



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:33 PM GMT

PDB ID : 3U02
Title : Crystal Structure of the tRNA modifier TiaS from *Pyrococcus furiosus*, Northeast Structural Genomics Consortium Target PfR225
Authors : Kuzin, A.; Lew, S.; Seetharaman, J.; Sahdev, S.; Xiao, R.; Ciccocanti, C.; Wang, H.; Everett, J.K.; Nair, R.; Acton, T.B.; Rost, B.; Montelione, G.T.; Tong, L.; Hunt, J.F.; Northeast Structural Genomics Consortium (NESG)
Deposited on : 2011-09-28
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

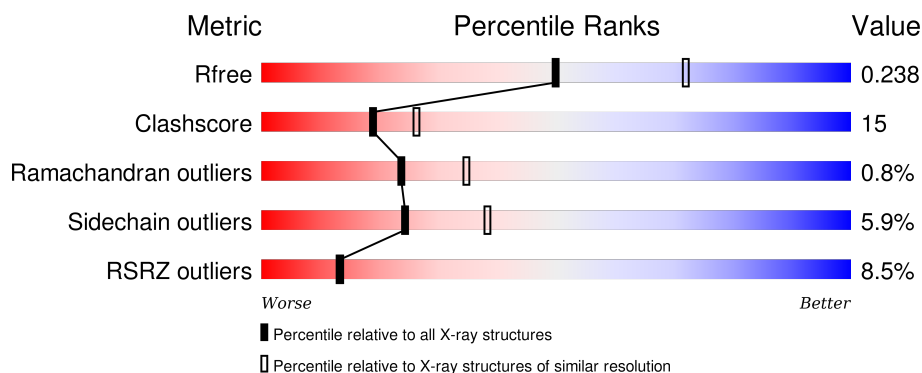
1 Overall quality at a glance ⓘ

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	252	<div> <div>7%</div> <div>69% 22% 5%</div> </div>
1	B	252	<div> <div>7%</div> <div>65% 27% . .</div> </div>
1	C	252	<div> <div>10%</div> <div>60% 26% . 10%</div> </div>
1	D	252	<div> <div>8%</div> <div>63% 25% 12%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	SO4	A	253	-	-	X	-
2	SO4	A	256	-	-	-	X
2	SO4	B	254	-	-	X	-

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 8052 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 Putative transcription-associated protein TFIIS.

Mol	Chain	Residues	Atoms							ZeroOcc	AltConf	Trace
1	A	240	Total	C	N	O	P	S	Se	0	0	0
			1952	1267	324	355	1	1	4			
1	B	244	Total	C	N	O	P	S	Se	0	1	0
			2011	1304	339	362	1	1	4			
1	C	226	Total	C	N	O	P	S	Se	0	1	0
			1857	1207	310	334	1	1	4			
1	D	223	Total	C	N	O	P	S	Se	0	1	0
			1836	1198	303	328	1	1	5			

- Molecule 2 is SULFATE ION (three-letter code: SO₄) (formula: O₄S).



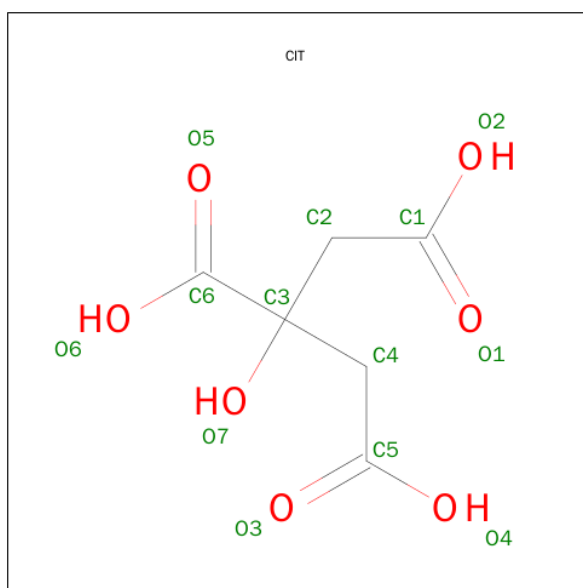
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		

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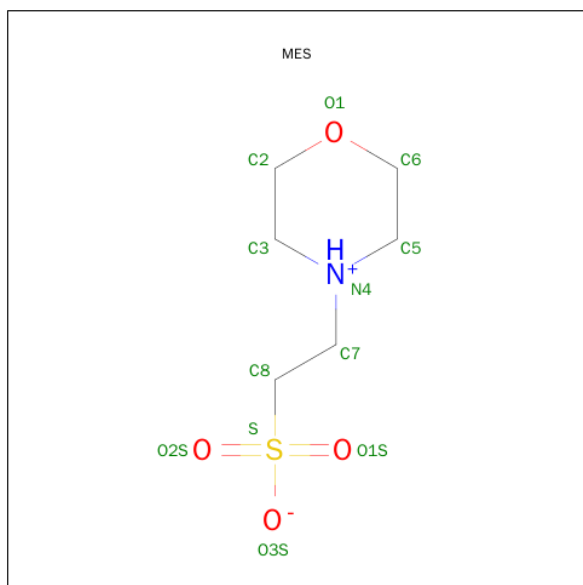
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total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

- Molecule 3 is CITRIC ACID (three-letter code: CIT) (formula: $C_6H_8O_7$).



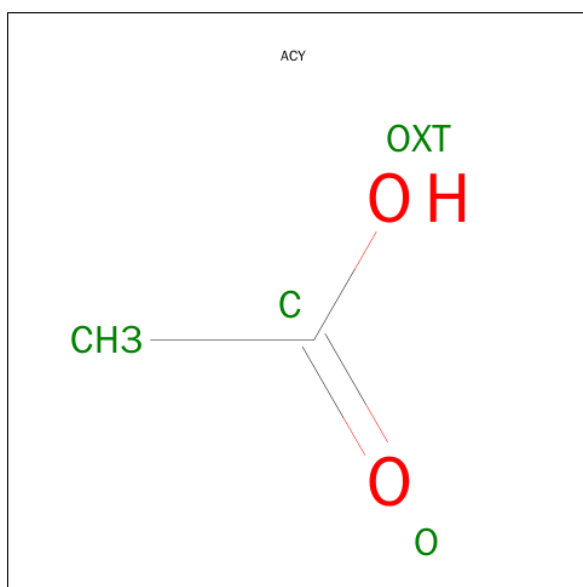
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			13	6	7		
3	B	1	Total	C	O	0	0
			13	6	7		
3	D	1	Total	C	O	0	0
			13	6	7		

- Molecule 4 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES) (formula: $C_6H_{13}NO_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

- Molecule 5 is ACETIC ACID (three-letter code: ACY) (formula: $C_2H_4O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	C	1	Total	C	O	0	0
			4	2	2		

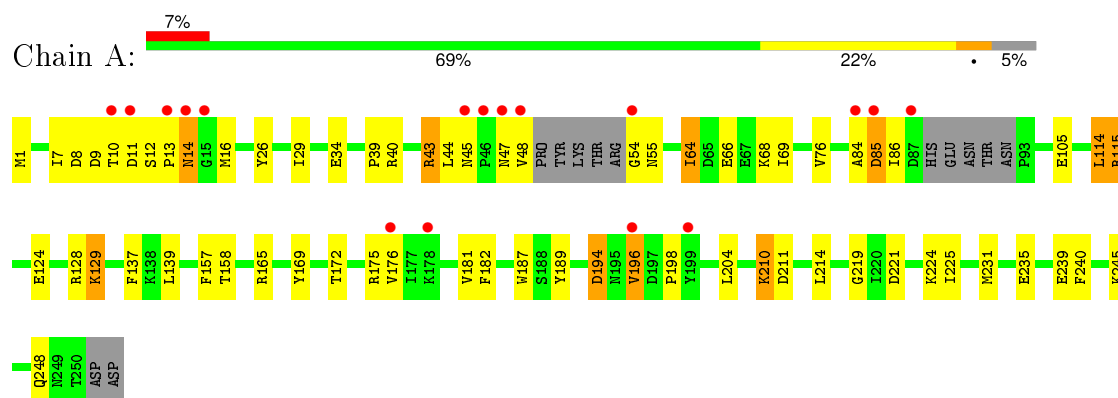
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	90	Total	O	0	0
			90	90		
6	B	81	Total	O	0	0
			81	81		
6	C	54	Total	O	0	0
			54	54		
6	D	45	Total	O	0	1
			46	46		

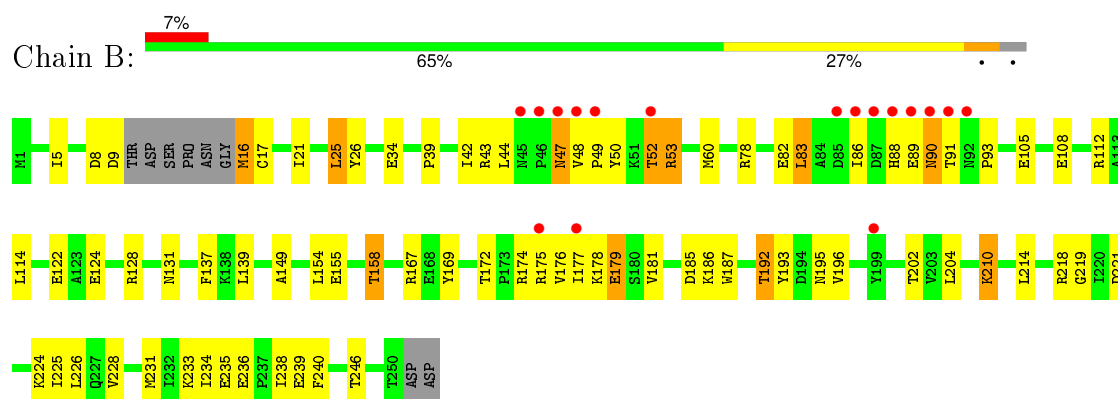
3 Residue-property plots

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.

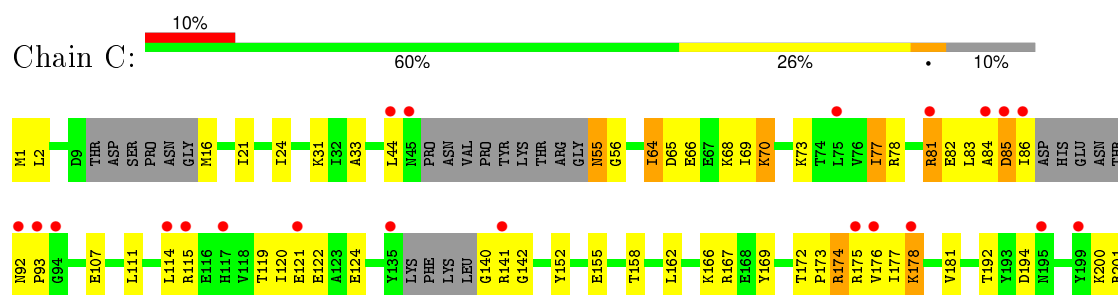
- Molecule 1: Putative transcription-associated protein TFIIS

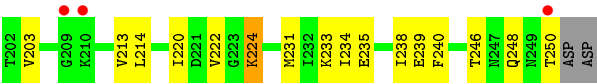


- Molecule 1: Putative transcription-associated protein TFIIS

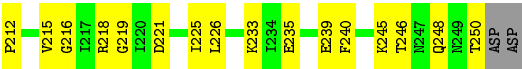
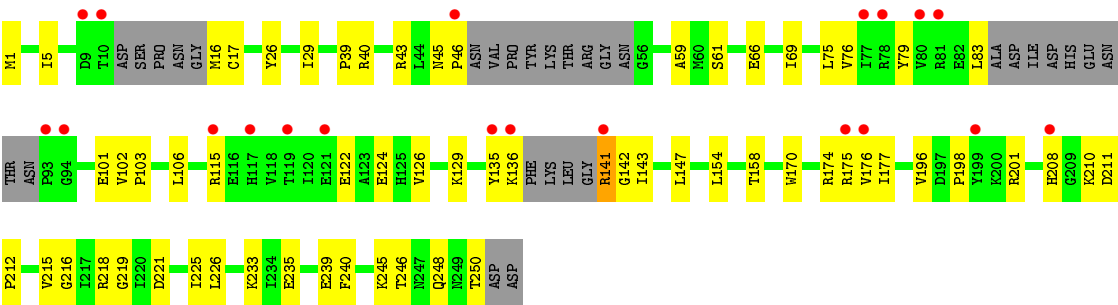


- Molecule 1: Putative transcription-associated protein TFIIS





● Molecule 1: Putative transcription-associated protein TFIIS



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	61.15Å 135.81Å 77.27Å 90.00° 93.76° 90.00°	Depositor
Resolution (Å)	43.90 – 2.40 45.38 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.5 (43.90-2.40) 99.5 (45.38-2.40)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.33 (at 2.39Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.1_743)	Depositor
R, R_{free}	0.179 , 0.242 0.172 , 0.238	Depositor DCC
R_{free} test set	2475 reflections (5.06%)	DCC
Wilson B-factor (Å ²)	33.5	Xtriage
Anisotropy	0.566	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 62.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 48946 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8052	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.80% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: TPO, ACY, SO4, MES, CIT

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	0.42	0/1979	0.56	0/2667
1	B	0.41	0/2042	0.56	0/2755
1	C	0.37	0/1883	0.54	0/2535
1	D	0.40	0/1863	0.55	0/2507
All	All	0.40	0/7767	0.55	0/10464

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1952	0	1988	52	0
1	B	2011	0	2049	72	0
1	C	1857	0	1897	71	0
1	D	1836	0	1885	49	0
2	A	20	0	0	4	0
2	B	25	0	0	3	0
2	C	10	0	0	0	0
2	D	15	0	0	1	0
3	A	13	0	5	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	13	0	5	1	0
3	D	13	0	5	2	0
4	A	12	0	13	3	0
5	C	4	0	3	0	0
6	A	90	0	0	4	0
6	B	81	0	0	4	0
6	C	54	0	0	4	0
6	D	46	0	0	1	0
All	All	8052	0	7850	239	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 15.

All (239) 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:C:201[B]:ARG:H	1:C:201[B]:ARG:HD3	1.26	1.01
1:C:16:MSE:HE3	1:C:167:ARG:HH12	1.27	0.96
1:C:201[B]:ARG:HH21	1:C:224:LYS:HE2	1.32	0.93
1:B:89:GLU:N	1:B:90:ASN:HA	1.87	0.88
1:B:90:ASN:ND2	1:B:90:ASN:O	2.08	0.85
1:B:181:VAL:HG21	1:B:231:MSE:HG2	1.59	0.84
1:B:25:LEU:HB3	1:B:60:MSE:HE1	1.60	0.82
1:C:181:VAL:HG21	1:C:231:MSE:HG2	1.64	0.80
1:B:43:ARG:HE	1:B:50:TYR:HB2	1.47	0.79
1:C:16:MSE:HE1	1:C:86:ILE:HG21	1.65	0.79
1:B:50:TYR:HD1	1:B:50:TYR:O	1.65	0.78
1:B:89:GLU:H	1:B:90:ASN:HA	1.45	0.78
1:C:16:MSE:HE3	1:C:167:ARG:NH1	2.00	0.77
1:A:210:LYS:H	1:A:210:LYS:HD3	1.49	0.76
1:B:8:ASP:HB3	6:B:326:HOH:O	1.88	0.72
1:C:31:LYS:NZ	6:C:308:HOH:O	2.21	0.72
1:A:124:GLU:O	1:A:128:ARG:HG3	1.90	0.72
1:A:169:TYR:O	1:A:172:THR:HB	1.91	0.71
1:A:175:ARG:HB3	1:A:235:GLU:HB2	1.74	0.70
1:A:8:ASP:OD2	1:A:10:THR:HG23	1.91	0.70
1:C:192:THR:HG23	1:C:194:ASP:OD2	1.92	0.69
1:D:141:ARG:CB	1:D:142:GLY:HA3	2.23	0.69
1:B:50:TYR:O	1:B:50:TYR:CD1	2.46	0.67
1:B:52:THR:OG1	1:B:53:ARG:HA	1.95	0.67
1:A:210:LYS:H	1:A:210:LYS:CD	2.08	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:124:GLU:O	1:B:128:ARG:HG3	1.95	0.66
1:C:73:LYS:O	1:C:77:ILE:HG12	1.96	0.66
1:B:48:VAL:CG1	1:B:49:PRO:HD2	2.25	0.66
1:D:1:MSE:HE1	1:D:66:GLU:HG3	1.77	0.65
1:C:16:MSE:CE	1:C:86:ILE:HG21	2.27	0.64
1:B:48:VAL:HG13	1:B:49:PRO:HD2	1.80	0.64
2:A:253:SO4:O1	6:A:344:HOH:O	2.12	0.64
1:D:141:ARG:HB2	1:D:142:GLY:HA3	1.80	0.63
1:D:122:GLU:O	1:D:126:VAL:HG23	1.98	0.63
1:C:92:ASN:HB2	1:C:93:PRO:HD3	1.81	0.63
1:A:10:THR:HG22	1:A:54:GLY:O	1.98	0.63
1:A:181:VAL:HG21	1:A:231:MSE:HG2	1.81	0.63
1:B:178:LYS:HD2	1:B:204:LEU:HD11	1.81	0.62
1:B:169:TYR:O	1:B:172:THR:HB	1.99	0.62
1:B:187:TRP:HD1	2:B:254:SO4:O1	1.82	0.61
1:B:155:GLU:OE2	1:D:198:PRO:HG2	1.99	0.61
1:B:86:ILE:HG21	1:B:167:ARG:HH12	1.66	0.61
1:D:40:ARG:NH2	1:D:61:SER:OG	2.30	0.61
1:B:185:ASP:OD1	1:B:224:LYS:NZ	2.22	0.60
1:A:10:THR:C	1:A:12:SER:H	2.03	0.60
1:C:21:ILE:HG12	1:C:83:LEU:HD23	1.82	0.60
1:C:201[B]:ARG:HG3	6:C:302:HOH:O	2.01	0.60
1:A:221:ASP:O	1:A:225:ILE:HG12	2.01	0.60
1:A:64:ILE:HD11	1:A:68:LYS:HB2	1.83	0.59
1:D:16[B]:MSE:HG3	1:D:17:CYS:H	1.67	0.59
1:C:231:MSE:HE2	6:C:301:HOH:O	2.03	0.59
1:C:111:LEU:O	1:C:115:ARG:HG3	2.02	0.59
1:A:43:ARG:HH12	1:A:48:VAL:HG12	1.68	0.59
1:A:210:LYS:N	1:A:210:LYS:HD3	2.17	0.59
1:C:140:GLY:C	1:C:142:GLY:H	2.06	0.59
1:A:66:GLU:HA	1:A:69:ILE:HD13	1.85	0.58
1:C:222:VAL:HG13	1:D:226:LEU:HD12	1.85	0.58
1:A:10:THR:CG2	1:A:54:GLY:O	2.51	0.58
1:C:158:THR:HG22	1:C:220:ILE:HG22	1.84	0.58
1:A:44:LEU:HA	1:A:114:LEU:HD12	1.84	0.58
1:A:231:MSE:HE2	6:A:300:HOH:O	2.02	0.58
1:C:158:THR:HG22	1:C:220:ILE:CG2	2.33	0.58
1:A:29:ILE:HD11	1:A:76:VAL:HG22	1.85	0.58
1:B:86:ILE:HG21	1:B:167:ARG:NH1	2.19	0.58
1:D:239:GLU:HG3	1:D:240:PHE:CD2	2.39	0.58
1:B:21:ILE:HG12	1:B:83:LEU:HD23	1.86	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:258:MES:H52	1:D:101:GLU:OE1	2.04	0.57
1:B:89:GLU:N	1:B:90:ASN:CA	2.63	0.57
1:C:178:LYS:HD3	1:C:178:LYS:C	2.25	0.57
1:A:224:LYS:HE2	2:A:255:SO4:O1	2.05	0.57
1:B:52:THR:HG21	1:B:53:ARG:NH2	2.20	0.57
1:D:175:ARG:HB3	1:D:235:GLU:HB2	1.88	0.56
1:B:219:GLY:HA3	1:B:225:ILE:HD11	1.85	0.56
1:B:43:ARG:NE	1:B:50:TYR:HB2	2.18	0.56
1:C:70:LYS:NZ	1:C:70:LYS:HB2	2.21	0.56
1:B:202:THR:HG21	1:B:228:VAL:HG21	1.88	0.56
1:B:16:MSE:HE2	1:B:210:LYS:HZ1	1.71	0.55
1:B:5:ILE:HD13	1:B:60:MSE:HE3	1.88	0.55
1:A:137:PHE:CE2	1:A:139:LEU:HA	2.42	0.55
1:C:201[B]:ARG:HH21	1:C:224:LYS:CE	2.11	0.54
1:B:47:ASN:HA	6:B:339:HOH:O	2.08	0.54
1:D:75:LEU:HD23	1:D:79:TYR:HD2	1.73	0.53
1:B:239:GLU:HG3	1:B:240:PHE:HD2	1.72	0.53
1:B:9:ASP:OD2	1:B:53:ARG:HG3	2.08	0.53
1:C:174:ARG:NH1	1:C:213:VAL:O	2.41	0.53
1:B:52:THR:HG23	1:B:53:ARG:HB2	1.90	0.53
1:D:158:THR:HG22	1:D:248:GLN:O	2.09	0.52
1:D:135:TYR:O	1:D:136:LYS:HB2	2.10	0.52
1:D:215:VAL:HG22	1:D:216:GLY:N	2.25	0.52
1:C:201[B]:ARG:NH2	1:C:224:LYS:HE2	2.14	0.52
1:D:141:ARG:HB2	1:D:142:GLY:CA	2.39	0.52
1:D:174:ARG:HG3	1:D:174:ARG:HH11	1.74	0.52
1:B:128:ARG:HD3	6:B:276:HOH:O	2.10	0.52
1:A:1:MSE:N	4:A:258:MES:H82	2.24	0.52
1:C:194:ASP:OD2	1:C:194:ASP:N	2.41	0.51
1:D:43:ARG:NH2	3:D:256:CIT:O1	2.43	0.51
1:D:141:ARG:N	1:D:141:ARG:HD3	2.26	0.51
1:C:1:MSE:SE	1:C:69:ILE:HD11	2.60	0.51
1:A:194:ASP:CG	1:A:194:ASP:O	2.48	0.51
1:C:64:ILE:HG13	1:C:68:LYS:HB2	1.91	0.51
1:C:78:ARG:O	1:C:81:ARG:HG3	2.11	0.51
1:A:64:ILE:HD11	1:A:68:LYS:CB	2.41	0.51
1:C:234:ILE:CD1	1:C:238:ILE:HD11	2.40	0.51
1:B:177:ILE:HG23	1:B:233:LYS:HB3	1.93	0.50
1:B:175:ARG:HB3	1:B:235:GLU:HB2	1.94	0.50
1:B:78[B]:ARG:O	1:B:82:GLU:HG3	2.11	0.50
1:C:33:ALA:HB2	1:C:64:ILE:HD13	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:187:TRP:HD1	2:A:253:SO4:O2	1.95	0.50
1:C:16:MSE:HE1	1:C:86:ILE:CG2	2.40	0.50
1:C:169:TYR:O	1:C:172:THR:HB	2.12	0.49
1:B:175:ARG:HD3	1:B:236:GLU:HB3	1.95	0.49
1:B:112:ARG:HH12	1:B:122:GLU:HG2	1.77	0.49
1:C:119:THR:OG1	1:C:122:GLU:HG3	2.11	0.49
1:C:84:ALA:O	1:C:86:ILE:N	2.38	0.49
1:A:9:ASP:OD2	1:A:12:SER:HB2	2.12	0.49
1:C:140:GLY:O	1:C:142:GLY:N	2.46	0.49
1:C:119:THR:HG1	1:C:122:GLU:HG3	1.78	0.49
1:D:170:TRP:CE2	1:D:212:PRO:HB3	2.47	0.49
1:B:89:GLU:O	1:B:89:GLU:HG3	2.12	0.49
1:C:120:ILE:O	1:C:124:GLU:HG2	2.13	0.49
1:B:9:ASP:C	1:B:93:PRO:HB3	2.32	0.49
1:D:106:LEU:HD11	1:D:147:LEU:HD22	1.95	0.49
1:B:158:THR:HG21	1:B:218:ARG:HD3	1.93	0.49
1:C:44:LEU:H	1:C:55:ASN:HB2	1.78	0.49
1:A:84:ALA:C	1:A:86:ILE:H	2.15	0.48
1:A:196:VAL:HG13	1:C:107:GLU:OE2	2.13	0.48
1:B:16:MSE:HB3	1:B:17:CYS:H	1.48	0.48
1:D:158:THR:OG1	1:D:219:GLY:O	2.30	0.48
1:A:47:ASN:O	1:A:48:VAL:HG22	2.13	0.48
1:C:16:MSE:SE	1:C:86:ILE:HG21	2.63	0.48
1:C:231:MSE:HE3	6:C:256:HOH:O	2.13	0.48
1:D:5:ILE:O	1:D:59:ALA:HA	2.13	0.48
1:A:13:PRO:O	1:A:14:ASN:HB2	2.13	0.48
1:D:246:THR:OG1	1:D:248:GLN:HG3	2.13	0.48
1:A:34:GLU:OE2	1:B:193:TYR:O	2.32	0.48
1:A:45:ASN:H	1:A:55:ASN:HD22	1.61	0.48
1:B:239:GLU:HG3	1:B:240:PHE:CD2	2.48	0.48
1:B:221:ASP:O	1:B:225:ILE:HG12	2.13	0.48
1:D:221:ASP:O	1:D:225:ILE:HG12	2.14	0.48
1:C:65:ASP:OD2	1:C:68:LYS:HG3	2.14	0.48
1:D:201:ARG:O	1:D:250:THR:HG21	2.13	0.48
1:D:239:GLU:HG3	1:D:240:PHE:HD2	1.79	0.47
1:D:1:MSE:CE	1:D:66:GLU:HG3	2.43	0.47
1:C:64:ILE:HD11	1:C:68:LYS:HD2	1.97	0.47
1:C:177:ILE:HB	1:C:233:LYS:HB3	1.97	0.47
1:B:25:LEU:CB	1:B:60:MSE:HE1	2.37	0.47
1:C:33:ALA:CB	1:C:64:ILE:HD13	2.45	0.47
1:D:40:ARG:HA	1:D:245:LYS:O	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:245:LYS:NZ	6:D:274:HOH:O	2.47	0.47
1:B:192:THR:O	1:B:195:ASN:ND2	2.46	0.47
1:B:50:TYR:CD1	1:B:50:TYR:C	2.88	0.46
1:B:174:ARG:HG3	1:B:174:ARG:HH11	1.80	0.46
1:C:81:ARG:HG3	1:C:82:GLU:N	2.31	0.46
1:D:177:ILE:HB	1:D:233:LYS:HB3	1.96	0.46
1:A:224:LYS:NZ	6:A:314:HOH:O	2.49	0.46
1:A:26:TYR:CE1	1:A:39:PRO:HD3	2.51	0.46
1:B:158:THR:HG22	1:B:246:THR:CG2	2.45	0.46
1:B:44:LEU:HA	1:B:114:LEU:HD22	1.98	0.46
1:B:26:TYR:CE1	1:B:39:PRO:HD3	2.51	0.46
1:C:174:ARG:HG2	1:C:214:LEU:HD12	1.97	0.46
1:A:86:ILE:HG22	1:A:86:ILE:O	2.15	0.46
1:C:172:THR:HG23	1:C:173:PRO:HD2	1.96	0.46
1:D:26:TYR:CE1	1:D:39:PRO:HD3	2.51	0.46
1:A:10:THR:C	1:A:12:SER:N	2.69	0.45
1:C:234:ILE:HD13	1:C:238:ILE:HD11	1.99	0.45
1:D:208:HIS:ND1	2:D:253:SO4:O4	2.27	0.45
1:A:157:PHE:CE1	1:A:245:LYS:HD2	2.52	0.45
1:A:45:ASN:O	1:A:55:ASN:ND2	2.49	0.45
1:C:92:ASN:HB2	1:C:93:PRO:CD	2.46	0.45
1:C:166:LYS:HD2	1:C:239:GLU:OE1	2.17	0.45
1:C:246:THR:HG23	1:C:248:GLN:H	1.80	0.45
1:D:1:MSE:SE	1:D:69:ILE:HD11	2.67	0.45
1:B:224:LYS:HE2	2:B:257:SO4:O2	2.17	0.45
1:A:189:TYR:HD1	2:A:256:SO4:O1	1.99	0.45
1:C:82:GLU:HA	1:C:85:ASP:HB3	1.99	0.45
1:C:64:ILE:HA	1:C:64:ILE:HD12	1.71	0.45
1:A:16:MSE:SE	1:A:211:ASP:HA	2.67	0.45
1:D:158:THR:HG21	1:D:218:ARG:HD2	1.98	0.44
1:A:194:ASP:OD1	1:A:194:ASP:O	2.36	0.44
1:B:179:GLU:H	1:B:179:GLU:CD	2.20	0.44
1:D:29:ILE:HD11	1:D:76:VAL:HG22	1.98	0.44
1:C:119:THR:HB	1:C:121:GLU:OE1	2.18	0.44
1:A:239:GLU:HG3	1:A:240:PHE:CD2	2.52	0.44
1:B:42:ILE:HG13	1:B:149:ALA:HA	1.99	0.44
1:D:106:LEU:CD1	1:D:147:LEU:HD22	2.46	0.44
1:D:124:GLU:H	1:D:124:GLU:HG2	1.70	0.44
1:C:44:LEU:H	1:C:55:ASN:CB	2.30	0.44
1:C:82:GLU:C	1:C:84:ALA:H	2.21	0.44
1:B:53:ARG:HB2	1:B:53:ARG:HE	1.48	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:77:ILE:H	1:C:77:ILE:HG12	1.56	0.43
1:A:1:MSE:H1	4:A:258:MES:H82	1.83	0.43
1:C:70:LYS:HB2	1:C:70:LYS:HZ2	1.83	0.43
3:B:258:CIT:C6	3:B:258:CIT:O2	2.66	0.43
1:B:174:ARG:HG3	1:B:174:ARG:NH1	2.33	0.43
1:B:231:MSE:HE3	6:B:261:HOH:O	2.18	0.43
1:B:43:ARG:HH21	1:B:50:TYR:H	1.66	0.43
1:D:102:VAL:HA	1:D:103:PRO:HD3	1.86	0.43
1:A:198:PRO:HG2	1:C:155:GLU:OE1	2.19	0.43
1:A:105:GLU:OE1	1:A:129:LYS:HD2	2.19	0.43
1:B:25:LEU:HD12	1:B:25:LEU:HA	1.81	0.43
1:C:44:LEU:HD12	1:C:44:LEU:N	2.34	0.43
1:A:115:ARG:NH1	6:A:307:HOH:O	2.52	0.43
1:A:158:THR:OG1	1:A:219:GLY:O	2.30	0.42
3:D:256:CIT:O7	3:D:256:CIT:O4	2.33	0.42
1:B:108:GLU:OE1	1:D:196:VAL:HG11	2.19	0.42
1:C:24:ILE:HG13	1:C:83:LEU:HD21	2.01	0.42
1:A:40:ARG:HA	1:A:245:LYS:O	2.19	0.42
1:A:165:ARG:HG2	1:A:214:LEU:HB2	2.00	0.42
1:B:16:MSE:HE2	1:B:210:LYS:NZ	2.34	0.42
1:D:79:TYR:O	1:D:83:LEU:HD12	2.20	0.42
1:B:219:GLY:CA	1:B:225:ILE:HD11	2.49	0.42
1:A:158:THR:HG22	1:A:248:GLN:O	2.20	0.42
1:D:129:LYS:HB3	1:D:129:LYS:HE2	1.76	0.42
1:A:196:VAL:HG13	1:C:107:GLU:CD	2.39	0.42
1:B:5:ILE:CD1	1:B:60:MSE:HE3	2.50	0.42
1:C:200:LYS:HD3	1:C:203:VAL:HG12	2.02	0.42
1:B:16:MSE:CE	1:B:210:LYS:HG3	2.50	0.41
1:D:45:ASN:HA	1:D:46:PRO:HD3	1.93	0.41
1:D:210:LYS:O	1:D:211:ASP:C	2.59	0.41
1:A:43:ARG:NH1	1:A:48:VAL:HG12	2.33	0.41
1:C:175:ARG:NH1	1:C:235:GLU:O	2.53	0.41
1:D:141:ARG:N	1:D:141:ARG:CD	2.83	0.41
1:C:66:GLU:O	1:C:69:ILE:HG12	2.20	0.41
1:D:174:ARG:HG3	1:D:174:ARG:NH1	2.34	0.41
1:C:114:LEU:HD22	1:C:152:TYR:HE1	1.85	0.41
1:B:187:TRP:CD1	2:B:254:SO4:O1	2.69	0.41
1:D:215:VAL:HG22	1:D:216:GLY:H	1.86	0.41
1:C:81:ARG:O	1:C:85:ASP:N	2.43	0.40
1:B:114:LEU:HD23	1:B:114:LEU:N	2.37	0.40
1:C:239:GLU:HG3	1:C:240:PHE:HD2	1.86	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:226:LEU:HD12	1:B:226:LEU:HA	1.92	0.40
1:D:115:ARG:HB3	1:D:115:ARG:HE	1.61	0.40
1:B:52:THR:CB	1:B:53:ARG:HA	2.51	0.40
1:B:234:ILE:HG22	1:B:236:GLU:O	2.21	0.40
1:C:140:GLY:C	1:C:142:GLY:N	2.72	0.40
1:D:75:LEU:C	1:D:75:LEU:HD23	2.42	0.40
1:B:137:PHE:CE2	1:B:139:LEU:HA	2.57	0.40
1:A:182:PHE:CD1	1:B:34:GLU:HG2	2.56	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	233/252 (92%)	219 (94%)	11 (5%)	3 (1%)	15	21
1	B	240/252 (95%)	230 (96%)	9 (4%)	1 (0%)	39	56
1	C	216/252 (86%)	208 (96%)	5 (2%)	3 (1%)	14	19
1	D	212/252 (84%)	205 (97%)	7 (3%)	0	100	100
All	All	901/1008 (89%)	862 (96%)	32 (4%)	7 (1%)	24	35

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	56	GLY
1	C	141	ARG
1	C	85	ASP
1	A	14	ASN
1	B	88	HIS
1	A	11	ASP
1	A	85	ASP

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	209/217 (96%)	197 (94%)	12 (6%)	25	40
1	B	215/217 (99%)	195 (91%)	20 (9%)	11	16
1	C	198/217 (91%)	186 (94%)	12 (6%)	23	36
1	D	197/217 (91%)	193 (98%)	4 (2%)	63	81
All	All	819/868 (94%)	771 (94%)	48 (6%)	24	38

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	ILE
1	A	43	ARG
1	A	64	ILE
1	A	85	ASP
1	A	114	LEU
1	A	115	ARG
1	A	129	LYS
1	A	176	VAL
1	A	194	ASP
1	A	196	VAL
1	A	204	LEU
1	A	210	LYS
1	B	16	MSE
1	B	25	LEU
1	B	47	ASN
1	B	52	THR
1	B	53	ARG
1	B	83	LEU
1	B	90	ASN
1	B	91	THR
1	B	105	GLU
1	B	131	ASN
1	B	154	LEU
1	B	158	THR

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Mol	Chain	Res	Type
1	B	176	VAL
1	B	179	GLU
1	B	186	LYS
1	B	192	THR
1	B	196	VAL
1	B	210	LYS
1	B	214	LEU
1	B	238	ILE
1	C	2	LEU
1	C	55	ASN
1	C	64	ILE
1	C	70	LYS
1	C	77	ILE
1	C	81	ARG
1	C	162	LEU
1	C	174	ARG
1	C	176	VAL
1	C	178	LYS
1	C	224	LYS
1	C	250	THR
1	D	141	ARG
1	D	143	ILE
1	D	154	LEU
1	D	176	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected

value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
1	TPO	A	18	1	8,10,11	1.16	0	7,14,16	1.27	1 (14%)
1	TPO	B	18	1	8,10,11	1.13	1 (12%)	7,14,16	1.32	0
1	TPO	C	18	1	8,10,11	1.08	0	7,14,16	1.51	2 (28%)
1	TPO	D	18	1	8,10,11	1.14	1 (12%)	7,14,16	1.46	1 (14%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	TPO	A	18	1	-	0/8/11/13	0/0/0/0
1	TPO	B	18	1	-	0/8/11/13	0/0/0/0
1	TPO	C	18	1	-	0/8/11/13	0/0/0/0
1	TPO	D	18	1	-	0/8/11/13	0/0/0/0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	18	TPO	P-O3P	2.00	1.61	1.54
1	B	18	TPO	P-O3P	2.11	1.62	1.54

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	18	TPO	CG2-CB-CA	-2.29	108.52	113.17
1	C	18	TPO	CG2-CB-CA	-2.04	109.02	113.17
1	A	18	TPO	C-CA-N	2.09	114.19	109.83
1	C	18	TPO	C-CA-N	2.49	115.03	109.83

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

19 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	253	-	4,4,4	0.08	0	6,6,6	0.10	0
2	SO4	A	254	-	4,4,4	0.17	0	6,6,6	0.28	0
2	SO4	A	255	-	4,4,4	0.20	0	6,6,6	0.39	0
2	SO4	A	256	-	4,4,4	0.19	0	6,6,6	0.16	0
3	CIT	A	257	-	3,12,12	1.41	0	3,17,17	1.71	1 (33%)
4	MES	A	258	-	11,12,12	1.55	2 (18%)	14,16,16	1.52	3 (21%)
2	SO4	B	253	-	4,4,4	0.17	0	6,6,6	0.06	0
2	SO4	B	254	-	4,4,4	0.10	0	6,6,6	0.14	0
2	SO4	B	255	-	4,4,4	0.16	0	6,6,6	0.10	0
2	SO4	B	256	-	4,4,4	0.21	0	6,6,6	0.19	0
2	SO4	B	257	-	4,4,4	0.19	0	6,6,6	0.10	0
3	CIT	B	258	-	3,12,12	1.36	0	3,17,17	2.52	2 (66%)
2	SO4	C	253	-	4,4,4	0.23	0	6,6,6	0.13	0
2	SO4	C	254	-	4,4,4	0.16	0	6,6,6	0.10	0
5	ACY	C	255	-	1,3,3	0.91	0	0,3,3	0.00	-
2	SO4	D	253	-	4,4,4	0.22	0	6,6,6	0.14	0
2	SO4	D	254	-	4,4,4	0.08	0	6,6,6	0.39	0
2	SO4	D	255	-	4,4,4	0.13	0	6,6,6	0.15	0
3	CIT	D	256	-	3,12,12	1.33	0	3,17,17	3.18	2 (66%)

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
2	SO4	A	253	-	-	0/0/0/0	0/0/0/0
2	SO4	A	254	-	-	0/0/0/0	0/0/0/0
2	SO4	A	255	-	-	0/0/0/0	0/0/0/0
2	SO4	A	256	-	-	0/0/0/0	0/0/0/0
3	CIT	A	257	-	-	0/6/16/16	0/0/0/0
4	MES	A	258	-	-	0/6/14/14	0/1/1/1
2	SO4	B	253	-	-	0/0/0/0	0/0/0/0
2	SO4	B	254	-	-	0/0/0/0	0/0/0/0
2	SO4	B	255	-	-	0/0/0/0	0/0/0/0
2	SO4	B	256	-	-	0/0/0/0	0/0/0/0
2	SO4	B	257	-	-	0/0/0/0	0/0/0/0
3	CIT	B	258	-	-	0/6/16/16	0/0/0/0
2	SO4	C	253	-	-	0/0/0/0	0/0/0/0
2	SO4	C	254	-	-	0/0/0/0	0/0/0/0
5	ACY	C	255	-	-	0/0/0/0	0/0/0/0
2	SO4	D	253	-	-	0/0/0/0	0/0/0/0
2	SO4	D	254	-	-	0/0/0/0	0/0/0/0
2	SO4	D	255	-	-	0/0/0/0	0/0/0/0
3	CIT	D	256	-	-	0/6/16/16	0/0/0/0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	258	MES	C5-N4	-3.25	1.38	1.46
4	A	258	MES	O3S-S	2.65	1.53	1.46

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	256	CIT	C3-C2-C1	-4.60	107.60	114.96
3	B	258	CIT	C3-C4-C5	-3.39	109.54	114.96
3	D	256	CIT	C3-C4-C5	-3.02	110.13	114.96
3	A	257	CIT	C3-C2-C1	-2.77	110.53	114.96
3	B	258	CIT	C3-C2-C1	-2.54	110.90	114.96
4	A	258	MES	C5-N4-C3	2.38	114.04	108.90
4	A	258	MES	C2-C3-N4	2.55	113.99	110.12
4	A	258	MES	O1S-S-C8	2.56	109.09	106.91

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

9 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	253	SO4	2	0
2	A	255	SO4	1	0
2	A	256	SO4	1	0
4	A	258	MES	3	0
2	B	254	SO4	2	0
2	B	257	SO4	1	0
3	B	258	CIT	1	0
2	D	253	SO4	1	0
3	D	256	CIT	2	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 ⓘ

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	235/252 (93%)	0.16	17 (7%)	18 18	17, 30, 82, 105	0
1	B	239/252 (94%)	0.39	17 (7%)	19 19	17, 32, 89, 123	0
1	C	221/252 (87%)	0.38	24 (10%)	7 7	19, 36, 79, 113	0
1	D	218/252 (86%)	0.28	20 (9%)	11 11	18, 36, 75, 105	0
All	All	913/1008 (90%)	0.30	78 (8%)	13 13	17, 34, 84, 123	0

All (78) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	48	VAL	6.1
1	A	13	PRO	6.0
1	A	48	VAL	5.8
1	D	93	PRO	5.1
1	B	90	ASN	5.1
1	B	88	HIS	5.1
1	A	84	ALA	5.0
1	A	14	ASN	4.8
1	D	199	TYR	4.8
1	A	178	LYS	4.8
1	C	84	ALA	4.6
1	D	10	THR	4.6
1	A	47	ASN	4.5
1	B	49	PRO	4.5
1	C	81	ARG	4.4
1	D	117	HIS	4.1
1	C	93	PRO	3.9
1	C	75	LEU	3.9
1	B	91	THR	3.8
1	D	175	ARG	3.7
1	C	114	LEU	3.6

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Mol	Chain	Res	Type	RSRZ
1	D	81	ARG	3.6
1	B	199	TYR	3.6
1	A	199	TYR	3.5
1	D	135	TYR	3.4
1	C	44	LEU	3.3
1	D	115	ARG	3.2
1	A	11	ASP	3.2
1	C	92	ASN	3.2
1	A	15	GLY	3.2
1	C	175	ARG	3.2
1	B	46	PRO	3.1
1	C	85	ASP	3.1
1	B	175	ARG	3.0
1	C	176	VAL	3.0
1	C	117	HIS	3.0
1	C	178	LYS	3.0
1	A	176	VAL	2.9
1	B	92	ASN	2.9
1	A	87	ASP	2.9
1	B	85	ASP	2.7
1	C	115	ARG	2.7
1	B	47	ASN	2.7
1	C	94	GLY	2.7
1	C	86	ILE	2.7
1	D	46	PRO	2.7
1	B	177	ILE	2.7
1	D	141	ARG	2.7
1	B	52	THR	2.7
1	B	86	ILE	2.7
1	A	196	VAL	2.6
1	C	209	GLY	2.6
1	C	210	LYS	2.5
1	D	77	ILE	2.5
1	B	45	ASN	2.4
1	C	45	ASN	2.4
1	A	46	PRO	2.4
1	D	136	LYS	2.4
1	C	199	TYR	2.4
1	C	195	ASN	2.3
1	A	45	ASN	2.3
1	D	94	GLY	2.2
1	A	54	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	10	THR	2.2
1	B	89	GLU	2.2
1	D	176	VAL	2.2
1	D	78	ARG	2.1
1	C	135	TYR	2.1
1	D	9	ASP	2.1
1	D	121	GLU	2.1
1	A	85	ASP	2.1
1	C	121	GLU	2.1
1	C	250	THR	2.1
1	D	208	HIS	2.0
1	D	80	VAL	2.0
1	D	119	THR	2.0
1	B	87	ASP	2.0
1	C	141	ARG	2.0

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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
1	TPO	B	18	11/12	0.89	0.14	-	23,29,93,93	0
1	TPO	A	18	11/12	0.94	0.13	-	23,31,74,93	0
1	TPO	D	18	11/12	0.93	0.15	-	30,36,137,138	0
1	TPO	C	18	11/12	0.93	0.15	-	31,35,92,143	0

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(\AA^2)	Q<0.9
2	SO4	A	256	5/5	0.86	0.25	3.56	71,82,88,98	0
3	CIT	A	257	13/13	0.93	0.16	1.88	30,40,52,53	0
3	CIT	D	256	13/13	0.90	0.20	1.48	50,61,68,75	0
2	SO4	B	255	5/5	0.92	0.18	1.46	78,86,87,90	0
2	SO4	A	254	5/5	0.96	0.17	0.59	45,54,62,62	0
2	SO4	C	253	5/5	0.95	0.21	0.56	75,79,88,88	0
4	MES	A	258	12/12	0.93	0.15	0.42	26,51,68,70	0
3	CIT	B	258	13/13	0.93	0.14	0.11	35,41,55,59	0
2	SO4	B	256	5/5	0.98	0.14	0.09	58,64,65,66	0
2	SO4	B	257	5/5	0.96	0.17	-0.13	53,64,68,82	0
2	SO4	D	254	5/5	0.96	0.15	-0.54	49,53,59,64	0
2	SO4	A	255	5/5	0.99	0.12	-0.56	31,41,45,46	0
2	SO4	D	255	5/5	0.94	0.13	-0.60	63,66,72,80	0
2	SO4	D	253	5/5	0.97	0.14	-0.75	59,67,70,71	0
2	SO4	C	254	5/5	0.88	0.18	-1.12	127,128,131,132	0
2	SO4	A	253	5/5	0.99	0.09	-1.83	30,32,40,41	0
2	SO4	B	254	5/5	0.98	0.11	-1.99	46,56,59,60	0
2	SO4	B	253	5/5	0.98	0.10	-2.10	38,56,64,64	0
5	ACY	C	255	4/4	0.92	0.35	-	32,38,39,47	0

6.5 Other polymers

There are no such residues in this entry.