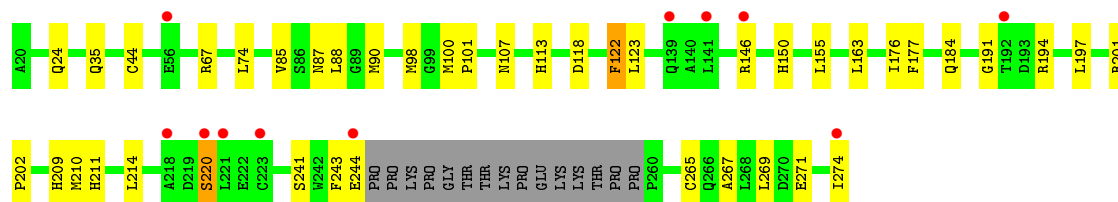
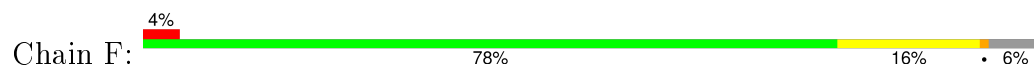




● Molecule 1: Penicillin-insensitive murein endopeptidase



● Molecule 1: Penicillin-insensitive murein endopeptidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	35.61Å 77.99Å 127.66Å 93.15° 95.93° 90.75°	Depositor
Resolution (Å)	20.00 – 2.40 19.95 – 2.40	Depositor EDS
% Data completeness (in resolution range)	98.2 (20.00-2.40) 96.6 (19.95-2.40)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	0.04	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.35 (at 2.41Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.234 , 0.266 0.235 , 0.268	Depositor DCC
R_{free} test set	2658 reflections (5.35%)	DCC
Wilson B-factor (Å ²)	25.9	Xtriage
Anisotropy	0.359	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 32.1	EDS
Estimated twinning fraction	0.036 for h,-k,-h-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 52369 reflections	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	11311	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 36.19 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 5.1911e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical 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

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, SO4

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	0.46	0/1851	0.64	0/2502
1	B	0.51	0/1912	0.66	0/2585
1	C	0.51	0/1928	0.66	0/2608
1	D	0.48	0/1928	0.68	0/2608
1	E	0.49	0/1928	0.68	0/2608
1	F	0.50	0/1920	0.65	0/2596
All	All	0.49	0/11467	0.66	0/15507

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

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	1813	0	1788	40	0
1	B	1871	0	1840	40	0
1	C	1885	0	1855	39	0
1	D	1885	0	1855	38	0
1	E	1885	0	1855	51	0
1	F	1878	0	1848	28	0
2	A	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	2	0	0	0	0
2	E	2	0	0	0	0
2	F	1	0	0	0	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0
3	C	5	0	0	0	0
3	D	5	0	0	0	0
3	E	5	0	0	0	0
3	F	5	0	0	0	0
4	A	6	0	0	1	0
4	B	12	0	0	3	0
4	C	9	0	0	1	0
4	D	8	0	0	0	0
4	E	12	0	0	1	0
4	F	8	0	0	0	0
All	All	11311	0	11041	227	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 10.

All (227) 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:B:146:ARG:N	1:B:146:ARG:HD3	1.58	1.09
1:B:146:ARG:H	1:B:146:ARG:CD	1.64	1.08
1:E:146:ARG:NH2	1:F:146:ARG:HH21	1.64	0.95
1:C:269:LEU:HD23	1:C:274:ILE:HD12	1.46	0.94
1:E:146:ARG:HH21	1:F:146:ARG:HH21	1.16	0.94
1:E:184:GLN:HE22	1:E:240:GLN:HE22	1.11	0.93
1:B:146:ARG:H	1:B:146:ARG:HD3	0.79	0.92
1:B:221:LEU:HB2	4:B:507:HOH:O	1.69	0.92
1:A:22:PRO:HB2	1:A:269:LEU:HD22	1.52	0.92
1:F:44:CYS:HG	1:F:265:CYS:HG	1.15	0.85
1:C:44:CYS:HG	1:C:265:CYS:HG	0.94	0.85
1:A:44:CYS:HG	1:A:265:CYS:HG	0.87	0.85
1:F:163:LEU:HG	1:F:210:MSE:HE1	1.61	0.82
1:B:66:ARG:HD2	4:B:509:HOH:O	1.80	0.80
1:E:24:GLN:NE2	1:E:107:ASN:HB2	1.99	0.77
1:A:177:PHE:HB2	1:A:211:HIS:HB3	1.68	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:123:LEU:HD21	1:B:210:MSE:HB3	1.67	0.75
1:D:123:LEU:HD21	1:D:210:MSE:HB3	1.67	0.75
1:B:44:CYS:HG	1:B:265:CYS:HG	0.90	0.74
1:D:88:LEU:HD12	1:D:90:MSE:HE2	1.69	0.74
1:A:127:LYS:HD2	1:C:221:LEU:HD13	1.68	0.73
1:F:123:LEU:HD21	1:F:210:MSE:HB3	1.69	0.71
1:F:194:ARG:HB2	1:F:197:LEU:HD12	1.72	0.71
1:B:88:LEU:HD12	1:B:90:MSE:HE2	1.73	0.70
1:B:163:LEU:HG	1:B:210:MSE:HE1	1.74	0.70
1:C:191:GLY:O	1:C:194:ARG:HD2	1.90	0.70
1:B:177:PHE:HB2	1:B:211:HIS:HB3	1.73	0.70
1:C:85:VAL:HA	1:C:90:MSE:HE3	1.72	0.70
1:C:35:GLN:OE1	1:C:220:SER:HA	1.91	0.70
1:E:85:VAL:HA	1:E:90:MSE:HE3	1.72	0.70
1:E:146:ARG:NH2	1:F:146:ARG:NH2	2.37	0.69
1:C:20:ALA:O	1:C:25:LYS:HE3	1.92	0.69
1:F:269:LEU:HD23	1:F:274:ILE:HD12	1.74	0.68
1:F:244:GLU:HA	1:F:244:GLU:OE2	1.93	0.68
1:A:191:GLY:O	1:A:194:ARG:HD2	1.93	0.68
1:A:104:GLY:C	1:A:114:GLN:HG2	2.14	0.67
1:E:150:HIS:CD2	1:F:150:HIS:CD2	2.83	0.67
1:C:269:LEU:CD2	1:C:274:ILE:HD12	2.25	0.66
1:E:190:ALA:HB3	1:E:194:ARG:NH1	2.12	0.65
1:E:90:MSE:HE1	1:E:166:LEU:CD1	2.27	0.65
1:B:146:ARG:NH1	1:B:152:VAL:HG22	2.12	0.64
1:C:137:ARG:HH11	1:C:137:ARG:HG3	1.61	0.64
1:A:146:ARG:HG3	1:A:146:ARG:HH11	1.63	0.63
1:E:194:ARG:HB2	1:E:197:LEU:HD12	1.80	0.62
1:E:244:GLU:HA	1:E:244:GLU:OE2	1.97	0.62
1:E:269:LEU:CD2	1:E:274:ILE:HD12	2.29	0.62
1:C:110:HIS:HE1	1:C:209:HIS:NE2	1.97	0.62
1:B:146:ARG:HH11	1:B:152:VAL:HG22	1.62	0.62
1:D:26:ILE:HB	1:D:274:ILE:OXT	2.00	0.61
1:E:149:LYS:HB2	1:E:150:HIS:CD2	2.36	0.61
1:A:170:ASP:HB3	1:A:173:VAL:HG23	1.82	0.61
1:D:163:LEU:HG	1:D:210:MSE:HE1	1.82	0.60
1:A:22:PRO:HB2	1:A:269:LEU:CD2	2.30	0.60
1:A:22:PRO:CB	1:A:269:LEU:HD22	2.31	0.60
1:A:146:ARG:NH1	4:A:603:HOH:O	2.35	0.59
1:B:85:VAL:HA	1:B:90:MSE:CE	2.33	0.58
1:E:112:SER:O	1:E:113:HIS:HB2	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:85:VAL:HA	1:D:90:MSE:CE	2.34	0.58
1:D:268:LEU:HB2	1:D:274:ILE:HD11	1.86	0.57
1:C:267:ALA:O	1:C:271:GLU:HG3	2.04	0.57
1:E:35:GLN:OE1	1:E:220:SER:HA	2.05	0.57
1:D:142:ASP:OD1	1:D:207:ARG:HD3	2.05	0.56
1:C:88:LEU:HD12	1:C:90:MSE:HE2	1.87	0.56
1:F:88:LEU:HD12	1:F:90:MSE:HE2	1.88	0.55
1:E:177:PHE:HB2	1:E:211:HIS:HB3	1.89	0.55
1:C:221:LEU:HB2	4:C:507:HOH:O	2.05	0.55
1:E:23:TRP:CH2	1:E:274:ILE:HD11	2.41	0.55
1:D:163:LEU:HG	1:D:210:MSE:CE	2.37	0.55
1:B:194:ARG:HB2	1:B:197:LEU:HD12	1.89	0.55
1:A:142:ASP:OD1	1:A:207:ARG:HD3	2.07	0.54
1:C:137:ARG:NH1	1:C:137:ARG:HG3	2.20	0.54
1:D:146:ARG:HG3	1:D:146:ARG:HH11	1.72	0.54
1:A:210:MSE:O	1:A:210:MSE:HG3	2.08	0.54
1:B:26:ILE:HD11	1:B:28:GLN:O	2.07	0.54
1:F:35:GLN:OE1	1:F:220:SER:HA	2.08	0.54
1:D:177:PHE:HB2	1:D:211:HIS:HB3	1.89	0.54
1:E:123:LEU:HD21	1:E:210:MSE:HB3	1.90	0.54
1:E:163:LEU:HG	1:E:210:MSE:HE1	1.88	0.54
1:F:113:HIS:HA	1:F:118:ASP:HB2	1.88	0.54
1:E:184:GLN:NE2	1:E:240:GLN:HE22	1.94	0.53
1:B:147:ASP:OD1	1:B:147:ASP:C	2.47	0.53
1:E:165:LYS:O	1:E:169:GLN:HG3	2.09	0.53
1:F:267:ALA:O	1:F:271:GLU:HG3	2.08	0.53
1:C:163:LEU:HG	1:C:210:MSE:HE1	1.89	0.53
1:D:113:HIS:HA	1:D:118:ASP:HB2	1.89	0.53
1:C:110:HIS:CE1	1:C:209:HIS:NE2	2.76	0.53
1:E:269:LEU:HD23	1:E:274:ILE:HD12	1.90	0.53
1:C:85:VAL:HA	1:C:90:MSE:CE	2.39	0.52
1:E:149:LYS:HB2	1:E:150:HIS:HD2	1.73	0.52
1:D:273:VAL:HG12	1:D:273:VAL:O	2.09	0.52
1:B:142:ASP:N	4:B:511:HOH:O	2.41	0.52
1:A:269:LEU:O	1:A:270:ASP:HB2	2.08	0.52
1:E:88:LEU:HB2	1:E:90:MSE:HE2	1.91	0.52
1:A:163:LEU:HG	1:A:210:MSE:HE1	1.91	0.51
1:A:123:LEU:HD21	1:A:210:MSE:HB3	1.93	0.51
1:A:176:ILE:HA	1:A:211:HIS:O	2.10	0.51
1:B:269:LEU:HD23	1:B:274:ILE:HD12	1.92	0.51
1:A:261:LEU:HD12	1:A:262:PRO:HD2	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:261:LEU:HD12	1:A:262:PRO:CD	2.41	0.51
1:D:85:VAL:HA	1:D:90:MSE:HE3	1.93	0.51
1:C:90:MSE:HE1	1:C:166:LEU:HD11	1.92	0.51
1:E:171:LYS:O	1:E:215:ARG:NH2	2.43	0.51
1:D:194:ARG:HB2	1:D:197:LEU:HD12	1.91	0.50
1:C:123:LEU:HD21	1:C:210:MSE:HB3	1.92	0.50
1:C:177:PHE:HB2	1:C:211:HIS:HB3	1.92	0.50
1:B:90:MSE:HE1	1:B:166:LEU:CD1	2.41	0.50
1:E:27:THR:HA	1:E:66:ARG:NH1	2.26	0.50
1:E:26:ILE:HG23	1:E:101:PRO:HB2	1.93	0.50
1:F:163:LEU:HG	1:F:210:MSE:CE	2.35	0.50
1:E:176:ILE:HA	1:E:211:HIS:O	2.12	0.50
1:E:142:ASP:OD1	1:E:207:ARG:HD3	2.12	0.50
1:D:27:THR:HA	1:D:66:ARG:NH1	2.27	0.49
1:A:268:LEU:C	1:A:270:ASP:H	2.14	0.49
1:B:85:VAL:HG13	1:B:90:MSE:HE3	1.93	0.49
1:D:269:LEU:HD23	1:D:274:ILE:HD12	1.95	0.49
1:E:90:MSE:HE1	1:E:166:LEU:HD12	1.93	0.48
1:D:92:THR:O	1:D:126:PRO:HD2	2.12	0.48
1:A:20:ALA:N	1:A:24:GLN:OE1	2.46	0.48
1:C:71:HIS:CE1	1:C:72:PRO:HD2	2.49	0.48
1:D:71:HIS:CD2	1:D:217:PRO:HG2	2.48	0.48
1:E:88:LEU:HD12	1:E:90:MSE:HE2	1.95	0.47
1:C:71:HIS:ND1	1:C:72:PRO:HD2	2.29	0.47
1:F:184:GLN:NE2	1:F:243:PHE:CE2	2.81	0.47
1:E:171:LYS:NZ	1:E:172:ASP:OD1	2.31	0.47
1:E:163:LEU:HG	1:E:210:MSE:CE	2.45	0.47
1:B:149:LYS:HA	1:B:243:PHE:CE1	2.50	0.47
1:A:201:ARG:HA	1:A:202:PRO:HD3	1.81	0.47
1:C:90:MSE:HE1	1:C:166:LEU:CD1	2.45	0.47
1:C:147:ASP:C	1:C:147:ASP:OD1	2.53	0.47
1:A:112:SER:O	1:A:113:HIS:HB2	2.15	0.47
1:E:113:HIS:HA	1:E:118:ASP:HB2	1.95	0.47
1:D:34:ALA:HA	1:D:46:VAL:O	2.15	0.47
1:E:88:LEU:CD1	1:E:90:MSE:HE2	2.44	0.47
1:E:131:THR:OG1	1:E:134:GLN:HG3	2.15	0.47
1:F:177:PHE:HB2	1:F:211:HIS:HB3	1.97	0.46
1:B:147:ASP:OD1	1:B:148:GLY:N	2.49	0.46
1:A:216:CYS:HB2	1:A:225:ASP:OD2	2.14	0.46
1:C:194:ARG:CZ	1:D:194:ARG:CZ	2.93	0.46
1:E:85:VAL:HA	1:E:90:MSE:CE	2.42	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:160:ILE:O	1:A:164:ILE:HG13	2.15	0.46
1:A:147:ASP:OD1	1:A:147:ASP:C	2.54	0.46
1:B:139:GLN:NE2	1:B:140:ALA:O	2.47	0.46
1:D:274:ILE:OXT	1:D:274:ILE:HG22	2.16	0.46
1:F:85:VAL:HG13	1:F:90:MSE:HE3	1.98	0.46
1:A:88:LEU:HD13	1:A:90:MSE:HE2	1.98	0.45
1:E:236:GLY:O	1:E:240:GLN:HG2	2.17	0.45
1:B:51:LEU:HD22	1:B:98:MSE:SE	2.66	0.45
1:B:145:SER:HB3	1:B:150:HIS:O	2.16	0.45
1:C:69:PHE:CE2	1:C:101:PRO:HG3	2.51	0.45
1:D:184:GLN:HB3	1:D:184:GLN:HE21	1.58	0.45
1:B:88:LEU:HD12	1:B:90:MSE:CE	2.46	0.45
1:E:190:ALA:HB3	1:E:194:ARG:HH12	1.82	0.45
1:D:269:LEU:HD13	1:F:87:ASN:HB3	1.99	0.45
1:C:201:ARG:HA	1:C:202:PRO:HD3	1.83	0.45
1:A:184:GLN:HB3	1:A:184:GLN:HE21	1.51	0.44
1:B:85:VAL:HA	1:B:90:MSE:HE2	1.98	0.44
1:F:191:GLY:O	1:F:194:ARG:HD2	2.17	0.44
1:F:176:ILE:HA	1:F:211:HIS:O	2.17	0.44
1:E:118:ASP:HB3	1:E:211:HIS:CE1	2.52	0.44
1:C:269:LEU:HD23	1:C:274:ILE:CD1	2.34	0.43
1:A:261:LEU:HD23	1:A:266:GLN:OE1	2.17	0.43
1:F:201:ARG:HA	1:F:202:PRO:HD3	1.82	0.43
1:D:274:ILE:CG2	1:D:274:ILE:OXT	2.65	0.43
1:A:127:LYS:HD2	1:C:221:LEU:CD1	2.45	0.43
1:F:100:MSE:HB3	1:F:101:PRO:HD2	2.01	0.43
1:D:124:GLN:CD	1:D:141:LEU:HD22	2.39	0.43
1:D:27:THR:HG22	1:D:66:ARG:CZ	2.48	0.43
1:A:179:ASN:CG	1:A:180:PRO:HD2	2.39	0.43
1:E:85:VAL:HG13	1:E:90:MSE:HB2	2.00	0.43
1:E:122:PHE:HZ	1:E:140:ALA:HA	1.83	0.43
1:E:157:LYS:HB3	1:E:159:GLU:OE1	2.18	0.43
1:F:122:PHE:HD2	1:F:209:HIS:CD2	2.35	0.43
1:A:261:LEU:HA	1:A:262:PRO:HD3	1.94	0.43
1:B:136:LEU:HA	1:B:136:LEU:HD12	1.86	0.43
1:D:210:MSE:O	1:D:210:MSE:HG3	2.19	0.43
1:B:85:VAL:HA	1:B:90:MSE:HE3	2.00	0.43
1:E:159:GLU:CD	1:E:159:GLU:H	2.22	0.43
1:F:74:LEU:HA	1:F:214:LEU:HD11	2.00	0.43
1:E:179:ASN:CG	1:E:180:PRO:HD2	2.39	0.42
1:A:175:ARG:HH11	1:A:175:ARG:HG3	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:184:GLN:HB3	1:B:184:GLN:HE21	1.62	0.42
1:C:259:PRO:HA	1:C:260:PRO:HD3	1.85	0.42
1:A:177:PHE:HD2	1:A:211:HIS:CD2	2.38	0.42
1:D:201:ARG:HA	1:D:202:PRO:HD3	1.80	0.42
1:F:67:ARG:O	1:F:98:MSE:HB2	2.20	0.42
1:A:67:ARG:O	1:A:98:MSE:HB2	2.19	0.42
1:E:146:ARG:HG2	1:E:146:ARG:H	1.45	0.42
1:A:163:LEU:HG	1:A:210:MSE:CE	2.50	0.42
1:B:92:THR:O	1:B:126:PRO:HD2	2.19	0.42
1:C:61:MSE:HE1	1:C:122:PHE:CD2	2.55	0.42
1:B:90:MSE:HE1	1:B:166:LEU:HD11	2.02	0.41
1:F:24:GLN:NE2	1:F:107:ASN:HB2	2.35	0.41
1:E:126:PRO:HB2	1:E:128:THR:O	2.20	0.41
1:A:150:HIS:CE1	1:B:149:LYS:HE3	2.55	0.41
1:C:179:ASN:CG	1:C:180:PRO:HD2	2.40	0.41
1:D:85:VAL:HG13	1:D:90:MSE:HE3	2.01	0.41
1:C:136:LEU:HD12	1:C:136:LEU:HA	1.91	0.41
1:B:42:ASN:OD1	1:B:261:LEU:HD12	2.20	0.41
1:D:238:GLU:O	1:D:241:SER:OG	2.26	0.41
1:F:155:LEU:HD23	1:F:155:LEU:HA	1.86	0.41
1:D:192:THR:O	1:D:192:THR:HG22	2.20	0.41
1:E:185:GLN:HB2	4:E:615:HOH:O	2.21	0.41
1:D:100:MSE:HB3	1:D:101:PRO:HD2	2.03	0.41
1:D:100:MSE:HE3	1:D:103:GLY:O	2.20	0.41
1:E:74:LEU:HA	1:E:214:LEU:HD11	2.03	0.41
1:C:163:LEU:HG	1:C:210:MSE:CE	2.50	0.41
1:D:144:VAL:HG22	1:D:181:ALA:HB3	2.02	0.41
1:B:94:LEU:HA	1:B:94:LEU:HD23	1.93	0.41
1:B:85:VAL:HG13	1:B:90:MSE:HB2	2.02	0.41
1:B:146:ARG:N	1:B:146:ARG:CD	2.42	0.41
1:A:104:GLY:O	1:A:114:GLN:HG2	2.21	0.41
1:C:184:GLN:HB3	1:C:184:GLN:HE21	1.70	0.41
1:B:163:LEU:HG	1:B:210:MSE:CE	2.46	0.41
1:D:36:SER:HB3	1:D:115:THR:HB	2.03	0.41
1:B:77:PHE:CZ	1:B:81:LEU:HD22	2.55	0.41
1:E:179:ASN:ND2	1:E:180:PRO:HD2	2.36	0.40
1:D:136:LEU:HA	1:D:136:LEU:HD12	1.81	0.40
1:D:124:GLN:OE1	1:D:141:LEU:HD22	2.22	0.40
1:C:122:PHE:CZ	1:C:140:ALA:HA	2.56	0.40
1:C:155:LEU:O	1:C:157:LYS:HE2	2.21	0.40
1:B:147:ASP:OD1	1:B:149:LYS:N	2.51	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:100:MSE:HB3	1:D:101:PRO:CD	2.52	0.40
1:A:113:HIS:NE2	1:A:211:HIS:HE1	2.18	0.40
1:C:206:HIS:HA	1:C:209:HIS:CE1	2.56	0.40
1:E:61:MSE:HE1	1:E:122:PHE:CD2	2.56	0.40
1:A:198:ARG:NH2	1:A:229:PRO:O	2.39	0.40
1:C:113:HIS:HA	1:C:118:ASP:HB2	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	226/255 (89%)	216 (96%)	8 (4%)	2 (1%)	21	30
1	B	235/255 (92%)	227 (97%)	7 (3%)	1 (0%)	39	56
1	C	237/255 (93%)	226 (95%)	11 (5%)	0	100	100
1	D	237/255 (93%)	224 (94%)	12 (5%)	1 (0%)	39	56
1	E	237/255 (93%)	225 (95%)	10 (4%)	2 (1%)	24	35
1	F	236/255 (92%)	225 (95%)	11 (5%)	0	100	100
All	All	1408/1530 (92%)	1343 (95%)	59 (4%)	6 (0%)	39	56

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	42	ASN
1	A	129	ARG
1	D	42	ASN
1	B	42	ASN
1	E	230	PRO
1	A	26	ILE

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	199/214 (93%)	192 (96%)	7 (4%)	43	64
1	B	205/214 (96%)	195 (95%)	10 (5%)	31	48
1	C	207/214 (97%)	201 (97%)	6 (3%)	50	71
1	D	207/214 (97%)	203 (98%)	4 (2%)	65	83
1	E	207/214 (97%)	202 (98%)	5 (2%)	57	76
1	F	206/214 (96%)	203 (98%)	3 (2%)	72	87
All	All	1231/1284 (96%)	1196 (97%)	35 (3%)	51	72

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	122	PHE
1	A	136	LEU
1	A	154	THR
1	A	184	GLN
1	A	220	SER
1	A	240	GLN
1	A	270	ASP
1	B	72	PRO
1	B	105	ARG
1	B	122	PHE
1	B	136	LEU
1	B	146	ARG
1	B	184	GLN
1	B	220	SER
1	B	225	ASP
1	B	240	GLN
1	B	264	SER
1	C	86	SER
1	C	122	PHE
1	C	128	THR
1	C	136	LEU
1	C	184	GLN

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Mol	Chain	Res	Type
1	C	220	SER
1	D	136	LEU
1	D	139	GLN
1	D	184	GLN
1	D	240	GLN
1	E	128	THR
1	E	146	ARG
1	E	154	THR
1	E	184	GLN
1	E	261	LEU
1	F	122	PHE
1	F	220	SER
1	F	241	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	169	GLN
1	A	184	GLN
1	A	185	GLN
1	B	139	GLN
1	B	169	GLN
1	B	185	GLN
1	B	240	GLN
1	C	124	GLN
1	C	205	GLN
1	D	184	GLN
1	D	185	GLN
1	E	57	HIS
1	E	65	GLN
1	E	184	GLN
1	F	169	GLN
1	F	185	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 15 ligands modelled in this entry, 9 are monoatomic - leaving 6 for Mogul analysis.

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	$\# Z > 2$
3	SO4	A	500	-	4,4,4	1.75	2 (50%)	6,6,6	1.05	0
3	SO4	B	501	-	4,4,4	1.79	1 (25%)	6,6,6	1.01	0
3	SO4	C	502	-	4,4,4	1.89	1 (25%)	6,6,6	0.84	0
3	SO4	D	503	-	4,4,4	1.84	3 (75%)	6,6,6	0.96	0
3	SO4	E	504	-	4,4,4	1.93	1 (25%)	6,6,6	0.95	0
3	SO4	F	505	-	4,4,4	1.68	1 (25%)	6,6,6	1.02	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
3	SO4	A	500	-	-	0/0/0/0	0/0/0/0
3	SO4	B	501	-	-	0/0/0/0	0/0/0/0
3	SO4	C	502	-	-	0/0/0/0	0/0/0/0
3	SO4	D	503	-	-	0/0/0/0	0/0/0/0
3	SO4	E	504	-	-	0/0/0/0	0/0/0/0
3	SO4	F	505	-	-	0/0/0/0	0/0/0/0

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	504	SO4	O1-S	-2.78	1.37	1.47
3	C	502	SO4	O3-S	-2.74	1.37	1.47
3	B	501	SO4	O1-S	-2.55	1.38	1.47
3	D	503	SO4	O1-S	-2.25	1.39	1.47
3	A	500	SO4	O3-S	-2.09	1.39	1.47
3	D	503	SO4	O3-S	-2.07	1.39	1.47
3	F	505	SO4	O3-S	-2.04	1.40	1.47
3	D	503	SO4	O4-S	-2.04	1.40	1.47
3	A	500	SO4	O1-S	-2.03	1.40	1.47

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	226/255 (88%)	0.43	15 (6%)	22 22	11, 29, 48, 57	0
1	B	233/255 (91%)	0.25	13 (5%)	28 28	7, 25, 48, 59	0
1	C	235/255 (92%)	0.07	6 (2%)	59 58	8, 23, 42, 54	0
1	D	235/255 (92%)	0.17	10 (4%)	39 40	6, 23, 43, 56	0
1	E	235/255 (92%)	0.25	13 (5%)	29 29	10, 24, 46, 57	0
1	F	234/255 (91%)	0.16	11 (4%)	35 36	7, 22, 46, 53	0
All	All	1398/1530 (91%)	0.22	68 (4%)	33 34	6, 25, 46, 59	0

All (68) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	141	LEU	11.4
1	D	141	LEU	10.3
1	A	141	LEU	7.8
1	A	20	ALA	6.6
1	E	140	ALA	6.0
1	B	141	LEU	5.9
1	B	274	ILE	5.4
1	D	140	ALA	5.3
1	A	231	SER	5.2
1	F	221	LEU	5.1
1	E	139	GLN	4.9
1	E	274	ILE	4.7
1	D	139	GLN	4.6
1	D	274	ILE	4.6
1	B	221	LEU	4.4
1	B	244	GLU	4.3
1	F	274	ILE	4.1
1	C	146	ARG	3.8
1	A	22	PRO	3.7

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Mol	Chain	Res	Type	RSRZ
1	C	274	ILE	3.7
1	F	141	LEU	3.6
1	A	139	GLN	3.6
1	D	138	PRO	3.6
1	A	105	ARG	3.5
1	E	192	THR	3.5
1	B	140	ALA	3.4
1	D	137	ARG	3.2
1	F	146	ARG	3.2
1	C	244	GLU	3.1
1	D	260	PRO	3.1
1	B	154	THR	3.1
1	C	260	PRO	3.0
1	A	25	LYS	3.0
1	E	146	ARG	3.0
1	A	24	GLN	3.0
1	B	139	GLN	2.9
1	F	244	GLU	2.9
1	B	220	SER	2.9
1	B	231	SER	2.9
1	F	192	THR	2.9
1	E	20	ALA	2.8
1	E	244	GLU	2.7
1	B	111	ALA	2.7
1	D	259	PRO	2.7
1	A	111	ALA	2.7
1	B	240	GLN	2.6
1	D	146	ARG	2.6
1	F	139	GLN	2.5
1	A	109	GLY	2.4
1	D	244	GLU	2.4
1	E	111	ALA	2.4
1	F	220	SER	2.3
1	B	273	VAL	2.3
1	A	270	ASP	2.3
1	A	41	SER	2.2
1	A	21	THR	2.2
1	C	20	ALA	2.2
1	E	272	HIS	2.2
1	F	218	ALA	2.2
1	A	140	ALA	2.1
1	E	107	ASN	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	105	ARG	2.1
1	A	27	THR	2.1
1	E	153	SER	2.1
1	F	223	CYS	2.0
1	E	218	ALA	2.0
1	F	56	GLU	2.0
1	C	221	LEU	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](#)

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
3	SO4	B	501	5/5	0.90	0.25	1.22	43,45,46,47	0
3	SO4	F	505	5/5	0.83	0.24	1.10	44,45,47,49	0
3	SO4	C	502	5/5	0.94	0.18	0.92	32,32,33,35	0
3	SO4	D	503	5/5	0.95	0.18	0.47	43,44,45,45	0
3	SO4	E	504	5/5	0.92	0.14	-0.46	39,40,42,43	0
2	ZN	A	601	1/1	0.77	0.14	-0.65	49,49,49,49	0
3	SO4	A	500	5/5	0.94	0.13	-1.17	43,43,44,46	0
2	ZN	E	603	1/1	0.92	0.07	-1.87	33,33,33,33	0
2	ZN	F	400	1/1	1.00	0.08	-2.88	18,18,18,18	0
2	ZN	C	400	1/1	1.00	0.04	-3.58	14,14,14,14	0
2	ZN	D	602	1/1	0.96	0.04	-3.63	27,27,27,27	0
2	ZN	E	400	1/1	1.00	0.05	-	16,16,16,16	0
2	ZN	B	400	1/1	0.99	0.04	-	15,15,15,15	0
2	ZN	A	400	1/1	0.98	0.06	-	25,25,25,25	0
2	ZN	D	400	1/1	1.00	0.05	-	11,11,11,11	0

6.5 Other polymers ⓘ

There are no such residues in this entry.