



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 8, 2021 – 01:04 PM BST

PDB ID : 7NZJ
Title : Structure of bsTrmB apo
Authors : Blersch, K.F.; Ficner, R.; Neumann, P.
Deposited on : 2021-03-24
Resolution : 1.98 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 : 2.23.1
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) : 2.23.1

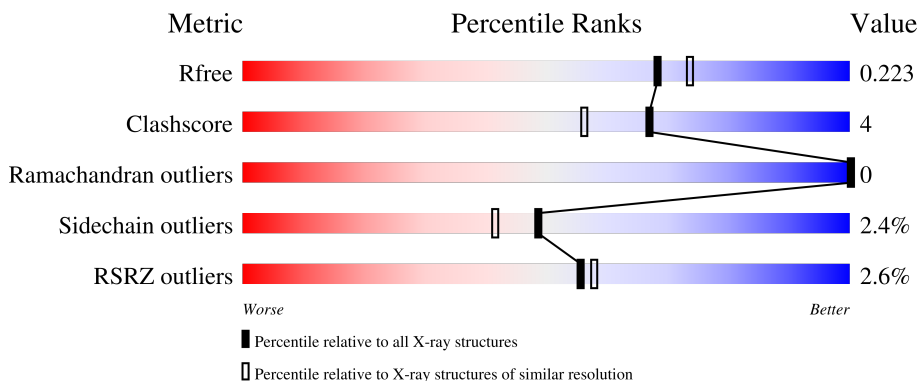
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.98 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	11647 (2.00-1.96)
Clashscore	141614	1014 (1.98-1.98)
Ramachandran outliers	138981	1006 (1.98-1.98)
Sidechain outliers	138945	1006 (1.98-1.98)
RSRZ outliers	127900	11410 (2.00-1.96)

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	213	 3% 87% 9% . .
1	B	213	 3% 84% 11% 5%
1	C	213	 2% 87% 8% . .
1	D	213	 3% 83% 12% 5%
1	E	213	 2% 85% 11% .

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Mol	Chain	Length	Quality of chain
1	F	213	 <p>A horizontal bar chart representing the quality of the chain. The bar is divided into four segments: a small red segment at the beginning labeled '2%', a large green segment labeled '85%', a yellow segment labeled '11%', and a small grey segment at the end. A small black dot is visible at the far right end of the bar.</p>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10868 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 tRNA (guanine-N(7)-)-methyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	206	Total 1667	C 1067	N 278	O 319	S 3	0	0	0
1	B	203	Total 1641	C 1050	N 274	O 314	S 3	0	0	0
1	C	205	Total 1654	C 1057	N 276	O 318	S 3	0	0	0
1	D	203	Total 1641	C 1050	N 274	O 314	S 3	0	0	0
1	E	206	Total 1668	C 1068	N 278	O 319	S 3	0	0	0
1	F	205	Total 1654	C 1057	N 276	O 318	S 3	0	0	0

- Molecule 2 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total 1	Na 1	0	0

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	C O	0	0
			6	3 3		
3	C	1	Total	C O	0	0
			6	3 3		
3	D	1	Total	C O	0	0
			6	3 3		
3	D	1	Total	C O	0	0
			6	3 3		
3	E	1	Total	C O	0	0
			6	3 3		

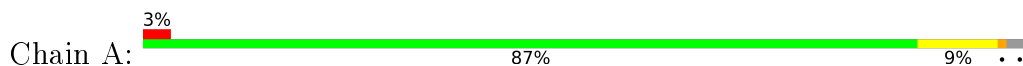
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	164	Total	O	0	0
			164	164		
4	B	144	Total	O	0	0
			144	144		
4	C	150	Total	O	0	0
			150	150		
4	D	146	Total	O	0	0
			146	146		
4	E	143	Total	O	0	0
			143	143		
4	F	165	Total	O	0	0
			165	165		

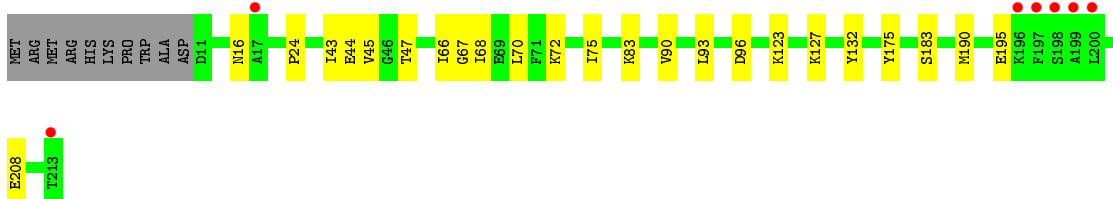
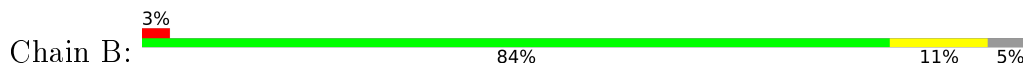
3 Residue-property plots

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.

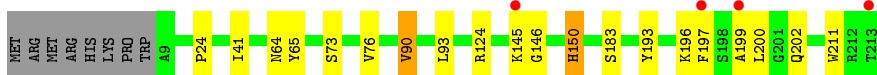
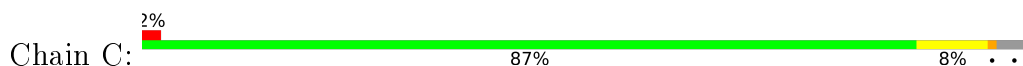
- Molecule 1: tRNA (guanine-N(7)-)-methyltransferase



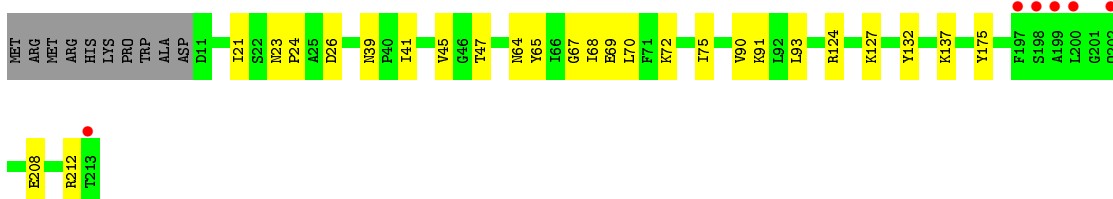
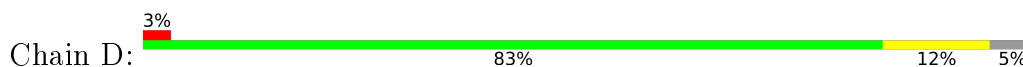
- Molecule 1: tRNA (guanine-N(7)-)-methyltransferase



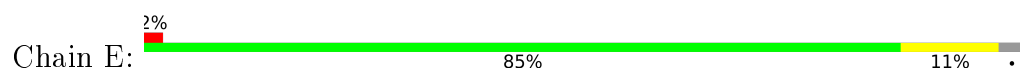
- Molecule 1: tRNA (guanine-N(7)-)-methyltransferase



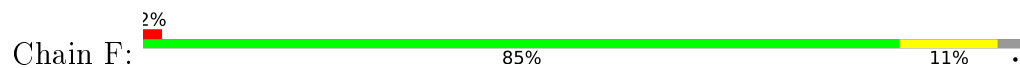
- Molecule 1: tRNA (guanine-N(7)-)-methyltransferase



- Molecule 1: tRNA (guanine-N(7)-)-methyltransferase



- Molecule 1: tRNA (guanine-N(7)-)-methyltransferase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	42.62Å 103.68Å 103.69Å 118.12° 97.91° 97.84°	Depositor
Resolution (Å)	44.53 – 1.98 44.49 – 1.98	Depositor EDS
% Data completeness (in resolution range)	91.7 (44.53-1.98) 91.4 (44.49-1.98)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.50 (at 1.98Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.187 , 0.223 0.193 , 0.223	Depositor DCC
R_{free} test set	5048 reflections (5.24%)	wwPDB-VP
Wilson B-factor (Å ²)	35.2	Xtriage
Anisotropy	0.552	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 36.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.458 for h,l,-h-k-l 0.458 for h,-h-k-l,k 0.010 for -h,-l,-k 0.010 for -h,-k,h+k+l 0.011 for -h,h+k+l,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	10868	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.23% 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: NA, 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	0.68	0/1707	0.71	0/2305
1	B	0.69	0/1679	0.71	0/2266
1	C	0.66	0/1692	0.71	0/2284
1	D	0.69	0/1679	0.70	0/2266
1	E	0.67	0/1708	0.72	0/2307
1	F	0.69	0/1692	0.72	0/2284
All	All	0.68	0/10157	0.71	0/13712

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1667	0	1620	10	0
1	B	1641	0	1605	14	0
1	C	1654	0	1614	11	0
1	D	1641	0	1605	16	0
1	E	1668	0	1624	15	0
1	F	1654	0	1614	13	0
2	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	6	0	8	0	0
3	C	6	0	8	1	0
3	D	12	0	16	0	0
3	E	6	0	8	0	0
4	A	164	0	0	0	0
4	B	144	0	0	1	0
4	C	150	0	0	3	0
4	D	146	0	0	2	0
4	E	143	0	0	2	0
4	F	165	0	0	2	0
All	All	10868	0	9722	76	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 4.

All (76) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:199:ALA:HB1	4:C:545:HOH:O	1.69	0.92
1:F:127:LYS:NZ	4:F:301:HOH:O	2.24	0.71
1:A:156:ARG:NH1	1:B:183:SER:OG	2.24	0.69
1:D:47:THR:HG21	1:D:67:GLY:HA3	1.77	0.67
1:A:38:ASP:OD1	1:E:125:HIS:ND1	2.26	0.66
1:D:212:ARG:NH1	4:D:502:HOH:O	2.26	0.64
1:E:24:PRO:HG3	1:E:93:LEU:HD22	1.85	0.59
1:D:23:ASN:ND2	1:D:26:ASP:OD1	2.34	0.58
1:B:47:THR:HG21	1:B:67:GLY:HA3	1.86	0.58
1:E:47:THR:HG21	1:E:67:GLY:HA3	1.85	0.57
1:A:86:GLU:O	1:A:88:GLN:NE2	2.37	0.56
1:F:58:LYS:HD2	1:F:85:SER:HB2	1.89	0.55
1:F:127:LYS:HD2	1:F:127:LYS:H	1.70	0.55
1:C:183:SER:HB3	3:C:301:GOL:H12	1.89	0.55
1:B:47:THR:HG21	1:B:67:GLY:CA	2.38	0.54
1:F:197:PHE:HE1	1:F:202:GLN:HB3	1.73	0.54
1:C:150:HIS:HD2	4:C:489:HOH:O	1.91	0.53
1:A:9:ALA:HB1	1:A:76:VAL:HG21	1.89	0.53
1:B:123:LYS:HE2	1:B:123:LYS:HA	1.89	0.53
1:D:137:LYS:NZ	1:F:84:ASP:O	2.42	0.53
1:F:68:ILE:HD13	1:F:93:LEU:HB2	1.91	0.53
1:C:146:GLY:O	1:C:211:TRP:HB2	2.08	0.52
1:A:197:PHE:HE1	1:A:202:GLN:HB2	1.75	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:30:LYS:NZ	4:E:508:HOH:O	2.43	0.50
1:D:21:ILE:HD13	1:D:91:LYS:HB3	1.92	0.50
1:C:193:TYR:HA	1:C:196:LYS:HE2	1.94	0.49
1:D:175:TYR:HB3	1:D:208:GLU:HB2	1.93	0.49
1:B:45:VAL:HA	1:B:68:ILE:HG12	1.95	0.49
1:C:65:TYR:HB2	1:C:90:VAL:HG13	1.94	0.49
1:C:73:SER:O	1:C:76:VAL:HG22	2.12	0.49
1:E:43:ILE:HA	1:E:66:ILE:O	2.12	0.49
1:E:181:HIS:CD2	1:E:198:SER:HB2	2.49	0.48
1:B:175:TYR:HB3	1:B:208:GLU:HB2	1.96	0.48
1:C:200:LEU:N	4:C:408:HOH:O	2.47	0.48
1:F:202:GLN:HG2	1:F:203:PRO:HD2	1.95	0.48
1:A:111:ARG:HD2	1:A:150:HIS:CD2	2.50	0.47
1:A:197:PHE:CE1	1:A:202:GLN:HB2	2.50	0.47
1:D:47:THR:HG21	1:D:67:GLY:CA	2.44	0.46
1:D:45:VAL:HA	1:D:68:ILE:HG12	1.97	0.46
1:E:53:ILE:HD11	1:E:65:TYR:HB3	1.98	0.46
1:E:47:THR:HG21	1:E:67:GLY:CA	2.46	0.45
1:F:24:PRO:HB3	1:F:93:LEU:HD22	1.98	0.45
1:A:111:ARG:HD3	1:A:148:SER:OG	2.16	0.45
1:F:83:LYS:O	1:F:85:SER:O	2.35	0.45
1:B:16:ASN:OD1	1:B:83:LYS:HD2	2.17	0.44
1:D:127:LYS:HA	1:D:132:TYR:CD2	2.52	0.44
1:E:16:ASN:OD1	1:E:83:LYS:HD2	2.17	0.44
1:A:65:TYR:O	1:A:90:VAL:HA	2.17	0.44
1:B:190:MET:HE2	1:B:195:GLU:HA	1.98	0.44
1:C:197:PHE:CE1	1:C:202:GLN:HB2	2.52	0.44
1:D:45:VAL:HG13	1:D:68:ILE:HD11	1.98	0.44
1:B:24:PRO:HB3	1:B:93:LEU:HD22	1.99	0.43
1:E:70:LEU:HD13	1:E:96:ASP:HB3	1.99	0.43
1:B:44:GLU:HB3	1:B:47:THR:HG22	1.99	0.43
1:F:43:ILE:HA	1:F:66:ILE:O	2.18	0.43
1:D:24:PRO:HB3	1:D:93:LEU:HD22	2.00	0.43
1:E:212:ARG:NH2	4:E:515:HOH:O	2.48	0.43
1:D:72:LYS:HA	1:D:75:ILE:HG22	2.01	0.42
1:A:43:ILE:HA	1:A:66:ILE:O	2.20	0.42
1:D:39:ASN:ND2	4:D:504:HOH:O	2.27	0.42
1:C:24:PRO:HB3	1:C:93:LEU:HD22	2.02	0.42
1:D:69:GLU:HB3	1:D:75:ILE:HD12	2.01	0.41
1:E:119:PRO:HA	1:E:158:LEU:HD13	2.02	0.41
1:E:111:ARG:HD2	1:E:150:HIS:CD2	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:65:TYR:O	1:D:90:VAL:HA	2.19	0.41
1:D:41:ILE:HA	1:D:64:ASN:O	2.21	0.41
1:F:9:ALA:N	1:F:72:LYS:HZ3	2.18	0.41
1:B:127:LYS:HA	1:B:132:TYR:CD2	2.55	0.41
1:B:43:ILE:HA	1:B:66:ILE:O	2.21	0.41
1:C:41:ILE:HA	1:C:64:ASN:O	2.21	0.41
1:F:23:ASN:ND2	1:F:26:ASP:OD1	2.44	0.40
1:E:170:GLY:HA2	4:F:310:HOH:O	2.20	0.40
1:B:72:LYS:HA	1:B:75:ILE:HG22	2.03	0.40
1:B:123:LYS:HD3	4:B:606:HOH:O	2.20	0.40
1:E:127:LYS:HA	1:E:132:TYR:CD2	2.56	0.40
1:F:197:PHE:CE1	1:F:202:GLN:HB3	2.56	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	204/213 (96%)	198 (97%)	6 (3%)	0	100	100
1	B	201/213 (94%)	198 (98%)	3 (2%)	0	100	100
1	C	203/213 (95%)	200 (98%)	3 (2%)	0	100	100
1	D	201/213 (94%)	197 (98%)	4 (2%)	0	100	100
1	E	204/213 (96%)	200 (98%)	4 (2%)	0	100	100
1	F	203/213 (95%)	198 (98%)	5 (2%)	0	100	100
All	All	1216/1278 (95%)	1191 (98%)	25 (2%)	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	179/187 (96%)	171 (96%)	8 (4%)	27	15
1	B	178/187 (95%)	175 (98%)	3 (2%)	60	53
1	C	179/187 (96%)	175 (98%)	4 (2%)	52	46
1	D	178/187 (95%)	176 (99%)	2 (1%)	73	70
1	E	180/187 (96%)	178 (99%)	2 (1%)	73	70
1	F	179/187 (96%)	172 (96%)	7 (4%)	32	19
All	All	1073/1122 (96%)	1047 (98%)	26 (2%)	49	41

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	TRP
1	A	70	LEU
1	A	74	VAL
1	A	90	VAL
1	A	123	LYS
1	A	197	PHE
1	A	209	VAL
1	A	213	THR
1	B	70	LEU
1	B	90	VAL
1	B	96	ASP
1	C	90	VAL
1	C	124	ARG
1	C	145	LYS
1	C	150	HIS
1	D	70	LEU
1	D	124	ARG
1	E	22	SER
1	E	111	ARG
1	F	70	LEU
1	F	71	PHE
1	F	74	VAL

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Mol	Chain	Res	Type
1	F	90	VAL
1	F	118	ASP
1	F	123	LYS
1	F	128	ARG

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

Mol	Chain	Res	Type
1	B	202	GLN
1	C	150	HIS
1	F	202	GLN

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 1 is monoatomic - leaving 5 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	$\# Z > 2$
3	GOL	B	501	-	5,5,5	0.07	0	5,5,5	0.17	0
3	GOL	D	401	-	5,5,5	0.19	0	5,5,5	0.55	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	GOL	D	402	-	5,5,5	0.12	0	5,5,5	0.21	0
3	GOL	C	301	-	5,5,5	0.14	0	5,5,5	0.32	0
3	GOL	E	401	-	5,5,5	0.13	0	5,5,5	0.37	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
3	GOL	B	501	-	-	2/4/4/4	-
3	GOL	D	401	-	-	2/4/4/4	-
3	GOL	D	402	-	-	0/4/4/4	-
3	GOL	C	301	-	-	0/4/4/4	-
3	GOL	E	401	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	401	GOL	O1-C1-C2-C3
3	E	401	GOL	C1-C2-C3-O3
3	B	501	GOL	O1-C1-C2-C3
3	B	501	GOL	O1-C1-C2-O2
3	D	401	GOL	O1-C1-C2-O2
3	E	401	GOL	O2-C2-C3-O3

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	301	GOL	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues

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	206/213 (96%)	-0.18	7 (3%) 45 48	30, 40, 70, 110	0
1	B	203/213 (95%)	-0.16	7 (3%) 45 48	31, 42, 66, 122	0
1	C	205/213 (96%)	-0.17	4 (1%) 65 66	29, 39, 67, 98	0
1	D	203/213 (95%)	-0.10	6 (2%) 50 52	30, 41, 64, 117	0
1	E	206/213 (96%)	-0.10	4 (1%) 66 68	30, 41, 74, 123	0
1	F	205/213 (96%)	-0.18	4 (1%) 65 66	30, 40, 68, 100	0
All	All	1228/1278 (96%)	-0.15	32 (2%) 56 58	29, 40, 72, 123	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	197	PHE	8.9
1	E	199	ALA	7.2
1	D	199	ALA	6.7
1	E	8	TRP	6.6
1	D	200	LEU	6.4
1	A	199	ALA	6.1
1	E	200	LEU	6.0
1	F	200	LEU	6.0
1	B	199	ALA	6.0
1	B	200	LEU	5.5
1	A	198	SER	5.3
1	A	213	THR	5.3
1	A	8	TRP	5.1
1	D	197	PHE	4.7
1	C	197	PHE	4.4
1	A	200	LEU	4.1
1	B	197	PHE	3.9
1	B	213	THR	3.7
1	F	199	ALA	3.7

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Mol	Chain	Res	Type	RSRZ
1	D	213	THR	3.7
1	C	199	ALA	3.6
1	A	197	PHE	3.5
1	F	9	ALA	3.4
1	C	213	THR	3.3
1	F	213	THR	3.2
1	B	196	LYS	2.7
1	D	198	SER	2.6
1	D	202	GLN	2.4
1	A	202	GLN	2.3
1	B	198	SER	2.3
1	B	17	ALA	2.2
1	C	145	LYS	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 monosaccharides 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(Å ²)	Q<0.9
3	GOL	E	401	6/6	0.83	0.21	53,58,61,62	0
2	NA	A	301	1/1	0.84	0.10	62,62,62,62	0
3	GOL	B	501	6/6	0.89	0.12	52,57,61,61	0
3	GOL	D	401	6/6	0.93	0.15	35,40,45,47	0
3	GOL	D	402	6/6	0.93	0.21	48,60,61,64	0
3	GOL	C	301	6/6	0.93	0.08	44,55,64,75	0

6.5 Other polymers [i](#)

There are no such residues in this entry.