



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 07:45 AM UTC

PDB ID : 6QTA / pdb_00006qta
Title : Crystal structure of Rea1-MIDAS/Rsa4-UBL complex from *Chaetomium thermophilum*
Authors : Ahmed, Y.L.; Thoms, M.; Hurt, E.; Sinning, I.
Deposited on : 2019-02-22
Resolution : 1.89 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

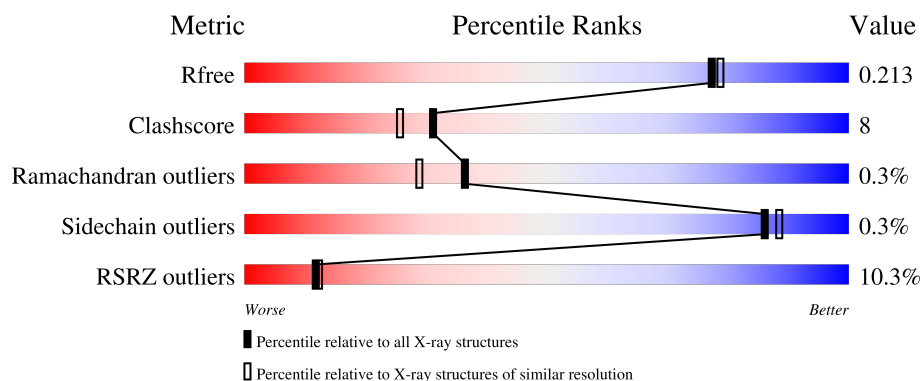
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.89 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	280	
2	B	108	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 3140 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 Midasin,Midasin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	258	Total	C	N	O	S	0	3	0
			2062	1303	369	381	9			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	4689	MET	-	initiating methionine	UNP G0SHE6
A	4734	GLY	-	linker	UNP G0SHE6
A	4735	SER	-	linker	UNP G0SHE6
A	4736	GLY	-	linker	UNP G0SHE6
A	4998	GLY	-	expression tag	UNP G0SHE6
A	4999	SER	-	expression tag	UNP G0SHE6
A	5000	HIS	-	expression tag	UNP G0SHE6
A	5001	HIS	-	expression tag	UNP G0SHE6
A	5002	HIS	-	expression tag	UNP G0SHE6
A	5003	HIS	-	expression tag	UNP G0SHE6
A	5004	HIS	-	expression tag	UNP G0SHE6
A	5005	HIS	-	expression tag	UNP G0SHE6

- Molecule 2 is a protein called Ribosome assembly protein 4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	101	Total	C	N	O	S	0	0	0
			801	504	135	160	2			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	21	MET	-	initiating methionine	UNP G0SC29
B	22	LYS	-	expression tag	UNP G0SC29
B	23	HIS	-	expression tag	UNP G0SC29
B	24	HIS	-	expression tag	UNP G0SC29

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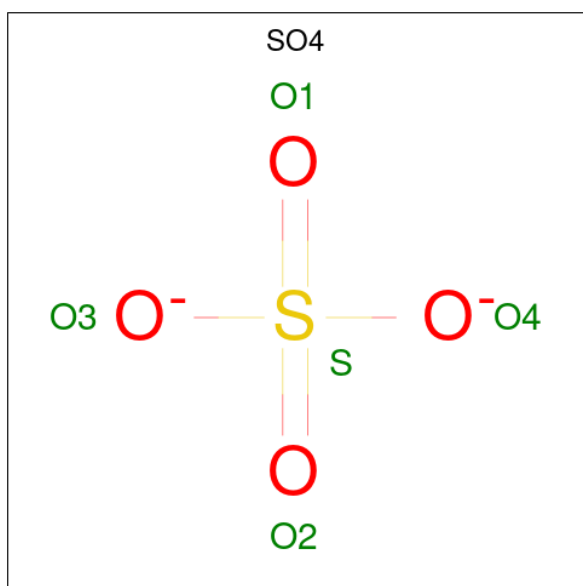
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Chain	Residue	Modelled	Actual	Comment	Reference
B	25	HIS	-	expression tag	UNP G0SC29
B	26	HIS	-	expression tag	UNP G0SC29
B	27	HIS	-	expression tag	UNP G0SC29
B	28	HIS	-	expression tag	UNP G0SC29
B	29	PRO	-	expression tag	UNP G0SC29
B	30	MET	VAL	conflict	UNP G0SC29

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

- Molecule 5 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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

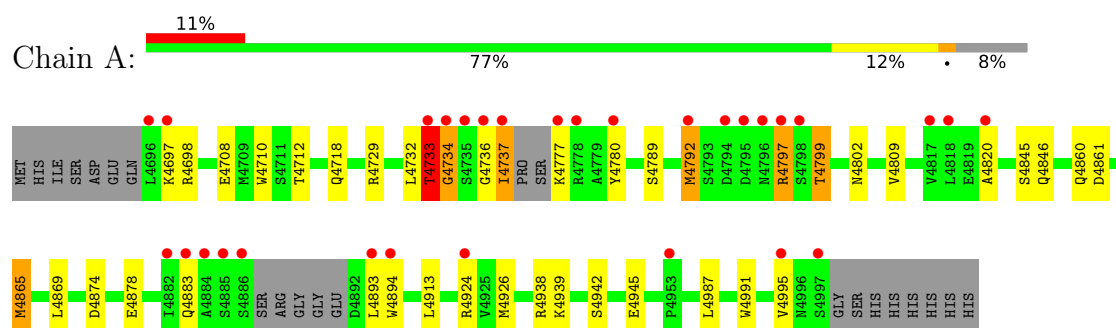
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	165	Total	O	0	0
			165	165		
6	B	54	Total	O	0	0
			54	54		

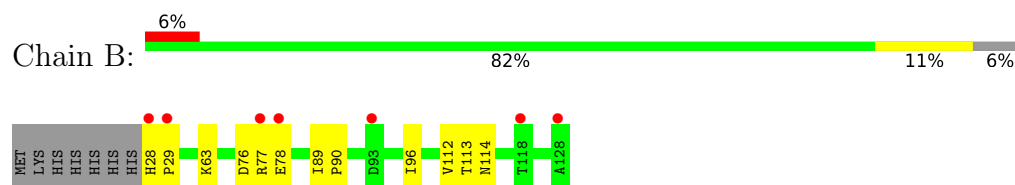
3 Residue-property plots [i](#)

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: Midasin,Midasin



- Molecule 2: Ribosome assembly protein 4



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	115.67Å 115.67Å 74.62Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.58 – 1.89 36.58 – 1.89	Depositor EDS
% Data completeness (in resolution range)	100.0 (36.58-1.89) 100.0 (36.58-1.89)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.20 (at 1.89Å)	Xtriage
Refinement program	PHENIX 1.12_2829	Depositor
R, R_{free}	0.176 , 0.211 0.178 , 0.213	Depositor DCC
R_{free} test set	2122 reflections (5.17%)	wwPDB-VP
Wilson B-factor (Å ²)	29.6	Xtriage
Anisotropy	0.249	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 58.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3140	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.59% 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: MG, SO4, GOL

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	0.87	2/2098 (0.1%)	0.87	6/2838 (0.2%)
2	B	0.59	0/817	0.66	0/1109
All	All	0.80	2/2915 (0.1%)	0.82	6/3947 (0.2%)

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	A	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	4792	MET	SD-CE	-8.03	1.59	1.79
1	A	4797	ARG	CZ-NH1	5.85	1.41	1.32

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	4924	ARG	CD-NE-CZ	7.31	134.63	124.40
1	A	4797	ARG	NE-CZ-NH2	-6.55	113.30	119.20
1	A	4733	THR	CA-C-N	6.49	128.36	121.48
1	A	4733	THR	C-N-CA	6.49	128.36	121.48
1	A	4734	GLY	N-CA-C	-6.41	101.75	111.10
1	A	4865	MET	CG-SD-CE	-5.67	88.43	100.90

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	4733	THR	Peptide

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2062	0	2067	37	0
2	B	801	0	777	10	0
3	A	1	0	0	0	0
4	A	15	0	0	1	0
5	A	42	0	54	6	0
6	A	165	0	0	6	1
6	B	54	0	0	3	0
All	All	3140	0	2898	48	1

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 8.

All (48) 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:A:4894:TRP:HZ3	6:A:5239:HOH:O	1.08	1.27
1:A:4799:THR:HG22	1:A:4802:ASN:H	1.39	0.86
1:A:4860:GLN:HG2	5:A:5111:GOL:H12	1.58	0.84
1:A:4861[B]:ASP:OD2	2:B:113:THR:OG1	1.98	0.80
1:A:4894:TRP:CZ3	6:A:5239:HOH:O	1.96	0.79
1:A:4846:GLN:H	5:A:5109:GOL:H11	1.47	0.79
1:A:4883:GLN:HG3	1:A:4893:LEU:HD21	1.68	0.75
1:A:4845:SER:HB2	5:A:5109:GOL:H31	1.70	0.73
1:A:4894:TRP:CH2	1:A:4926:MET:HB2	2.24	0.73
2:B:28:HIS:HB3	2:B:29:PRO:CD	2.24	0.67
1:A:4874:ASP:O	1:A:4878:GLU:HG3	1.95	0.66
2:B:28:HIS:N	6:B:202:HOH:O	2.31	0.63
1:A:4869:LEU:HD22	1:A:4913:LEU:HD22	1.82	0.62
1:A:4718:GLN:OE1	6:A:5201:HOH:O	2.16	0.61
1:A:4733:THR:HA	1:A:4737:ILE:HG23	1.84	0.59
4:A:5102:SO4:O4	5:A:5106:GOL:O3	2.22	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:4732:LEU:O	1:A:4737:ILE:HG22	2.06	0.56
1:A:4697:LYS:HG3	6:A:5252:HOH:O	2.08	0.52
1:A:4894:TRP:CH2	1:A:4926:MET:CB	2.93	0.51
1:A:4789:SER:O	1:A:4792:MET:HG2	2.10	0.51
1:A:4736:GLY:O	1:A:4777:LYS:N	2.45	0.50
1:A:4697:LYS:HE2	1:A:4697:LYS:HA	1.93	0.50
1:A:4780:TYR:CE2	1:A:4820:ALA:HB3	2.46	0.49
2:B:28:HIS:HB3	2:B:29:PRO:HD3	1.92	0.49
5:A:5107:GOL:H31	6:A:5255:HOH:O	2.12	0.49
1:A:4991:TRP:CZ2	1:A:4995:VAL:HG21	2.48	0.49
1:A:4894:TRP:CH2	1:A:4991:TRP:CH2	3.02	0.48
1:A:4894:TRP:HH2	1:A:4926:MET:HB2	1.74	0.48
1:A:4708:GLU:O	1:A:4712[B]:THR:HG22	2.14	0.47
2:B:63:LYS:NZ	6:B:201:HOH:O	2.16	0.47
2:B:89:ILE:HD11	2:B:96:ILE:HD12	1.96	0.47
2:B:90:PRO:HD2	2:B:112:VAL:HG21	1.96	0.47
2:B:114:ASN:HB2	6:B:250:HOH:O	2.16	0.45
1:A:4797:ARG:HD2	6:A:5298:HOH:O	2.15	0.45
1:A:4729:ARG:O	1:A:4733:THR:HB	2.17	0.45
1:A:4991:TRP:CE2	1:A:4995:VAL:HG21	2.53	0.44
1:A:4697:LYS:HD3	1:A:4698:ARG:H	1.82	0.44
1:A:4869:LEU:HD22	1:A:4913:LEU:CD2	2.47	0.44
1:A:4938:ARG:HG3	1:A:4939:LYS:N	2.33	0.44
1:A:4710:TRP:CZ3	1:A:4809:VAL:HG22	2.53	0.43
1:A:4987:LEU:HD23	1:A:4987:LEU:C	2.44	0.43
2:B:76:ASP:OD2	2:B:77:ARG:CZ	2.67	0.43
1:A:4732:LEU:C	1:A:4734:GLY:H	2.26	0.42
1:A:4942:SER:HB3	1:A:4945:GLU:OE1	2.18	0.42
2:B:78:GLU:HG3	2:B:78:GLU:O	2.20	0.42
1:A:4865:MET:HE2	1:A:4865:MET:HB2	1.70	0.42
1:A:4860:GLN:CD	5:A:5111:GOL:H31	2.45	0.41
1:A:4732:LEU:O	1:A:4734:GLY:O	2.39	0.41

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:A:5336:HOH:O	6:A:5362:HOH:O[7_555]	2.13	0.07

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	255/280 (91%)	249 (98%)	5 (2%)	1 (0%)	30	22
2	B	99/108 (92%)	98 (99%)	1 (1%)	0	100	100
All	All	354/388 (91%)	347 (98%)	6 (2%)	1 (0%)	36	29

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	4799	THR

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	224/240 (93%)	223 (100%)	1 (0%)	84	87
2	B	90/97 (93%)	90 (100%)	0	100	100
All	All	314/337 (93%)	313 (100%)	1 (0%)	86	88

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4737	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	4714	GLN
1	A	4802	ASN
2	B	127	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 11 ligands modelled in this entry, 1 is monoatomic - leaving 10 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
4	SO4	A	5103	-	4,4,4	0.25	0	6,6,6	0.21	0
5	GOL	A	5110	-	5,5,5	1.39	1 (20%)	5,5,5	1.19	0
4	SO4	A	5104	-	4,4,4	0.20	0	6,6,6	0.54	0
4	SO4	A	5102	-	4,4,4	0.47	0	6,6,6	0.48	0
5	GOL	A	5108	-	5,5,5	1.36	0	5,5,5	0.71	0
5	GOL	A	5109	-	5,5,5	1.07	0	5,5,5	1.01	0
5	GOL	A	5105	-	5,5,5	1.01	0	5,5,5	2.06	2 (40%)
5	GOL	A	5111	-	5,5,5	1.67	2 (40%)	5,5,5	0.75	0
5	GOL	A	5106	-	5,5,5	1.48	1 (20%)	5,5,5	0.77	0
5	GOL	A	5107	-	5,5,5	1.28	0	5,5,5	0.89	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	GOL	A	5110	-	-	2/4/4/4	-
5	GOL	A	5108	-	-	2/4/4/4	-
5	GOL	A	5109	-	-	4/4/4/4	-
5	GOL	A	5105	-	-	2/4/4/4	-
5	GOL	A	5111	-	-	3/4/4/4	-
5	GOL	A	5106	-	-	1/4/4/4	-
5	GOL	A	5107	-	-	4/4/4/4	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	5106	GOL	O2-C2	-2.61	1.35	1.43
5	A	5111	GOL	C1-C2	2.52	1.61	1.51
5	A	5110	GOL	C3-C2	2.47	1.61	1.51
5	A	5111	GOL	C3-C2	2.29	1.60	1.51

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	5105	GOL	O2-C2-C3	3.21	122.45	109.18
5	A	5105	GOL	O2-C2-C1	3.18	122.37	109.18

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	5107	GOL	O1-C1-C2-C3
5	A	5107	GOL	C1-C2-C3-O3
5	A	5110	GOL	O1-C1-C2-C3
5	A	5108	GOL	O1-C1-C2-O2
5	A	5110	GOL	O1-C1-C2-O2
5	A	5111	GOL	O2-C2-C3-O3
5	A	5108	GOL	O1-C1-C2-C3
5	A	5109	GOL	O1-C1-C2-C3
5	A	5109	GOL	C1-C2-C3-O3
5	A	5111	GOL	C1-C2-C3-O3

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Mol	Chain	Res	Type	Atoms
5	A	5105	GOL	O2-C2-C3-O3
5	A	5107	GOL	O1-C1-C2-O2
5	A	5109	GOL	O2-C2-C3-O3
5	A	5107	GOL	O2-C2-C3-O3
5	A	5106	GOL	O2-C2-C3-O3
5	A	5111	GOL	O1-C1-C2-O2
5	A	5105	GOL	O1-C1-C2-O2
5	A	5109	GOL	O1-C1-C2-O2

There are no ring outliers.

5 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	5102	SO4	1	0
5	A	5109	GOL	2	0
5	A	5111	GOL	2	0
5	A	5106	GOL	1	0
5	A	5107	GOL	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

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	258/280 (92%)	0.17	30 (11%) 9 10	15, 30, 84, 113	3 (1%)
2	B	101/108 (93%)	0.45	7 (6%) 23 24	26, 45, 84, 118	0
All	All	359/388 (92%)	0.25	37 (10%) 12 12	15, 36, 85, 118	3 (0%)

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	4696	LEU	6.0
1	A	4894	TRP	5.2
1	A	4737	ILE	5.0
1	A	4995	VAL	4.6
1	A	4734	GLY	4.5
1	A	4736	GLY	4.0
1	A	4780	TYR	3.7
2	B	28	HIS	3.6
1	A	4797	ARG	3.6
1	A	4792	MET	3.5
2	B	128	ALA	3.5
1	A	4882	ILE	3.3
1	A	4997	SER	3.3
1	A	4795	ASP	3.3
1	A	4886	SER	3.1
1	A	4777	LYS	3.0
1	A	4798	SER	3.0
1	A	4884	ALA	3.0
2	B	29	PRO	2.9
1	A	4697	LYS	2.9
1	A	4735	SER	2.8
1	A	4817	VAL	2.8
1	A	4820	ALA	2.8
1	A	4796	ASN	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	4953	PRO	2.6
1	A	4818	LEU	2.5
1	A	4893	LEU	2.5
1	A	4733	THR	2.5
1	A	4794	ASP	2.4
1	A	4885	SER	2.4
1	A	4778	ARG	2.3
2	B	118	THR	2.3
2	B	78	GLU	2.3
2	B	93	ASP	2.3
1	A	4883	GLN	2.1
1	A	4924	ARG	2.1
2	B	77	ARG	2.1

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	5106	6/6	0.74	0.17	48,59,70,71	0
5	GOL	A	5109	6/6	0.78	0.16	73,80,84,87	0
5	GOL	A	5107	6/6	0.82	0.16	73,78,80,81	0
4	SO4	A	5104	5/5	0.85	0.11	72,79,82,83	0
4	SO4	A	5103	5/5	0.85	0.10	86,87,89,90	0
5	GOL	A	5110	6/6	0.85	0.16	58,71,72,72	0
5	GOL	A	5111	6/6	0.87	0.16	58,66,68,71	0
5	GOL	A	5108	6/6	0.88	0.15	65,74,78,87	0
4	SO4	A	5102	5/5	0.91	0.10	53,63,75,81	0
5	GOL	A	5105	6/6	0.97	0.07	32,34,36,38	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	MG	A	5101	1/1	0.99	0.06	25,25,25,25	0

6.5 Other polymers [i](#)

There are no such residues in this entry.