



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:51 PM UTC

PDB ID : 3IP3 / pdb_00003ip3
Title : Structure of putative oxidoreductase (TM_0425) from *Thermotoga maritima*
Authors : Ramagopal, U.A.; Morano, C.; Burley, S.K.; Almo, S.C.; New York SGX
Research Center for Structural Genomics (NYSGXRC)
Deposited on : 2009-08-16
Resolution : 2.14 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

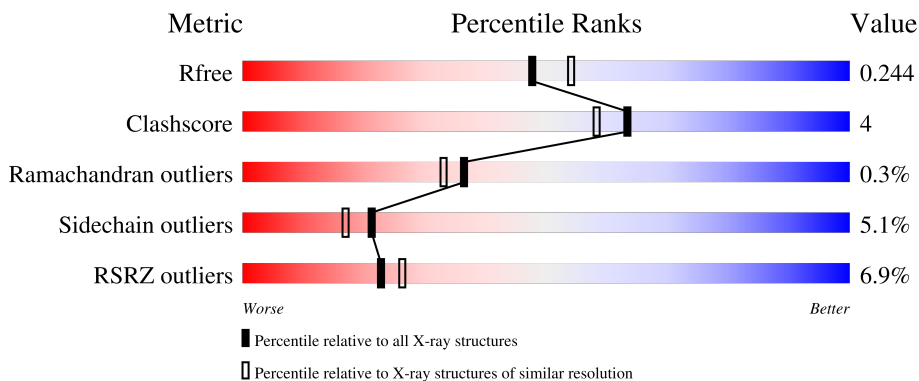
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.14 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.



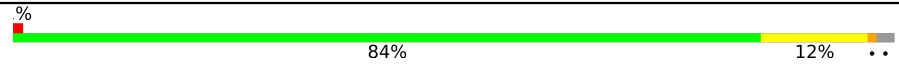

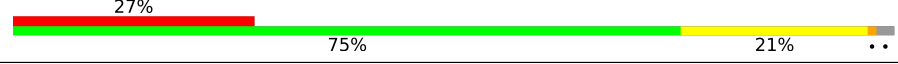
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3689 (2.16-2.12)
Clashscore	190562	3812 (2.16-2.12)
Ramachandran outliers	187476	3773 (2.16-2.12)
Sidechain outliers	187428	3772 (2.16-2.12)
RSRZ outliers	180081	3691 (2.16-2.12)

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	337	 86% 10% ..
1	B	337	 90% 8% ..
1	C	337	 85% 11% ..
1	D	337	 87% 10% ..
1	E	337	 9% 83% 13% ..

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Mol	Chain	Length	Quality of chain
1	F	337	 <p>% 12% 84% 12% ..</p>
1	G	337	 <p>12% 81% 15% ..</p>
1	H	337	 <p>27% 75% 21% ..</p>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 21415 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 Oxidoreductase, putative.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	328	Total 2632	C 1686	N 453	O 484	S 4	Se 5	0	0	0
1	B	331	Total 2657	C 1700	N 459	O 489	S 4	Se 5	0	0	0
1	C	330	Total 2647	C 1694	N 456	O 488	S 4	Se 5	0	0	0
1	D	328	Total 2640	C 1691	N 454	O 486	S 4	Se 5	0	1	0
1	E	328	Total 2634	C 1687	N 454	O 484	S 4	Se 5	0	0	0
1	F	330	Total 2647	C 1694	N 456	O 488	S 4	Se 5	0	0	0
1	G	327	Total 2628	C 1684	N 453	O 482	S 4	Se 5	0	0	0
1	H	329	Total 2641	C 1691	N 455	O 486	S 4	Se 5	0	0	0

There are 80 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MSE	-	expression tag	UNP Q9WYQ6
A	1	SER	-	expression tag	UNP Q9WYQ6
A	329	GLU	-	expression tag	UNP Q9WYQ6
A	330	GLY	-	expression tag	UNP Q9WYQ6
A	331	HIS	-	expression tag	UNP Q9WYQ6
A	332	HIS	-	expression tag	UNP Q9WYQ6
A	333	HIS	-	expression tag	UNP Q9WYQ6
A	334	HIS	-	expression tag	UNP Q9WYQ6
A	335	HIS	-	expression tag	UNP Q9WYQ6
A	336	HIS	-	expression tag	UNP Q9WYQ6
B	0	MSE	-	expression tag	UNP Q9WYQ6
B	1	SER	-	expression tag	UNP Q9WYQ6
B	329	GLU	-	expression tag	UNP Q9WYQ6

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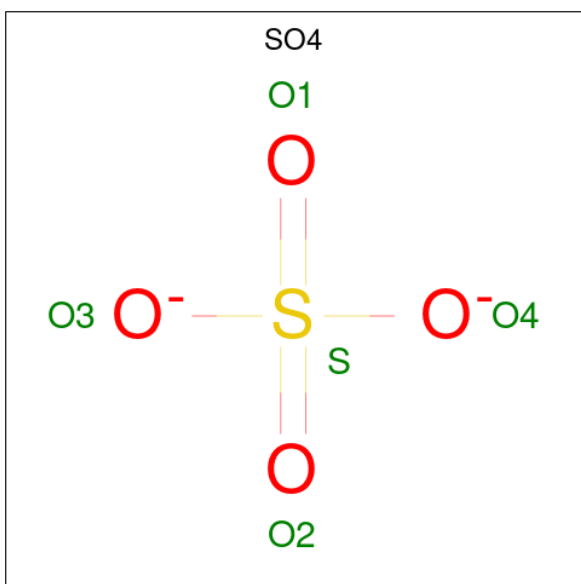
Chain	Residue	Modelled	Actual	Comment	Reference
B	330	GLY	-	expression tag	UNP Q9WYQ6
B	331	HIS	-	expression tag	UNP Q9WYQ6
B	332	HIS	-	expression tag	UNP Q9WYQ6
B	333	HIS	-	expression tag	UNP Q9WYQ6
B	334	HIS	-	expression tag	UNP Q9WYQ6
B	335	HIS	-	expression tag	UNP Q9WYQ6
B	336	HIS	-	expression tag	UNP Q9WYQ6
C	0	MSE	-	expression tag	UNP Q9WYQ6
C	1	SER	-	expression tag	UNP Q9WYQ6
C	329	GLU	-	expression tag	UNP Q9WYQ6
C	330	GLY	-	expression tag	UNP Q9WYQ6
C	331	HIS	-	expression tag	UNP Q9WYQ6
C	332	HIS	-	expression tag	UNP Q9WYQ6
C	333	HIS	-	expression tag	UNP Q9WYQ6
C	334	HIS	-	expression tag	UNP Q9WYQ6
C	335	HIS	-	expression tag	UNP Q9WYQ6
C	336	HIS	-	expression tag	UNP Q9WYQ6
D	0	MSE	-	expression tag	UNP Q9WYQ6
D	1	SER	-	expression tag	UNP Q9WYQ6
D	329	GLU	-	expression tag	UNP Q9WYQ6
D	330	GLY	-	expression tag	UNP Q9WYQ6
D	331	HIS	-	expression tag	UNP Q9WYQ6
D	332	HIS	-	expression tag	UNP Q9WYQ6
D	333	HIS	-	expression tag	UNP Q9WYQ6
D	334	HIS	-	expression tag	UNP Q9WYQ6
D	335	HIS	-	expression tag	UNP Q9WYQ6
D	336	HIS	-	expression tag	UNP Q9WYQ6
E	0	MSE	-	expression tag	UNP Q9WYQ6
E	1	SER	-	expression tag	UNP Q9WYQ6
E	329	GLU	-	expression tag	UNP Q9WYQ6
E	330	GLY	-	expression tag	UNP Q9WYQ6
E	331	HIS	-	expression tag	UNP Q9WYQ6
E	332	HIS	-	expression tag	UNP Q9WYQ6
E	333	HIS	-	expression tag	UNP Q9WYQ6
E	334	HIS	-	expression tag	UNP Q9WYQ6
E	335	HIS	-	expression tag	UNP Q9WYQ6
E	336	HIS	-	expression tag	UNP Q9WYQ6
F	0	MSE	-	expression tag	UNP Q9WYQ6
F	1	SER	-	expression tag	UNP Q9WYQ6
F	329	GLU	-	expression tag	UNP Q9WYQ6
F	330	GLY	-	expression tag	UNP Q9WYQ6
F	331	HIS	-	expression tag	UNP Q9WYQ6

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Chain	Residue	Modelled	Actual	Comment	Reference
F	332	HIS	-	expression tag	UNP Q9WYQ6
F	333	HIS	-	expression tag	UNP Q9WYQ6
F	334	HIS	-	expression tag	UNP Q9WYQ6
F	335	HIS	-	expression tag	UNP Q9WYQ6
F	336	HIS	-	expression tag	UNP Q9WYQ6
G	0	MSE	-	expression tag	UNP Q9WYQ6
G	1	SER	-	expression tag	UNP Q9WYQ6
G	329	GLU	-	expression tag	UNP Q9WYQ6
G	330	GLY	-	expression tag	UNP Q9WYQ6
G	331	HIS	-	expression tag	UNP Q9WYQ6
G	332	HIS	-	expression tag	UNP Q9WYQ6
G	333	HIS	-	expression tag	UNP Q9WYQ6
G	334	HIS	-	expression tag	UNP Q9WYQ6
G	335	HIS	-	expression tag	UNP Q9WYQ6
G	336	HIS	-	expression tag	UNP Q9WYQ6
H	0	MSE	-	expression tag	UNP Q9WYQ6
H	1	SER	-	expression tag	UNP Q9WYQ6
H	329	GLU	-	expression tag	UNP Q9WYQ6
H	330	GLY	-	expression tag	UNP Q9WYQ6
H	331	HIS	-	expression tag	UNP Q9WYQ6
H	332	HIS	-	expression tag	UNP Q9WYQ6
H	333	HIS	-	expression tag	UNP Q9WYQ6
H	334	HIS	-	expression tag	UNP Q9WYQ6
H	335	HIS	-	expression tag	UNP Q9WYQ6
H	336	HIS	-	expression tag	UNP Q9WYQ6

- 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	B	1	Total	O	S	0	0
			5	4	1		
2	F	1	Total	O	S	0	0
			5	4	1		
2	F	1	Total	O	S	0	0
			5	4	1		

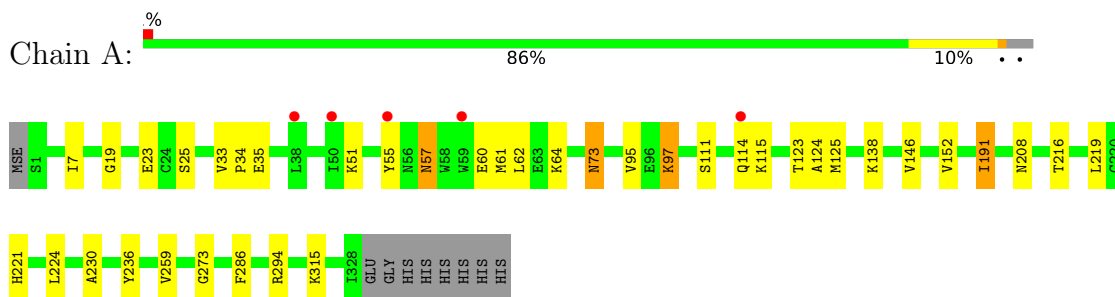
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	59	Total	O	0	0
			59	59		
3	B	67	Total	O	0	0
			67	67		
3	C	18	Total	O	0	0
			18	18		
3	D	30	Total	O	0	0
			30	30		
3	E	19	Total	O	0	0
			19	19		
3	F	59	Total	O	0	0
			59	59		
3	G	8	Total	O	0	0
			8	8		
3	H	4	Total	O	0	0
			4	4		

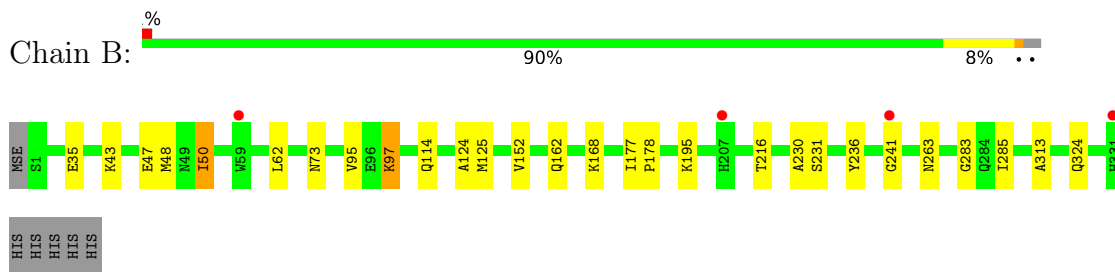
3 Residue-property plots [i](#)

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.

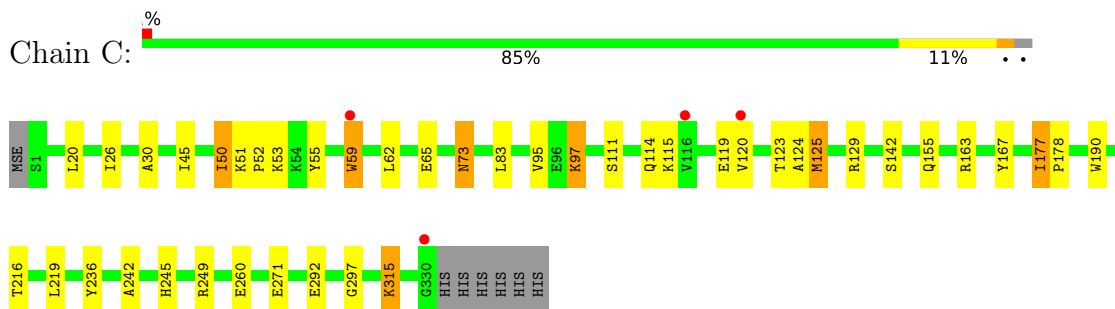
- Molecule 1: Oxidoreductase, putative



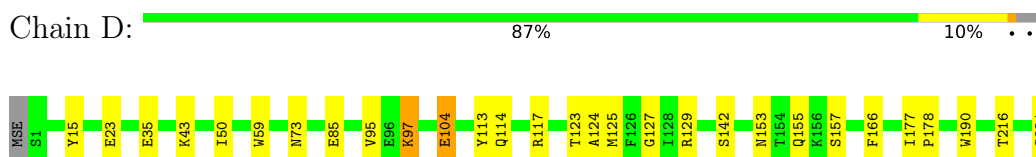
- Molecule 1: Oxidoreductase, putative

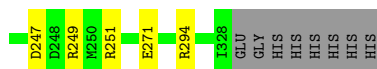


- Molecule 1: Oxidoreductase, putative

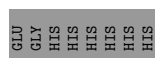
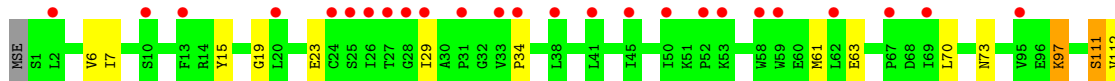
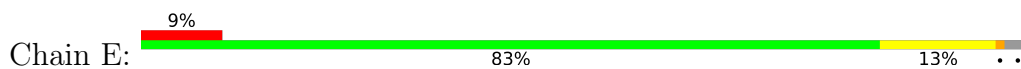


- Molecule 1: Oxidoreductase, putative

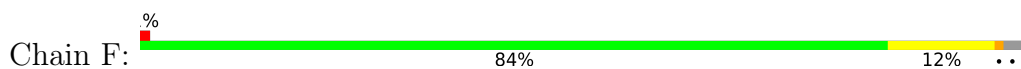




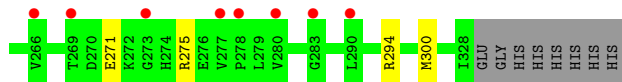
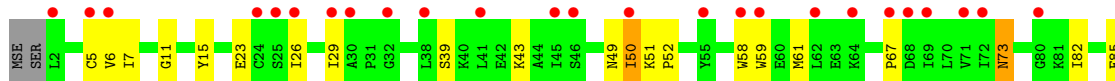
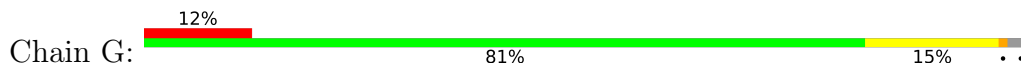
- Molecule 1: Oxidoreductase, putative



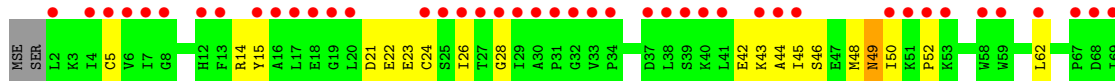
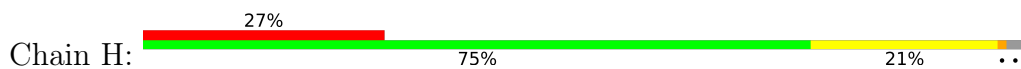
- Molecule 1: Oxidoreductase, putative

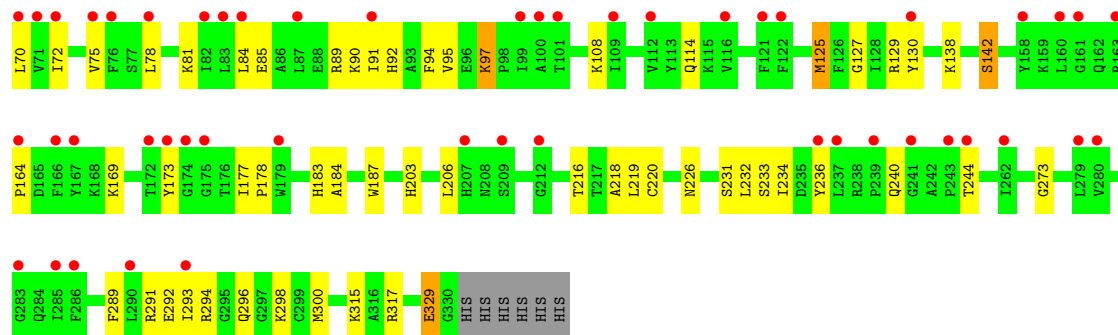


- Molecule 1: Oxidoreductase, putative



- Molecule 1: Oxidoreductase, putative





4 Data and refinement statistics i

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	118.01Å 94.18Å 142.85Å 90.00° 91.91° 90.00°	Depositor
Resolution (Å)	50.00 – 2.14 50.00 – 2.15	Depositor EDS
% Data completeness (in resolution range)	99.5 (50.00-2.14) 99.4 (50.00-2.15)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.17 (at 2.14Å)	Xtrriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.204 , 0.242 0.208 , 0.244	Depositor DCC
R_{free} test set	8541 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	42.2	Xtrriage
Anisotropy	0.026	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 26.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.016 for h,-k,-l	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	21415	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.73% 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: 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	0.80	1/2683 (0.0%)	0.86	0/3615
1	B	0.92	1/2709 (0.0%)	0.88	2/3649 (0.1%)
1	C	0.71	2/2698 (0.1%)	0.88	5/3634 (0.1%)
1	D	0.72	1/2694 (0.0%)	0.86	0/3629
1	E	0.68	1/2685 (0.0%)	0.87	0/3617
1	F	0.79	1/2698 (0.0%)	0.88	0/3634
1	G	0.62	2/2679 (0.1%)	0.84	0/3609
1	H	0.62	1/2692 (0.0%)	0.87	0/3626
All	All	0.74	10/21538 (0.0%)	0.87	7/29013 (0.0%)

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	177	ILE	CA-CB	7.55	1.58	1.54
1	B	125	MSE	SE-CE	-6.68	1.75	1.95
1	D	125	MSE	SE-CE	-5.71	1.78	1.95
1	F	125	MSE	SE-CE	-5.69	1.78	1.95
1	H	125	MSE	SE-CE	-5.57	1.78	1.95
1	G	125	MSE	SE-CE	-5.57	1.78	1.95
1	E	125	MSE	SE-CE	-5.55	1.78	1.95
1	A	125	MSE	SE-CE	-5.55	1.78	1.95
1	C	125	MSE	SE-CE	-5.33	1.79	1.95
1	G	125	MSE	CG-SE	-5.12	1.80	1.95

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	242	ALA	CA-C-N	7.45	127.81	119.32
1	C	242	ALA	C-N-CA	7.45	127.81	119.32
1	C	59	TRP	CA-CB-CG	5.66	124.35	113.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	285	ILE	N-CA-C	5.59	116.33	110.62
1	B	283	GLY	N-CA-C	-5.49	105.23	112.81
1	C	30	ALA	CA-C-N	-5.14	115.43	120.52
1	C	30	ALA	C-N-CA	-5.14	115.43	120.52

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	2632	0	2659	21	0
1	B	2657	0	2682	11	0
1	C	2647	0	2675	21	0
1	D	2640	0	2672	18	0
1	E	2634	0	2666	21	0
1	F	2647	0	2675	19	0
1	G	2628	0	2658	27	0
1	H	2641	0	2667	42	0
2	A	5	0	0	0	0
2	B	10	0	0	0	0
2	F	10	0	0	0	0
3	A	59	0	0	0	0
3	B	67	0	0	1	0
3	C	18	0	0	0	0
3	D	30	0	0	0	0
3	E	19	0	0	0	0
3	F	59	0	0	1	0
3	G	8	0	0	1	0
3	H	4	0	0	0	0
All	All	21415	0	21354	171	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 (171) 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:256:ARG:HH11	1:G:256:ARG:HG3	1.28	0.98
1:H:129:ARG:HD2	1:H:300:MSE:HE3	1.68	0.73
1:F:131:ARG:HD3	3:F:365:HOH:O	1.89	0.72
1:G:256:ARG:HG3	1:G:256:ARG:NH1	2.03	0.71
1:H:291:ARG:HG2	1:H:296:GLN:HB2	1.74	0.70
1:C:111:SER:O	1:C:114:GLN:HG2	1.94	0.68
1:C:45:ILE:HB	1:C:50:ILE:CD1	2.24	0.67
1:E:130:TYR:CZ	1:E:300:MSE:HE2	2.29	0.66
1:E:23:GLU:HB2	1:E:294:ARG:NH2	2.11	0.66
1:A:57:ASN:ND2	1:A:60:GLU:H	1.95	0.65
1:E:113:TYR:O	1:E:117:ARG:HB3	1.97	0.64
1:F:138:LYS:O	1:F:142:SER:HB2	1.98	0.64
1:D:59:TRP:CH2	1:D:85[A]:GLU:HG2	2.34	0.63
1:H:177:ILE:HB	1:H:178:PRO:HD3	1.80	0.63
1:C:45:ILE:HB	1:C:50:ILE:HD13	1.82	0.62
1:H:220:CYS:HB2	1:H:232:LEU:HB3	1.81	0.61
1:G:113:TYR:O	1:G:117:ARG:HB3	2.01	0.61
1:E:130:TYR:CE2	1:E:300:MSE:HE2	2.37	0.60
1:E:7:ILE:HG12	1:E:61:MSE:HE1	1.84	0.59
1:G:241:GLY:O	1:H:273:GLY:HA2	2.04	0.58
1:B:195:LYS:HE2	3:G:338:HOH:O	2.03	0.58
1:A:55:TYR:HE2	1:A:64:LYS:HD2	1.66	0.58
1:H:294:ARG:HB3	1:H:296:GLN:HE22	1.70	0.57
1:E:292:GLU:OE2	1:E:298:LYS:HA	2.05	0.57
1:H:97:LYS:HE3	1:H:97:LYS:O	2.05	0.56
1:E:97:LYS:HD2	1:E:182:ILE:CG2	2.35	0.56
1:H:22:GLU:C	1:H:24:CYS:H	2.14	0.56
1:C:260:GLU:OE2	1:D:249:ARG:NH1	2.29	0.56
1:H:15:TYR:CE1	1:H:127:GLY:HA3	2.41	0.56
1:D:104:GLU:H	1:D:104:GLU:CD	2.12	0.56
1:D:23:GLU:HB2	1:D:294:ARG:NH2	2.22	0.55
1:E:315:LYS:HG2	1:E:328:ILE:HG12	1.88	0.55
1:F:315:LYS:HG3	1:F:326:VAL:HG11	1.90	0.54
1:B:35:GLU:H	1:B:35:GLU:CD	2.16	0.54
1:G:23:GLU:HB2	1:G:294:ARG:NH2	2.23	0.54
1:A:57:ASN:C	1:A:57:ASN:HD22	2.17	0.53
1:D:15:TYR:CZ	1:D:127:GLY:HA3	2.44	0.53
1:G:5:CYS:HB2	1:G:67:PRO:HG3	1.90	0.53
1:H:89:ARG:O	1:H:91:ILE:HG13	2.08	0.53
1:H:240:GLN:HE21	1:H:240:GLN:HA	1.74	0.52
1:H:72:ILE:HB	1:H:95:VAL:HG23	1.92	0.52
1:E:6:VAL:HB	1:E:29:ILE:HG22	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:129:ARG:HD2	1:E:300:MSE:HE3	1.92	0.52
1:H:92:HIS:CD2	1:H:293:ILE:HD12	2.45	0.52
1:H:43:LYS:N	1:H:43:LYS:HD2	2.25	0.51
1:G:39:SER:O	1:G:43:LYS:HD3	2.10	0.51
1:G:15:TYR:CE1	1:G:127:GLY:HA3	2.46	0.51
1:C:55:TYR:OH	1:C:65:GLU:OE2	2.23	0.50
1:D:153:ASN:HD21	1:D:251:ARG:HH11	1.58	0.50
1:F:114:GLN:NE2	1:F:115:LYS:HE3	2.26	0.50
1:H:84:LEU:HD13	1:H:108:LYS:HB3	1.94	0.50
1:G:130:TYR:CZ	1:G:300:MSE:HE2	2.47	0.50
1:C:219:LEU:HD11	1:D:231:SER:HB3	1.94	0.50
1:A:152:VAL:O	1:A:230:ALA:HA	2.12	0.49
1:F:302:THR:HB	1:F:303:PRO:HD2	1.94	0.49
1:H:97:LYS:HD3	1:H:183:HIS:CE1	2.47	0.49
1:B:152:VAL:O	1:B:230:ALA:HA	2.12	0.49
1:E:7:ILE:HG13	1:E:70:LEU:HD11	1.93	0.49
1:F:114:GLN:HE21	1:F:115:LYS:HG3	1.77	0.49
1:F:152:VAL:O	1:F:230:ALA:HA	2.13	0.49
1:D:155:GLN:HB3	1:D:249:ARG:HB2	1.95	0.49
1:G:6:VAL:HB	1:G:29:ILE:HG22	1.95	0.49
1:H:26:ILE:HG22	1:H:28:GLY:H	1.78	0.48
1:H:85:GLU:OE1	1:H:89:ARG:HD2	2.12	0.48
1:H:169:LYS:O	1:H:173:TYR:HB3	2.13	0.48
1:G:7:ILE:HG12	1:G:61:MSE:HE1	1.95	0.48
1:G:152:VAL:O	1:G:230:ALA:HA	2.13	0.48
1:A:57:ASN:HD21	1:A:60:GLU:H	1.62	0.48
1:C:155:GLN:HB3	1:C:249:ARG:HB2	1.96	0.48
1:F:37:ASP:OD1	1:F:39:SER:OG	2.32	0.48
1:E:19:GLY:HA3	1:E:286:PHE:HB3	1.95	0.48
1:F:23:GLU:HB2	1:F:294:ARG:NH2	2.29	0.47
1:A:138:LYS:HA	1:A:191:ILE:CD1	2.45	0.47
1:F:315:LYS:HG2	1:F:328:ILE:HG12	1.96	0.47
1:A:111:SER:O	1:A:114:GLN:HG2	2.15	0.47
1:G:50:ILE:HG23	1:G:52:PRO:HD3	1.97	0.47
1:G:97:LYS:HD2	1:G:182:ILE:CG2	2.45	0.46
1:A:19:GLY:HA3	1:A:286:PHE:HB3	1.97	0.46
1:B:48:MSE:HB2	1:B:50:ILE:HD13	1.96	0.46
1:F:57:ASN:HD22	1:F:60:GLU:CD	2.23	0.46
1:D:113:TYR:O	1:D:117:ARG:HB3	2.16	0.46
1:G:58:TRP:HH2	1:G:85:GLU:HG3	1.81	0.46
1:G:151:LEU:HD21	1:H:233:SER:OG	2.14	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:168:LYS:HE3	3:B:370:HOH:O	2.16	0.45
1:C:95:VAL:O	1:C:124:ALA:HA	2.16	0.45
1:G:59:TRP:CZ3	1:G:89:ARG:NH2	2.84	0.45
1:E:138:LYS:HG2	1:E:190:TRP:CZ2	2.51	0.45
1:G:256:ARG:NH1	1:G:256:ARG:CG	2.75	0.45
1:D:97:LYS:H	1:D:97:LYS:HG3	1.56	0.45
1:G:95:VAL:O	1:G:124:ALA:HA	2.16	0.45
1:D:177:ILE:HB	1:D:178:PRO:HD3	1.99	0.45
1:A:97:LYS:H	1:A:97:LYS:HG3	1.57	0.45
1:D:95:VAL:O	1:D:124:ALA:HA	2.16	0.45
1:E:15:TYR:CZ	1:E:127:GLY:HA3	2.51	0.45
1:H:45:ILE:HG13	1:H:46:SER:N	2.31	0.45
1:A:273:GLY:HA2	1:B:241:GLY:O	2.17	0.44
1:D:153:ASN:ND2	1:D:251:ARG:HH11	2.14	0.44
1:H:138:LYS:O	1:H:142:SER:HB3	2.17	0.44
1:A:138:LYS:HA	1:A:191:ILE:HD11	1.99	0.44
1:C:163:ARG:HD2	1:C:167:TYR:CE2	2.53	0.44
1:D:166:PHE:CD1	1:D:166:PHE:C	2.96	0.44
1:F:114:GLN:HE21	1:F:115:LYS:HE3	1.83	0.44
1:F:114:GLN:HA	1:F:117:ARG:NH1	2.33	0.44
1:H:240:GLN:HA	1:H:240:GLN:NE2	2.32	0.44
1:F:281:GLU:HG3	1:F:282:LYS:N	2.31	0.44
1:A:73:ASN:CG	1:A:73:ASN:O	2.61	0.43
1:C:177:ILE:HB	1:C:178:PRO:HD3	2.00	0.43
1:E:97:LYS:H	1:E:97:LYS:HG3	1.67	0.43
1:F:315:LYS:HG3	1:F:326:VAL:CG1	2.47	0.43
1:G:219:LEU:HD11	1:H:231:SER:HB3	2.00	0.43
1:A:35:GLU:OE2	1:A:35:GLU:N	2.35	0.43
1:B:97:LYS:H	1:B:97:LYS:HG3	1.59	0.43
1:B:178:PRO:HA	1:B:313:ALA:HB1	2.00	0.43
1:E:289:PHE:O	1:E:292:GLU:HB3	2.18	0.43
1:H:218:ALA:HB3	1:H:234:ILE:HB	2.00	0.43
1:C:83:LEU:HD22	1:C:95:VAL:HG11	2.01	0.43
1:F:73:ASN:O	1:F:73:ASN:CG	2.62	0.43
1:A:7:ILE:HG12	1:A:61:MSE:HE1	2.01	0.43
1:E:177:ILE:HB	1:E:178:PRO:HD3	2.01	0.43
1:F:42:GLU:HA	1:F:45:ILE:HG12	2.00	0.43
1:H:94:PHE:CE2	1:H:125:MSE:HG2	2.54	0.43
1:H:130:TYR:CZ	1:H:300:MSE:HE2	2.53	0.43
1:A:219:LEU:HD11	1:B:231:SER:HB3	2.01	0.43
1:B:95:VAL:O	1:B:124:ALA:HA	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:73:ASN:O	1:C:73:ASN:CG	2.62	0.43
1:A:208:ASN:HA	1:A:216:THR:HG21	2.00	0.43
1:G:88:GLU:C	1:G:90:LYS:H	2.26	0.43
1:H:21:ASP:H	1:H:24:CYS:HB2	1.84	0.43
1:C:97:LYS:H	1:C:97:LYS:HG3	1.48	0.43
1:C:111:SER:O	1:C:114:GLN:CG	2.64	0.43
1:H:48:MSE:O	1:H:49:ASN:HB2	2.19	0.42
1:H:78:LEU:HA	1:H:81:LYS:HD3	2.00	0.42
1:C:315:LYS:HA	1:C:315:LYS:HD2	1.81	0.42
1:G:7:ILE:HD12	1:G:82:ILE:HG21	2.01	0.42
1:A:23:GLU:HB2	1:A:294:ARG:NH2	2.34	0.42
1:E:141:VAL:HG11	1:E:149:ILE:HD11	2.02	0.42
1:E:219:LEU:HD11	1:F:231:SER:HB3	2.01	0.42
1:H:5:CYS:O	1:H:70:LEU:HD12	2.19	0.42
1:E:111:SER:O	1:E:114:GLN:HG2	2.20	0.42
1:G:11:GLY:HA3	1:G:73:ASN:HB2	2.02	0.42
1:G:26:ILE:O	1:G:52:PRO:HB3	2.20	0.42
1:C:114:GLN:HE21	1:C:115:LYS:NZ	2.17	0.42
1:H:329:GLU:OE2	1:H:329:GLU:HA	2.20	0.41
1:C:129:ARG:HD3	1:C:190:TRP:CD2	2.55	0.41
1:C:292:GLU:HA	1:C:297:GLY:O	2.20	0.41
1:H:23:GLU:HB2	1:H:294:ARG:NH2	2.35	0.41
1:C:45:ILE:HD12	1:C:52:PRO:HG2	2.03	0.41
1:H:84:LEU:HD22	1:H:108:LYS:HD3	2.02	0.41
1:H:184:ALA:HA	1:H:187:TRP:CD2	2.55	0.41
1:G:231:SER:HB3	1:H:219:LEU:HD21	2.02	0.41
1:H:289:PHE:O	1:H:293:ILE:HG12	2.20	0.41
1:D:104:GLU:CD	1:D:104:GLU:N	2.78	0.41
1:A:221:HIS:CD2	1:A:221:HIS:C	2.98	0.41
1:C:20:LEU:HD11	1:C:26:ILE:HD11	2.02	0.41
1:D:35:GLU:HG2	1:D:35:GLU:O	2.20	0.41
1:D:129:ARG:HD3	1:D:190:TRP:CD2	2.55	0.41
1:D:157:SER:HB3	1:D:247:ASP:HB3	2.03	0.41
1:H:22:GLU:C	1:H:24:CYS:N	2.79	0.41
1:E:152:VAL:O	1:E:230:ALA:HA	2.21	0.41
1:C:73:ASN:HB3	1:C:125:MSE:HE1	2.04	0.40
1:F:20:LEU:HD12	1:F:48:MSE:HE1	2.03	0.40
1:H:178:PRO:HG2	1:H:317:ARG:HD3	2.03	0.40
1:H:184:ALA:HA	1:H:187:TRP:CE3	2.56	0.40
1:A:95:VAL:O	1:A:124:ALA:HA	2.22	0.40
1:H:42:GLU:C	1:H:44:ALA:H	2.29	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:292:GLU:OE2	1:H:298:LYS:HA	2.22	0.40
1:A:146:VAL:CG1	1:A:259:VAL:HG23	2.52	0.40
1:A:57:ASN:ND2	1:A:57:ASN:C	2.78	0.40
1:G:50:ILE:HD13	1:G:51:LYS:N	2.36	0.40
1:B:177:ILE:HB	1:B:178:PRO:HD3	2.02	0.40
1:G:177:ILE:HB	1:G:178:PRO:HD3	2.02	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	326/337 (97%)	313 (96%)	12 (4%)	1 (0%)	36	33
1	B	329/337 (98%)	318 (97%)	11 (3%)	0	100	100
1	C	328/337 (97%)	315 (96%)	13 (4%)	0	100	100
1	D	327/337 (97%)	316 (97%)	11 (3%)	0	100	100
1	E	326/337 (97%)	309 (95%)	16 (5%)	1 (0%)	36	33
1	F	328/337 (97%)	319 (97%)	9 (3%)	0	100	100
1	G	325/337 (96%)	304 (94%)	20 (6%)	1 (0%)	36	33
1	H	327/337 (97%)	297 (91%)	25 (8%)	5 (2%)	8	3
All	All	2616/2696 (97%)	2491 (95%)	117 (4%)	8 (0%)	36	33

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	H	49	ASN
1	E	34	PRO
1	G	49	ASN
1	H	90	LYS

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Mol	Chain	Res	Type
1	H	164	PRO
1	H	52	PRO
1	A	34	PRO
1	H	75	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	286/289 (99%)	273 (96%)	13 (4%)	24	21
1	B	289/289 (100%)	277 (96%)	12 (4%)	26	23
1	C	288/289 (100%)	272 (94%)	16 (6%)	19	14
1	D	288/289 (100%)	277 (96%)	11 (4%)	29	28
1	E	287/289 (99%)	272 (95%)	15 (5%)	21	16
1	F	288/289 (100%)	265 (92%)	23 (8%)	11	7
1	G	286/289 (99%)	273 (96%)	13 (4%)	24	21
1	H	287/289 (99%)	273 (95%)	14 (5%)	22	18
All	All	2299/2312 (99%)	2182 (95%)	117 (5%)	21	17

All (117) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	SER
1	A	33	VAL
1	A	51	LYS
1	A	57	ASN
1	A	62	LEU
1	A	73	ASN
1	A	97	LYS
1	A	115	LYS
1	A	123	THR
1	A	191	ILE
1	A	224	LEU

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Mol	Chain	Res	Type
1	A	236	TYR
1	A	315	LYS
1	B	43	LYS
1	B	47	GLU
1	B	50	ILE
1	B	62	LEU
1	B	73	ASN
1	B	97	LYS
1	B	114	GLN
1	B	162	GLN
1	B	216	THR
1	B	236	TYR
1	B	263	ASN
1	B	324	GLN
1	C	50	ILE
1	C	51	LYS
1	C	53	LYS
1	C	59	TRP
1	C	62	LEU
1	C	73	ASN
1	C	97	LYS
1	C	119	GLU
1	C	120	VAL
1	C	123	THR
1	C	142	SER
1	C	216	THR
1	C	236	TYR
1	C	245	HIS
1	C	271	GLU
1	C	315	LYS
1	D	43	LYS
1	D	50	ILE
1	D	73	ASN
1	D	97	LYS
1	D	104	GLU
1	D	114	GLN
1	D	123	THR
1	D	142	SER
1	D	216	THR
1	D	236	TYR
1	D	271	GLU
1	E	63	GLU

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Mol	Chain	Res	Type
1	E	73	ASN
1	E	97	LYS
1	E	111	SER
1	E	112	VAL
1	E	117	ARG
1	E	119	GLU
1	E	120	VAL
1	E	123	THR
1	E	142	SER
1	E	191	ILE
1	E	236	TYR
1	E	263	ASN
1	E	281	GLU
1	E	315	LYS
1	F	22	GLU
1	F	39	SER
1	F	49	ASN
1	F	50	ILE
1	F	51	LYS
1	F	53	LYS
1	F	62	LEU
1	F	64	LYS
1	F	73	ASN
1	F	97	LYS
1	F	104	GLU
1	F	111	SER
1	F	114	GLN
1	F	120	VAL
1	F	123	THR
1	F	142	SER
1	F	148	GLU
1	F	209	SER
1	F	216	THR
1	F	236	TYR
1	F	271	GLU
1	F	315	LYS
1	F	324	GLN
1	G	50	ILE
1	G	73	ASN
1	G	97	LYS
1	G	104	GLU
1	G	111	SER

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Mol	Chain	Res	Type
1	G	114	GLN
1	G	123	THR
1	G	195	LYS
1	G	216	THR
1	G	236	TYR
1	G	256	ARG
1	G	271	GLU
1	G	275	ARG
1	H	14	ARG
1	H	50	ILE
1	H	62	LEU
1	H	97	LYS
1	H	114	GLN
1	H	142	SER
1	H	203	HIS
1	H	206	LEU
1	H	216	THR
1	H	226	ASN
1	H	236	TYR
1	H	244	THR
1	H	315	LYS
1	H	329	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	57	ASN
1	A	114	GLN
1	A	118	ASN
1	A	240	GLN
1	A	245	HIS
1	B	118	ASN
1	B	189	HIS
1	C	114	GLN
1	C	245	HIS
1	D	118	ASN
1	D	153	ASN
1	F	57	ASN
1	F	114	GLN
1	F	240	GLN
1	F	274	HIS
1	G	240	GLN

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Mol	Chain	Res	Type
1	G	245	HIS
1	H	12	HIS
1	H	49	ASN
1	H	73	ASN
1	H	114	GLN
1	H	226	ASN
1	H	240	GLN
1	H	296	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](#)

5 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	F	337	-	4,4,4	0.24	0	6,6,6	0.07	0
2	SO4	B	338	-	4,4,4	0.24	0	6,6,6	0.12	0
2	SO4	A	337	-	4,4,4	0.26	0	6,6,6	0.12	0
2	SO4	F	338	-	4,4,4	0.24	0	6,6,6	0.12	0
2	SO4	B	337	-	4,4,4	0.24	0	6,6,6	0.14	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	323/337 (95%)	0.03	5 (1%) 72 76	27, 41, 68, 80	0
1	B	326/337 (96%)	-0.14	4 (1%) 76 80	26, 35, 54, 70	0
1	C	325/337 (96%)	0.36	4 (1%) 76 80	35, 50, 65, 74	0
1	D	323/337 (95%)	0.03	1 (0%) 90 92	25, 45, 59, 70	1 (0%)
1	E	323/337 (95%)	0.60	29 (8%) 15 17	32, 53, 89, 96	0
1	F	325/337 (96%)	-0.14	2 (0%) 85 88	28, 38, 53, 66	0
1	G	322/337 (95%)	0.91	41 (12%) 8 9	33, 65, 104, 109	0
1	H	324/337 (96%)	1.49	92 (28%) 1 1	36, 77, 120, 122	0
All	All	2591/2696 (96%)	0.39	178 (6%) 23 26	25, 47, 94, 122	1 (0%)

All (178) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	34	PRO	5.6
1	H	26	ILE	5.0
1	H	33	VAL	5.0
1	H	29	ILE	5.0
1	H	2	LEU	5.0
1	H	28	GLY	4.8
1	H	20	LEU	4.6
1	H	50	ILE	4.5
1	G	50	ILE	4.5
1	H	27	THR	4.3
1	G	6	VAL	4.2
1	H	45	ILE	4.2
1	H	4	ILE	4.1
1	C	59	TRP	4.1
1	H	32	GLY	4.1
1	H	13	PHE	4.0

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Mol	Chain	Res	Type	RSRZ
1	H	69	ILE	4.0
1	H	5	CYS	4.0
1	E	59	TRP	3.8
1	G	38	LEU	3.8
1	H	59	TRP	3.8
1	H	41	LEU	3.8
1	H	38	LEU	3.7
1	H	72	ILE	3.6
1	H	16	ALA	3.6
1	H	15	TYR	3.6
1	H	166	PHE	3.5
1	G	145	ALA	3.5
1	H	83	LEU	3.5
1	G	2	LEU	3.5
1	H	8	GLY	3.5
1	E	120	VAL	3.5
1	H	6	VAL	3.5
1	H	241	GLY	3.4
1	G	24	CYS	3.4
1	H	71	VAL	3.4
1	H	91	ILE	3.4
1	H	70	LEU	3.4
1	H	12	HIS	3.4
1	H	24	CYS	3.3
1	H	58	TRP	3.3
1	E	25	SER	3.2
1	H	84	LEU	3.2
1	B	207	HIS	3.2
1	G	29	ILE	3.2
1	H	99	ILE	3.2
1	G	58	TRP	3.1
1	H	101	THR	3.1
1	H	172	THR	3.1
1	G	59	TRP	3.1
1	H	158	TYR	3.1
1	H	39	SER	3.1
1	H	62	LEU	3.1
1	H	243	PRO	3.0
1	E	50	ILE	3.0
1	E	27	THR	3.0
1	C	120	VAL	3.0
1	E	34	PRO	3.0

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Mol	Chain	Res	Type	RSRZ
1	H	19	GLY	3.0
1	H	161	GLY	3.0
1	C	116	VAL	2.9
1	G	45	ILE	2.9
1	H	109	ILE	2.9
1	E	62	LEU	2.9
1	E	29	ILE	2.9
1	H	44	ALA	2.9
1	E	28	GLY	2.9
1	G	41	LEU	2.8
1	H	121	PHE	2.8
1	H	279	LEU	2.8
1	H	67	PRO	2.8
1	H	52	PRO	2.8
1	G	116	VAL	2.8
1	E	26	ILE	2.8
1	G	26	ILE	2.8
1	E	289	PHE	2.7
1	H	100	ALA	2.7
1	E	2	LEU	2.7
1	B	59	TRP	2.7
1	G	62	LEU	2.7
1	E	67	PRO	2.7
1	H	285	ILE	2.7
1	G	290	LEU	2.7
1	H	130	TYR	2.7
1	H	207	HIS	2.7
1	E	45	ILE	2.7
1	H	31	PRO	2.6
1	G	120	VAL	2.6
1	G	280	VAL	2.6
1	H	51	LYS	2.6
1	A	59	TRP	2.6
1	H	179	TRP	2.6
1	B	331	HIS	2.6
1	H	78	LEU	2.6
1	H	237	LEU	2.6
1	H	293	ILE	2.6
1	G	32	GLY	2.6
1	G	273	GLY	2.6
1	H	174	GLY	2.6
1	H	25	SER	2.6

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Mol	Chain	Res	Type	RSRZ
1	H	280	VAL	2.6
1	F	117	ARG	2.6
1	H	17	LEU	2.6
1	H	160	LEU	2.6
1	H	7	ILE	2.5
1	H	167	TYR	2.5
1	H	283	GLY	2.5
1	E	24	CYS	2.5
1	H	87	LEU	2.5
1	H	122	PHE	2.5
1	H	286	PHE	2.5
1	G	69	ILE	2.5
1	H	82	ILE	2.5
1	H	236	TYR	2.5
1	H	40	LYS	2.5
1	G	283	GLY	2.5
1	G	71	VAL	2.4
1	G	5	CYS	2.4
1	E	52	PRO	2.4
1	E	41	LEU	2.4
1	H	290	LEU	2.4
1	E	69	ILE	2.4
1	H	75	VAL	2.4
1	A	55	TYR	2.4
1	E	38	LEU	2.4
1	G	151	LEU	2.4
1	G	278	PRO	2.4
1	A	38	LEU	2.4
1	A	50	ILE	2.4
1	G	277	VAL	2.4
1	H	262	ILE	2.4
1	H	173	TYR	2.3
1	G	30	ALA	2.3
1	H	239	PRO	2.3
1	G	266	VAL	2.3
1	H	18	GLU	2.3
1	E	33	VAL	2.3
1	H	244	THR	2.3
1	G	68	ASP	2.3
1	H	53	LYS	2.3
1	H	212	GLY	2.3
1	F	1	SER	2.2

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Mol	Chain	Res	Type	RSRZ
1	H	116	VAL	2.2
1	E	13	PHE	2.2
1	H	76	PHE	2.2
1	H	68	ASP	2.2
1	H	30	ALA	2.2
1	E	290	LEU	2.2
1	G	158	TYR	2.2
1	E	293	ILE	2.2
1	D	241	GLY	2.2
1	G	25	SER	2.2
1	G	109	ILE	2.1
1	G	144	GLY	2.1
1	H	163	ARG	2.1
1	H	209	SER	2.1
1	H	164	PRO	2.1
1	H	112	VAL	2.1
1	H	37	ASP	2.1
1	G	46	SER	2.1
1	A	114	GLN	2.1
1	E	31	PRO	2.1
1	G	269	THR	2.1
1	G	72	ILE	2.1
1	E	95	VAL	2.1
1	E	58	TRP	2.1
1	G	245	HIS	2.0
1	B	241	GLY	2.0
1	G	55	TYR	2.0
1	H	175	GLY	2.0
1	E	10	SER	2.0
1	E	20	LEU	2.0
1	E	53	LYS	2.0
1	G	64	LYS	2.0
1	H	43	LYS	2.0
1	G	67	PRO	2.0
1	C	330	GLY	2.0
1	G	80	GLY	2.0

6.2 Non-standard residues in protein, DNA, RNA chains

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(\AA^2)	Q<0.9
2	SO4	F	337	5/5	0.82	0.11	102,102,102,102	0
2	SO4	B	338	5/5	0.83	0.12	108,108,108,109	0
2	SO4	A	337	5/5	0.84	0.10	105,105,106,106	0
2	SO4	B	337	5/5	0.92	0.10	88,88,89,89	0
2	SO4	F	338	5/5	0.92	0.10	96,96,96,96	0

6.5 Other polymers [i](#)

There are no such residues in this entry.