



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 03:13 AM UTC

PDB ID : 1Y1T / pdb_00001y1t
Title : Crystal Structure of the Uridine Phosphorylase from Salmonella Typhimurium at 1.77Å Resolution
Authors : Gabdoulkhakov, A.G.; Dontsova, M.V.; Kachalova, G.S.; Betzel, C.; Ealick, S.E.; Mikhailov, A.M.
Deposited on : 2004-11-19
Resolution : 1.77 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

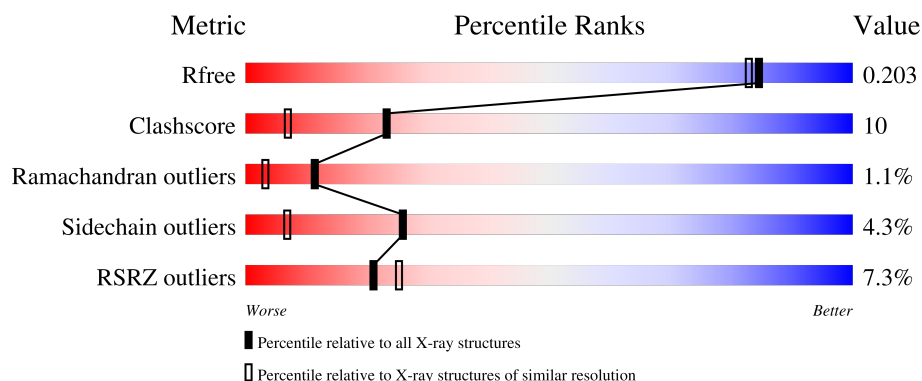
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.77 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	1365 (1.78-1.78)
Clashscore	190562	1395 (1.78-1.78)
Ramachandran outliers	187476	1382 (1.78-1.78)
Sidechain outliers	187428	1382 (1.78-1.78)
RSRZ outliers	180081	1365 (1.78-1.78)

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	253	9% 74% 21% • •
1	F	253	4% 67% 16% • 14%

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	GOL	A	5001	-	X	-	-

2 Entry composition [i](#)

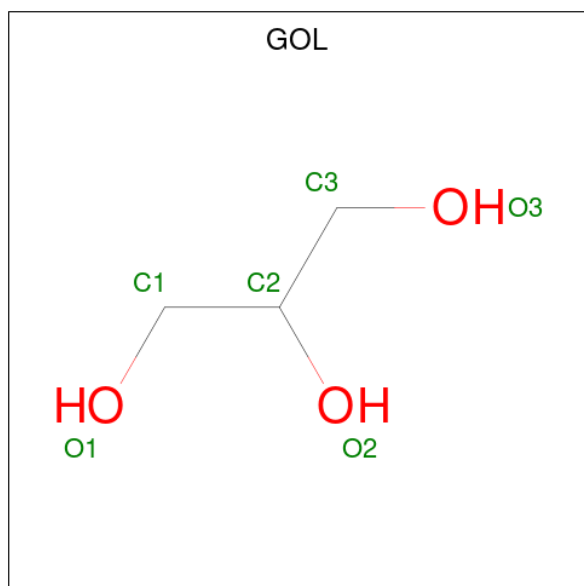
There are 4 unique types of molecules in this entry. The entry contains 3773 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Uridine phosphorylase.

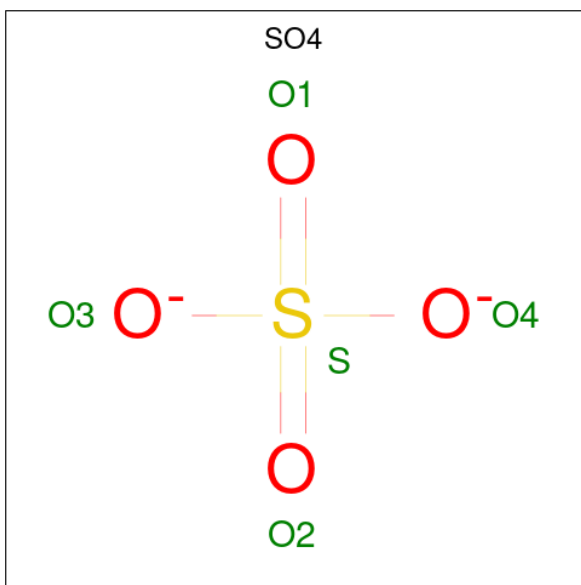
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	250	Total	C	N	O	S	0	0	0
			1876	1174	330	360	12			
1	F	218	Total	C	N	O	S	0	0	0
			1607	1009	278	309	11			

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O_4S).



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

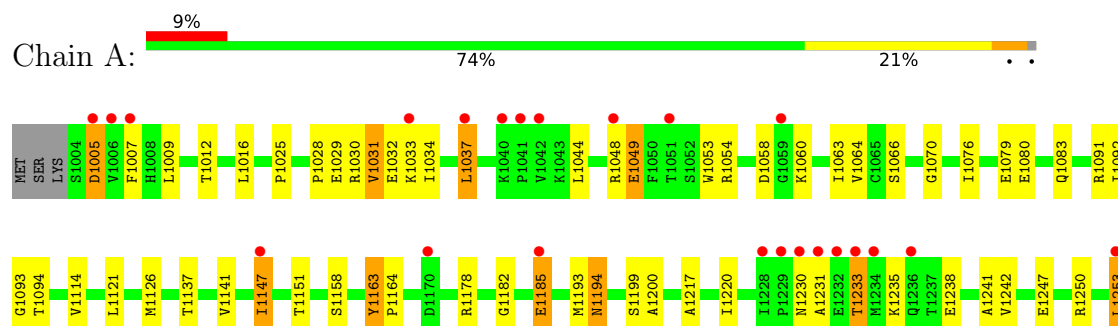
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	153	Total	O	0	0
			153	153		
4	F	126	Total	O	0	0
			126	126		

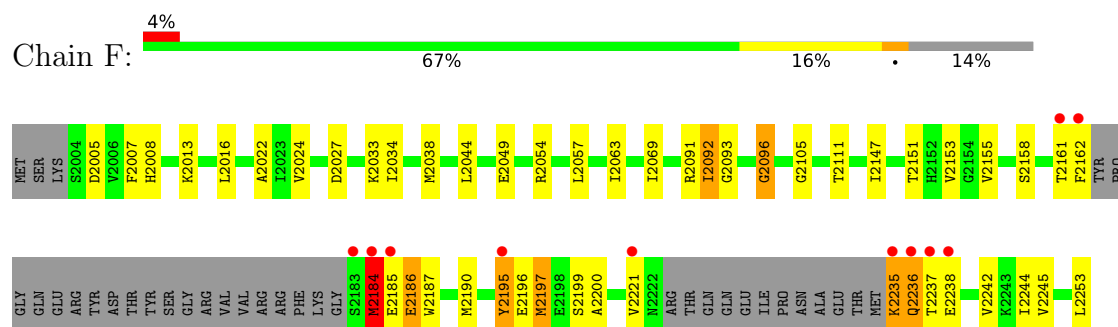
3 Residue-property plots [i](#)

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: Uridine phosphorylase



• Molecule 1: Uridine phosphorylase



4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	150.82Å 150.82Å 47.84Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	25.40 – 1.77 25.40 – 1.77	Depositor EDS
% Data completeness (in resolution range)	97.3 (25.40-1.77) 97.4 (25.40-1.77)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.45 (at 1.77Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.184 , 0.205 0.183 , 0.203	Depositor DCC
R_{free} test set	1925 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	18.9	Xtriage
Anisotropy	0.478	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 55.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.027 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3773	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.16% 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, 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	0.45	0/1906	0.94	8/2584 (0.3%)
1	F	0.53	1/1629 (0.1%)	0.98	11/2208 (0.5%)
All	All	0.49	1/3535 (0.0%)	0.96	19/4792 (0.4%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	2184	MET	C-N	11.80	1.49	1.34

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	2184	MET	O-C-N	7.71	132.85	122.59
1	F	2253	LEU	CA-C-O	-7.24	108.50	120.80
1	F	2007	PHE	N-CA-C	7.21	119.14	111.28
1	F	2185	GLU	N-CA-C	-6.82	103.87	111.71
1	A	1253	LEU	CA-C-O	-6.76	109.30	120.80
1	A	1199	SER	N-CA-C	6.56	118.43	111.28
1	A	1070	GLY	N-CA-C	5.44	119.54	112.68
1	A	1114	VAL	N-CA-C	-5.42	101.73	108.89
1	A	1151	THR	N-CA-C	5.42	118.31	109.59
1	F	2008	HIS	N-CA-C	5.41	120.24	113.43
1	F	2096	GLY	N-CA-C	-5.40	103.28	112.51
1	A	1007	PHE	N-CA-C	5.34	119.42	113.01
1	F	2155	VAL	N-CA-C	5.31	116.94	108.87
1	F	2199	SER	N-CA-C	5.21	116.96	111.28
1	F	2027	ASP	CA-C-N	5.12	124.74	119.05
1	F	2027	ASP	C-N-CA	5.12	124.74	119.05
1	A	1193	MET	N-CA-C	5.10	117.96	111.69
1	A	1031	VAL	N-CA-C	5.05	115.20	110.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	2151	THR	N-CA-C	5.04	117.71	109.59

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1876	0	1887	44	0
1	F	1607	0	1627	32	0
2	A	6	0	4	0	0
3	F	5	0	0	0	0
4	A	153	0	0	4	0
4	F	126	0	0	2	0
All	All	3773	0	3518	72	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 10.

All (72) 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:1092:ILE:HD11	1:A:1241:ALA:HB1	1.53	0.90
1:F:2235:LYS:HB2	1:F:2235:LYS:HZ3	1.52	0.74
1:A:1094:THR:HG21	1:A:1238:GLU:HG2	1.69	0.74
1:F:2187:TRP:CD1	1:F:2195:TYR:HH	2.11	0.68
1:F:2158:SER:HB3	1:F:2200:ALA:HB2	1.75	0.68
1:F:2069:ILE:HG23	1:F:2197:MET:HE1	1.76	0.68
1:F:2038:MET:HG2	1:F:2057:LEU:HD21	1.77	0.66
1:A:1079:GLU:O	1:A:1083:GLN:HG3	1.97	0.65
1:A:1182:GLY:O	1:A:1185:GLU:HG3	1.97	0.64
1:A:1060:LYS:HD2	1:A:1253:LEU:HB3	1.80	0.62
1:A:1048:ARG:HG2	1:A:1048:ARG:HH11	1.65	0.62
1:A:1094:THR:CG2	1:A:1238:GLU:HG2	2.29	0.61
1:A:1194:ASN:HD22	1:A:1194:ASN:N	1.98	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1032:GLU:HG3	1:A:1053:TRP:CH2	2.38	0.59
1:F:2069:ILE:CG2	1:F:2197:MET:HE1	2.33	0.58
1:A:1049:GLU:HG2	1:F:2049:GLU:CD	2.29	0.58
1:A:1048:ARG:HG3	4:A:3090:HOH:O	2.04	0.57
1:A:1049:GLU:HG2	1:F:2049:GLU:HG3	1.87	0.57
1:A:1030:ARG:HH22	1:A:1093:GLY:HA2	1.70	0.57
1:A:1220:ILE:HB	1:A:1233:THR:O	2.04	0.56
1:A:1158:SER:HB3	1:A:1200:ALA:HB2	1.87	0.56
1:A:1030:ARG:NH2	1:A:1093:GLY:HA2	2.22	0.55
1:A:1092:ILE:HD11	1:A:1241:ALA:CB	2.33	0.55
1:A:1054:ARG:HG3	1:A:1063:ILE:HD13	1.91	0.52
1:A:1030:ARG:HA	1:A:1033:LYS:HD2	1.90	0.52
1:F:2092:ILE:C	1:F:2092:ILE:HD13	2.34	0.51
1:F:2147:ILE:HD11	1:F:2244:ILE:HD11	1.93	0.51
1:F:2033:LYS:HE3	4:F:3145:HOH:O	2.09	0.51
1:F:2044:LEU:HD11	1:F:2054:ARG:HB2	1.93	0.51
1:A:1005:ASP:OD2	1:A:1012:THR:HA	2.10	0.51
1:A:1044:LEU:HD11	1:A:1054:ARG:HB2	1.93	0.50
1:A:1121:LEU:HD21	1:A:1126:MET:HE2	1.92	0.50
1:F:2186:GLU:HG3	1:F:2190:MET:HE3	1.94	0.50
1:A:1034:ILE:HG12	1:A:1242:VAL:HG13	1.93	0.50
1:F:2187:TRP:HD1	1:F:2195:TYR:HH	1.58	0.50
1:A:1058:ASP:OD2	1:A:1250:ARG:HG3	2.13	0.49
1:A:1076:ILE:HD11	1:F:2162:PHE:HD1	1.77	0.49
1:A:1137:THR:O	1:A:1141:VAL:HG23	2.13	0.49
1:F:2092:ILE:HD13	1:F:2093:GLY:N	2.27	0.49
1:F:2013:LYS:HG2	4:F:3173:HOH:O	2.13	0.49
1:A:1049:GLU:HG2	1:F:2049:GLU:CG	2.45	0.47
1:F:2105:GLY:HA2	1:F:2237:THR:HG22	1.96	0.47
1:F:2005:ASP:OD2	1:F:2013:LYS:HD2	2.15	0.47
1:A:1230:ASN:O	1:A:1233:THR:N	2.48	0.46
1:F:2242:VAL:O	1:F:2245:VAL:HG12	2.15	0.46
1:F:2111:THR:HG23	1:F:2153:VAL:HG12	1.98	0.45
1:A:1009:LEU:HG	1:A:1080:GLU:OE1	2.17	0.45
1:A:1185:GLU:HG2	4:A:3042:HOH:O	2.17	0.45
1:A:1238:GLU:O	1:A:1242:VAL:HG23	2.18	0.44
1:F:2096:GLY:HA2	1:F:2221:VAL:O	2.16	0.44
1:A:1194:ASN:N	1:A:1194:ASN:ND2	2.65	0.44
1:F:2024:VAL:O	1:F:2091:ARG:HG3	2.18	0.44
1:A:1031:VAL:HG13	1:A:1064:VAL:HG12	2.00	0.44
1:A:1025:PRO:O	1:A:1066:SER:HA	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1054:ARG:HG3	1:A:1063:ILE:CD1	2.48	0.43
1:A:1016:LEU:HG	1:A:1063:ILE:HG13	2.00	0.43
1:F:2238:GLU:O	1:F:2242:VAL:HG23	2.19	0.43
1:A:1178:ARG:HD2	4:A:3209:HOH:O	2.18	0.42
1:F:2022:ALA:HA	1:F:2063:ILE:O	2.19	0.42
1:F:2034:ILE:HG12	1:F:2242:VAL:HG13	2.00	0.42
1:F:2162:PHE:HD2	1:F:2162:PHE:O	2.02	0.42
1:A:1247:GLU:HB3	4:A:3213:HOH:O	2.19	0.42
1:A:1163:TYR:HB2	1:A:1164:PRO:CD	2.50	0.42
1:A:1091:ARG:HH11	1:A:1091:ARG:HG3	1.86	0.41
1:A:1147:ILE:O	1:A:1147:ILE:HG23	2.20	0.41
1:F:2016:LEU:HD22	1:F:2063:ILE:CD1	2.50	0.41
1:F:2235:LYS:HB2	1:F:2235:LYS:NZ	2.29	0.41
1:F:2236:GLN:HE21	1:F:2236:GLN:HB2	1.65	0.41
1:A:1093:GLY:O	1:A:1217:ALA:HA	2.21	0.41
1:F:2186:GLU:OE2	1:F:2190:MET:HE1	2.20	0.41
1:A:1037:LEU:HD22	1:A:1037:LEU:HA	1.92	0.40
1:A:1028:PRO:HG3	1:A:1049:GLU:OE2	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	248/253 (98%)	234 (94%)	10 (4%)	4 (2%)	7	1
1	F	212/253 (84%)	210 (99%)	1 (0%)	1 (0%)	24	11
All	All	460/506 (91%)	444 (96%)	11 (2%)	5 (1%)	11	3

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1235	LYS
1	F	2184	MET
1	A	1163	TYR
1	A	1231	ALA
1	A	1233	THR

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	199/202 (98%)	192 (96%)	7 (4%)	32	11
1	F	171/202 (85%)	162 (95%)	9 (5%)	20	4
All	All	370/404 (92%)	354 (96%)	16 (4%)	26	7

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1005	ASP
1	A	1029	GLU
1	A	1037	LEU
1	A	1049	GLU
1	A	1147	ILE
1	A	1185	GLU
1	A	1194	ASN
1	F	2092	ILE
1	F	2161	THR
1	F	2184	MET
1	F	2186	GLU
1	F	2195	TYR
1	F	2196	GLU
1	F	2197	MET
1	F	2235	LYS
1	F	2236	GLN

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

Mol	Chain	Res	Type
1	A	1083	GLN
1	A	1103	ASN
1	A	1122	HIS
1	A	1166	GLN
1	A	1194	ASN
1	A	1209	GLN
1	A	1225	GLN
1	A	1226	GLN
1	F	2236	GLN
1	F	2240	HIS

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](#)

2 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	GOL	A	5001	-	5,5,5	4.71	5 (100%)	5,5,5	6.00	3 (60%)
3	SO4	F	4001	-	4,4,4	0.56	0	6,6,6	0.18	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
2	GOL	A	5001	-	-	2/4/4/4	-

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	5001	GOL	C3-C2	-7.79	1.22	1.51
2	A	5001	GOL	O1-C1	5.00	1.63	1.42
2	A	5001	GOL	O3-C3	3.49	1.57	1.42
2	A	5001	GOL	C1-C2	-2.69	1.41	1.51
2	A	5001	GOL	O2-C2	-2.44	1.36	1.43

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	5001	GOL	O3-C3-C2	10.88	159.34	110.38
2	A	5001	GOL	O2-C2-C3	7.13	138.68	109.18
2	A	5001	GOL	O1-C1-C2	3.26	125.04	110.38

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	5001	GOL	O1-C1-C2-C3
2	A	5001	GOL	C1-C2-C3-O3

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	250/253 (98%)	0.46	23 (9%) 14 16	11, 21, 47, 78	0
1	F	218/253 (86%)	0.17	11 (5%) 34 40	10, 19, 41, 92	0
All	All	468/506 (92%)	0.32	34 (7%) 21 25	10, 20, 46, 92	0

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	2162	PHE	5.1
1	A	1233	THR	4.5
1	A	1231	ALA	4.4
1	A	1007	PHE	4.3
1	F	2184	MET	4.0
1	F	2236	GLN	3.9
1	A	1170	ASP	3.8
1	A	1232	GLU	3.8
1	F	2195	TYR	3.6
1	F	2237	THR	3.2
1	A	1228	ILE	3.0
1	A	1234	MET	3.0
1	A	1253	LEU	3.0
1	A	1147	ILE	2.9
1	A	1042	VAL	2.8
1	F	2185	GLU	2.7
1	A	1040	LYS	2.7
1	F	2183	SER	2.6
1	F	2221	VAL	2.6
1	A	1006	VAL	2.5
1	A	1041	PRO	2.5
1	A	1059	GLY	2.5
1	F	2161	THR	2.5
1	A	1033	LYS	2.5

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Mol	Chain	Res	Type	RSRZ
1	F	2235	LYS	2.5
1	A	1048	ARG	2.5
1	A	1005	ASP	2.3
1	A	1229	PRO	2.3
1	A	1230	ASN	2.2
1	A	1185	GLU	2.2
1	A	1051	THR	2.2
1	A	1037	LEU	2.1
1	A	1236	GLN	2.1
1	F	2238	GLU	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
2	GOL	A	5001	6/6	0.71	0.18	20,20,20,20	0
3	SO4	F	4001	5/5	0.87	0.28	20,20,20,20	0

6.5 Other polymers [i](#)

There are no such residues in this entry.