



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:53 PM GMT

PDB ID : 4QAM
Title : Crystal Structure of the RPGR RCC1-like domain in complex with the RPGR-interacting domain of RPGRIP1
Authors : Remans, K.; Buerger, M.; Vetter, I.R.; Wittinghofer, A.
Deposited on : 2014-05-05
Resolution : 1.83 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
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

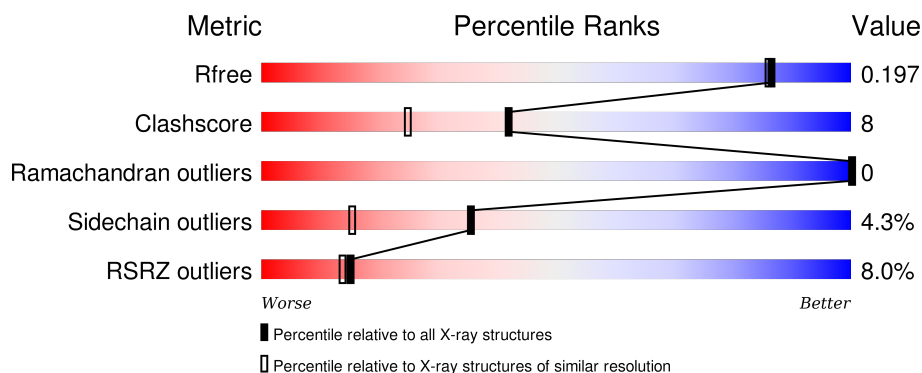
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.83 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}	91344	2634 (1.86-1.82)
Clashscore	102246	2862 (1.86-1.82)
Ramachandran outliers	100387	2831 (1.86-1.82)
Sidechain outliers	100360	2832 (1.86-1.82)
RSRZ outliers	91569	2639 (1.86-1.82)

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. 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	400	
2	B	201	

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
4	GOL	A	502	-	-	-	X
4	GOL	A	504	-	-	-	X
4	GOL	A	505	-	-	-	X
4	GOL	B	1301	-	-	-	X
4	GOL	B	1302	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4654 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 X-linked retinitis pigmentosa GTPase regulator.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	364	2824	1793	482	531	18	0	11	0

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	195	VAL	ILE	ENGINEERED MUTATION	UNP Q92834
A	393	TRP	-	EXPRESSION TAG	UNP Q92834
A	394	SER	-	EXPRESSION TAG	UNP Q92834
A	395	HIS	-	EXPRESSION TAG	UNP Q92834
A	396	PRO	-	EXPRESSION TAG	UNP Q92834
A	397	GLN	-	EXPRESSION TAG	UNP Q92834
A	398	PHE	-	EXPRESSION TAG	UNP Q92834
A	399	GLU	-	EXPRESSION TAG	UNP Q92834
A	400	LYS	-	EXPRESSION TAG	UNP Q92834

- Molecule 2 is a protein called X-linked retinitis pigmentosa GTPase regulator-interacting protein 1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
2	B	180	1464	930	233	294	2	5	0	2	0

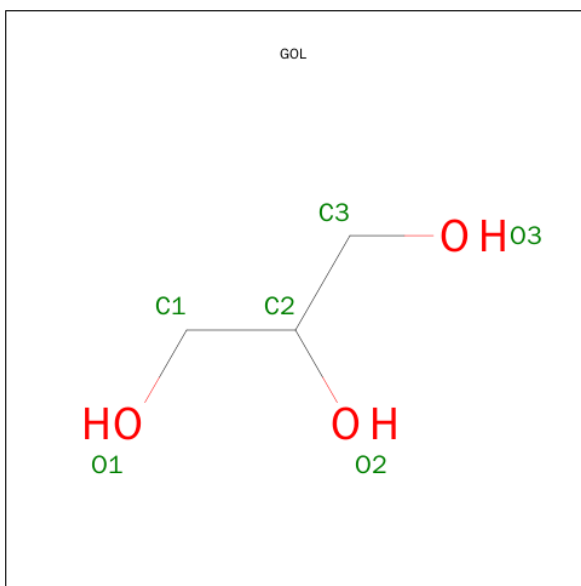
There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	1086	GLY	-	EXPRESSION TAG	UNP Q96KN7
B	1087	ALA	-	EXPRESSION TAG	UNP Q96KN7
B	1088	MSE	-	EXPRESSION TAG	UNP Q96KN7
B	1089	GLY	-	EXPRESSION TAG	UNP Q96KN7
B	1090	SER	-	EXPRESSION TAG	UNP Q96KN7

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Mg	0	0
			2	2		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	239	Total	O	0	0
			239	239		

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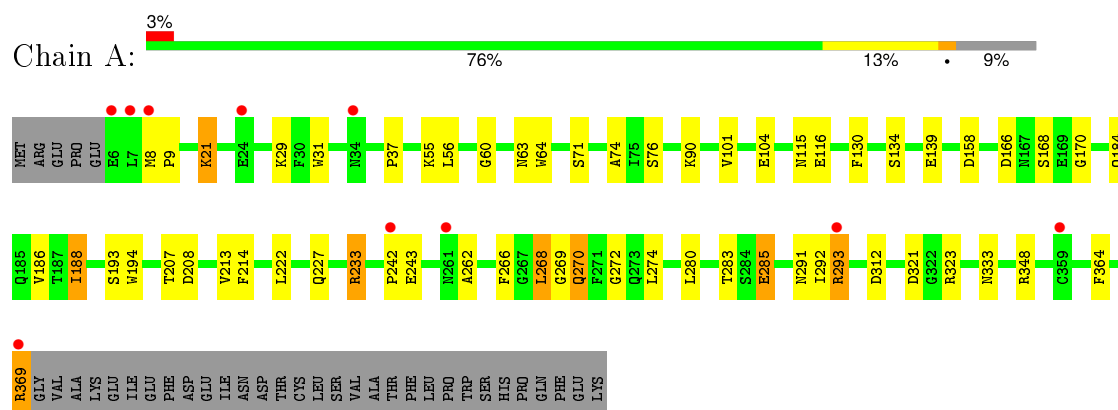
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	83	Total	O	0	0
			83	83		

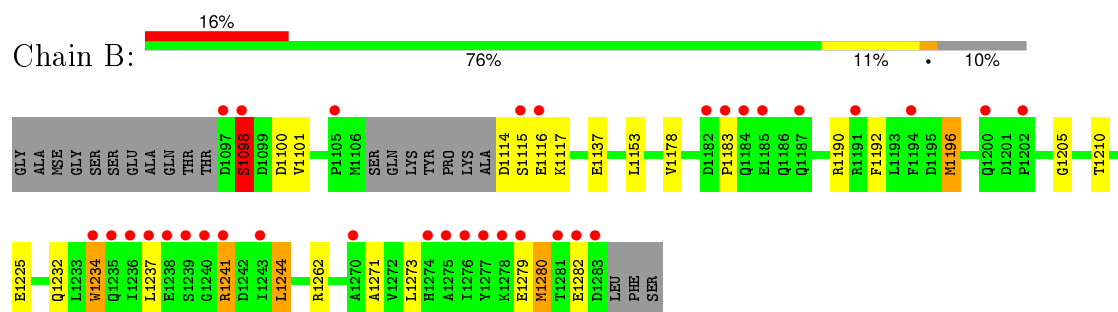
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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: X-linked retinitis pigmentosa GTPase regulator



- Molecule 2: X-linked retinitis pigmentosa GTPase regulator-interacting protein 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	63.12Å 63.80Å 161.44Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.23 – 1.83 43.23 – 1.83	Depositor EDS
% Data completeness (in resolution range)	100.0 (43.23-1.83) 100.0 (43.23-1.83)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.93 (at 1.83Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.162 , 0.197 0.163 , 0.197	Depositor DCC
R_{free} test set	2917 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	25.3	Xtriage
Anisotropy	0.166	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 44.8	EDS
Estimated twinning fraction	0.023 for k,h,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	2 of 58344 reflections (0.003%)	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	4654	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.68% 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, 0.375 respectively for untwinned datasets, and 0.333, 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, MG

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	1.23	5/2918 (0.2%)	1.10	13/3952 (0.3%)
2	B	1.08	1/1493 (0.1%)	1.10	3/2010 (0.1%)
All	All	1.18	6/4411 (0.1%)	1.10	16/5962 (0.3%)

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
2	B	0	1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	194	TRP	CD2-CE2	6.97	1.49	1.41
1	A	134	SER	CB-OG	-6.76	1.33	1.42
2	B	1234	TRP	CD2-CE2	5.71	1.48	1.41
1	A	31	TRP	CD2-CE2	5.64	1.48	1.41
1	A	64	TRP	CD2-CE2	5.40	1.47	1.41
1	A	272	GLY	N-CA	5.01	1.53	1.46

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	233	ARG	NE-CZ-NH1	10.46	125.53	120.30
1	A	233	ARG	NE-CZ-NH2	-9.27	115.66	120.30
2	B	1262	ARG	NE-CZ-NH1	-8.26	116.17	120.30
1	A	312	ASP	CB-CG-OD1	7.29	124.86	118.30
1	A	55	LYS	CD-CE-NZ	-6.91	95.81	111.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	312	ASP	CB-CG-OD2	-6.48	112.47	118.30
1	A	323	ARG	NE-CZ-NH1	-6.42	117.09	120.30
1	A	233	ARG	CG-CD-NE	-5.86	99.50	111.80
1	A	208	ASP	CB-CG-OD1	5.83	123.54	118.30
1	A	158	ASP	CB-CG-OD1	5.58	123.33	118.30
1	A	166	ASP	CB-CG-OD2	-5.42	113.42	118.30
1	A	321	ASP	CB-CG-OD1	5.19	122.97	118.30
2	B	1153	LEU	CB-CG-CD1	-5.17	102.21	111.00
2	B	1190	ARG	NE-CZ-NH1	5.11	122.86	120.30
1	A	268	LEU	CB-CG-CD1	-5.02	102.46	111.00
1	A	186	VAL	N-CA-C	-5.01	97.47	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	1098	SER	Peptide

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	2824	0	2785	44	0
2	B	1464	0	1426	24	0
3	A	2	0	0	0	0
4	A	30	0	40	4	0
4	B	12	0	16	4	0
5	A	239	0	0	14	1
5	B	83	0	0	3	0
All	All	4654	0	4267	72	1

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 8.

All (72) 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:293:ARG:HG3	1:A:293:ARG:HH11	1.15	1.09
4:A:505:GOL:H31	5:A:838:HOH:O	1.57	1.01
1:A:291[B]:ASN:ND2	5:A:834:HOH:O	1.93	0.99
1:A:333:ASN:ND2	5:A:756:HOH:O	1.95	0.96
1:A:115:ASN:HB2	5:A:722:HOH:O	1.69	0.93
1:A:270[B]:GLN:NE2	5:A:776:HOH:O	2.07	0.86
1:A:293:ARG:HG3	1:A:293:ARG:NH1	1.87	0.86
2:B:1232:GLN:HE21	2:B:1234:TRP:HB3	1.40	0.84
1:A:369:ARG:NH1	1:A:369:ARG:HB2	1.95	0.81
1:A:188:ILE:CD1	5:A:782:HOH:O	2.28	0.80
1:A:369:ARG:HH11	1:A:369:ARG:HB2	1.48	0.79
1:A:188:ILE:HD12	5:A:782:HOH:O	1.81	0.79
1:A:213[A]:VAL:HG21	1:A:222:LEU:O	1.86	0.75
2:B:1244:LEU:HD12	2:B:1244:LEU:H	1.50	0.75
1:A:242:PRO:HD2	1:A:243:GLU:OE1	1.90	0.71
1:A:8[B]:MET:CE	1:A:29:LYS:HZ1	2.04	0.71
1:A:90:LYS:HE3	1:A:104:GLU:OE2	1.91	0.70
1:A:8[B]:MET:HE2	1:A:29:LYS:HZ1	1.57	0.70
1:A:293:ARG:HH11	1:A:293:ARG:CG	2.01	0.69
2:B:1241:ARG:HB2	2:B:1241:ARG:CZ	2.23	0.69
2:B:1116:GLU:HG2	2:B:1183:PRO:HD3	1.74	0.68
1:A:262:ALA:HB2	1:A:293:ARG:HE	1.60	0.67
1:A:8[B]:MET:CE	1:A:29:LYS:NZ	2.58	0.67
2:B:1114:ASP:HB2	2:B:1117[A]:LYS:HG2	1.77	0.65
2:B:1232:GLN:NE2	2:B:1234:TRP:HB3	2.11	0.65
1:A:8[B]:MET:HE2	1:A:29:LYS:NZ	2.11	0.65
2:B:1098:SER:C	2:B:1100:ASP:H	2.00	0.64
2:B:1225:GLU:H	4:B:1302:GOL:H11	1.63	0.63
2:B:1137:GLU:HG2	5:B:1432:HOH:O	2.00	0.62
4:A:506:GOL:H31	5:A:839:HOH:O	2.01	0.60
2:B:1114:ASP:HB3	2:B:1116:GLU:H	1.68	0.57
1:A:293:ARG:HD3	5:A:668:HOH:O	2.04	0.56
1:A:9:PRO:HG3	5:A:835:HOH:O	2.05	0.56
2:B:1241:ARG:HB2	2:B:1241:ARG:NH1	2.23	0.54
2:B:1210:THR:HG21	2:B:1225:GLU:HG2	1.90	0.53
1:A:168:SER:O	1:A:233:ARG:NH2	2.42	0.53
1:A:101[A]:VAL:HG23	1:A:130:PHE:CE1	2.44	0.52
1:A:74:ALA:HB2	5:A:667:HOH:O	2.09	0.52
2:B:1244:LEU:CD1	2:B:1244:LEU:H	2.22	0.52
1:A:56:LEU:HD22	1:A:101[B]:VAL:HG11	1.92	0.51
2:B:1225:GLU:H	4:B:1302:GOL:C1	2.24	0.51
1:A:227:GLN:CD	1:A:227:GLN:H	2.14	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:1196:MSE:O	2:B:1205:GLY:HA2	2.12	0.49
2:B:1115:SER:HB3	2:B:1271:ALA:HB1	1.94	0.49
1:A:116:GLU:OE1	4:A:505:GOL:H2	2.13	0.49
1:A:21:LYS:HB2	1:A:60:GLY:HA2	1.96	0.48
1:A:227:GLN:H	1:A:227:GLN:NE2	2.12	0.48
2:B:1192:PHE:CZ	2:B:1196:MSE:HE1	2.49	0.47
2:B:1098:SER:C	2:B:1100:ASP:N	2.67	0.47
2:B:1192:PHE:CE2	2:B:1196:MSE:HE1	2.51	0.45
2:B:1098:SER:CA	2:B:1100:ASP:H	2.29	0.45
2:B:1116:GLU:CG	2:B:1183:PRO:HD3	2.46	0.44
1:A:348[B]:ARG:NH2	5:A:699:HOH:O	2.51	0.43
1:A:8[B]:MET:HE3	1:A:29:LYS:NZ	2.32	0.43
1:A:63:ASN:OD1	4:A:504:GOL:O1	2.36	0.43
1:A:193:SER:HB3	1:A:207:THR:HG22	1.99	0.43
2:B:1279:GLU:OE1	2:B:1280:MSE:HE3	2.18	0.43
1:A:269:GLY:HA3	1:A:274:LEU:HG	2.01	0.43
1:A:37:PRO:HG2	1:A:364:PHE:CD1	2.53	0.43
4:B:1301:GOL:C2	5:B:1483:HOH:O	2.67	0.43
1:A:170:GLY:HA3	1:A:233:ARG:HA	2.00	0.42
1:A:266:PHE:HA	1:A:283:THR:HG23	2.01	0.42
2:B:1117[A]:LYS:HD2	2:B:1178:VAL:HG11	2.01	0.42
1:A:21:LYS:O	1:A:76:SER:HA	2.19	0.42
1:A:90:LYS:CE	1:A:104:GLU:OE2	2.64	0.42
1:A:184:GLN:HB2	5:A:781:HOH:O	2.20	0.42
1:A:262:ALA:CB	1:A:293:ARG:HE	2.31	0.41
2:B:1196:MSE:HB2	2:B:1196:MSE:HE2	1.51	0.41
1:A:280:LEU:HA	1:A:280:LEU:HD12	1.89	0.41
1:A:285[B]:GLU:OE2	5:A:779:HOH:O	2.22	0.41
4:B:1301:GOL:H11	5:B:1483:HOH:O	2.20	0.40

All (1) 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)
5:A:689:HOH:O	5:A:829:HOH:O[4_445]	2.10	0.10

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	373/400 (93%)	365 (98%)	8 (2%)	0	100	100
2	B	178/201 (89%)	170 (96%)	8 (4%)	0	100	100
All	All	551/601 (92%)	535 (97%)	16 (3%)	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	312/334 (93%)	299 (96%)	13 (4%)	36	16
2	B	166/174 (95%)	157 (95%)	9 (5%)	27	9
All	All	478/508 (94%)	456 (95%)	22 (5%)	35	13

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	21	LYS
1	A	71	SER
1	A	139	GLU
1	A	188	ILE
1	A	214	PHE
1	A	268	LEU
1	A	270[A]	GLN
1	A	270[B]	GLN

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Mol	Chain	Res	Type
1	A	285[A]	GLU
1	A	285[B]	GLU
1	A	292	ILE
1	A	293	ARG
1	A	369	ARG
2	B	1098	SER
2	B	1101	VAL
2	B	1196	MSE
2	B	1237	LEU
2	B	1241	ARG
2	B	1244	LEU
2	B	1273	LEU
2	B	1280	MSE
2	B	1282	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	324	HIS
1	A	333	ASN
1	A	337	HIS
1	A	368	HIS
2	B	1232	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 2 are monoatomic - leaving 7 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$
4	GOL	A	502	-	5,5,5	0.35	0	5,5,5	1.23	0
4	GOL	A	503	-	5,5,5	1.43	1 (20%)	5,5,5	1.54	1 (20%)
4	GOL	A	504	-	5,5,5	1.31	1 (20%)	5,5,5	0.97	0
4	GOL	A	505	-	5,5,5	0.60	0	5,5,5	1.56	2 (40%)
4	GOL	A	506	-	5,5,5	0.44	0	5,5,5	0.73	0
4	GOL	B	1301	-	5,5,5	0.26	0	5,5,5	1.35	2 (40%)
4	GOL	B	1302	-	5,5,5	0.51	0	5,5,5	1.41	1 (20%)

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
4	GOL	A	502	-	-	0/4/4/4	0/0/0/0
4	GOL	A	503	-	-	0/4/4/4	0/0/0/0
4	GOL	A	504	-	-	0/4/4/4	0/0/0/0
4	GOL	A	505	-	-	0/4/4/4	0/0/0/0
4	GOL	A	506	-	-	0/4/4/4	0/0/0/0
4	GOL	B	1301	-	-	0/4/4/4	0/0/0/0
4	GOL	B	1302	-	-	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)
4	A	503	GOL	O2-C2	2.35	1.50	1.43
4	A	504	GOL	O3-C3	2.43	1.52	1.42

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1302	GOL	O2-C2-C3	-2.86	95.52	108.65
4	B	1301	GOL	O3-C3-C2	-2.15	99.78	110.18

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1301	GOL	O1-C1-C2	-2.09	100.06	110.18
4	A	505	GOL	O3-C3-C2	-2.04	100.31	110.18
4	A	505	GOL	O1-C1-C2	-2.03	100.32	110.18
4	A	503	GOL	C3-C2-C1	2.02	119.05	111.12

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	504	GOL	1	0
4	A	505	GOL	2	0
4	A	506	GOL	1	0
4	B	1301	GOL	2	0
4	B	1302	GOL	2	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	364/400 (91%)	-0.00	10 (2%) 58 54	16, 25, 50, 83	0
2	B	175/201 (87%)	0.73	33 (18%) 2 1	16, 32, 76, 107	0
All	All	539/601 (89%)	0.23	43 (7%) 15 14	16, 27, 63, 107	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	7	LEU	8.0
2	B	1283	ASP	7.5
2	B	1098	SER	6.5
2	B	1277	TYR	6.4
2	B	1234	TRP	5.2
2	B	1274	HIS	5.1
2	B	1282	GLU	4.8
2	B	1097	ASP	4.8
2	B	1281	THR	4.7
2	B	1241	ARG	4.2
2	B	1275	ALA	4.1
2	B	1115	SER	3.9
2	B	1278	LYS	3.8
1	A	369	ARG	3.8
1	A	242	PRO	3.7
2	B	1182	ASP	3.6
2	B	1270	ALA	3.6
2	B	1184	GLN	3.5
2	B	1183	PRO	3.5
2	B	1105	PRO	3.4
2	B	1238	GLU	3.4
2	B	1200	GLN	3.4
2	B	1239	SER	3.3
2	B	1243	ILE	3.3

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Mol	Chain	Res	Type	RSRZ
2	B	1194	PHE	3.2
2	B	1237	LEU	3.2
1	A	293	ARG	3.2
2	B	1235	GLN	3.1
2	B	1187	GLN	3.0
1	A	6	GLU	3.0
2	B	1240	GLY	2.9
2	B	1276	ILE	2.9
2	B	1279	GLU	2.7
2	B	1236	ILE	2.5
2	B	1185	GLU	2.4
2	B	1116	GLU	2.3
1	A	261	ASN	2.2
2	B	1202	PRO	2.2
2	B	1191	ARG	2.2
1	A	34	ASN	2.2
1	A	8[A]	MET	2.1
1	A	24	GLU	2.1
1	A	359[A]	CYS	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. 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	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
4	GOL	B	1301	6/6	0.72	0.26	6.93	46,67,69,70	0
4	GOL	A	502	6/6	0.92	0.16	6.86	20,24,26,31	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	GOL	B	1302	6/6	0.77	0.15	5.08	42,45,53,60	0
4	GOL	A	505	6/6	0.69	0.17	3.36	47,52,56,59	0
4	GOL	A	504	6/6	0.93	0.14	2.91	22,37,38,41	0
4	GOL	A	503	6/6	0.94	0.11	0.12	16,22,26,30	0
3	MG	A	507	1/1	0.99	0.10	-1.56	28,28,28,28	1
3	MG	A	501	1/1	0.93	0.15	-	49,49,49,49	0
4	GOL	A	506	6/6	0.86	0.24	-	45,56,59,64	0

6.5 Other polymers

There are no such residues in this entry.