



Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 02:34 pm BST

PDB ID : 3TIV
Title : Crystal structure of subunit B mutant N157A of the A1AO ATP synthase
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Deposited on : 2011-08-22
Resolution : 1.75 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

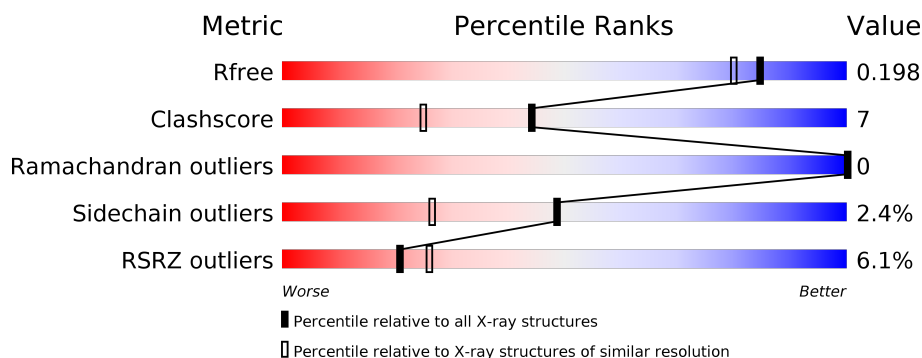
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.75 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}	130704	2340 (1.76-1.76)
Clashscore	141614	2466 (1.76-1.76)
Ramachandran outliers	138981	2437 (1.76-1.76)
Sidechain outliers	138945	2437 (1.76-1.76)
RSRZ outliers	127900	2298 (1.76-1.76)

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 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	460	<div> <div>6%</div> <div> <div></div> <div>85%</div> <div>10%</div> <div>5%</div> </div> </div>
1	B	460	<div> <div>5%</div> <div> <div></div> <div>81%</div> <div>12%</div> <div>7%</div> </div> </div>

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 7957 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 V-type ATP synthase beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	437	Total	C	N	O	S	0	14	0
			3435	2175	597	652	11			
1	B	430	Total	C	N	O	S	0	14	0
			3398	2156	590	641	11			

There are 4 discrepancies between the modelled and reference sequences:

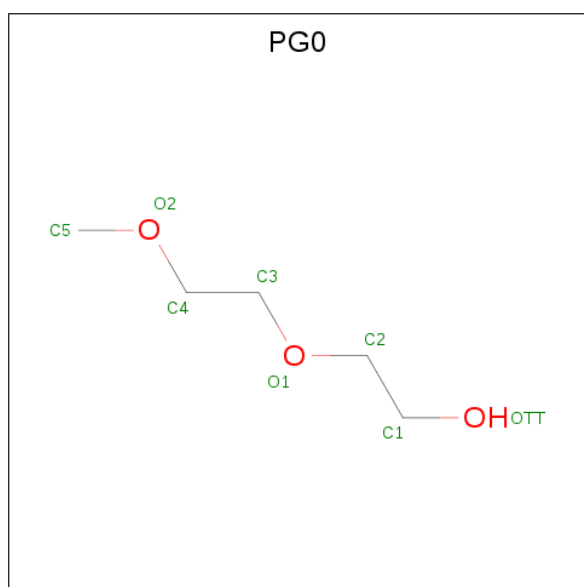
Chain	Residue	Modelled	Actual	Comment	Reference
A	2	VAL	ALA	SEE REMARK 999	UNP Q60187
A	157	ALA	ASN	ENGINEERED MUTATION	UNP Q60187
B	2	VAL	ALA	SEE REMARK 999	UNP Q60187
B	157	ALA	ASN	ENGINEERED MUTATION	UNP Q60187

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



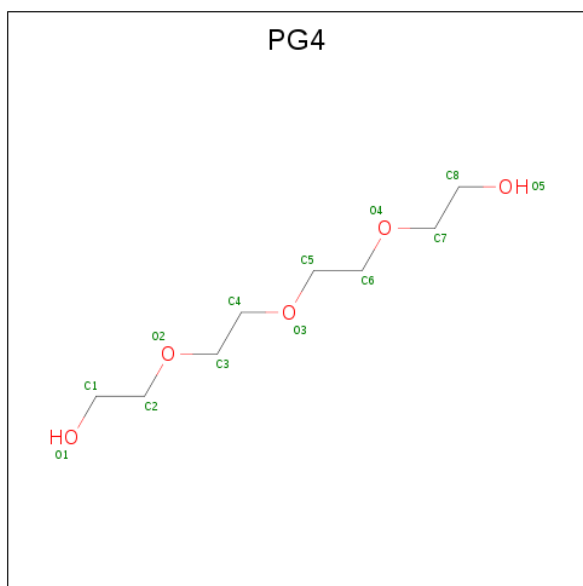
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 6 3 3	0	0
2	A	1	Total C O 6 3 3	0	0
2	A	1	Total C O 6 3 3	0	0
2	A	1	Total C O 6 3 3	0	0
2	A	1	Total C O 6 3 3	0	0
2	A	1	Total C O 6 3 3	0	0
2	A	1	Total C O 6 3 3	0	0
2	A	1	Total C O 6 3 3	0	0
2	B	1	Total C O 6 3 3	0	0
2	B	1	Total C O 6 3 3	0	0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

- Molecule 3 is 2-(2-METHOXYETHOXY)ETHANOL (three-letter code: PG0) (formula: $C_5H_{12}O_3$).



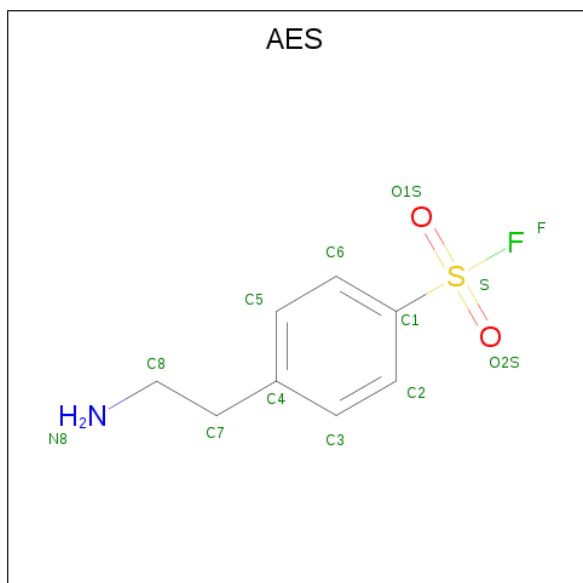
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			8	5	3		

- Molecule 4 is TETRAETHYLENE GLYCOL (three-letter code: PG4) (formula: $C_8H_{18}O_5$).



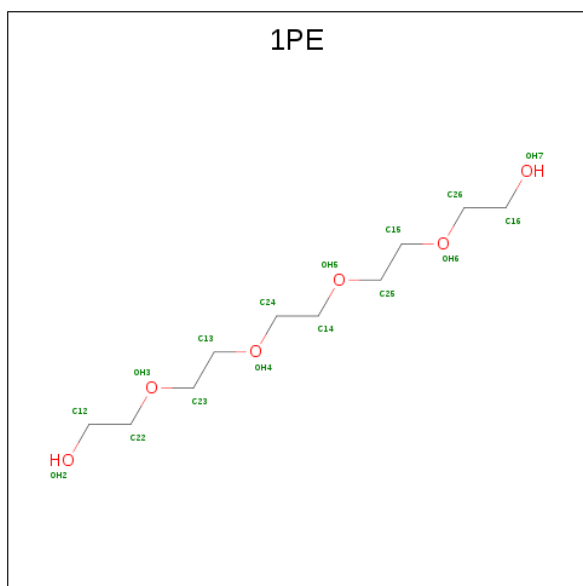
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			13	8	5		

- Molecule 5 is 4-(2-AMINOETHYL)BENZENESULFONYL FLUORIDE (three-letter code: AES) (formula: $C_8H_{10}FNO_2S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	B	1	Total	C	F	N	O	S	
			13	8	1	1	2	1	

- Molecule 6 is PENTAETHYLENE GLYCOL (three-letter code: 1PE) (formula: $C_{10}H_{22}O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	B	1	Total	C	O		
			16	10	6	0	0

- Molecule 7 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	B	1	Total	Cl		
			1	1	0	0

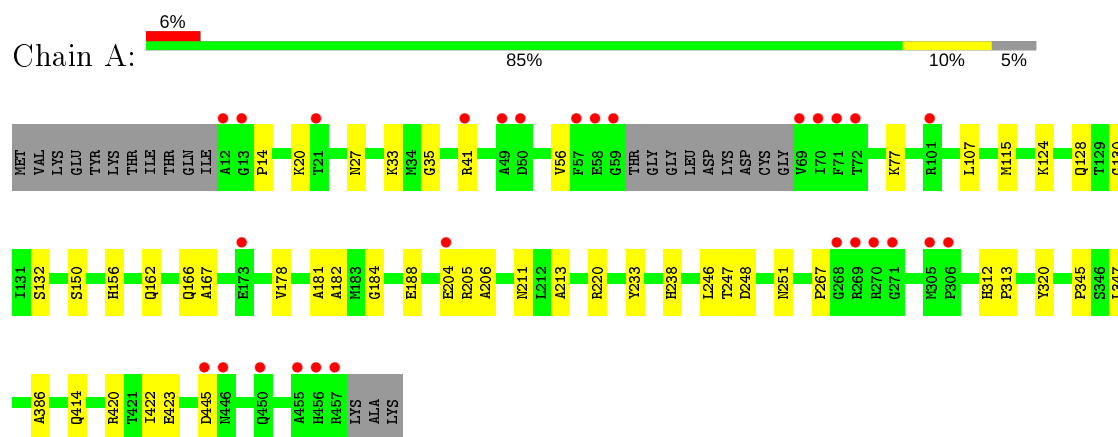
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	611	Total	O		
			611	611	0	0
8	B	396	Total	O		
			396	396	0	0

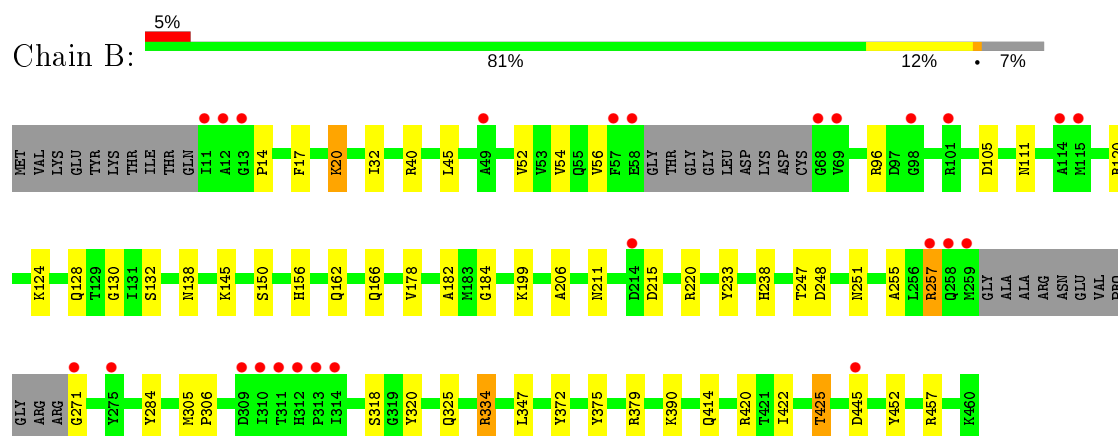
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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: V-type ATP synthase beta chain



• Molecule 1: V-type ATP synthase beta chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	73.72Å 95.77Å 130.16Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.46 – 1.75 19.46 – 1.75	Depositor EDS
% Data completeness (in resolution range)	99.7 (19.46-1.75) 99.7 (19.46-1.75)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.13 (at 1.76Å)	Xtriage
Refinement program	REFMAC 5.5.0072	Depositor
R, R_{free}	0.167 , 0.199 0.167 , 0.198	Depositor DCC
R_{free} test set	4664 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	16.6	Xtriage
Anisotropy	0.038	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 59.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7957	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.08% 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, AES, CL, 1PE, PG4, PG0

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	0.43	0/3541	0.59	0/4798
1	B	0.45	0/3502	0.61	0/4740
All	All	0.44	0/7043	0.60	0/9538

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

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	3435	0	3472	41	0
1	B	3398	0	3448	56	0
2	A	48	0	64	4	0
2	B	18	0	24	1	0
3	A	8	0	12	1	0
4	B	13	0	18	0	0
5	B	13	0	10	3	0
6	B	16	0	22	0	0
7	B	1	0	0	0	0
8	A	611	0	0	4	0
8	B	396	0	0	7	0
All	All	7957	0	7070	92	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 7.

All (92) 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:B:20:LYS:HE2	5:B:1474:AES:O2S	1.47	1.13
1:B:40[B]:ARG:HH11	1:B:40[B]:ARG:HG2	0.95	1.07
1:B:40[B]:ARG:HG2	1:B:40[B]:ARG:NH1	1.76	0.94
1:B:20:LYS:HE2	5:B:1474:AES:S	2.12	0.89
1:B:334[A]:ARG:HH11	1:B:334[A]:ARG:CG	1.86	0.89
1:B:40[B]:ARG:HH11	1:B:40[B]:ARG:CG	1.85	0.88
1:B:20:LYS:CE	5:B:1474:AES:O2S	2.23	0.86
1:A:20:LYS:HE2	2:A:468:GOL:H2	1.57	0.86
1:B:184:GLY:H	1:B:211:ASN:HD22	1.21	0.85
1:B:257:ARG:HH11	1:B:257:ARG:HG3	1.43	0.82
1:A:184:GLY:HA3	1:A:251[B]:ASN:HD21	1.44	0.81
1:B:17:PHE:H	1:B:111:ASN:HD21	1.31	0.78
1:B:334[A]:ARG:HH11	1:B:334[A]:ARG:HG2	1.49	0.76
1:A:35:GLY:HA2	1:A:77[A]:LYS:NZ	2.01	0.74
1:A:184:GLY:HA3	1:A:251[B]:ASN:ND2	2.02	0.73
1:A:35:GLY:HA2	1:A:77[A]:LYS:HZ1	1.55	0.71
1:A:312:HIS:HE1	8:A:558:HOH:O	1.74	0.70
1:A:184:GLY:H	1:A:211:ASN:HD22	1.37	0.70
1:A:213:ALA:HB3	2:A:465:GOL:H12	1.77	0.67
1:B:132:SER:H	1:B:414:GLN:HE22	1.43	0.66
1:B:220:ARG:HD3	1:B:251[A]:ASN:OD1	1.96	0.66
1:A:132:SER:H	1:A:414:GLN:HE22	1.45	0.65
1:B:334[A]:ARG:NH1	1:B:334[A]:ARG:CG	2.58	0.62
1:B:199:LYS:NZ	8:B:587:HOH:O	2.33	0.62
1:B:220:ARG:HD2	1:B:255:ALA:HB2	1.85	0.59
1:B:425:THR:CG2	8:B:693:HOH:O	2.50	0.58
1:A:35:GLY:CA	1:A:77[A]:LYS:NZ	2.66	0.58
1:A:20:LYS:HE2	2:A:468:GOL:C2	2.32	0.57
1:A:20:LYS:CE	2:A:468:GOL:H2	2.32	0.57
1:B:184:GLY:H	1:B:211:ASN:ND2	1.97	0.57
1:A:267:PRO:HB2	1:B:334[B]:ARG:HD2	1.86	0.57
1:A:115:MET:H	3:A:6108:PG0:H53	1.70	0.56
1:B:20:LYS:HB2	1:B:20:LYS:HZ2	1.71	0.56
1:A:347:LEU:HD22	1:B:347:LEU:HD22	1.88	0.55
1:B:14:PRO:HG2	1:B:56:VAL:CG1	2.37	0.54
1:B:162:GLN:HE21	1:B:166:GLN:HE22	1.55	0.54
1:B:257:ARG:NH1	1:B:257:ARG:HG3	2.19	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:452:TYR:O	1:B:457[B]:ARG:HD2	2.09	0.52
1:A:162:GLN:HE21	1:A:166:GLN:HE22	1.58	0.52
1:B:150:SER:OG	1:B:156:HIS:HD2	1.94	0.50
1:A:184:GLY:H	1:A:211:ASN:ND2	2.09	0.50
1:B:40[B]:ARG:NH1	1:B:40[B]:ARG:CG	2.54	0.50
1:A:188:GLU:H	1:A:188:GLU:CD	2.15	0.50
1:B:257:ARG:HD3	1:B:271:GLY:O	2.12	0.50
1:A:77[B]:LYS:HD2	1:A:107:LEU:HB3	1.93	0.50
1:B:375:TYR:CZ	1:B:379:ARG:HD2	2.47	0.49
1:B:238:HIS:HD2	8:B:487:HOH:O	1.96	0.49
1:B:17:PHE:H	1:B:111:ASN:ND2	2.04	0.48
1:B:32:ILE:HD11	1:B:54:VAL:HG11	1.95	0.48
1:A:14:PRO:HD2	1:A:56:VAL:HG23	1.96	0.48
1:B:233:TYR:OH	1:B:238:HIS:HE1	1.95	0.48
1:B:247:THR:HA	1:B:248[B]:ASP:HA	1.67	0.47
1:B:96:ARG:NH2	1:B:215:ASP:OD1	2.46	0.47
1:A:128:GLN:NE2	1:A:422:ILE:H	2.13	0.47
1:A:14:PRO:HD2	1:A:56:VAL:CG2	2.45	0.47
1:B:178:VAL:O	1:B:206:ALA:HA	2.15	0.47
1:B:334[A]:ARG:HH11	1:B:334[A]:ARG:HG3	1.76	0.46
1:B:425:THR:HG23	8:B:693:HOH:O	2.14	0.46
1:A:386:ALA:HB1	1:B:14:PRO:HB3	1.98	0.46
1:B:372:TYR:CD1	2:B:461:GOL:H31	2.51	0.46
1:A:251[A]:ASN:ND2	8:A:777:HOH:O	2.50	0.45
1:A:167:ALA:O	1:A:205[B]:ARG:NH2	2.47	0.45
1:A:182:ALA:HA	1:A:247:THR:O	2.17	0.45
1:B:425:THR:HG21	8:B:693:HOH:O	2.16	0.45
1:B:156:HIS:HE1	8:B:509:HOH:O	2.01	0.44
1:A:220:ARG:HD3	1:A:251[B]:ASN:HD22	1.82	0.43
1:B:128:GLN:NE2	1:B:422:ILE:H	2.16	0.43
1:A:124:LYS:NZ	8:A:614:HOH:O	2.51	0.43
1:B:132:SER:H	1:B:414:GLN:NE2	2.12	0.43
1:B:150:SER:OG	1:B:156:HIS:CD2	2.71	0.43
1:A:35:GLY:CA	1:A:77[A]:LYS:HZ1	2.24	0.43
1:A:247:THR:HG22	8:A:774:HOH:O	2.19	0.43
1:B:257:ARG:HH11	1:B:257:ARG:CG	2.22	0.42
1:A:345:PRO:HG3	1:B:318:SER:HB2	2.01	0.42
1:B:182:ALA:HA	1:B:247:THR:O	2.19	0.42
1:B:14:PRO:HG2	1:B:56:VAL:HG13	2.00	0.42
1:B:305:MET:HA	1:B:306:PRO:HD3	1.94	0.42
1:A:247:THR:HA	1:A:248[B]:ASP:HA	1.73	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:27:ASN:HA	1:A:27:ASN:HD22	1.76	0.41
1:A:267:PRO:HG2	1:A:313:PRO:HB2	2.03	0.41
1:A:267:PRO:HB2	1:B:334[B]:ARG:CD	2.48	0.41
1:A:130:GLY:HA2	1:A:420:ARG:O	2.21	0.41
1:B:120:ARG:NH2	8:B:644:HOH:O	2.53	0.41
1:B:145:LYS:HD3	1:B:284:TYR:O	2.21	0.41
1:A:423:GLU:CD	1:A:423:GLU:H	2.24	0.41
1:A:178:VAL:O	1:A:206:ALA:HA	2.20	0.41
1:A:181:ALA:HB3	1:A:246:LEU:HD23	2.02	0.41
1:A:233:TYR:OH	1:A:238:HIS:HE1	2.03	0.41
1:B:130:GLY:HA2	1:B:420:ARG:O	2.20	0.41
1:B:20:LYS:HE3	1:B:52:VAL:HG22	2.02	0.40
1:A:150:SER:OG	1:A:156:HIS:HD2	2.05	0.40
1:B:184:GLY:N	1:B:211:ASN:HD22	2.03	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	447/460 (97%)	435 (97%)	12 (3%)	0	100	100
1	B	438/460 (95%)	426 (97%)	12 (3%)	0	100	100
All	All	885/920 (96%)	861 (97%)	24 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	369/374 (99%)	363 (98%)	6 (2%)	62	45
1	B	365/374 (98%)	350 (96%)	15 (4%)	30	10
All	All	734/748 (98%)	713 (97%)	21 (3%)	49	19

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

Mol	Chain	Res	Type
1	A	33	LYS
1	A	41	ARG
1	A	204[A]	GLU
1	A	204[B]	GLU
1	A	320	TYR
1	A	445	ASP
1	B	20	LYS
1	B	45	LEU
1	B	105	ASP
1	B	124	LYS
1	B	138[A]	ASN
1	B	138[B]	ASN
1	B	257	ARG
1	B	320	TYR
1	B	325[A]	GLN
1	B	325[B]	GLN
1	B	334[A]	ARG
1	B	334[B]	ARG
1	B	390	LYS
1	B	425	THR
1	B	445	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	27	ASN
1	A	116	ASN
1	A	128	GLN
1	A	156	HIS
1	A	166	GLN
1	A	211	ASN

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Mol	Chain	Res	Type
1	A	238	HIS
1	A	299	GLN
1	A	312	HIS
1	A	414	GLN
1	A	417	ASN
1	A	431	GLN
1	B	55	GLN
1	B	111	ASN
1	B	128	GLN
1	B	156	HIS
1	B	166	GLN
1	B	211	ASN
1	B	238	HIS
1	B	299	GLN
1	B	414	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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 16 ligands modelled in this entry, 1 is monoatomic - leaving 15 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
2	GOL	A	468	-	5,5,5	0.39	0	5,5,5	0.35	0
2	GOL	A	467	-	5,5,5	0.34	0	5,5,5	0.26	0
2	GOL	B	461	-	5,5,5	0.33	0	5,5,5	0.50	0
5	AES	B	1474	-	11,13,13	0.68	0	16,18,18	1.46	1 (6%)
2	GOL	A	464	-	5,5,5	0.34	0	5,5,5	0.34	0
2	GOL	A	463	-	5,5,5	0.35	0	5,5,5	0.30	0
6	1PE	B	463	-	15,15,15	0.50	0	14,14,14	0.57	0
2	GOL	B	464	-	5,5,5	0.35	0	5,5,5	0.19	0
4	PG4	B	6112	-	12,12,12	0.49	0	11,11,11	0.50	0
2	GOL	A	465	-	5,5,5	0.35	0	5,5,5	0.57	0
2	GOL	A	461	-	5,5,5	0.33	0	5,5,5	0.30	0
3	PG0	A	6108	-	7,7,7	0.52	0	6,6,6	0.52	0
2	GOL	A	466	-	5,5,5	0.34	0	5,5,5	0.42	0
2	GOL	A	462	-	5,5,5	0.31	0	5,5,5	0.41	0
2	GOL	B	462	-	5,5,5	0.45	0	5,5,5	0.24	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	468	-	-	2/4/4/4	-
2	GOL	A	467	-	-	4/4/4/4	-
2	GOL	B	461	-	-	2/4/4/4	-
5	AES	B	1474	-	-	9/9/9/9	0/1/1/1
2	GOL	A	464	-	-	4/4/4/4	-
2	GOL	A	463	-	-	4/4/4/4	-
6	1PE	B	463	-	-	4/13/13/13	-
2	GOL	B	464	-	-	0/4/4/4	-
4	PG4	B	6112	-	-	3/10/10/10	-
2	GOL	A	465	-	-	2/4/4/4	-
2	GOL	A	461	-	-	0/4/4/4	-
3	PG0	A	6108	-	-	4/5/5/5	-
2	GOL	A	466	-	-	2/4/4/4	-
2	GOL	A	462	-	-	4/4/4/4	-
2	GOL	B	462	-	-	0/4/4/4	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	1474	AES	O2S-S-O1S	-4.48	105.09	119.07

There are no chirality outliers.

All (44) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	467	GOL	O1-C1-C2-C3
2	A	467	GOL	C1-C2-C3-O3
2	B	461	GOL	O1-C1-C2-C3
5	B	1474	AES	C6-C1-S-F
5	B	1474	AES	C4-C7-C8-N8
2	A	464	GOL	O1-C1-C2-C3
2	A	463	GOL	O1-C1-C2-C3
2	A	463	GOL	C1-C2-C3-O3
2	A	465	GOL	O1-C1-C2-O2
2	A	462	GOL	O1-C1-C2-O2
3	A	6108	PG0	O1-C3-C4-O2
2	A	463	GOL	O1-C1-C2-O2
6	B	463	1PE	OH7-C16-C26-OH6
4	B	6112	PG4	O1-C1-C2-O2
2	A	468	GOL	C1-C2-C3-O3
2	A	464	GOL	C1-C2-C3-O3
2	A	465	GOL	O1-C1-C2-C3
2	A	466	GOL	C1-C2-C3-O3
2	A	462	GOL	O1-C1-C2-C3
2	A	462	GOL	C1-C2-C3-O3
3	A	6108	PG0	OTT-C1-C2-O1
2	A	464	GOL	O1-C1-C2-O2
2	A	463	GOL	O2-C2-C3-O3
2	A	466	GOL	O2-C2-C3-O3
2	A	468	GOL	O2-C2-C3-O3
2	B	461	GOL	O1-C1-C2-O2
5	B	1474	AES	C2-C1-S-O1S
5	B	1474	AES	C2-C1-S-O2S
5	B	1474	AES	C6-C1-S-O1S
5	B	1474	AES	C6-C1-S-O2S
5	B	1474	AES	C5-C4-C7-C8
2	A	467	GOL	O1-C1-C2-O2
2	A	462	GOL	O2-C2-C3-O3
5	B	1474	AES	C3-C4-C7-C8
3	A	6108	PG0	C3-C4-O2-C5

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Mol	Chain	Res	Type	Atoms
5	B	1474	AES	C2-C1-S-F
2	A	467	GOL	O2-C2-C3-O3
2	A	464	GOL	O2-C2-C3-O3
6	B	463	1PE	C15-C25-OH5-C14
4	B	6112	PG4	C8-C7-O4-C6
6	B	463	1PE	C25-C15-OH6-C26
6	B	463	1PE	C16-C26-OH6-C15
3	A	6108	PG0	C4-C3-O1-C2
4	B	6112	PG4	C5-C6-O4-C7

There are no ring outliers.

5 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	468	GOL	3	0
2	B	461	GOL	1	0
5	B	1474	AES	3	0
2	A	465	GOL	1	0
3	A	6108	PG0	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	437/460 (95%)	0.12	28 (6%)	19 25	8, 16, 31, 44	0
1	B	430/460 (93%)	0.14	25 (5%)	23 28	7, 15, 30, 42	0
All	All	867/920 (94%)	0.13	53 (6%)	21 26	7, 16, 30, 44	0

All (53) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	69	VAL	12.5
1	B	11	ILE	10.4
1	B	68	GLY	8.5
1	A	12	ALA	6.6
1	B	69	VAL	6.4
1	B	12	ALA	6.2
1	B	258	GLN	6.2
1	B	310	ILE	6.1
1	A	270	ARG	5.9
1	B	311	THR	5.7
1	B	58	GLU	5.0
1	B	312	HIS	4.8
1	A	58	GLU	4.7
1	A	70	ILE	4.6
1	A	269	ARG	3.9
1	A	41	ARG	3.7
1	A	57	PHE	3.5
1	B	114	ALA	3.5
1	B	271	GLY	3.3
1	A	71	PHE	3.2
1	A	455	ALA	2.9
1	B	275	TYR	2.9
1	B	214	ASP	2.9
1	A	446	ASN	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	309	ASP	2.7
1	B	115	MET	2.7
1	A	204[A]	GLU	2.6
1	B	13	GLY	2.6
1	A	59	GLY	2.5
1	B	314	ILE	2.4
1	B	57	PHE	2.4
1	B	259	MET	2.3
1	A	173	GLU	2.3
1	B	101	ARG	2.2
1	A	21	THR	2.2
1	A	50	ASP	2.2
1	A	457	ARG	2.2
1	A	306	PRO	2.2
1	B	313	PRO	2.2
1	A	445	ASP	2.2
1	A	49	ALA	2.1
1	A	268	GLY	2.1
1	A	271	GLY	2.1
1	A	72	THR	2.1
1	B	49	ALA	2.1
1	A	456	HIS	2.1
1	B	257	ARG	2.1
1	A	13	GLY	2.1
1	A	450	GLN	2.0
1	A	305	MET	2.0
1	A	101	ARG	2.0
1	B	445	ASP	2.0
1	B	98	GLY	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands

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(Å ²)	Q<0.9
2	GOL	A	463	6/6	0.46	0.27	62,62,62,62	0
2	GOL	A	467	6/6	0.67	0.22	45,47,47,48	0
2	GOL	A	464	6/6	0.68	0.23	35,42,43,43	0
3	PG0	A	6108	8/8	0.69	0.24	70,70,70,70	0
2	GOL	A	468	6/6	0.75	0.30	53,54,54,55	0
2	GOL	A	466	6/6	0.77	0.29	32,39,41,42	0
2	GOL	A	462	6/6	0.83	0.14	24,31,33,35	0
2	GOL	A	465	6/6	0.84	0.25	25,31,32,33	0
2	GOL	A	461	6/6	0.85	0.16	19,25,27,27	0
6	1PE	B	463	16/16	0.85	0.17	35,37,46,47	0
2	GOL	B	464	6/6	0.86	0.17	56,57,58,58	0
2	GOL	B	461	6/6	0.89	0.18	17,28,29,30	0
4	PG4	B	6112	13/13	0.90	0.14	41,43,45,46	0
5	AES	B	1474	13/13	0.91	0.15	20,22,26,30	13
2	GOL	B	462	6/6	0.92	0.12	15,21,24,28	0
7	CL	B	465	1/1	0.98	0.04	20,20,20,20	0

6.5 Other polymers

There are no such residues in this entry.