



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 11:26 AM UTC

PDB ID : 3KCM / pdb_00003kcm
Title : The crystal structure of thioredoxin protein from *Geobacter metallireducens*
Authors : Zhang, Z.; Burley, S.K.; Swaminathan, S.; New York SGX Research Center
for Structural Genomics (NYSGXRC)
Deposited on : 2009-10-21
Resolution : 2.45 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

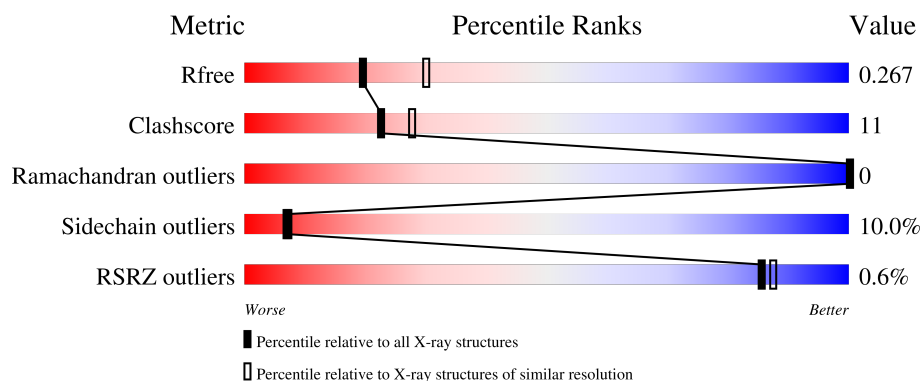
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.45 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	1190 (2.46-2.46)
Clashscore	190562	1229 (2.46-2.46)
Ramachandran outliers	187476	1218 (2.46-2.46)
Sidechain outliers	187428	1218 (2.46-2.46)
RSRZ outliers	180081	1190 (2.46-2.46)

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	154	 63% 24% 10%
1	B	154	 2% 66% 21% 5% 8%
1	C	154	 60% 23% 6% 10%
1	D	154	 71% 18% 8%
1	E	154	 70% 19% 8%

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Mol	Chain	Length	Quality of chain
1	F	154	62% 23% 5% 10%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6720 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 Thioredoxin family protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	139	Total	C	N	O	S	Se	0	0	0
			1081	698	179	196	3	5			
1	B	141	Total	C	N	O	S	Se	0	0	0
			1088	704	180	196	3	5			
1	C	139	Total	C	N	O	S	Se	0	0	0
			1075	695	177	195	3	5			
1	D	142	Total	C	N	O	S	Se	0	0	0
			1091	705	181	197	3	5			
1	E	141	Total	C	N	O	S	Se	0	0	0
			1094	706	184	196	3	5			
1	F	138	Total	C	N	O	S	Se	0	0	0
			1063	687	176	192	3	5			

There are 66 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	25	MSE	-	expression tag	UNP Q39SU7
A	26	SER	-	expression tag	UNP Q39SU7
A	27	LEU	-	expression tag	UNP Q39SU7
A	171	GLU	-	expression tag	UNP Q39SU7
A	172	GLY	-	expression tag	UNP Q39SU7
A	173	HIS	-	expression tag	UNP Q39SU7
A	174	HIS	-	expression tag	UNP Q39SU7
A	175	HIS	-	expression tag	UNP Q39SU7
A	176	HIS	-	expression tag	UNP Q39SU7
A	177	HIS	-	expression tag	UNP Q39SU7
A	178	HIS	-	expression tag	UNP Q39SU7
B	25	MSE	-	expression tag	UNP Q39SU7
B	26	SER	-	expression tag	UNP Q39SU7
B	27	LEU	-	expression tag	UNP Q39SU7
B	171	GLU	-	expression tag	UNP Q39SU7
B	172	GLY	-	expression tag	UNP Q39SU7
B	173	HIS	-	expression tag	UNP Q39SU7

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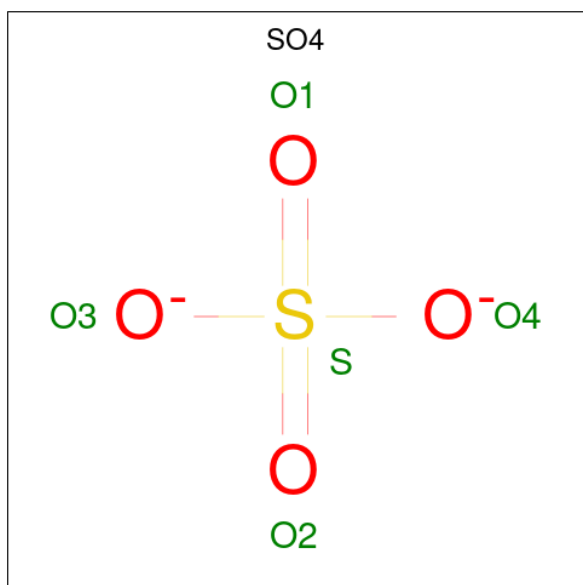
Chain	Residue	Modelled	Actual	Comment	Reference
B	174	HIS	-	expression tag	UNP Q39SU7
B	175	HIS	-	expression tag	UNP Q39SU7
B	176	HIS	-	expression tag	UNP Q39SU7
B	177	HIS	-	expression tag	UNP Q39SU7
B	178	HIS	-	expression tag	UNP Q39SU7
C	25	MSE	-	expression tag	UNP Q39SU7
C	26	SER	-	expression tag	UNP Q39SU7
C	27	LEU	-	expression tag	UNP Q39SU7
C	171	GLU	-	expression tag	UNP Q39SU7
C	172	GLY	-	expression tag	UNP Q39SU7
C	173	HIS	-	expression tag	UNP Q39SU7
C	174	HIS	-	expression tag	UNP Q39SU7
C	175	HIS	-	expression tag	UNP Q39SU7
C	176	HIS	-	expression tag	UNP Q39SU7
C	177	HIS	-	expression tag	UNP Q39SU7
C	178	HIS	-	expression tag	UNP Q39SU7
D	25	MSE	-	expression tag	UNP Q39SU7
D	26	SER	-	expression tag	UNP Q39SU7
D	27	LEU	-	expression tag	UNP Q39SU7
D	171	GLU	-	expression tag	UNP Q39SU7
D	172	GLY	-	expression tag	UNP Q39SU7
D	173	HIS	-	expression tag	UNP Q39SU7
D	174	HIS	-	expression tag	UNP Q39SU7
D	175	HIS	-	expression tag	UNP Q39SU7
D	176	HIS	-	expression tag	UNP Q39SU7
D	177	HIS	-	expression tag	UNP Q39SU7
D	178	HIS	-	expression tag	UNP Q39SU7
E	25	MSE	-	expression tag	UNP Q39SU7
E	26	SER	-	expression tag	UNP Q39SU7
E	27	LEU	-	expression tag	UNP Q39SU7
E	171	GLU	-	expression tag	UNP Q39SU7
E	172	GLY	-	expression tag	UNP Q39SU7
E	173	HIS	-	expression tag	UNP Q39SU7
E	174	HIS	-	expression tag	UNP Q39SU7
E	175	HIS	-	expression tag	UNP Q39SU7
E	176	HIS	-	expression tag	UNP Q39SU7
E	177	HIS	-	expression tag	UNP Q39SU7
E	178	HIS	-	expression tag	UNP Q39SU7
F	25	MSE	-	expression tag	UNP Q39SU7
F	26	SER	-	expression tag	UNP Q39SU7
F	27	LEU	-	expression tag	UNP Q39SU7
F	171	GLU	-	expression tag	UNP Q39SU7

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Chain	Residue	Modelled	Actual	Comment	Reference
F	172	GLY	-	expression tag	UNP Q39SU7
F	173	HIS	-	expression tag	UNP Q39SU7
F	174	HIS	-	expression tag	UNP Q39SU7
F	175	HIS	-	expression tag	UNP Q39SU7
F	176	HIS	-	expression tag	UNP Q39SU7
F	177	HIS	-	expression tag	UNP Q39SU7
F	178	HIS	-	expression tag	UNP Q39SU7

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



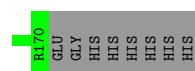
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	E	1	Total	O	S	0	0
			5	4	1		
2	F	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is water.

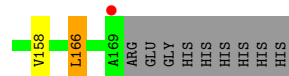
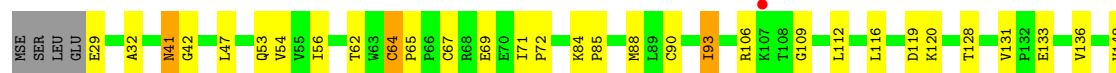
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	39	Total 39	O 39	0	0
3	B	28	Total 28	O 28	0	0
3	C	27	Total 27	O 27	0	0
3	D	37	Total 37	O 37	0	0
3	E	39	Total 39	O 39	0	0
3	F	28	Total 28	O 28	0	0

- Molecule 1: Thioredoxin family protein

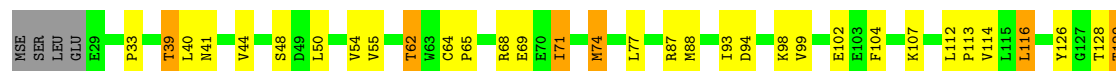




- Molecule 1: Thioredoxin family protein



- Molecule 1: Thioredoxin family protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	51.99Å 74.56Å 82.41Å 100.66° 109.05° 99.07°	Depositor
Resolution (Å)	75.36 – 2.45 75.36 – 2.45	Depositor EDS
% Data completeness (in resolution range)	94.4 (75.36-2.45) 94.4 (75.36-2.45)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.77 (at 2.45Å)	Xtriage
Refinement program	REFMAC 5.5.0072	Depositor
R, R_{free}	0.214 , 0.270 0.214 , 0.267	Depositor DCC
R_{free} test set	1953 reflections (4.75%)	wwPDB-VP
Wilson B-factor (Å ²)	37.0	Xtriage
Anisotropy	0.050	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 42.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	6720	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.82% 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 ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: 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	1.14	1/1101 (0.1%)	1.18	3/1487 (0.2%)
1	B	1.12	1/1108 (0.1%)	1.25	3/1498 (0.2%)
1	C	1.03	1/1095 (0.1%)	1.20	3/1480 (0.2%)
1	D	1.17	0/1111	1.27	8/1501 (0.5%)
1	E	1.20	5/1114 (0.4%)	1.19	4/1503 (0.3%)
1	F	1.07	2/1083 (0.2%)	1.19	5/1464 (0.3%)
All	All	1.13	10/6612 (0.2%)	1.22	26/8933 (0.3%)

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	71	ILE	CA-CB	7.72	1.58	1.54
1	E	53	GLN	C-O	-6.24	1.15	1.23
1	B	113	PRO	C-O	-6.03	1.16	1.24
1	E	47	LEU	C-O	-5.78	1.17	1.24
1	E	158	VAL	CA-CB	5.51	1.60	1.54
1	A	51	LYS	C-O	-5.49	1.17	1.23
1	F	71	ILE	CA-CB	5.29	1.60	1.54
1	E	71	ILE	CA-C	5.17	1.57	1.52
1	E	65	PRO	CA-C	5.09	1.56	1.52
1	F	99	VAL	C-O	-5.00	1.18	1.24

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	155	HIS	CA-C-N	7.46	127.00	119.24
1	A	155	HIS	C-N-CA	7.46	127.00	119.24
1	D	30	ASN	N-CA-C	7.28	119.97	109.48
1	B	155	HIS	CA-C-N	7.12	128.73	119.84
1	B	155	HIS	C-N-CA	7.12	128.73	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	65	PRO	CA-C-N	-7.02	111.56	119.28
1	D	65	PRO	C-N-CA	-7.02	111.56	119.28
1	F	65	PRO	CA-C-N	-6.45	112.35	119.19
1	F	65	PRO	C-N-CA	-6.45	112.35	119.19
1	E	32	ALA	CA-C-N	-5.94	113.41	119.83
1	E	32	ALA	C-N-CA	-5.94	113.41	119.83
1	C	65	PRO	N-CA-C	5.83	117.81	110.70
1	C	70	GLU	CA-C-N	5.79	124.96	120.33
1	C	70	GLU	C-N-CA	5.79	124.96	120.33
1	E	65	PRO	N-CA-C	5.78	117.75	110.70
1	D	84	LYS	CA-C-N	5.57	126.80	119.84
1	D	84	LYS	C-N-CA	5.57	126.80	119.84
1	E	42	GLY	N-CA-C	5.56	123.30	115.30
1	B	112	LEU	N-CA-C	5.53	113.20	108.22
1	A	78	ASN	N-CA-C	5.38	116.83	111.07
1	D	47	LEU	N-CA-C	5.36	117.20	111.36
1	F	65	PRO	N-CA-C	5.32	117.19	110.70
1	F	116	LEU	N-CA-C	5.18	117.68	109.24
1	F	129	THR	N-CA-C	-5.17	107.64	114.31
1	D	155	HIS	CA-C-N	5.14	125.18	119.32
1	D	155	HIS	C-N-CA	5.14	125.18	119.32

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	1081	0	1081	29	0
1	B	1088	0	1083	20	0
1	C	1075	0	1072	29	0
1	D	1091	0	1091	19	0
1	E	1094	0	1107	16	0
1	F	1063	0	1057	27	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	5	0	0	1	0
2	D	5	0	0	0	0
2	E	5	0	0	0	0
2	F	5	0	0	0	0
3	A	39	0	0	3	0
3	B	28	0	0	1	0
3	C	27	0	0	2	0
3	D	37	0	0	0	0
3	E	39	0	0	3	0
3	F	28	0	0	1	0
All	All	6720	0	6491	137	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 11.

All (137) 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:D:88:MSE:HE1	1:D:112:LEU:HD22	1.28	1.09
1:D:88:MSE:HE3	1:D:90:CYS:SG	2.03	0.98
1:D:88:MSE:CE	1:D:112:LEU:HD22	1.96	0.96
1:B:95:GLU:HG3	3:B:231:HOH:O	1.73	0.88
1:F:139:ARG:HE	1:F:139:ARG:H	1.22	0.88
1:A:39:THR:HG21	3:A:204:HOH:O	1.72	0.87
1:A:88:MSE:HE3	1:A:90:CYS:SG	2.16	0.85
1:C:139:ARG:HG2	3:C:247:HOH:O	1.76	0.85
1:C:28:GLU:OE1	1:C:146:LYS:HB2	1.80	0.81
1:A:88:MSE:CE	1:A:112:LEU:HD22	2.14	0.78
1:F:62:THR:HB	1:F:94:ASP:OD1	1.88	0.74
1:A:88:MSE:HE1	1:A:112:LEU:HD22	1.68	0.73
1:A:165:GLU:HG3	3:A:197:HOH:O	1.88	0.73
1:B:138:ASP:HB3	1:B:144:LEU:HD23	1.72	0.72
1:C:62:THR:HB	1:C:94:ASP:OD1	1.90	0.71
1:E:88:MSE:CE	1:E:112:LEU:HD22	2.21	0.71
1:C:95:GLU:HB2	2:C:6:SO4:O4	1.91	0.71
1:C:138:ASP:HB3	1:C:144:LEU:HD22	1.74	0.70
1:F:138:ASP:HB2	1:F:139:ARG:HH21	1.59	0.68
1:E:88:MSE:HE1	1:E:112:LEU:HD22	1.75	0.67
1:F:71:ILE:HA	1:F:74:MSE:HE2	1.77	0.67
1:D:88:MSE:HE1	1:D:112:LEU:CD2	2.14	0.67
1:C:84:LYS:HB3	1:C:85:PRO:HD2	1.78	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:133:GLU:HG2	1:B:148:VAL:HG12	1.78	0.64
1:C:163:ASN:HA	1:C:166:LEU:HD12	1.80	0.63
1:F:98:LYS:HG2	1:F:102:GLU:OE2	1.99	0.63
1:C:84:LYS:HE3	1:C:163:ASN:OD1	1.99	0.62
1:D:88:MSE:CE	1:D:90:CYS:SG	2.86	0.62
1:F:55:VAL:HA	1:F:87:ARG:O	2.01	0.60
1:A:152:GLU:OE1	1:A:152:GLU:HA	2.02	0.60
1:C:28:GLU:OE1	1:C:146:LYS:CB	2.50	0.59
1:C:77:LEU:HD13	1:C:153:TRP:O	2.02	0.59
1:F:93:ILE:HD13	1:F:131:VAL:HG11	1.84	0.59
1:C:74:MSE:CE	1:C:90:CYS:SG	2.90	0.59
1:B:74:MSE:HG2	1:B:88:MSE:SE	2.53	0.58
1:F:39:THR:HG23	1:F:41:ASN:H	1.67	0.58
1:F:133:GLU:HG3	1:F:148:VAL:HG12	1.85	0.58
1:A:88:MSE:CE	1:A:90:CYS:SG	2.91	0.58
1:A:154:ASP:OD2	1:A:155:HIS:CD2	2.57	0.58
1:F:88:MSE:HE3	1:F:113:PRO:CD	2.36	0.56
1:A:147:VAL:HG11	1:A:151:MSE:HG3	1.88	0.55
1:F:132:PRO:HD2	1:F:149:GLY:HA2	1.89	0.55
1:A:39:THR:HG22	1:A:40:LEU:N	2.22	0.55
1:C:98:LYS:O	1:C:102:GLU:HG2	2.07	0.54
1:B:48:SER:C	1:B:50:LEU:H	2.15	0.54
1:C:84:LYS:HB3	1:C:85:PRO:CD	2.38	0.54
1:E:128:THR:HA	1:E:133:GLU:OE2	2.08	0.53
1:A:31:PRO:HA	1:A:142:VAL:HA	1.91	0.53
1:C:88:MSE:CE	1:C:112:LEU:HD22	2.39	0.53
1:B:89:LEU:HD22	1:B:115:LEU:HD11	1.91	0.53
1:F:138:ASP:HB2	1:F:139:ARG:NH2	2.22	0.53
1:A:39:THR:CG2	3:A:204:HOH:O	2.44	0.52
1:B:93:ILE:O	1:B:120:LYS:NZ	2.42	0.52
1:D:37:LEU:HD22	1:D:116:LEU:O	2.09	0.52
1:F:128:THR:HG22	1:F:133:GLU:CD	2.34	0.52
1:C:152:GLU:HA	1:C:152:GLU:OE1	2.10	0.52
1:C:34:ASP:OD2	1:C:48:SER:OG	2.28	0.52
1:E:72:PRO:HD2	3:E:197:HOH:O	2.11	0.51
1:A:129:THR:HG23	1:D:87:ARG:NH2	2.26	0.50
1:E:54:VAL:HG13	1:E:54:VAL:O	2.11	0.50
1:A:155:HIS:O	1:A:159:ILE:HG12	2.12	0.50
1:F:151:MSE:HE3	3:F:211:HOH:O	2.11	0.50
1:D:88:MSE:HE2	1:D:113:PRO:HD2	1.93	0.50
1:F:40:LEU:HG	1:F:114:VAL:HB	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:34:ASP:OD2	1:B:48:SER:OG	2.28	0.50
1:F:139:ARG:H	1:F:139:ARG:NE	1.99	0.50
1:B:130:GLY:O	1:B:133:GLU:HG3	2.13	0.49
1:F:68:ARG:HG2	1:F:104:PHE:CZ	2.47	0.49
1:C:88:MSE:HE3	1:C:112:LEU:HD22	1.94	0.49
1:E:54:VAL:HG12	1:E:166:LEU:HD11	1.94	0.49
1:E:84:LYS:HB3	1:E:85:PRO:HD2	1.95	0.48
1:B:163:ASN:O	1:B:164:ASN:C	2.53	0.48
1:D:112:LEU:HB3	1:D:113:PRO:HD2	1.96	0.48
1:A:136:VAL:HB	1:A:145:LYS:HB3	1.96	0.48
1:F:88:MSE:HE3	1:F:113:PRO:HD2	1.96	0.48
1:D:67:CYS:SG	1:D:132:PRO:HG3	2.54	0.47
1:E:88:MSE:HE3	1:E:90:CYS:SG	2.54	0.47
1:E:109:GLY:HA2	3:E:186:HOH:O	2.14	0.47
1:A:39:THR:HG22	1:A:41:ASN:H	1.79	0.47
1:D:62:THR:HB	1:D:94:ASP:OD1	2.14	0.47
1:C:65:PRO:HB2	1:C:66:PRO:HD3	1.97	0.47
1:F:68:ARG:HG2	1:F:104:PHE:HZ	1.80	0.47
1:C:147:VAL:HG11	1:C:151:MSE:HG3	1.97	0.47
1:B:74:MSE:HE3	1:B:90:CYS:SG	2.55	0.46
1:A:32:ALA:HB2	1:A:143:ILE:HG13	1.96	0.46
1:B:47:LEU:O	1:B:50:LEU:HB2	2.15	0.46
1:E:64:CYS:O	1:E:67:CYS:HB2	2.15	0.46
1:A:39:THR:CG2	1:A:40:LEU:N	2.79	0.46
1:B:54:VAL:HG12	1:B:166:LEU:HD23	1.98	0.45
1:E:119:ASP:C	1:E:120:LYS:HG2	2.41	0.45
1:B:119:ASP:C	1:B:120:LYS:HG2	2.41	0.45
1:D:73:SER:HB2	1:D:154:ASP:OD2	2.17	0.45
1:A:104:PHE:O	1:A:108:THR:HG23	2.17	0.45
1:A:28:GLU:O	1:A:29:GLU:HB2	2.18	0.44
1:B:29:GLU:OE1	1:B:29:GLU:HA	2.18	0.44
1:E:56:ILE:HD13	1:E:136:VAL:HG22	2.00	0.44
1:B:86:PHE:C	1:B:86:PHE:CD2	2.95	0.44
1:E:54:VAL:CG1	1:E:166:LEU:HD11	2.48	0.44
1:C:58:ASN:ND2	1:C:74:MSE:HE1	2.32	0.44
1:F:39:THR:CG2	1:F:41:ASN:H	2.29	0.44
1:F:71:ILE:HG12	1:F:112:LEU:HD11	1.98	0.44
1:C:111:THR:O	1:C:112:LEU:HD23	2.18	0.43
1:D:86:PHE:CD2	1:D:86:PHE:C	2.96	0.43
1:A:88:MSE:HE1	1:A:112:LEU:CD2	2.42	0.43
1:E:93:ILE:O	1:E:120:LYS:NZ	2.44	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:128:THR:HA	1:F:133:GLU:OE2	2.18	0.43
1:A:65:PRO:HB2	1:A:66:PRO:HD3	2.00	0.43
1:F:128:THR:HG22	1:F:133:GLU:OE2	2.18	0.43
1:B:43:GLU:OE2	1:F:133:GLU:OE1	2.37	0.43
1:F:33:PRO:HD2	1:F:126:TYR:CZ	2.54	0.43
1:B:152:GLU:HB2	1:B:155:HIS:CD2	2.53	0.42
1:D:34:ASP:OD2	1:D:48:SER:OG	2.35	0.42
1:A:34:ASP:OD2	1:A:48:SER:OG	2.32	0.42
1:E:41:ASN:ND2	3:E:185:HOH:O	2.52	0.42
1:F:71:ILE:HA	1:F:74:MSE:CE	2.47	0.42
1:A:119:ASP:C	1:A:120:LYS:HG2	2.44	0.42
1:D:54:VAL:CG2	1:D:166:LEU:HD12	2.49	0.42
1:C:112:LEU:O	1:C:114:VAL:HG23	2.20	0.42
1:A:155:HIS:HB3	1:A:156:PRO:CD	2.51	0.41
1:C:90:CYS:SG	1:C:112:LEU:HD13	2.61	0.41
1:C:138:ASP:HB3	1:C:144:LEU:CD2	2.48	0.41
1:D:37:LEU:HD11	1:D:122:VAL:HG21	2.02	0.41
1:B:78:ASN:OD1	1:B:88:MSE:HE2	2.20	0.41
1:D:99:VAL:O	1:D:103:GLU:HG2	2.20	0.41
1:E:56:ILE:CD1	1:E:136:VAL:HG22	2.50	0.41
1:A:32:ALA:HA	1:A:33:PRO:HD2	1.96	0.41
1:B:128:THR:HG22	1:B:133:GLU:OE2	2.21	0.41
1:C:103:GLU:O	1:C:107:LYS:HD3	2.20	0.41
1:C:119:ASP:OD2	1:C:121:ARG:HB3	2.20	0.41
1:A:48:SER:C	1:A:50:LEU:H	2.29	0.41
1:C:62:THR:HB	1:C:94:ASP:CG	2.45	0.41
1:C:33:PRO:HD2	1:C:126:TYR:CZ	2.56	0.41
1:C:73:SER:HB3	3:C:235:HOH:O	2.19	0.41
1:A:64:CYS:O	1:A:68:ARG:HG3	2.21	0.40
1:F:77:LEU:HD22	1:F:153:TRP:HB3	2.01	0.40
1:A:133:GLU:OE1	1:D:43:GLU:OE2	2.39	0.40
1:D:155:HIS:HA	1:D:156:PRO:HD3	1.98	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	137/154 (89%)	129 (94%)	8 (6%)	0	100	100
1	B	139/154 (90%)	132 (95%)	7 (5%)	0	100	100
1	C	137/154 (89%)	128 (93%)	9 (7%)	0	100	100
1	D	140/154 (91%)	134 (96%)	6 (4%)	0	100	100
1	E	139/154 (90%)	135 (97%)	4 (3%)	0	100	100
1	F	136/154 (88%)	130 (96%)	6 (4%)	0	100	100
All	All	828/924 (90%)	788 (95%)	40 (5%)	0	100	100

There are no Ramachandran outliers to report.

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	117/126 (93%)	109 (93%)	8 (7%)	14	20
1	B	116/126 (92%)	102 (88%)	14 (12%)	5	4
1	C	116/126 (92%)	100 (86%)	16 (14%)	3	3
1	D	117/126 (93%)	110 (94%)	7 (6%)	17	26
1	E	119/126 (94%)	108 (91%)	11 (9%)	8	9
1	F	114/126 (90%)	100 (88%)	14 (12%)	4	4
All	All	699/756 (92%)	629 (90%)	70 (10%)	7	7

All (70) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	LEU
1	A	56	ILE
1	A	64	CYS

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Mol	Chain	Res	Type
1	A	74	MSE
1	A	95	GLU
1	A	129	THR
1	A	152	GLU
1	A	165	GLU
1	B	39	THR
1	B	41	ASN
1	B	54	VAL
1	B	56	ILE
1	B	64	CYS
1	B	69	GLU
1	B	74	MSE
1	B	77	LEU
1	B	86	PHE
1	B	93	ILE
1	B	94	ASP
1	B	136	VAL
1	B	144	LEU
1	B	152	GLU
1	C	28	GLU
1	C	41	ASN
1	C	44	VAL
1	C	53	GLN
1	C	62	THR
1	C	77	LEU
1	C	84	LYS
1	C	86	PHE
1	C	94	ASP
1	C	106	ARG
1	C	107	LYS
1	C	125	LEU
1	C	129	THR
1	C	139	ARG
1	C	144	LEU
1	C	152	GLU
1	D	46	LYS
1	D	50	LEU
1	D	62	THR
1	D	64	CYS
1	D	103	GLU
1	D	144	LEU
1	D	166	LEU

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Mol	Chain	Res	Type
1	E	29	GLU
1	E	41	ASN
1	E	62	THR
1	E	64	CYS
1	E	69	GLU
1	E	93	ILE
1	E	106	ARG
1	E	116	LEU
1	E	131	VAL
1	E	148	VAL
1	E	166	LEU
1	F	39	THR
1	F	44	VAL
1	F	48	SER
1	F	50	LEU
1	F	54	VAL
1	F	62	THR
1	F	64	CYS
1	F	69	GLU
1	F	74	MSE
1	F	107	LYS
1	F	116	LEU
1	F	129	THR
1	F	139	ARG
1	F	144	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	38	ASN
1	A	155	HIS
1	B	155	HIS
1	C	38	ASN
1	C	41	ASN
1	C	53	GLN
1	C	155	HIS
1	D	155	HIS
1	E	30	ASN
1	E	41	ASN

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](#)

6 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	SO4	A	1	-	4,4,4	0.31	0	6,6,6	0.52	0
2	SO4	B	2	-	4,4,4	0.29	0	6,6,6	0.13	0
2	SO4	D	3	-	4,4,4	0.27	0	6,6,6	0.52	0
2	SO4	C	6	-	4,4,4	0.16	0	6,6,6	0.20	0
2	SO4	F	5	-	4,4,4	0.30	0	6,6,6	0.33	0
2	SO4	E	4	-	4,4,4	0.28	0	6,6,6	0.51	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.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	6	SO4	1	0

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	134/154 (87%)	-0.27	0 100 100	12, 21, 36, 49	0
1	B	136/154 (88%)	-0.20	3 (2%) 62 64	12, 22, 40, 47	0
1	C	134/154 (87%)	-0.27	0 100 100	12, 21, 38, 51	0
1	D	137/154 (88%)	-0.47	0 100 100	12, 20, 33, 39	0
1	E	136/154 (88%)	-0.26	2 (1%) 72 74	13, 21, 39, 46	0
1	F	133/154 (86%)	-0.30	0 100 100	13, 23, 40, 47	0
All	All	810/924 (87%)	-0.30	5 (0%) 85 87	12, 21, 39, 51	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	28	GLU	2.8
1	E	169	ALA	2.7
1	B	27	LEU	2.6
1	E	107	LYS	2.4
1	B	167	SER	2.2

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(Å ²)	Q<0.9
2	SO4	E	4	5/5	0.96	0.08	24,25,26,27	0
2	SO4	D	3	5/5	0.97	0.06	26,26,29,29	0
2	SO4	C	6	5/5	0.98	0.10	27,28,28,31	0
2	SO4	F	5	5/5	0.98	0.06	24,25,25,26	0
2	SO4	A	1	5/5	0.99	0.06	19,20,22,25	0
2	SO4	B	2	5/5	0.99	0.06	22,24,25,25	0

6.5 Other polymers [i](#)

There are no such residues in this entry.