



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 02:20 PM UTC

PDB ID : 2FK6 / pdb_00002fk6
Title : Crystal Structure of RNase Z/tRNA(Thr) complex
Authors : Li de la Sierra-Gallay, I.; Mathy, N.; Pellegrini, O.; Condon, C.
Deposited on : 2006-01-04
Resolution : 2.90 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

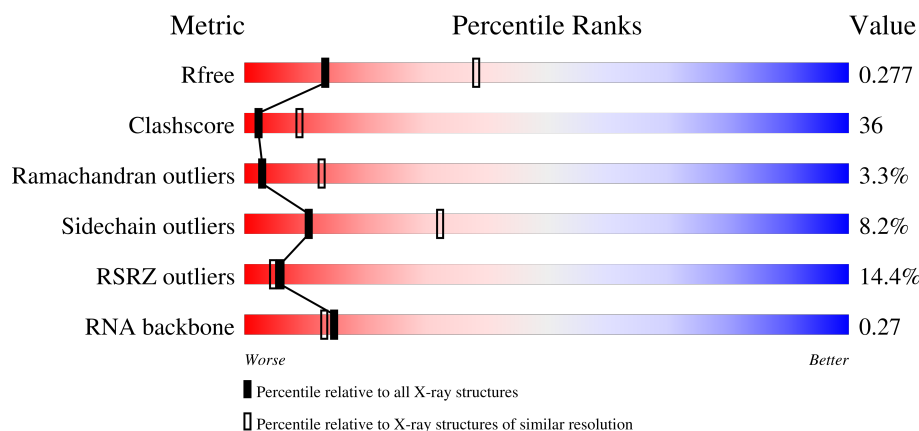
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)
RNA backbone	3983	1120 (3.10-2.70)

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	R	79	
2	A	320	

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
5	GOL	A	501	-	X	-	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 3550 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 RNA chain called TRNA(THR).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	R	53	Total	C	N	O	P	0	0	0
			1125	504	202	368	51			

- Molecule 2 is a protein called RIBONUCLEASE Z.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	307	Total	C	N	O	S	0	0	0
			2395	1524	408	456	7			

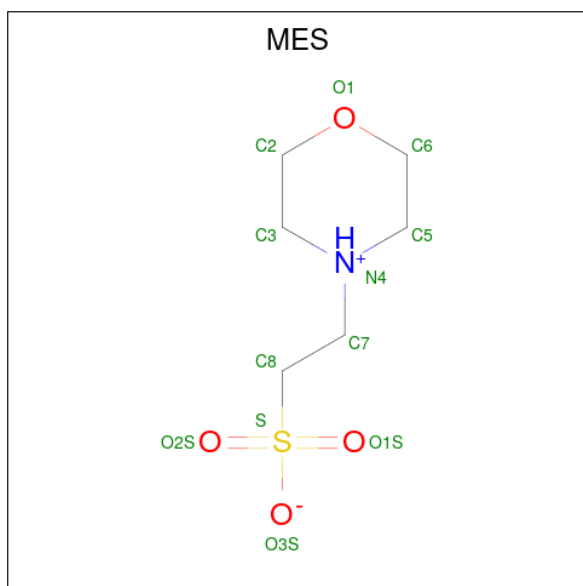
There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	46	MET	ILE	engineered mutation	UNP P54548
A	65	ALA	HIS	engineered mutation	UNP P54548
A	228	MET	LEU	engineered mutation	UNP P54548
A	308	LYS	-	expression tag	UNP P54548
A	309	LEU	-	expression tag	UNP P54548
A	310	ALA	-	expression tag	UNP P54548
A	311	ALA	-	expression tag	UNP P54548
A	312	ALA	-	expression tag	UNP P54548
A	313	LEU	-	expression tag	UNP P54548
A	314	GLU	-	expression tag	UNP P54548
A	315	HIS	-	expression tag	UNP P54548
A	316	HIS	-	expression tag	UNP P54548
A	317	HIS	-	expression tag	UNP P54548
A	318	HIS	-	expression tag	UNP P54548
A	319	HIS	-	expression tag	UNP P54548
A	320	HIS	-	expression tag	UNP P54548

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

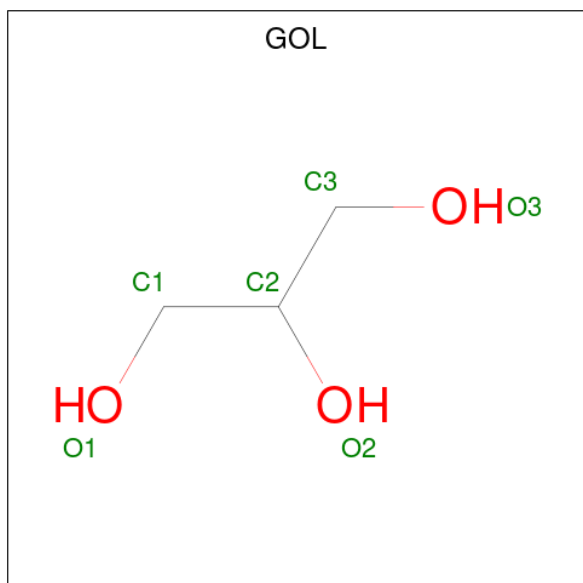
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Zn	0	0
			1	1		

- Molecule 4 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (CCD ID: 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 GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			6	3	3		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	R	1	Total	O	0	0
			1	1		
6	A	10	Total	O	0	0
			10	10		

- Molecule 1: tRNA(THR)



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	173.54Å 42.40Å 110.43Å 90.00° 121.42° 90.00°	Depositor
Resolution (Å)	47.12 – 2.90 47.12 – 2.90	Depositor EDS
% Data completeness (in resolution range)	95.9 (47.12-2.90) 96.0 (47.12-2.90)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.53 (at 2.91Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.250 , 0.299 0.237 , 0.277	Depositor DCC
R_{free} test set	747 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	57.6	Xtriage
Anisotropy	0.726	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 72.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	3550	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.88% 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: GOL, MES, ZN

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	R	0.40	0/1256	1.17	27/1955 (1.4%)
2	A	0.59	0/2441	1.06	10/3301 (0.3%)
All	All	0.53	0/3697	1.10	37/5256 (0.7%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	R	0	1

There are no bond length outliers.

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	162	VAL	N-CA-C	-13.91	99.05	112.96
1	R	72	C	C4'-C3'-O3'	9.31	126.97	113.00
1	R	67	G	C4'-C3'-O3'	8.25	125.37	113.00
1	R	68	G	C4'-C3'-O3'	8.24	125.35	113.00
1	R	66	U	C2'-C3'-O3'	7.54	125.01	113.70
1	R	71	G	C2'-C3'-O3'	7.49	120.73	109.50
1	R	70	A	C4'-C3'-O3'	7.32	123.98	113.00
1	R	66	U	C4'-C3'-O3'	7.06	123.59	113.00
1	R	71	G	C4'-C3'-O3'	6.65	119.38	109.40
2	A	276	GLY	N-CA-C	-6.58	107.10	114.40
1	R	69	A	C2'-C3'-O3'	6.57	123.56	113.70
2	A	217	LYS	N-CA-C	-6.51	103.87	110.97
1	R	67	G	C2'-C3'-O3'	6.46	123.39	113.70
1	R	64	C	C4'-C3'-O3'	6.33	122.50	113.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	R	63	G	C2'-C3'-O3'	6.18	122.97	113.70
1	R	7	A	N9-C1'-C2'	6.16	121.24	112.00
1	R	69	A	C4'-C3'-O3'	6.13	122.20	113.00
2	A	58	LYS	N-CA-C	6.06	118.26	108.32
2	A	20	THR	N-CA-C	6.00	123.59	110.80
1	R	22	G	C2'-C3'-O3'	6.00	118.50	109.50
1	R	70	A	C2'-C3'-O3'	5.88	122.52	113.70
1	R	68	G	C4'-C3'-C2'	5.84	108.44	102.60
1	R	7	A	C2'-C3'-O3'	-5.76	105.06	113.70
1	R	69	A	C4'-C3'-C2'	5.60	108.20	102.60
2	A	32	SER	N-CA-C	5.60	115.94	108.38
1	R	67	G	C4'-C3'-C2'	5.54	108.14	102.60
1	R	70	A	C4'-C3'-C2'	5.48	108.08	102.60
1	R	64	C	C4'-C3'-C2'	5.44	108.04	102.60
1	R	65	U	C4'-C3'-C2'	5.37	107.97	102.60
2	A	145	PHE	N-CA-C	5.35	117.81	109.52
2	A	176	LYS	N-CA-C	-5.32	105.56	111.36
1	R	66	U	C4'-C3'-C2'	5.27	107.87	102.60
1	R	10	G	N9-C1'-C2'	5.27	119.90	112.00
2	A	120	ILE	N-CA-C	5.22	120.20	109.34
1	R	63	G	C4'-C3'-C2'	5.21	107.81	102.60
2	A	131	PHE	N-CA-C	5.15	117.64	109.50
1	R	62	C	C4'-C3'-C2'	5.11	107.71	102.60

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	R	71	G	Sidechain

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	R	1125	0	576	66	0
2	A	2395	0	2411	173	0
3	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	12	0	13	2	0
5	A	6	0	4	0	0
6	A	10	0	0	0	0
6	R	1	0	0	0	0
All	All	3550	0	3004	233	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 36.

All (233) 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)
2:A:240:LYS:H	2:A:240:LYS:HD3	1.24	0.99
1:R:9:A:H62	1:R:23:A:H62	1.19	0.89
2:A:177:ILE:HG23	2:A:193:GLY:HA2	1.54	0.88
2:A:190:ILE:HD13	2:A:190:ILE:H	1.40	0.81
2:A:249:THR:H	2:A:252:GLN:HE21	1.28	0.81
1:R:9:A:N6	1:R:23:A:H62	1.80	0.80
2:A:283:GLN:O	2:A:287:VAL:HG23	1.82	0.79
1:R:66:U:C2'	1:R:67:G:H5''	2.12	0.78
2:A:204:ARG:HD2	2:A:306:ARG:HA	1.64	0.78
2:A:190:ILE:HD13	2:A:190:ILE:N	2.01	0.75
2:A:185:LEU:HB2	2:A:189:ARG:HB3	1.68	0.75
2:A:63:HIS:CE1	2:A:211:ASP:HB2	2.22	0.75
1:R:13:C:H2'	1:R:14:A:H5''	1.69	0.74
2:A:198:GLU:OE1	2:A:199:PRO:HD2	1.87	0.74
1:R:65:U:C2'	1:R:66:U:H5''	2.17	0.74
2:A:191:ILE:HD13	2:A:191:ILE:H	1.53	0.73
2:A:283:GLN:HE21	2:A:295:ALA:H	1.34	0.73
1:R:71:G:O2'	1:R:72:C:O5'	2.07	0.72
2:A:194:ASN:HA	2:A:197:LEU:HD12	1.72	0.71
2:A:249:THR:H	2:A:252:GLN:NE2	1.88	0.71
1:R:61:C:O2'	1:R:62:C:H5'	1.91	0.70
1:R:66:U:H2'	1:R:67:G:H8	1.57	0.70
1:R:64:C:O2'	1:R:65:U:H5'	1.92	0.70
2:A:278:ALA:HA	2:A:281:GLU:OE1	1.92	0.69
2:A:185:LEU:HD12	2:A:189:ARG:HB3	1.74	0.68
1:R:7:A:H5'	1:R:8:U:OP2	1.93	0.68
2:A:201:LYS:HE2	2:A:306:ARG:HB3	1.74	0.68
2:A:4:LEU:C	2:A:4:LEU:HD23	2.19	0.67
2:A:9:GLY:O	2:A:269:HIS:HA	1.94	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:54:ARG:HD2	2:A:82:GLN:O	1.94	0.67
2:A:87:GLU:HA	2:A:114:PRO:HG2	1.77	0.67
1:R:61:C:H5''	2:A:31:ARG:HH21	1.60	0.67
1:R:65:U:O2'	1:R:66:U:H5''	1.94	0.67
2:A:35:LEU:HD11	2:A:46:MET:CE	2.26	0.66
1:R:4:U:H5'	2:A:54:ARG:NH1	2.11	0.66
2:A:275:GLN:HB2	2:A:277:ASP:OD1	1.96	0.66
1:R:66:U:H2'	1:R:67:G:H5''	1.75	0.66
2:A:96:ILE:HG22	2:A:119:GLU:HG2	1.78	0.66
2:A:236:LYS:HG2	2:A:251:GLU:HB2	1.76	0.66
2:A:251:GLU:O	2:A:255:VAL:HG23	1.97	0.65
1:R:60:C:H5''	1:R:61:C:H5	1.60	0.65
1:R:64:C:H2'	1:R:65:U:C6	2.33	0.64
2:A:191:ILE:HD13	2:A:191:ILE:N	2.12	0.64
2:A:173:VAL:O	2:A:177:ILE:HG13	1.97	0.64
1:R:69:A:O2'	1:R:70:A:O5'	2.17	0.63
2:A:185:LEU:HD12	2:A:189:ARG:HG2	1.81	0.63
1:R:65:U:H2'	1:R:66:U:H5''	1.80	0.63
1:R:59:G:H8	1:R:59:G:H5'	1.64	0.62
2:A:275:GLN:HB2	2:A:277:ASP:CG	2.23	0.62
2:A:2:GLU:HG3	2:A:303:ASN:HD22	1.65	0.62
1:R:71:G:O2'	1:R:72:C:C6	2.52	0.62
1:R:63:G:O2'	1:R:64:C:H5'	1.99	0.62
1:R:58:A:H4'	1:R:59:G:OP1	2.00	0.62
2:A:131:PHE:CE1	2:A:151:GLU:HB2	2.35	0.62
2:A:156:GLY:HA3	2:A:198:GLU:O	1.99	0.62
2:A:161:ASP:O	2:A:165:GLU:HG3	2.00	0.62
1:R:18:G:H2'	1:R:57:G:N2	2.14	0.61
2:A:35:LEU:HD11	2:A:46:MET:HE1	1.83	0.61
1:R:2:C:O2'	1:R:3:U:H5'	2.01	0.61
2:A:12:ILE:N	2:A:12:ILE:HD13	2.15	0.61
1:R:71:G:O2'	1:R:72:C:H6	1.85	0.60
2:A:52:LYS:H	2:A:55:LYS:HD3	1.66	0.60
2:A:154:VAL:HG21	2:A:201:LYS:HD3	1.85	0.59
2:A:171:GLY:O	2:A:174:TYR:HB2	2.01	0.59
2:A:1:MET:HA	2:A:25:LYS:O	2.02	0.59
2:A:240:LYS:H	2:A:240:LYS:CD	2.07	0.59
2:A:183:VAL:HG23	2:A:191:ILE:HD11	1.85	0.58
2:A:238:ASP:HA	2:A:240:LYS:NZ	2.18	0.58
2:A:240:LYS:HD3	2:A:240:LYS:N	2.07	0.58
2:A:209:SER:HB3	2:A:218:LEU:HD21	1.86	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:43:GLN:O	2:A:46:MET:HB2	2.04	0.58
1:R:61:C:H5'	2:A:31:ARG:NH2	2.19	0.58
1:R:70:A:H2'	1:R:71:G:C8	2.39	0.57
2:A:172:PRO:O	2:A:174:TYR:N	2.37	0.57
2:A:97:LYS:HG2	2:A:101:GLU:OE1	2.05	0.57
2:A:102:THR:O	2:A:106:VAL:HG22	2.04	0.56
2:A:214:VAL:HG13	2:A:252:GLN:HB3	1.87	0.56
2:A:83:GLY:O	2:A:84:GLY:O	2.24	0.56
2:A:172:PRO:C	2:A:174:TYR:H	2.14	0.55
2:A:201:LYS:CE	2:A:306:ARG:HB3	2.36	0.55
2:A:30:ARG:NH2	2:A:128:ASP:OD2	2.39	0.55
2:A:236:LYS:HG3	2:A:285:GLU:OE1	2.07	0.55
1:R:9:A:H62	1:R:23:A:N6	1.98	0.54
2:A:8:THR:HG21	2:A:297:TYR:N	2.23	0.54
2:A:170:PRO:HA	2:A:174:TYR:CZ	2.42	0.54
2:A:172:PRO:C	2:A:174:TYR:N	2.64	0.54
1:R:64:C:H2'	1:R:65:U:H6	1.72	0.54
2:A:283:GLN:NE2	2:A:295:ALA:H	2.04	0.54
2:A:97:LYS:HA	2:A:117:ILE:HG21	1.88	0.54
2:A:87:GLU:HG3	2:A:114:PRO:HG2	1.89	0.53
2:A:222:ALA:O	2:A:225:CYS:HB3	2.08	0.53
2:A:52:LYS:HB2	2:A:55:LYS:HD3	1.89	0.53
2:A:80:SER:OG	2:A:112:THR:HG23	2.08	0.53
2:A:223:ARG:HG3	2:A:224:ASP:OD2	2.07	0.53
1:R:7:A:C5'	1:R:8:U:OP2	2.57	0.53
1:R:61:C:C2'	1:R:62:C:H5'	2.39	0.53
2:A:205:SER:O	2:A:226:ASP:HB2	2.09	0.53
2:A:249:THR:O	2:A:250:THR:C	2.50	0.52
2:A:113:TYR:HB2	2:A:114:PRO:HD2	1.91	0.52
2:A:176:LYS:N	2:A:176:LYS:HE2	2.24	0.52
2:A:209:SER:CB	2:A:218:LEU:HD21	2.40	0.52
1:R:66:U:C2	1:R:67:G:C8	2.98	0.52
1:R:8:U:O4'	1:R:48:C:O2'	2.28	0.52
2:A:2:GLU:HG3	2:A:303:ASN:ND2	2.25	0.52
2:A:170:PRO:HA	2:A:174:TYR:CE1	2.45	0.52
2:A:142:VAL:O	2:A:143:GLU:HB2	2.10	0.51
1:R:22:G:O2'	1:R:23:A:P	2.69	0.51
2:A:176:LYS:HE2	2:A:176:LYS:CA	2.40	0.51
2:A:27:LEU:O	2:A:31:ARG:HA	2.10	0.51
2:A:100:ILE:HG21	2:A:117:ILE:HD11	1.93	0.50
2:A:6:LEU:O	2:A:45:GLN:HG3	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:97:LYS:O	2:A:101:GLU:HB2	2.11	0.50
1:R:64:C:O2'	1:R:65:U:C5'	2.58	0.50
2:A:86:ASP:CG	4:A:801:MES:H32	2.35	0.50
2:A:4:LEU:C	2:A:4:LEU:CD2	2.84	0.49
2:A:5:PHE:CD2	2:A:22:VAL:HG22	2.46	0.49
2:A:5:PHE:CZ	2:A:266:ILE:HG21	2.46	0.49
2:A:66:GLY:HA3	2:A:70:TYR:CE2	2.47	0.49
2:A:162:VAL:O	2:A:166:MET:HG3	2.11	0.49
2:A:132:ILE:HG23	2:A:150:GLN:HB3	1.93	0.49
2:A:223:ARG:C	2:A:225:CYS:H	2.21	0.49
2:A:54:ARG:NE	2:A:84:GLY:H	2.11	0.49
2:A:175:GLN:HG3	2:A:176:LYS:HE3	1.93	0.49
2:A:249:THR:N	2:A:252:GLN:NE2	2.59	0.49
2:A:30:ARG:HH11	2:A:57:GLU:CD	2.20	0.49
1:R:65:U:O2'	1:R:66:U:C5'	2.60	0.49
2:A:190:ILE:N	2:A:190:ILE:CD1	2.71	0.49
2:A:238:ASP:HA	2:A:240:LYS:HZ1	1.77	0.48
1:R:48:C:O4'	1:R:48:C:P	2.71	0.48
1:R:66:U:O2'	1:R:67:G:H5''	2.13	0.48
2:A:94:LYS:HA	2:A:119:GLU:OE1	2.14	0.48
2:A:182:THR:HG22	2:A:192:ASN:CG	2.38	0.48
1:R:66:U:C2	1:R:67:G:N7	2.81	0.48
2:A:185:LEU:HD12	2:A:189:ARG:CB	2.42	0.48
1:R:1:G:O2'	1:R:2:C:H5'	2.14	0.48
2:A:154:VAL:HG23	2:A:201:LYS:HB3	1.96	0.47
2:A:228:MET:HE1	2:A:256:THR:HG22	1.97	0.47
2:A:136:VAL:O	2:A:145:PHE:HA	2.14	0.47
2:A:32:SER:HB3	4:A:801:MES:H62	1.96	0.47
1:R:60:C:H5''	1:R:61:C:C5	2.46	0.47
2:A:236:LYS:O	2:A:239:ARG:HG2	2.14	0.47
1:R:66:U:N3	1:R:67:G:N7	2.61	0.47
2:A:213:ARG:O	2:A:214:VAL:C	2.55	0.47
1:R:71:G:O2'	1:R:72:C:C5'	2.63	0.47
2:A:265:LEU:HD12	2:A:266:ILE:N	2.30	0.47
1:R:4:U:C5'	2:A:54:ARG:NH1	2.78	0.47
2:A:249:THR:HG23	2:A:252:GLN:NE2	2.30	0.47
2:A:136:VAL:CG1	2:A:137:SER:N	2.78	0.47
2:A:185:LEU:HD11	2:A:191:ILE:HD12	1.96	0.47
2:A:96:ILE:CG2	2:A:119:GLU:HG2	2.43	0.47
1:R:56:C:O4'	2:A:160:ALA:HB2	2.15	0.46
1:R:49:A:H2'	1:R:50:G:C8	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:1:MET:HB3	2:A:204:ARG:HH21	1.80	0.46
2:A:185:LEU:HD12	2:A:189:ARG:CG	2.43	0.46
2:A:209:SER:HA	2:A:218:LEU:HD11	1.97	0.46
2:A:36:PHE:HZ	2:A:149:VAL:HG21	1.81	0.46
1:R:3:U:C2	1:R:71:G:N2	2.83	0.46
2:A:30:ARG:NH1	2:A:57:GLU:OE2	2.49	0.46
2:A:42:THR:HG22	2:A:46:MET:HE3	1.96	0.46
2:A:173:VAL:O	2:A:173:VAL:HG22	2.16	0.46
2:A:176:LYS:HE2	2:A:176:LYS:HA	1.97	0.46
2:A:57:GLU:O	2:A:88:LEU:HD12	2.16	0.46
2:A:175:GLN:HG3	2:A:176:LYS:CE	2.46	0.46
2:A:78:SER:O	2:A:82:GLN:HG3	2.17	0.45
2:A:161:ASP:C	2:A:161:ASP:OD1	2.59	0.45
2:A:234:PHE:O	2:A:249:THR:HA	2.17	0.45
2:A:275:GLN:C	2:A:277:ASP:N	2.74	0.45
1:R:8:U:C4'	1:R:48:C:O2'	2.64	0.45
1:R:53:G:O2'	1:R:54:U:H5'	2.17	0.45
1:R:49:A:H2'	1:R:50:G:H8	1.81	0.45
2:A:41:ALA:O	2:A:44:HIS:HB2	2.16	0.45
2:A:138:VAL:HB	2:A:212:THR:HB	1.98	0.45
2:A:238:ASP:OD1	2:A:238:ASP:N	2.49	0.45
2:A:72:LEU:HB2	2:A:73:PRO:HD3	1.98	0.45
2:A:154:VAL:HG22	2:A:201:LYS:HG2	1.98	0.45
2:A:249:THR:O	2:A:252:GLN:HB2	2.17	0.45
1:R:18:G:O2'	1:R:60:C:N4	2.49	0.45
2:A:192:ASN:O	2:A:193:GLY:C	2.59	0.45
2:A:52:LYS:HB2	2:A:55:LYS:CD	2.46	0.45
2:A:88:LEU:HD12	2:A:89:THR:N	2.31	0.45
1:R:68:G:O2'	1:R:69:A:H5'	2.17	0.45
2:A:46:MET:C	2:A:48:HIS:H	2.24	0.44
1:R:21:A:O2'	1:R:22:G:P	2.75	0.44
1:R:58:A:O2'	1:R:60:C:OP2	2.30	0.44
2:A:275:GLN:C	2:A:277:ASP:H	2.26	0.44
1:R:65:U:HO2'	1:R:66:U:H5''	1.80	0.44
1:R:71:G:O2'	1:R:72:C:P	2.76	0.44
1:R:52:G:O2'	1:R:53:G:H5'	2.18	0.44
2:A:245:TYR:O	2:A:246:TYR:HB2	2.18	0.44
2:A:69:VAL:C	2:A:71:GLY:H	2.25	0.43
1:R:66:U:H2'	1:R:67:G:C8	2.44	0.43
2:A:192:ASN:HB3	2:A:195:ASP:OD2	2.19	0.43
1:R:22:G:O2'	1:R:23:A:OP1	2.36	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:121:GLU:O	2:A:122:GLU:C	2.61	0.43
2:A:268:THR:OG1	2:A:269:HIS:N	2.52	0.43
2:A:248:SER:HA	2:A:252:GLN:HE22	1.83	0.43
2:A:12:ILE:HD13	2:A:12:ILE:H	1.84	0.43
2:A:72:LEU:O	2:A:73:PRO:C	2.61	0.43
2:A:230:HIS:ND1	2:A:231:GLU:O	2.52	0.43
2:A:300:LEU:HD12	2:A:301:GLU:N	2.34	0.43
1:R:70:A:H2'	1:R:71:G:H8	1.84	0.42
2:A:128:ASP:C	2:A:128:ASP:OD1	2.62	0.42
2:A:131:PHE:CD1	2:A:151:GLU:HA	2.54	0.42
2:A:222:ALA:CB	2:A:228:MET:HE3	2.49	0.42
2:A:230:HIS:HD1	2:A:231:GLU:N	2.16	0.42
1:R:3:U:H2'	1:R:4:U:C6	2.55	0.42
2:A:1:MET:HB3	2:A:204:ARG:NH2	2.34	0.42
2:A:204:ARG:HD2	2:A:306:ARG:CA	2.43	0.42
2:A:65:ALA:O	2:A:68:HIS:N	2.43	0.42
2:A:69:VAL:C	2:A:71:GLY:N	2.77	0.42
2:A:281:GLU:O	2:A:285:GLU:HG3	2.19	0.42
1:R:4:U:H4'	2:A:54:ARG:HH12	1.84	0.42
2:A:159:LYS:HD3	2:A:198:GLU:HG2	2.02	0.42
2:A:249:THR:O	2:A:252:GLN:N	2.51	0.42
2:A:63:HIS:CE1	2:A:211:ASP:CB	3.00	0.42
2:A:199:PRO:HA	2:A:200:PRO:HD3	1.92	0.41
2:A:254:ALA:HB1	2:A:289:VAL:HB	2.02	0.41
2:A:63:HIS:CE1	2:A:68:HIS:CE1	3.08	0.41
1:R:13:C:H2'	1:R:14:A:C5'	2.44	0.41
2:A:273:ARG:HG3	2:A:274:TYR:CD2	2.56	0.41
2:A:248:SER:HA	2:A:252:GLN:NE2	2.34	0.41
2:A:111:LEU:C	2:A:113:TYR:H	2.29	0.41
2:A:136:VAL:HG12	2:A:137:SER:N	2.36	0.41
1:R:6:C:O2	1:R:6:C:H2'	2.21	0.41
1:R:14:A:C5	1:R:22:G:C2	3.08	0.41
2:A:100:ILE:CG2	2:A:117:ILE:HD11	2.49	0.41
2:A:185:LEU:CD1	2:A:189:ARG:HB3	2.49	0.41
2:A:243:TYR:CD1	2:A:243:TYR:C	2.99	0.40
2:A:30:ARG:O	2:A:31:ARG:HB2	2.21	0.40
2:A:106:VAL:HG23	2:A:107:THR:HG23	2.03	0.40
2:A:208:PHE:CD1	2:A:208:PHE:N	2.90	0.40
1:R:63:G:C2'	1:R:64:C:H5'	2.52	0.40
2:A:260:ALA:O	2:A:261:ARG:HB2	2.20	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	
2	A	305/320 (95%)	258 (85%)	37 (12%)	10 (3%)	3	13

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	A	21	SER
2	A	84	GLY
2	A	20	THR
2	A	66	GLY
2	A	122	GLU
2	A	173	VAL
2	A	238	ASP
2	A	37	ASP
2	A	53	PRO
2	A	193	GLY

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	
2	A	257/267 (96%)	236 (92%)	21 (8%)	10	32

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	A	12	ILE

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Mol	Chain	Res	Type
2	A	20	THR
2	A	53	PRO
2	A	101	GLU
2	A	104	LEU
2	A	111	LEU
2	A	115	LEU
2	A	134	THR
2	A	158	LEU
2	A	189	ARG
2	A	190	ILE
2	A	191	ILE
2	A	204	ARG
2	A	214	VAL
2	A	216	ASP
2	A	225	CYS
2	A	231	GLU
2	A	240	LYS
2	A	284	LYS
2	A	300	LEU
2	A	306	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
2	A	18	ASN
2	A	63	HIS
2	A	68	HIS
2	A	194	ASN
2	A	252	GLN
2	A	269	HIS
2	A	283	GLN
2	A	303	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	R	51/79 (64%)	24 (47%)	10 (19%)

All (24) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	R	7	A
1	R	8	U
1	R	9	A
1	R	10	G
1	R	14	A
1	R	16	C
1	R	18	G
1	R	19	G
1	R	20	U
1	R	22	G
1	R	23	A
1	R	25	C
1	R	48	C
1	R	49	A
1	R	59	G
1	R	61	C
1	R	66	U
1	R	67	G
1	R	68	G
1	R	69	A
1	R	70	A
1	R	71	G
1	R	72	C
1	R	73	U

All (10) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	R	18	G
1	R	19	G
1	R	21	A
1	R	22	G
1	R	58	A
1	R	68	G
1	R	69	A
1	R	70	A
1	R	71	G
1	R	72	C

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](#)

Of 3 ligands modelled in this entry, 1 is monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	GOL	A	501	-	5,5,5	4.73	5 (100%)	5,5,5	6.14	3 (60%)
4	MES	A	801	-	12,12,12	9.31	8 (66%)	15,16,16	2.24	5 (33%)

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
5	GOL	A	501	-	-	2/4/4/4	-
4	MES	A	801	-	-	0/6/14/14	0/1/1/1

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	801	MES	C8-S	-24.45	1.43	1.77
4	A	801	MES	O2S-S	12.68	1.80	1.45
4	A	801	MES	O1S-S	12.45	1.80	1.45
4	A	801	MES	O3S-S	8.74	1.80	1.47
5	A	501	GOL	C3-C2	-7.98	1.21	1.51
4	A	801	MES	C7-C8	-4.75	1.39	1.52
5	A	501	GOL	O1-C1	4.50	1.61	1.42
5	A	501	GOL	O3-C3	3.31	1.56	1.42
5	A	501	GOL	O2-C2	-3.00	1.34	1.43
4	A	801	MES	C3-C2	-2.91	1.39	1.50
5	A	501	GOL	C1-C2	-2.87	1.40	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	801	MES	C5-C6	-2.51	1.41	1.50
4	A	801	MES	C7-N4	-2.41	1.41	1.47

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	501	GOL	O3-C3-C2	11.13	160.50	110.38
5	A	501	GOL	O2-C2-C3	7.23	139.10	109.18
4	A	801	MES	O3S-S-C8	5.17	116.11	106.00
4	A	801	MES	O1S-S-C8	3.78	112.44	106.73
5	A	501	GOL	O1-C1-C2	3.44	125.84	110.38
4	A	801	MES	O2S-S-O1S	-2.56	105.48	113.82
4	A	801	MES	O3S-S-O2S	-2.55	105.03	111.40
4	A	801	MES	O2S-S-C8	2.27	110.16	106.73

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	501	GOL	C1-C2-C3-O3
5	A	501	GOL	O1-C1-C2-O2

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	801	MES	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	R	53/79 (67%)	2.38	32 (60%) 0 0	53, 97, 156, 179	0
2	A	307/320 (95%)	0.67	20 (6%) 25 20	33, 50, 84, 98	0
All	All	360/399 (90%)	0.93	52 (14%) 6 5	33, 52, 113, 179	0

All (52) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	R	8	U	6.1
1	R	9	A	5.9
1	R	46	G	4.4
1	R	10	G	3.9
1	R	73	U	3.8
2	A	175	GLN	3.6
1	R	25	C	3.6
2	A	63	HIS	3.5
1	R	22	G	3.5
1	R	71	G	3.4
1	R	18	G	3.4
1	R	59	G	3.3
1	R	49	A	3.3
1	R	60	C	3.3
2	A	307	GLY	3.3
2	A	263	LYS	3.3
2	A	234	PHE	3.1
1	R	50	G	3.1
1	R	69	A	3.0
1	R	70	A	3.0
2	A	157	SER	2.9
1	R	11	C	2.9
1	R	13	C	2.9
2	A	260	ALA	2.9

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Mol	Chain	Res	Type	RSRZ
2	A	139	ILE	2.9
1	R	24	G	2.8
1	R	16	C	2.8
1	R	1	G	2.7
2	A	211	ASP	2.7
1	R	15	G	2.7
2	A	262	ALA	2.7
2	A	289	VAL	2.6
1	R	58	A	2.5
1	R	17	A	2.5
2	A	140	HIS	2.5
2	A	85	GLU	2.4
1	R	21	A	2.4
1	R	20	U	2.4
1	R	7	A	2.3
2	A	241	LEU	2.3
2	A	83	GLY	2.3
1	R	53	G	2.3
1	R	12	U	2.2
2	A	86	ASP	2.2
2	A	248	SER	2.1
2	A	120	ILE	2.1
1	R	48	C	2.1
2	A	19	VAL	2.1
1	R	51	C	2.1
1	R	4	U	2.1
2	A	188	GLY	2.0
1	R	68	G	2.0

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

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

6.3 Carbohydrates [i](#)

There are no 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($\text{\AA}^2$)	Q<0.9
5	GOL	A	501	6/6	0.91	0.12	69,71,72,73	0
4	MES	A	801	12/12	0.92	0.13	84,86,89,89	0
3	ZN	A	401	1/1	0.98	0.11	99,99,99,99	0

6.5 Other polymers [i](#)

There are no such residues in this entry.