



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 09:15 AM GMT

PDB ID : 2NQR
Title : MoeA D142N
Authors : Nicolas, J.; Xiang, S.; Schindelin, H.; Rajagopalan, K.V.
Deposited on : 2006-10-31
Resolution : 2.20 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

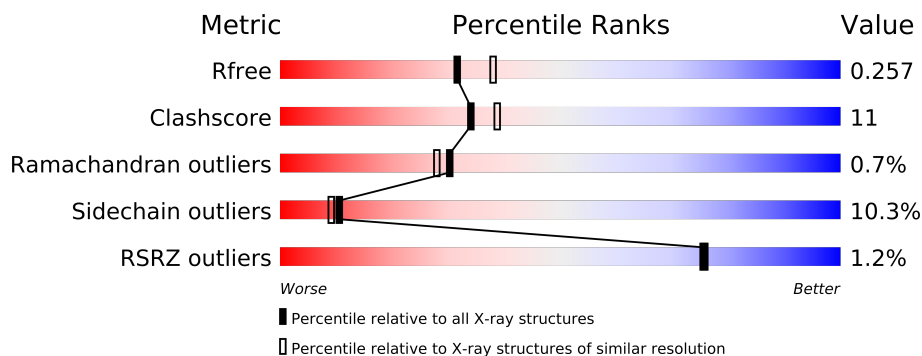
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.20 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}	66092	2938 (2.20-2.20)
Clashscore	79885	3751 (2.20-2.20)
Ramachandran outliers	78287	3681 (2.20-2.20)
Sidechain outliers	78261	3682 (2.20-2.20)
RSRZ outliers	66119	2939 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	411	
1	B	411	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
2	GOL	A	901	-	X
2	GOL	A	904	-	X
2	GOL	A	905	-	X
2	GOL	A	906	-	X
2	GOL	A	907	-	X
2	GOL	A	910	-	X

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Mol	Type	Chain	Res	Geometry	Electron density
2	GOL	A	911	-	X
2	GOL	A	914	-	X
2	GOL	A	917	-	X
2	GOL	A	918	-	X
2	GOL	B	903	-	X
2	GOL	B	908	-	X
2	GOL	B	909	-	X
2	GOL	B	913	-	X
2	GOL	B	916	-	X

2 Entry composition i

There are 3 unique types of molecules in this entry. The entry contains 7031 atoms, of which 0 are hydrogen and 0 are deuterium.

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.

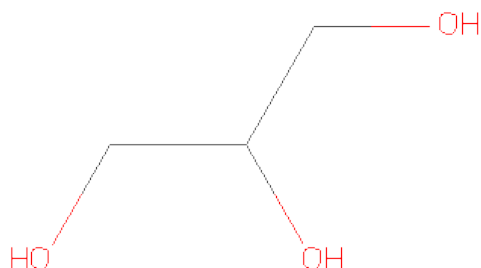
- Molecule 1 is a protein called Molybdopterin biosynthesis protein moeA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	404	Total	C	N	O	S	0	0	0
			3044	1920	533	578	13			
1	B	403	Total	C	N	O	S	0	0	0
			3040	1918	532	577	13			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	142	ASN	ASP	ENGINEERED	UNP P12281
B	142	ASN	ASP	ENGINEERED	UNP P12281

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



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

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

- Molecule 3 is water.

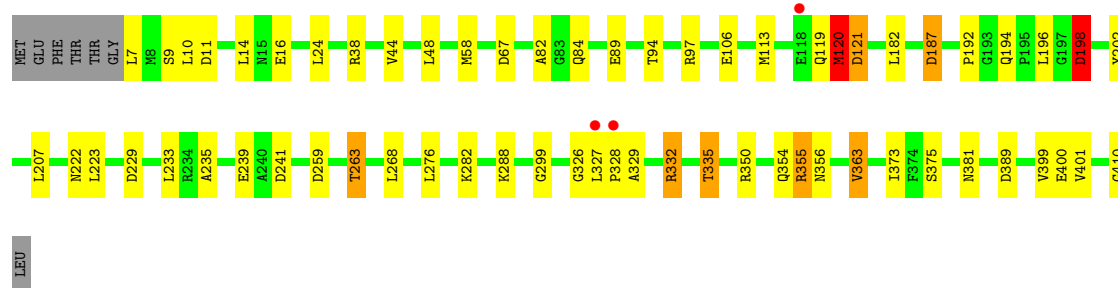
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	438	Total O 438 438	0	0
3	B	401	Total O 401 401	0	0

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

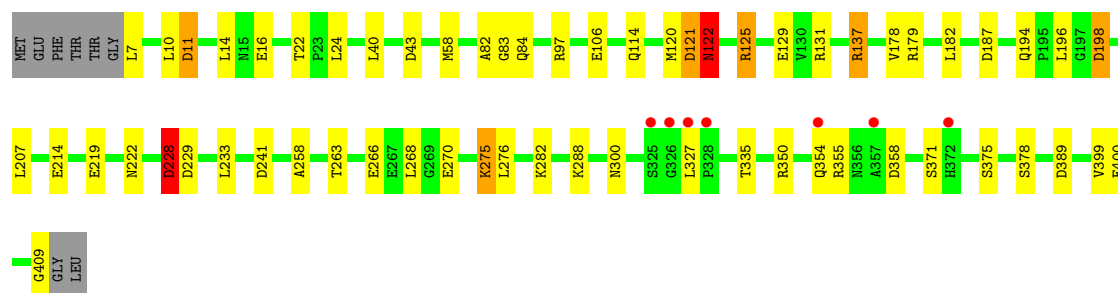
- Molecule 1: Molybdopterin biosynthesis protein moeA

Chain A: 



- Molecule 1: Molybdopterin biosynthesis protein moeA

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	75.28Å 100.08Å 102.48Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.64 – 2.20 39.01 – 2.20	Depositor EDS
% Data completeness (in resolution range)	98.8 (45.64-2.20) 98.8 (39.01-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.27 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.162 , 0.253 0.172 , 0.257	Depositor DCC
R_{free} test set	1981 reflections (5.28%)	DCC
Wilson B-factor (Å ²)	20.2	Xtriage
Anisotropy	0.030	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 44.4	EDS
Estimated twinning fraction	0.014 for -h,l,k	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 39535 reflections (0.003%)	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7031	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.51% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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	0.89	0/3103	1.04	12/4217 (0.3%)
1	B	0.95	5/3099 (0.2%)	1.03	12/4212 (0.3%)
All	All	0.92	5/6202 (0.1%)	1.04	24/8429 (0.3%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	214	GLU	CD-OE1	8.27	1.34	1.25
1	B	275	LYS	CD-CE	6.32	1.67	1.51
1	B	214	GLU	CG-CD	5.97	1.60	1.51
1	B	258	ALA	CA-CB	5.08	1.63	1.52
1	B	122	ASN	CB-CG	5.02	1.62	1.51

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	228	ASP	CB-CG-OD1	9.43	126.79	118.30
1	A	389	ASP	CB-CG-OD2	9.02	126.41	118.30
1	B	11	ASP	CB-CG-OD2	7.98	125.48	118.30
1	B	275	LYS	CD-CE-NZ	7.82	129.69	111.70
1	B	198	ASP	CB-CG-OD2	7.73	125.26	118.30
1	A	332	ARG	NE-CZ-NH1	7.69	124.14	120.30
1	B	121	ASP	CB-CG-OD2	7.68	125.21	118.30
1	A	259	ASP	CB-CG-OD1	7.46	125.01	118.30
1	A	363	VAL	CB-CA-C	-7.31	97.50	111.40
1	A	241	ASP	CB-CG-OD2	7.11	124.70	118.30
1	A	198	ASP	CB-CG-OD2	6.97	124.57	118.30
1	A	187	ASP	CB-CG-OD2	6.88	124.49	118.30
1	B	43	ASP	CB-CG-OD2	6.73	124.36	118.30
1	B	229	ASP	CB-CG-OD2	6.73	124.35	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	11	ASP	CB-CG-OD1	-6.67	112.30	118.30
1	B	358	ASP	CB-CG-OD2	6.57	124.21	118.30
1	B	409	GLY	CA-C-O	-6.48	108.94	120.60
1	A	67	ASP	CB-CG-OD2	6.41	124.07	118.30
1	A	121	ASP	CB-CG-OD2	6.20	123.88	118.30
1	B	241	ASP	CB-CG-OD2	5.77	123.49	118.30
1	A	120	MET	CA-CB-CG	5.51	122.66	113.30
1	B	214	GLU	CA-CB-CG	5.48	125.47	113.40
1	A	38	ARG	NE-CZ-NH2	5.08	122.84	120.30
1	A	335	THR	CB-CA-C	-5.05	97.96	111.60

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3044	0	0	29	0
1	B	3040	0	0	28	0
2	A	78	0	103	29	1
2	B	30	0	40	10	0
3	A	438	0	0	17	3
3	B	401	0	0	13	4
All	All	7031	0	143	68	5

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 11.

All (68) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:A:901:GOL:H32	1:B:84:GLN:NE2	1.47	1.26
1:A:187:ASP:OD2	2:A:902:GOL:H2	1.44	1.15
2:B:909:GOL:H32	3:B:987:HOH:O	1.52	1.08
1:B:7:LEU:N	3:B:1050:HOH:O	1.82	1.07

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:187:ASP:OD1	2:A:901:GOL:H2	1.54	1.07
2:A:902:GOL:H31	3:A:1007:HOH:O	1.59	1.00
2:A:901:GOL:C3	1:B:84:GLN:NE2	2.23	1.00
1:B:228:ASP:OD2	2:B:908:GOL:H32	1.68	0.94
2:B:909:GOL:C3	3:B:987:HOH:O	2.10	0.91
2:A:902:GOL:C3	3:A:1007:HOH:O	2.17	0.89
1:A:375:SER:CB	2:A:914:GOL:H12	2.14	0.78
1:B:120:MET:CB	1:B:122:ASN:ND2	2.49	0.76
1:B:378:SER:O	2:B:909:GOL:H2	1.86	0.75
2:A:901:GOL:H11	3:A:1332:HOH:O	1.88	0.72
1:A:239:GLU:OE2	2:A:907:GOL:H2	1.89	0.72
1:B:378:SER:O	3:B:987:HOH:O	2.09	0.71
1:B:389:ASP:OD2	3:B:1197:HOH:O	2.09	0.70
1:A:187:ASP:OD1	2:A:901:GOL:C2	2.38	0.70
2:A:917:GOL:C1	3:A:1132:HOH:O	2.39	0.70
1:B:187:ASP:OD2	2:B:908:GOL:H11	1.92	0.69
2:A:917:GOL:H11	3:A:1132:HOH:O	1.91	0.69
2:A:902:GOL:O2	1:B:82:ALA:O	2.13	0.67
1:B:266:GLU:O	3:B:1222:HOH:O	2.12	0.65
2:A:901:GOL:C1	3:A:1332:HOH:O	2.41	0.64
1:B:228:ASP:CB	3:B:1306:HOH:O	2.45	0.64
2:B:903:GOL:O1	3:B:1132:HOH:O	2.06	0.64
1:A:187:ASP:CG	2:A:902:GOL:H2	2.19	0.61
1:A:355:ARG:NH2	1:B:400:GLU:OE2	2.33	0.61
1:A:400:GLU:OE2	1:B:355:ARG:NH2	2.35	0.59
1:B:198:ASP:N	3:B:1169:HOH:O	2.36	0.58
1:A:375:SER:CA	2:A:914:GOL:H12	2.34	0.57
2:A:905:GOL:H32	3:A:943:HOH:O	2.02	0.57
1:B:228:ASP:OD2	2:B:908:GOL:C3	2.48	0.57
2:A:917:GOL:H12	3:A:1132:HOH:O	2.04	0.57
1:B:270:GLU:CB	3:B:1172:HOH:O	2.54	0.56
1:A:7:LEU:N	3:A:1039:HOH:O	2.42	0.52
1:A:7:LEU:N	3:A:1079:HOH:O	2.44	0.50
1:A:44:VAL:CA	2:A:911:GOL:H32	2.42	0.50
1:A:235:ALA:O	2:A:907:GOL:H32	2.12	0.48
1:A:239:GLU:OE2	2:A:907:GOL:C1	2.62	0.48
1:B:178:VAL:N	2:B:903:GOL:HO3	2.11	0.48
3:A:1346:HOH:O	1:B:375:SER:CB	2.62	0.47
1:A:229:ASP:CB	2:A:918:GOL:H32	2.45	0.46
2:A:902:GOL:H12	2:A:906:GOL:H12	1.97	0.46
1:A:263:THR:CG2	3:A:1342:HOH:O	2.63	0.46
1:B:121:ASP:O	1:B:122:ASN:CB	2.63	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:120:MET:O	1:A:121:ASP:CB	2.64	0.45
1:B:300:ASN:ND2	1:B:371:SER:CB	2.80	0.45
1:B:137:ARG:NH1	3:B:1206:HOH:O	2.51	0.44
1:A:410:GLY:O	3:A:1275:HOH:O	2.21	0.44
1:A:239:GLU:OE2	2:A:907:GOL:C2	2.64	0.44
1:A:198:ASP:O	3:A:1208:HOH:O	2.21	0.43
1:A:326:GLY:CA	3:A:965:HOH:O	2.66	0.43
1:B:300:ASN:ND2	1:B:371:SER:O	2.52	0.42
1:A:356:ASN:OD1	1:A:356:ASN:C	2.57	0.42
1:A:82:ALA:C	2:B:908:GOL:O2	2.52	0.42
1:A:194:GLN:O	1:A:202:TYR:OH	2.38	0.42
1:A:223:LEU:N	1:A:223:LEU:CD1	2.83	0.41
1:B:187:ASP:CG	2:B:908:GOL:H11	2.40	0.41
1:B:129:GLU:CG	3:B:1201:HOH:O	2.68	0.41
1:A:192:PRO:CD	3:A:976:HOH:O	2.67	0.41
1:A:187:ASP:OD1	2:A:901:GOL:C3	2.68	0.41
1:B:125:ARG:CG	3:B:1007:HOH:O	2.69	0.41
1:B:179:ARG:NH2	1:B:219:GLU:OE1	2.54	0.41
2:A:906:GOL:O1	3:A:1200:HOH:O	2.06	0.41
2:A:901:GOL:O1	1:B:83:GLY:O	2.37	0.41
1:A:381:ASN:ND2	2:A:912:GOL:H11	2.36	0.40
1:A:299:GLY:O	2:A:904:GOL:H31	2.21	0.40

All (5) 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	Distance(Å)	Clash(Å)
2:A:911:GOL:O2	3:A:1261:HOH:O[4_465]	1.78	0.42
3:B:981:HOH:O	3:B:1316:HOH:O[1_545]	1.93	0.27
3:B:943:HOH:O	3:B:1243:HOH:O[3_645]	2.06	0.14
3:A:1073:HOH:O	3:B:1269:HOH:O[3_545]	2.16	0.04
3:A:1029:HOH:O	3:B:1027:HOH:O[3_545]	2.18	0.02

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	402/411 (98%)	390 (97%)	8 (2%)	4 (1%)	22	18
1	B	401/411 (98%)	387 (96%)	12 (3%)	2 (0%)	38	38
All	All	803/822 (98%)	777 (97%)	20 (2%)	6 (1%)	30	28

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	198	ASP
1	A	282	LYS
1	A	328	PRO
1	B	122	ASN
1	B	282	LYS
1	A	329	ALA

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	325/331 (98%)	290 (89%)	35 (11%)	9	8
1	B	325/331 (98%)	293 (90%)	32 (10%)	12	10
All	All	650/662 (98%)	583 (90%)	67 (10%)	10	9

All (67) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	SER
1	A	10	LEU
1	A	11	ASP
1	A	14	LEU
1	A	16	GLU
1	A	24	LEU
1	A	48	LEU
1	A	58	MET
1	A	84	GLN
1	A	89	GLU
1	A	94	THR

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Mol	Chain	Res	Type
1	A	97	ARG
1	A	106	GLU
1	A	113	MET
1	A	119	GLN
1	A	120	MET
1	A	182	LEU
1	A	196	LEU
1	A	207	LEU
1	A	222	ASN
1	A	233	LEU
1	A	263	THR
1	A	268	LEU
1	A	276	LEU
1	A	288	LYS
1	A	327	LEU
1	A	332	ARG
1	A	335	THR
1	A	350	ARG
1	A	354	GLN
1	A	355	ARG
1	A	363	VAL
1	A	373	ILE
1	A	399	VAL
1	A	401	VAL
1	B	10	LEU
1	B	11	ASP
1	B	14	LEU
1	B	16	GLU
1	B	22	THR
1	B	24	LEU
1	B	40	LEU
1	B	58	MET
1	B	97	ARG
1	B	106	GLU
1	B	114	GLN
1	B	122	ASN
1	B	125	ARG
1	B	131	ARG
1	B	137	ARG
1	B	182	LEU
1	B	194	GLN
1	B	196	LEU

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Mol	Chain	Res	Type
1	B	207	LEU
1	B	222	ASN
1	B	228	ASP
1	B	233	LEU
1	B	263	THR
1	B	268	LEU
1	B	275	LYS
1	B	276	LEU
1	B	288	LYS
1	B	327	LEU
1	B	335	THR
1	B	350	ARG
1	B	354	GLN
1	B	399	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

18 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	901	-	5,5,5	1.78	1 (20%)	5,5,5	0.91	0
2	GOL	A	902	-	5,5,5	1.96	1 (20%)	5,5,5	0.92	0
2	GOL	A	904	-	5,5,5	0.48	0	5,5,5	1.02	0
2	GOL	A	905	-	5,5,5	0.96	0	5,5,5	2.08	2 (40%)
2	GOL	A	906	-	5,5,5	0.43	0	5,5,5	0.73	0
2	GOL	A	907	-	5,5,5	0.71	0	5,5,5	1.84	2 (40%)
2	GOL	A	910	-	5,5,5	0.33	0	5,5,5	0.49	0
2	GOL	A	911	-	5,5,5	0.53	0	5,5,5	1.42	1 (20%)
2	GOL	A	912	-	5,5,5	0.33	0	5,5,5	0.18	0
2	GOL	A	914	-	5,5,5	0.38	0	5,5,5	1.46	1 (20%)
2	GOL	A	915	-	5,5,5	0.31	0	5,5,5	0.30	0
2	GOL	A	917	-	5,5,5	0.46	0	5,5,5	0.56	0
2	GOL	A	918	-	5,5,5	0.46	0	5,5,5	0.83	0
2	GOL	B	903	-	5,5,5	0.52	0	5,5,5	0.57	0
2	GOL	B	908	1	5,5,5	0.60	0	5,5,5	0.99	0
2	GOL	B	909	-	5,5,5	0.69	0	5,5,5	0.38	0
2	GOL	B	913	-	5,5,5	0.63	0	5,5,5	1.32	0
2	GOL	B	916	-	5,5,5	0.44	0	5,5,5	0.34	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	901	-	-	0/4/4/4	0/0/0/0
2	GOL	A	902	-	-	0/4/4/4	0/0/0/0
2	GOL	A	904	-	-	0/4/4/4	0/0/0/0
2	GOL	A	905	-	-	0/4/4/4	0/0/0/0
2	GOL	A	906	-	-	0/4/4/4	0/0/0/0
2	GOL	A	907	-	-	0/4/4/4	0/0/0/0
2	GOL	A	910	-	-	0/4/4/4	0/0/0/0
2	GOL	A	911	-	-	0/4/4/4	0/0/0/0
2	GOL	A	912	-	-	0/4/4/4	0/0/0/0
2	GOL	A	914	-	-	0/4/4/4	0/0/0/0
2	GOL	A	915	-	-	0/4/4/4	0/0/0/0
2	GOL	A	917	-	-	0/4/4/4	0/0/0/0
2	GOL	A	918	-	-	0/4/4/4	0/0/0/0
2	GOL	B	903	-	-	0/4/4/4	0/0/0/0
2	GOL	B	908	1	-	0/4/4/4	0/0/0/0
2	GOL	B	909	-	-	0/4/4/4	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GOL	B	913	-	-	0/4/4/4	0/0/0/0
2	GOL	B	916	-	-	0/4/4/4	0/0/0/0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	902	GOL	O2-C2	-4.28	1.30	1.43
2	A	901	GOL	C3-C2	-3.69	1.37	1.52

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	905	GOL	O3-C3-C2	3.69	127.73	109.71
2	A	907	GOL	O1-C1-C2	2.93	124.02	109.71
2	A	914	GOL	O1-C1-C2	2.82	123.47	109.71
2	A	911	GOL	O3-C3-C2	2.60	122.41	109.71
2	A	905	GOL	O2-C2-C3	2.49	119.55	108.22
2	A	907	GOL	O2-C2-C1	2.19	118.18	108.22

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers

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	404/411 (98%)	-0.56	3 (0%) 84 86	11, 20, 35, 57	0
1	B	403/411 (98%)	-0.51	7 (1%) 67 68	9, 20, 38, 63	0
All	All	807/822 (98%)	-0.54	10 (1%) 75 76	9, 20, 36, 63	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	327	LEU	5.9
1	A	327	LEU	4.4
1	A	328	PRO	3.9
1	B	357	ALA	3.8
1	B	326	GLY	3.5
1	B	328	PRO	3.1
1	B	372	HIS	2.7
1	B	325	SER	2.4
1	B	354	GLN	2.4
1	A	118	GLU	2.1

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	GOL	A	910	6/6	0.28	29.44	44,45,45,47	0
2	GOL	A	917	6/6	0.24	18.63	40,48,50,52	0
2	GOL	A	914	6/6	0.26	13.94	36,41,43,47	0
2	GOL	A	906	6/6	0.15	10.35	38,43,44,46	0
2	GOL	A	918	6/6	0.18	7.38	31,38,41,42	0
2	GOL	B	903	6/6	0.19	6.82	30,36,37,39	0
2	GOL	A	905	6/6	0.17	6.42	24,28,31,33	0
2	GOL	B	913	6/6	0.19	5.78	21,33,36,39	0
2	GOL	A	911	6/6	0.17	5.51	37,41,43,46	0
2	GOL	B	916	6/6	0.27	4.19	55,56,57,58	0
2	GOL	B	909	6/6	0.15	4.17	46,47,48,49	0
2	GOL	A	904	6/6	0.12	3.36	15,25,27,34	0
2	GOL	A	901	6/6	0.21	3.34	7,19,23,24	0
2	GOL	B	908	6/6	0.28	2.95	39,40,40,44	0
2	GOL	A	907	6/6	0.17	2.27	26,30,32,35	0
2	GOL	A	902	6/6	0.15	1.99	14,16,25,25	0
2	GOL	A	915	6/6	0.28	1.19	49,49,51,51	0
2	GOL	A	912	6/6	0.14	0.69	43,45,46,48	0

6.5 Other polymers ⓘ

There are no such residues in this entry.