



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 02:15 AM UTC

PDB ID : 3MW6 / pdb_00003mw6
Title : Crystal structure of NMB1681 from Neisseria meningitidis MC58, a FinO-like RNA chaperone
Authors : Tan, K.; Zhou, M.; Duggan, E.; Abdullah, J.; Joachimiak, A.; Midwest Center for Structural Genomics (MCSG)
Deposited on : 2010-05-05
Resolution : 2.21 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

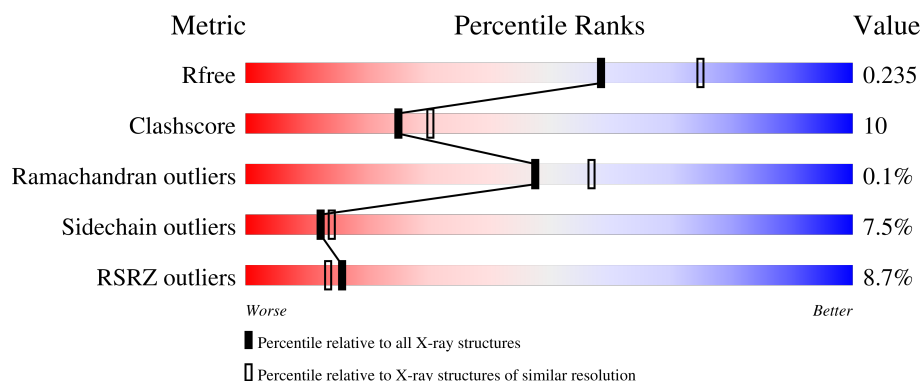
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.21 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	144	5% 70% 12% • 15%
1	B	144	16% 40% 24% • 35%
1	C	144	5% 62% 13% • 22%
1	D	144	7% 65% 14% • 18%
1	E	144	4% 66% 15% • 17%

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Mol	Chain	Length	Quality of chain
1	F	144	3% 66% 11% • 19%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5597 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 uncharacterized protein NMB1681.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	122	Total	C	N	O	S	Se	0	3	0
			981	620	186	172	1	2			
1	B	94	Total	C	N	O	S	Se	0	1	0
			769	493	144	130	1	1			
1	C	112	Total	C	N	O	S	Se	0	0	0
			884	563	164	154	1	2			
1	D	118	Total	C	N	O	S	Se	0	0	0
			931	590	172	165	1	3			
1	E	119	Total	C	N	O	S	Se	0	0	0
			939	594	174	167	1	3			
1	F	116	Total	C	N	O	S	Se	0	0	0
			917	582	169	162	1	3			

There are 18 discrepancies between the modelled and reference sequences:

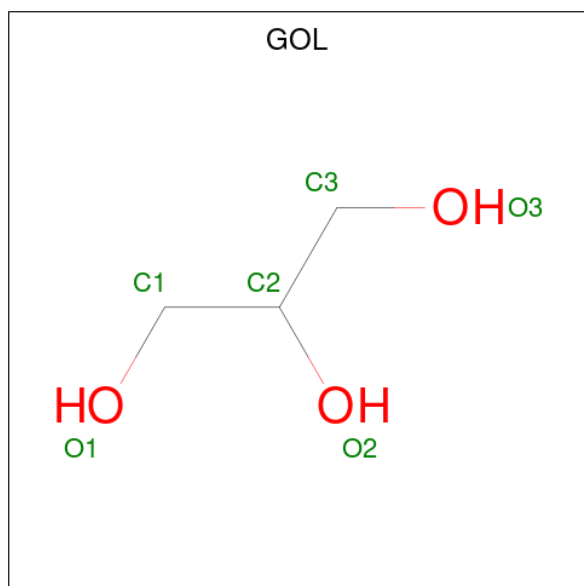
Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	UNP Q9JY98
A	-1	ASN	-	expression tag	UNP Q9JY98
A	0	ALA	-	expression tag	UNP Q9JY98
B	-2	SER	-	expression tag	UNP Q9JY98
B	-1	ASN	-	expression tag	UNP Q9JY98
B	0	ALA	-	expression tag	UNP Q9JY98
C	-2	SER	-	expression tag	UNP Q9JY98
C	-1	ASN	-	expression tag	UNP Q9JY98
C	0	ALA	-	expression tag	UNP Q9JY98
D	-2	SER	-	expression tag	UNP Q9JY98
D	-1	ASN	-	expression tag	UNP Q9JY98
D	0	ALA	-	expression tag	UNP Q9JY98
E	-2	SER	-	expression tag	UNP Q9JY98
E	-1	ASN	-	expression tag	UNP Q9JY98
E	0	ALA	-	expression tag	UNP Q9JY98
F	-2	SER	-	expression tag	UNP Q9JY98
F	-1	ASN	-	expression tag	UNP Q9JY98

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Chain	Residue	Modelled	Actual	Comment	Reference
F	0	ALA	-	expression tag	UNP Q9JY98

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



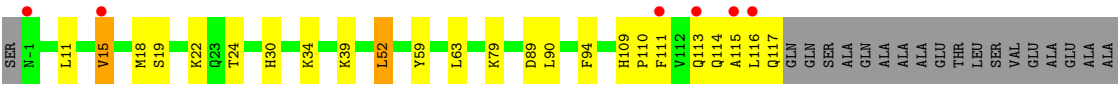
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	C	O	0	0
			6	3	3		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	33	Total	O	0	0
			33	33		
3	B	11	Total	O	0	0
			11	11		
3	C	27	Total	O	0	0
			27	27		
3	D	29	Total	O	0	0
			29	29		
3	E	26	Total	O	0	0
			26	26		
3	F	44	Total	O	0	0
			44	44		

GLU
ALA
GLU
ALA
ALA
GLU
SER
SER
ALA
ALA
GLU

● Molecule 1: uncharacterized protein NMB1681



GLU
SER
SER
SER
ALA
ALA
GLU

● Molecule 1: uncharacterized protein NMB1681



ALA
GLU

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	126.48Å 76.53Å 122.72Å 90.00° 116.84° 90.00°	Depositor
Resolution (Å)	42.89 – 2.21 42.89 – 2.21	Depositor EDS
% Data completeness (in resolution range)	93.2 (42.89-2.21) 95.8 (42.89-2.21)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.60 (at 2.20Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.5_2)	Depositor
R, R_{free}	0.191 , 0.235 0.192 , 0.235	Depositor DCC
R_{free} test set	2556 reflections (4.86%)	wwPDB-VP
Wilson B-factor (Å ²)	41.3	Xtriage
Anisotropy	0.544	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 56.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5597	wwPDB-VP
Average B, all atoms (Å ²)	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.72% 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: 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	0.49	0/1004	0.76	0/1345
1	B	0.43	0/787	0.80	0/1055
1	C	0.50	0/898	0.78	0/1205
1	D	0.54	0/944	0.77	0/1265
1	E	0.52	0/952	0.81	0/1276
1	F	0.51	0/931	0.77	0/1249
All	All	0.50	0/5516	0.78	0/7395

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	981	0	1018	15	0
1	B	769	0	794	27	0
1	C	884	0	913	18	0
1	D	931	0	959	17	0
1	E	939	0	965	21	0
1	F	917	0	946	15	0
2	C	6	0	8	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	33	0	0	0	0
3	B	11	0	0	0	0
3	C	27	0	0	0	0
3	D	29	0	0	0	0
3	E	26	0	0	1	0
3	F	44	0	0	1	0
All	All	5597	0	5603	107	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 (107) 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:E:79:LYS:HD3	1:E:116:LEU:HD11	1.58	0.86
1:C:5:THR:HG23	1:C:6:ALA:H	1.41	0.84
1:B:48:ILE:HD11	1:B:90:LEU:HG	1.66	0.77
1:A:5:THR:HG22	1:A:7:LEU:H	1.51	0.74
1:A:110:PRO:HG2	1:B:110:PRO:HG2	1.68	0.73
1:E:79:LYS:CD	1:E:116:LEU:HD11	2.18	0.72
1:C:20:LYS:O	1:C:24:THR:HG23	1.90	0.72
1:C:115:ALA:C	1:C:116:LEU:HD13	2.18	0.69
1:C:15:VAL:HA	1:C:18:MSE:HE3	1.73	0.69
1:A:3:GLN:HG3	1:A:4:GLU:H	1.58	0.68
1:C:24:THR:HG22	1:C:63:LEU:HD22	1.76	0.68
1:B:99:THR:OG1	1:B:102:GLU:CG	2.42	0.67
1:D:23:GLN:HA	1:D:26:MSE:HE3	1.77	0.66
1:A:13:SER:O	1:A:17:THR:HG23	1.95	0.66
1:F:89:ASP:HB2	3:F:154:HOH:O	1.96	0.65
1:F:2:THR:HG22	1:F:5:THR:HG23	1.80	0.64
1:F:103:GLN:O	1:F:107:GLN:HG3	2.00	0.61
1:C:24:THR:CG2	1:C:63:LEU:HD22	2.30	0.61
1:B:99:THR:OG1	1:B:102:GLU:HG3	2.03	0.59
1:F:2:THR:HG22	1:F:5:THR:H	1.67	0.58
1:A:5:THR:HG22	1:A:7:LEU:N	2.17	0.58
1:B:55:ALA:O	1:B:56:LEU:HD23	2.04	0.57
1:D:22:LYS:HD2	1:D:26:MSE:HE2	1.85	0.57
1:F:2:THR:HG23	1:F:4:GLU:H	1.70	0.57
1:C:109:HIS:CE1	1:C:111:PHE:CD2	2.92	0.56
1:E:110:PRO:HA	1:E:113:GLN:HE21	1.71	0.56
1:E:116:LEU:HD12	1:E:116:LEU:N	2.21	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:20:LYS:O	1:B:24:THR:HG23	2.06	0.56
1:D:14:ALA:O	1:D:17:THR:HB	2.06	0.56
1:E:115:ALA:HB3	1:E:116:LEU:HD12	1.87	0.56
1:E:30:HIS:HD2	1:E:59:TYR:OH	1.88	0.55
1:B:99:THR:OG1	1:B:102:GLU:HG2	2.06	0.54
1:A:113:GLN:HE21	1:A:113:GLN:HA	1.72	0.53
1:B:93:ARG:HA	1:C:116:LEU:HD12	1.92	0.52
1:E:114:GLN:HG3	1:E:115:ALA:N	2.25	0.51
1:A:3:GLN:HG3	1:A:4:GLU:N	2.24	0.51
1:B:44:LEU:O	1:B:87:ARG:HD3	2.10	0.51
1:B:35:TYR:CD2	1:B:90:LEU:HD22	2.46	0.51
1:D:110:PRO:HA	1:D:113:GLN:OE1	2.11	0.51
1:E:113:GLN:HA	1:E:117:GLN:HG2	1.93	0.51
1:B:23:GLN:O	1:B:27:ILE:HD13	2.11	0.50
1:D:18:MSE:HG3	1:D:26:MSE:HE1	1.93	0.50
1:F:113:GLN:HE21	1:F:113:GLN:N	2.10	0.49
1:A:113:GLN:HA	1:A:113:GLN:NE2	2.28	0.49
1:B:52:LEU:HD13	1:B:90:LEU:HD21	1.94	0.49
1:D:49:ASP:HB3	1:D:68:LEU:CD2	2.43	0.49
1:E:113:GLN:O	1:E:117:GLN:HG2	2.13	0.48
1:B:70:ASN:O	1:B:73:ARG:HG2	2.13	0.48
1:B:27:ILE:HG22	1:B:67:VAL:HG21	1.96	0.48
1:B:27:ILE:HG13	1:B:59:TYR:CD1	2.48	0.48
1:E:24:THR:OG1	1:E:63:LEU:HD22	2.14	0.48
1:D:11:LEU:C	1:D:11:LEU:HD23	2.39	0.48
1:D:12:LYS:O	1:D:16:GLN:HG3	2.14	0.48
1:C:110:PRO:HD3	1:D:109:HIS:NE2	2.29	0.47
1:E:19:SER:HB3	1:E:22:LYS:HB2	1.97	0.47
1:A:3:GLN:CG	1:A:4:GLU:H	2.23	0.47
1:C:49:ASP:OD2	1:C:50:GLN:N	2.48	0.47
1:F:109:HIS:CE1	1:F:111:PHE:CD2	3.02	0.47
1:A:118:GLN:O	1:A:122:GLN:CG	2.63	0.47
1:F:2:THR:CG2	1:F:5:THR:HG23	2.44	0.46
1:C:49:ASP:HB3	1:C:68:LEU:CD2	2.45	0.46
1:C:7:LEU:H	1:C:7:LEU:HG	1.52	0.46
1:F:82:ALA:HB1	1:F:107:GLN:HG2	1.98	0.46
1:A:23:GLN:O	1:A:27:ILE:HG13	2.14	0.46
1:D:7:LEU:O	1:D:8:GLY:C	2.58	0.46
1:E:89:ASP:OD1	1:E:89:ASP:C	2.58	0.46
1:D:34:LYS:NZ	1:D:34:LYS:HB2	2.31	0.46
1:D:17:THR:HG22	1:D:18:MSE:HE2	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:34:LYS:O	1:E:34:LYS:HG2	2.15	0.45
1:B:36:ASP:HB2	1:B:92:ASN:HD22	1.81	0.45
1:B:53:ILE:HD13	1:B:61:ALA:HB1	1.98	0.45
1:C:5:THR:HG23	1:C:6:ALA:N	2.21	0.45
1:A:99:THR:OG1	1:A:102:GLU:HG3	2.17	0.45
1:E:89:ASP:HB2	3:E:167:HOH:O	2.17	0.44
1:A:118:GLN:O	1:A:122:GLN:HG2	2.17	0.44
1:E:52:LEU:HD13	1:E:90:LEU:HD21	1.99	0.44
1:C:110:PRO:HG2	1:D:110:PRO:HG2	1.99	0.44
1:C:101:GLU:O	1:C:105:ILE:HG13	2.18	0.44
1:C:116:LEU:HD13	1:C:116:LEU:N	2.33	0.44
1:E:109:HIS:CE1	1:E:111:PHE:CD2	3.06	0.43
1:E:11:LEU:O	1:E:15:VAL:HG13	2.19	0.43
1:A:5:THR:HG22	1:A:6:ALA:N	2.32	0.43
1:A:109:HIS:CE1	1:A:111:PHE:HD1	2.36	0.43
1:B:28:ALA:HB2	1:B:67:VAL:HG13	2.01	0.43
1:F:115:ALA:C	1:F:116:LEU:HG	2.43	0.43
1:D:116:LEU:HD13	1:E:94:PHE:CZ	2.54	0.42
1:B:56:LEU:HB3	1:B:59:TYR:HD2	1.84	0.42
1:B:23:GLN:C	1:B:27:ILE:HD13	2.44	0.42
1:B:24:THR:HG22	1:B:63:LEU:HB3	2.00	0.42
1:E:110:PRO:HG2	1:F:110:PRO:HG2	2.01	0.42
1:B:52:LEU:HD12	1:B:52:LEU:HA	1.87	0.42
1:C:27:ILE:HG12	1:C:59:TYR:CG	2.54	0.42
1:E:18:MSE:HE2	1:E:22:LYS:HG2	2.01	0.42
1:F:109:HIS:HE1	1:F:111:PHE:CD2	2.38	0.42
1:F:86:LYS:HA	1:F:86:LYS:HD2	1.82	0.42
1:B:50:GLN:HA	1:B:50:GLN:OE1	2.19	0.42
1:D:113:GLN:O	1:D:117:GLN:HB2	2.19	0.42
1:F:72:CYS:HA	1:F:77:TYR:CD2	2.55	0.41
1:B:48:ILE:CD1	1:B:90:LEU:HG	2.44	0.41
1:C:109:HIS:HE1	1:C:111:PHE:CD2	2.37	0.41
1:B:21:LYS:N	1:B:21:LYS:HD2	2.36	0.41
1:B:31:ILE:C	1:B:33:GLY:N	2.78	0.41
1:F:114:GLN:H	1:F:114:GLN:HG2	1.60	0.41
1:B:49:ASP:OD2	1:B:50:GLN:N	2.50	0.40
1:D:30:HIS:CD2	1:D:30:HIS:C	3.00	0.40
1:D:78:LEU:HD13	1:D:78:LEU:HA	1.93	0.40
1:E:110:PRO:HA	1:E:113:GLN:NE2	2.35	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	123/144 (85%)	122 (99%)	1 (1%)	0	100	100
1	B	93/144 (65%)	89 (96%)	4 (4%)	0	100	100
1	C	110/144 (76%)	108 (98%)	2 (2%)	0	100	100
1	D	116/144 (81%)	115 (99%)	0	1 (1%)	14	14
1	E	117/144 (81%)	114 (97%)	3 (3%)	0	100	100
1	F	114/144 (79%)	111 (97%)	3 (3%)	0	100	100
All	All	673/864 (78%)	659 (98%)	13 (2%)	1 (0%)	48	57

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	8	GLY

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	99/108 (92%)	92 (93%)	7 (7%)	13	16
1	B	78/108 (72%)	69 (88%)	9 (12%)	5	5
1	C	89/108 (82%)	84 (94%)	5 (6%)	19	24
1	D	94/108 (87%)	86 (92%)	8 (8%)	10	11
1	E	95/108 (88%)	92 (97%)	3 (3%)	34	47
1	F	93/108 (86%)	84 (90%)	9 (10%)	8	8

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	548/648 (85%)	507 (92%)	41 (8%)	12	14

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	GLU
1	A	7	LEU
1	A	36	ASP
1	A	97	GLU
1	A	107	GLN
1	A	116	LEU
1	A	122	GLN
1	B	26	MSE
1	B	46	LEU
1	B	52	LEU
1	B	58	GLN
1	B	66	ARG
1	B	73	ARG
1	B	81	LEU
1	B	97	GLU
1	B	102	GLU
1	C	7	LEU
1	C	21	LYS
1	C	24	THR
1	C	39	LYS
1	C	116	LEU
1	D	5	THR
1	D	7	LEU
1	D	11	LEU
1	D	13	SER
1	D	34	LYS
1	D	40	ARG
1	D	52	LEU
1	D	78	LEU
1	E	15	VAL
1	E	39	LYS
1	E	52	LEU
1	F	1	MSE
1	F	2	THR
1	F	40	ARG
1	F	48	ILE
1	F	52	LEU

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Mol	Chain	Res	Type
1	F	86	LYS
1	F	113	GLN
1	F	114	GLN
1	F	116	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	113	GLN
1	A	117	GLN
1	B	58	GLN
1	C	114	GLN
1	D	23	GLN
1	D	30	HIS
1	D	70	ASN
1	D	107	GLN
1	E	30	HIS
1	E	113	GLN
1	E	114	GLN
1	F	107	GLN
1	F	113	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

1 ligand is 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	C	142	-	5,5,5	0.40	0	5,5,5	0.35	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	C	142	-	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

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	120/144 (83%)	0.13	7 (5%)	29 25	30, 49, 106, 152	3 (2%)
1	B	93/144 (64%)	1.10	23 (24%)	2 1	25, 73, 134, 169	1 (1%)
1	C	110/144 (76%)	0.22	7 (6%)	25 22	34, 53, 112, 148	0
1	D	115/144 (79%)	0.30	10 (8%)	16 13	30, 49, 89, 130	0
1	E	116/144 (80%)	0.20	6 (5%)	33 29	29, 54, 90, 116	0
1	F	113/144 (78%)	0.06	5 (4%)	39 36	32, 48, 87, 108	0
All	All	667/864 (77%)	0.31	58 (8%)	16 13	25, 52, 110, 169	4 (0%)

All (58) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	6	ALA	6.2
1	B	63	LEU	5.7
1	A	124	ALA	5.4
1	F	116	LEU	4.4
1	C	5	THR	4.3
1	F	111	PHE	4.3
1	B	23	GLN	4.3
1	A	121	ALA	4.3
1	C	116	LEU	4.2
1	F	115	ALA	3.9
1	B	22	LYS	3.9
1	C	7	LEU	3.6
1	D	116	LEU	3.6
1	E	113	GLN	3.5
1	D	0	ALA	3.5
1	A	123	ALA	3.3
1	B	57	PRO	3.3
1	D	111	PHE	3.3
1	B	65	ALA	3.2

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Mol	Chain	Res	Type	RSRZ
1	C	113	GLN	3.2
1	B	55	ALA	3.2
1	B	62	ALA	3.1
1	B	53	ILE	3.0
1	D	5	THR	3.0
1	D	113	GLN	2.8
1	A	120	SER	2.8
1	F	3	GLN	2.8
1	E	15	VAL	2.8
1	C	114	GLN	2.8
1	B	46	LEU	2.8
1	E	116	LEU	2.8
1	A	5	THR	2.7
1	B	20	LYS	2.7
1	B	67	VAL	2.7
1	E	115	ALA	2.6
1	D	115	ALA	2.6
1	A	116	LEU	2.6
1	D	9	ALA	2.6
1	B	113	GLN	2.5
1	C	115	ALA	2.5
1	E	-1	ASN	2.5
1	D	117	GLN	2.5
1	B	54	ALA	2.4
1	B	69	ALA	2.4
1	B	61	ALA	2.3
1	E	111	PHE	2.3
1	F	5	THR	2.3
1	D	114	GLN	2.3
1	A	3	GLN	2.3
1	B	50	GLN	2.3
1	D	10	ALA	2.2
1	B	52	LEU	2.2
1	B	27	ILE	2.2
1	B	31	ILE	2.2
1	B	40	ARG	2.1
1	B	64	ILE	2.1
1	B	35	TYR	2.1
1	B	21	LYS	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	C	142	6/6	0.82	0.27	97,97,98,99	0

6.5 Other polymers [i](#)

There are no such residues in this entry.