



Full wwPDB X-ray Structure Validation Report ⓘ

May 14, 2020 – 01:43 pm BST

PDB ID : 4G2V
Title : Structure complex of LGN binding with FRMPD1
Authors : Shang, Y.; Pan, Z.; Wen, W.; Wang, W.; Zhang, M.
Deposited on : 2012-07-13
Resolution : 2.40 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

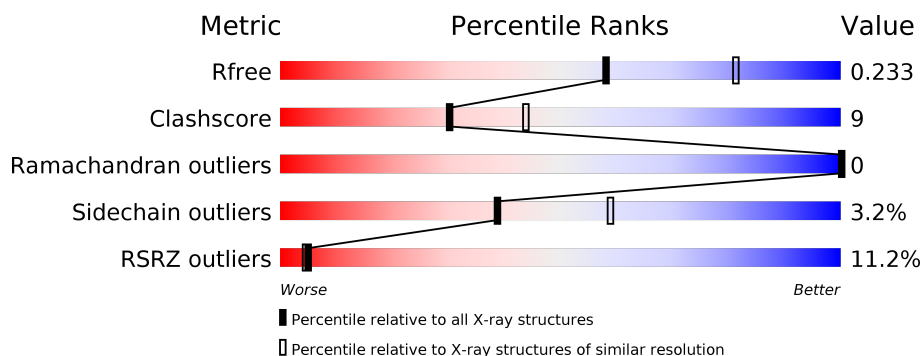
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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	340	
2	B	38	

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
3	GOL	A	403	-	-	X	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 2732 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 G-protein-signaling modulator 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	321	Total	C	N	O	S	0	1	0
			2456	1536	436	476	8			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	11	GLY	-	EXPRESSION TAG	UNP Q8VDU0
A	12	PRO	-	EXPRESSION TAG	UNP Q8VDU0
A	13	GLY	-	EXPRESSION TAG	UNP Q8VDU0
A	14	SER	-	EXPRESSION TAG	UNP Q8VDU0

- Molecule 2 is a protein called peptide from FERM and PDZ domain-containing protein 1.

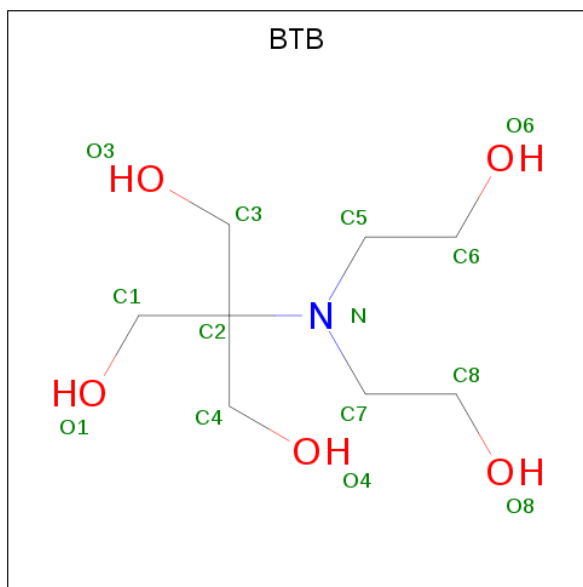
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	17	Total	C	N	O	S	0	0	0
			129	78	18	30	3			

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



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

- Molecule 4 is 2-[BIS-(2-HYDROXY-ETHYL)-AMINO]-2-HYDROXYMETHYL-PROPAN E-1,3-DIOL (three-letter code: BTB) (formula: C₈H₁₉NO₅).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 5 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Cl	0	0
			1	1		

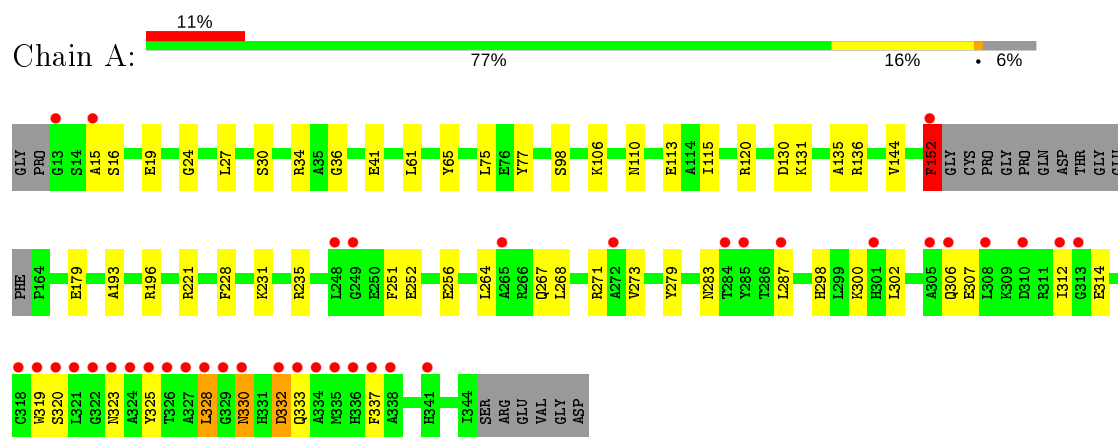
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	78	Total	O	0	0
			78	78		
6	B	6	Total	O	0	0
			6	6		

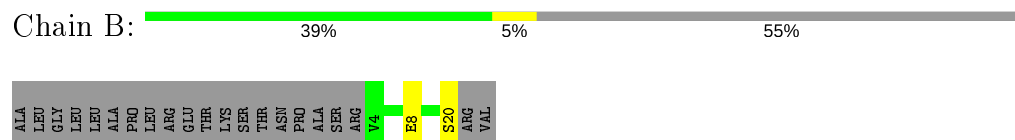
3 Residue-property plots [i](#)

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: G-protein-signaling modulator 2



- Molecule 2: peptide from FERM and PDZ domain-containing protein 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	93.83 Å 93.83 Å 172.86 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	41.23 – 2.40 41.23 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.8 (41.23-2.40) 99.9 (41.23-2.40)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.59 (at 2.39 Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.1_743)	Depositor
R, R_{free}	0.190 , 0.232 0.188 , 0.233	Depositor DCC
R_{free} test set	930 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å ²)	38.7	Xtriage
Anisotropy	0.602	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 66.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	2732	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 6.74% 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, CL, BTB

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	0.26	0/2501	0.43	2/3378 (0.1%)
2	B	0.28	0/129	0.50	0/172
All	All	0.26	0/2630	0.43	2/3550 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	152	PHE	CG-CD1-CE1	-5.72	114.51	120.80
1	A	152	PHE	CB-CG-CD1	5.34	124.54	120.80

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	152	PHE	Sidechain

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	2456	0	2322	46	4
2	B	129	0	117	1	1
3	A	42	0	56	14	0
3	B	6	0	8	1	0
4	A	14	0	19	1	0
5	A	1	0	0	1	0
6	A	78	0	0	2	1
6	B	6	0	0	0	0
All	All	2732	0	2522	48	5

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 9.

All (48) 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:106:LYS:HZ2	3:A:405:GOL:H31	1.34	0.91
1:A:235:ARG:HD2	1:A:273:VAL:HG11	1.55	0.85
1:A:312:ILE:HD12	1:A:312:ILE:H	1.48	0.78
1:A:106:LYS:NZ	3:A:405:GOL:H31	1.99	0.77
1:A:319:TRP:O	1:A:323:ASN:ND2	2.19	0.76
1:A:193:ALA:HB2	3:A:403:GOL:H31	1.68	0.75
1:A:330:ASN:OD1	1:A:333:GLN:HB2	1.89	0.72
1:A:120[B]:ARG:NH1	5:A:409:CL:CL	2.63	0.69
1:A:179:GLU:OE1	6:A:567:HOH:O	2.10	0.68
3:A:402:GOL:H31	3:B:101:GOL:H32	1.78	0.65
1:A:136:ARG:HG3	3:A:403:GOL:H11	1.81	0.62
1:A:235:ARG:CD	1:A:273:VAL:HG11	2.28	0.62
1:A:235:ARG:HD2	1:A:273:VAL:CG1	2.29	0.59
1:A:332:ASP:N	1:A:332:ASP:OD1	2.34	0.59
1:A:271:ARG:NH1	1:A:307:GLU:OE2	2.19	0.57
1:A:130:ASP:OD2	3:A:402:GOL:O2	2.24	0.55
1:A:136:ARG:CG	3:A:403:GOL:H11	2.35	0.55
1:A:325:TYR:HA	1:A:328:LEU:HD11	1.89	0.54
1:A:135:ALA:CB	3:A:403:GOL:H12	2.39	0.53
1:A:231:LYS:HB2	1:A:268:LEU:HD21	1.91	0.52
1:A:325:TYR:O	1:A:328:LEU:HD12	2.11	0.51
1:A:228:PHE:HB2	3:A:406:GOL:H11	1.94	0.50
1:A:231:LYS:CB	1:A:268:LEU:HD21	2.43	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:302:LEU:O	1:A:306:GLN:HG3	2.12	0.49
1:A:27:LEU:HA	1:A:30:SER:HB3	1.95	0.48
1:A:136:ARG:HA	3:A:403:GOL:H11	1.95	0.47
1:A:34:ARG:NH2	6:A:541:HOH:O	2.48	0.47
1:A:306:GLN:HG2	1:A:314:GLU:OE2	2.15	0.46
1:A:61:LEU:HD13	1:A:77:TYR:CE1	2.52	0.45
1:A:279:TYR:O	1:A:283:ASN:ND2	2.50	0.45
1:A:136:ARG:HG3	3:A:403:GOL:C1	2.46	0.45
1:A:251:PHE:CE2	1:A:287:LEU:HB3	2.51	0.44
1:A:196:ARG:HA	2:B:8:GLU:HG3	2.00	0.43
4:A:408:BTB:H71	4:A:408:BTB:H42	1.82	0.43
1:A:41:GLU:OE1	1:A:65:TYR:OH	2.29	0.43
1:A:98:SER:OG	1:A:120[A]:ARG:HB3	2.18	0.43
1:A:325:TYR:HE2	1:A:337:PHE:CE2	2.37	0.43
1:A:330:ASN:O	1:A:330:ASN:CG	2.58	0.43
1:A:252:GLU:O	1:A:256:GLU:HG2	2.19	0.42
1:A:298:HIS:CE1	1:A:320:SER:HB3	2.54	0.42
1:A:228:PHE:CB	3:A:406:GOL:H11	2.50	0.42
1:A:131:LYS:HA	1:A:131:LYS:HD3	1.91	0.42
1:A:228:PHE:CD2	3:A:406:GOL:H11	2.55	0.42
1:A:264:LEU:HA	1:A:267:GLN:HB2	2.02	0.41
1:A:115:ILE:HG13	1:A:144:VAL:HG11	2.02	0.41
1:A:136:ARG:CA	3:A:403:GOL:H11	2.51	0.40
1:A:15:ALA:O	1:A:19:GLU:HG3	2.21	0.40
1:A:24:GLY:HA2	1:A:36:GLY:HA2	2.04	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	Interatomic distance (Å)	Clash overlap (Å)
1:A:152:PHE:CD1	1:A:152:PHE:CE1[9_554]	1.40	0.80
1:A:152:PHE:CG	1:A:152:PHE:CD1[9_554]	1.41	0.79
1:A:152:PHE:CG	1:A:152:PHE:CE1[9_554]	1.70	0.50
6:A:563:HOH:O	6:A:572:HOH:O[6_555]	2.12	0.08
1:A:300:LYS:NZ	2:B:20:SER:O[8_545]	2.17	0.03

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	318/340 (94%)	315 (99%)	3 (1%)	0	100	100
2	B	15/38 (40%)	15 (100%)	0	0	100	100
All	All	333/378 (88%)	330 (99%)	3 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	238/266 (90%)	230 (97%)	8 (3%)	37	56
2	B	15/34 (44%)	15 (100%)	0	100	100
All	All	253/300 (84%)	245 (97%)	8 (3%)	39	59

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	SER
1	A	75	LEU
1	A	110	ASN
1	A	113	GLU
1	A	221	ARG
1	A	328	LEU
1	A	330	ASN
1	A	332	ASP

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

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 10 ligands modelled in this entry, 1 is monoatomic - leaving 9 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	A	405	-	5,5,5	0.38	0	5,5,5	0.14	0
3	GOL	A	404	-	5,5,5	0.37	0	5,5,5	0.33	0
3	GOL	A	403	-	5,5,5	0.37	0	5,5,5	0.56	0
4	BTB	A	408	-	13,13,13	1.50	4 (30%)	7,16,16	1.90	3 (42%)
3	GOL	A	407	-	5,5,5	0.38	0	5,5,5	0.28	0
3	GOL	B	101	-	5,5,5	0.37	0	5,5,5	0.28	0
3	GOL	A	402	-	5,5,5	0.43	0	5,5,5	0.20	0
3	GOL	A	401	-	5,5,5	0.39	0	5,5,5	0.30	0
3	GOL	A	406	-	5,5,5	0.36	0	5,5,5	0.31	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	A	405	-	-	2/4/4/4	-
3	GOL	A	404	-	-	2/4/4/4	-
3	GOL	A	403	-	-	2/4/4/4	-
4	BTB	A	408	-	-	6/21/21/21	-
3	GOL	A	407	-	-	2/4/4/4	-
3	GOL	B	101	-	-	3/4/4/4	-
3	GOL	A	402	-	-	2/4/4/4	-
3	GOL	A	401	-	-	0/4/4/4	-
3	GOL	A	406	-	-	2/4/4/4	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	408	BTB	C5-N	-2.65	1.44	1.48
4	A	408	BTB	C3-C2	-2.27	1.50	1.53
4	A	408	BTB	C4-C2	-2.21	1.50	1.53
4	A	408	BTB	C1-C2	-2.06	1.50	1.53

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	408	BTB	O3-C3-C2	2.69	118.79	111.44
4	A	408	BTB	O1-C1-C2	2.27	117.65	111.44
4	A	408	BTB	O4-C4-C2	2.14	117.30	111.44

There are no chirality outliers.

All (21) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	404	GOL	O1-C1-C2-C3
3	A	403	GOL	C1-C2-C3-O3
4	A	408	BTB	O1-C1-C2-C3
4	A	408	BTB	O1-C1-C2-C4
4	A	408	BTB	C4-C2-C3-O3
3	A	407	GOL	O1-C1-C2-O2
3	A	407	GOL	O1-C1-C2-C3
3	B	101	GOL	O1-C1-C2-C3
3	B	101	GOL	O1-C1-C2-O2
3	A	402	GOL	O1-C1-C2-C3
3	A	406	GOL	O1-C1-C2-C3

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Mol	Chain	Res	Type	Atoms
3	A	404	GOL	O1-C1-C2-O2
3	A	402	GOL	O1-C1-C2-O2
3	A	403	GOL	O2-C2-C3-O3
3	A	405	GOL	O1-C1-C2-C3
4	A	408	BTB	C1-C2-C3-O3
4	A	408	BTB	O1-C1-C2-N
4	A	408	BTB	N-C2-C3-O3
3	A	405	GOL	O1-C1-C2-O2
3	A	406	GOL	O1-C1-C2-O2
3	B	101	GOL	C1-C2-C3-O3

There are no ring outliers.

6 monomers are involved in 15 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	405	GOL	2	0
3	A	403	GOL	7	0
4	A	408	BTB	1	0
3	B	101	GOL	1	0
3	A	402	GOL	2	0
3	A	406	GOL	3	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	321/340 (94%)	0.32	38 (11%) 4 4	29, 53, 113, 134	0
2	B	17/38 (44%)	-0.04	0 100 100	33, 48, 99, 105	0
All	All	338/378 (89%)	0.31	38 (11%) 5 4	29, 53, 113, 134	0

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	337	PHE	6.9
1	A	335	MET	5.4
1	A	334	ALA	5.1
1	A	329	GLY	5.0
1	A	328	LEU	4.2
1	A	287	LEU	4.1
1	A	308	LEU	3.9
1	A	321	LEU	3.7
1	A	338	ALA	3.6
1	A	327	ALA	3.6
1	A	323	ASN	3.5
1	A	318	CYS	3.3
1	A	324	ALA	3.3
1	A	152	PHE	3.2
1	A	332	ASP	3.1
1	A	326	THR	3.1
1	A	336	HIS	3.1
1	A	284	THR	3.1
1	A	13	GLY	3.0
1	A	333	GLN	3.0
1	A	305	ALA	2.9
1	A	249	GLY	2.9
1	A	265	ALA	2.9
1	A	341	HIS	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	272	ALA	2.6
1	A	306	GLN	2.5
1	A	325	TYR	2.4
1	A	310	ASP	2.4
1	A	319	TRP	2.4
1	A	301	HIS	2.4
1	A	320	SER	2.3
1	A	313	GLY	2.2
1	A	322	GLY	2.2
1	A	285	TYR	2.1
1	A	330	ASN	2.1
1	A	15	ALA	2.0
1	A	248	LEU	2.0
1	A	312	ILE	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 carbohydrates 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	A	405	6/6	0.81	0.31	71,80,84,86	0
3	GOL	A	402	6/6	0.81	0.29	47,61,63,66	0
3	GOL	A	406	6/6	0.84	0.23	45,64,68,72	0
3	GOL	A	407	6/6	0.86	0.25	67,75,83,85	0
5	CL	A	409	1/1	0.88	0.11	90,90,90,90	0
3	GOL	B	101	6/6	0.90	0.14	48,69,75,77	0
3	GOL	A	404	6/6	0.92	0.14	52,63,76,86	0
4	BTB	A	408	14/14	0.93	0.15	45,61,77,84	0
3	GOL	A	403	6/6	0.93	0.38	6,17,21,21	6

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	GOL	A	401	6/6	0.96	0.12	31,46,61,70	0

6.5 Other polymers [i](#)

There are no such residues in this entry.