



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 11:16 AM GMT

PDB ID : 3OG6
Title : The crystal structure of human interferon lambda 1 complexed with its high affinity receptor in space group P212121
Authors : Miknis, Z.J.; Magracheva, E.; Lei, W.; Zdanov, A.; Kotenko, S.V.; Wlodawer, A.
Deposited on : 2010-08-16
Resolution : 2.10 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

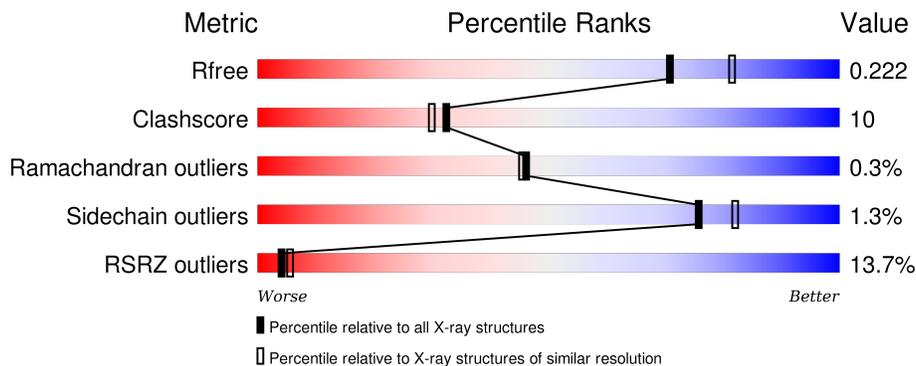
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	3939 (2.10-2.10)
Clashscore	102246	4460 (2.10-2.10)
Ramachandran outliers	100387	4413 (2.10-2.10)
Sidechain outliers	100360	4414 (2.10-2.10)
RSRZ outliers	91569	3948 (2.10-2.10)

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	196	
2	B	226	

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	183	-	-	-	X
3	GOL	B	225	-	-	-	X
3	GOL	B	226	-	-	-	X
4	NAG	B	301	-	-	-	X

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 3082 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 Interleukin-29.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	143	1122	714	203	202	3	0	0	0

There are 17 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-14	HIS	-	EXPRESSION TAG	UNP Q8IU54
A	-13	HIS	-	EXPRESSION TAG	UNP Q8IU54
A	-12	HIS	-	EXPRESSION TAG	UNP Q8IU54
A	-11	HIS	-	EXPRESSION TAG	UNP Q8IU54
A	-10	HIS	-	EXPRESSION TAG	UNP Q8IU54
A	-9	HIS	-	EXPRESSION TAG	UNP Q8IU54
A	-8	ASP	-	EXPRESSION TAG	UNP Q8IU54
A	-7	TYR	-	EXPRESSION TAG	UNP Q8IU54
A	-6	LYS	-	EXPRESSION TAG	UNP Q8IU54
A	-5	ASP	-	EXPRESSION TAG	UNP Q8IU54
A	-4	ASP	-	EXPRESSION TAG	UNP Q8IU54
A	-3	ASP	-	EXPRESSION TAG	UNP Q8IU54
A	-2	ASP	-	EXPRESSION TAG	UNP Q8IU54
A	-1	LYS	-	EXPRESSION TAG	UNP Q8IU54
A	0	ALA	-	EXPRESSION TAG	UNP Q8IU54
A	1	ASP	GLY	ENGINEERED MUTATION	UNP Q8IU54
A	10	PRO	THR	ENGINEERED MUTATION	UNP Q8IU54

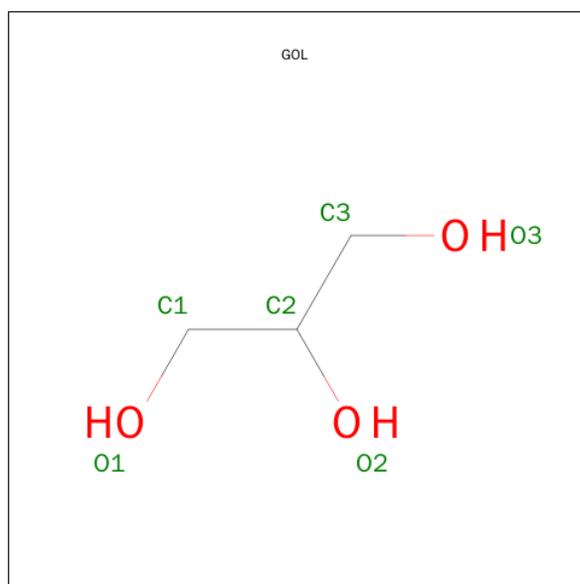
- Molecule 2 is a protein called Interleukin 28 receptor, alpha (Interferon, lambda receptor).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	200	1603	1039	262	293	9	0	1	0

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	207	ALA	-	EXPRESSION TAG	UNP Q5VTX7
B	208	SER	-	EXPRESSION TAG	UNP Q5VTX7
B	209	GLY	-	EXPRESSION TAG	UNP Q5VTX7
B	210	SER	-	EXPRESSION TAG	UNP Q5VTX7
B	211	SER	-	EXPRESSION TAG	UNP Q5VTX7
B	212	GLY	-	EXPRESSION TAG	UNP Q5VTX7
B	213	GLY	-	EXPRESSION TAG	UNP Q5VTX7
B	214	SER	-	EXPRESSION TAG	UNP Q5VTX7
B	215	SER	-	EXPRESSION TAG	UNP Q5VTX7
B	216	GLY	-	EXPRESSION TAG	UNP Q5VTX7
B	217	THR	-	EXPRESSION TAG	UNP Q5VTX7
B	218	SER	-	EXPRESSION TAG	UNP Q5VTX7
B	219	HIS	-	EXPRESSION TAG	UNP Q5VTX7
B	220	HIS	-	EXPRESSION TAG	UNP Q5VTX7
B	221	HIS	-	EXPRESSION TAG	UNP Q5VTX7
B	222	HIS	-	EXPRESSION TAG	UNP Q5VTX7
B	223	HIS	-	EXPRESSION TAG	UNP Q5VTX7
B	224	HIS	-	EXPRESSION TAG	UNP Q5VTX7

- 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	B	1	Total	C	O	0	0
			6	3	3		

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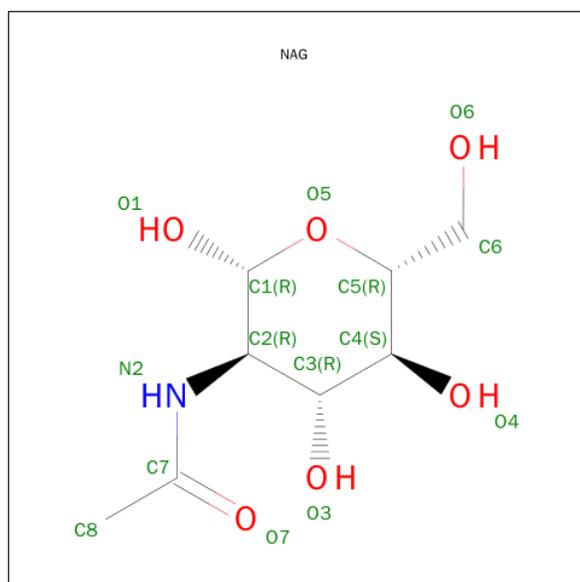
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
3	B	1	6	3	3	0	0

- Molecule 4 is a polymer of unknown type called SUGAR (3-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
4	B	3	39	22	2	15	0	0

- Molecule 5 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



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

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
6	A	114	114	114	0	0
6	B	152	152	152	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	64.96Å 85.79Å 116.51Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.39 – 2.10 29.39 – 2.10	Depositor EDS
% Data completeness (in resolution range)	90.0 (29.39-2.10) 90.0 (29.39-2.10)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.37 (at 2.10Å)	Xtrriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, R_{free}	0.184 , 0.228 0.177 , 0.222	Depositor DCC
R_{free} test set	1072 reflections (3.03%)	DCC
Wilson B-factor (Å ²)	37.5	Xtrriage
Anisotropy	0.399	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 66.4	EDS
Estimated twinning fraction	No twinning to report.	Xtrriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Outliers	0 of 37762 reflections	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3082	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.55% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, BMA, NAG

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	0.99	1/1146 (0.1%)	0.91	2/1555 (0.1%)
2	B	0.78	1/1656 (0.1%)	0.79	0/2263
All	All	0.87	2/2802 (0.1%)	0.84	2/3818 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	39	GLU	CG-CD	5.46	1.60	1.51
2	B	16	ASN	CB-CG	5.05	1.62	1.51

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	35	ARG	NE-CZ-NH1	5.16	122.88	120.30
1	A	133	ARG	NE-CZ-NH1	-5.10	117.75	120.30

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	1122	0	1130	16	1
2	B	1603	0	1586	44	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	12	0	16	2	0
3	B	12	0	16	3	0
4	B	39	0	34	0	0
5	B	28	0	26	0	0
6	A	114	0	0	1	0
6	B	152	0	0	4	1
All	All	3082	0	2808	59	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 10.

All (59) 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)
2:B:49:ARG:HD3	6:B:338:HOH:O	1.69	0.92
2:B:79:ARG:HD3	2:B:91:TRP:CE3	2.14	0.83
2:B:144:LYS:HD3	2:B:172:GLU:OE1	1.88	0.73
2:B:6:PRO:HG3	2:B:90:PRO:HD2	1.70	0.72
2:B:35:TYR:CE1	2:B:82:THR:HG22	2.24	0.72
1:A:159:THR:O	1:A:163:LYS:HE2	1.90	0.72
2:B:82:THR:OG1	2:B:89:SER:HB3	1.95	0.66
2:B:120:SER:HB3	2:B:163:GLN:OE1	1.97	0.64
2:B:79:ARG:HD3	2:B:91:TRP:CD2	2.33	0.64
2:B:202:VAL:HB	2:B:203:PRO:HD3	1.80	0.64
1:A:129:HIS:HB2	6:A:267:HOH:O	1.99	0.62
2:B:38:ALA:HB3	2:B:79:ARG:HD2	1.82	0.62
2:B:35:TYR:HE1	2:B:82:THR:HG22	1.67	0.60
2:B:48:TRP:CZ3	2:B:79:ARG:HG2	2.37	0.59
1:A:76:LEU:HD21	1:A:105:ILE:HG22	1.85	0.57
1:A:32:LYS:HD2	2:B:101:PHE:CE2	2.40	0.57
1:A:97:GLN:HB2	1:A:98:PRO:HD3	1.86	0.57
2:B:86:SER:O	2:B:87:SER:HB2	2.07	0.55
2:B:16:ASN:O	2:B:17:PHE:HB2	2.07	0.54
2:B:166:LEU:HD13	2:B:200:LEU:HD22	1.91	0.53
1:A:33:LYS:HE2	3:A:183:GOL:H31	1.91	0.52
2:B:156:THR:CG2	2:B:157:PRO:HD2	2.40	0.52
1:A:61:ARG:HD3	1:A:164:TYR:CZ	2.45	0.51
2:B:150:LYS:NZ	3:B:226:GOL:H2	2.25	0.51
1:A:47:TRP:CE3	1:A:47:TRP:O	2.63	0.51
2:B:79:ARG:HD3	2:B:91:TRP:CZ3	2.47	0.49
2:B:47[A]:ARG:HD3	6:B:368:HOH:O	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:83:VAL:HG12	2:B:88:LYS:HG2	1.94	0.48
2:B:27:LEU:HD12	2:B:27:LEU:H	1.79	0.47
2:B:41:SER:HB3	2:B:49:ARG:HE	1.80	0.47
2:B:48:TRP:CE3	2:B:79:ARG:HG2	2.50	0.47
1:A:76:LEU:HD21	1:A:105:ILE:CG2	2.45	0.46
2:B:16:ASN:HB2	6:B:239:HOH:O	2.16	0.46
1:A:33:LYS:HE2	3:A:183:GOL:C3	2.45	0.46
2:B:8:GLN:OE1	2:B:27:LEU:HG	2.15	0.46
1:A:157:LEU:C	1:A:157:LEU:HD23	2.36	0.46
2:B:156:THR:HG23	2:B:157:PRO:HD2	1.98	0.45
2:B:168:PRO:HA	2:B:202:VAL:CG2	2.46	0.45
2:B:29:ASN:N	2:B:30:PRO:HD3	2.32	0.45
2:B:13:LEU:HD12	2:B:13:LEU:C	2.37	0.45
2:B:53:GLU:HG3	6:B:344:HOH:O	2.16	0.45
2:B:84:SER:HB2	2:B:85:PRO:HD2	1.99	0.45
2:B:71:ASP:CG	3:B:225:GOL:H32	2.37	0.44
2:B:7:PRO:HG3	2:B:82:THR:HG23	1.98	0.44
2:B:84:SER:HB2	2:B:85:PRO:CD	2.47	0.44
1:A:47:TRP:O	1:A:47:TRP:HE3	2.01	0.44
2:B:47[A]:ARG:HB2	2:B:47[A]:ARG:CZ	2.48	0.44
1:A:97:GLN:N	1:A:98:PRO:CD	2.81	0.43
1:A:137:ALA:HB3	1:A:138:PRO:HD3	1.99	0.43
2:B:202:VAL:CB	2:B:203:PRO:HD3	2.49	0.42
2:B:71:ASP:OD2	3:B:225:GOL:H32	2.19	0.42
2:B:174:HIS:CE1	2:B:200:LEU:HD12	2.55	0.42
2:B:156:THR:CG2	2:B:157:PRO:CD	2.96	0.42
1:A:68:ARG:N	1:A:69:PRO:CD	2.83	0.41
2:B:61:LEU:C	2:B:61:LEU:HD12	2.40	0.41
2:B:4:LEU:HG	2:B:87:SER:HB3	2.02	0.41
2:B:200:LEU:HA	2:B:200:LEU:HD23	1.82	0.41
1:A:136:GLU:HB3	1:A:140:LYS:HD3	2.01	0.41
2:B:68:LYS:O	2:B:69:LYS:HB2	2.21	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)
1:A:96:ASP:OD1	6:B:283:HOH:O[4_466]	2.16	0.04

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	139/196 (71%)	137 (99%)	2 (1%)	0	100	100
2	B	199/226 (88%)	183 (92%)	15 (8%)	1 (0%)	34	30
All	All	338/422 (80%)	320 (95%)	17 (5%)	1 (0%)	46	45

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	30	PRO

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	120/170 (71%)	118 (98%)	2 (2%)	68	74
2	B	182/201 (90%)	180 (99%)	2 (1%)	80	85
All	All	302/371 (81%)	298 (99%)	4 (1%)	76	82

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	26	GLN
1	A	133	ARG
2	B	79	ARG
2	B	135	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](#)

3 carbohydrates 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
4	NAG	B	301	2,4	14,14,15	0.62	0	15,19,21	0.83	0
4	NAG	B	302	4	14,14,15	0.54	0	15,19,21	0.58	0
4	BMA	B	303	4	11,11,12	0.68	0	14,15,17	0.62	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
4	NAG	B	301	2,4	-	0/6/23/26	0/1/1/1
4	NAG	B	302	4	-	0/6/23/26	0/1/1/1
4	BMA	B	303	4	-	0/2/19/22	0/1/1/1

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.

5.6 Ligand geometry [i](#)

6 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
3	GOL	A	182	-	5,5,5	0.41	0	5,5,5	0.52	0
3	GOL	A	183	-	5,5,5	0.41	0	5,5,5	0.20	0
3	GOL	B	225	-	5,5,5	0.43	0	5,5,5	0.95	0
3	GOL	B	226	-	5,5,5	0.22	0	5,5,5	0.83	0
5	NAG	B	300	2	14,14,15	0.55	0	15,19,21	0.59	0
5	NAG	B	304	2	14,14,15	0.96	1 (7%)	15,19,21	1.83	2 (13%)

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	182	-	-	0/4/4/4	0/0/0/0
3	GOL	A	183	-	-	0/4/4/4	0/0/0/0
3	GOL	B	225	-	-	0/4/4/4	0/0/0/0
3	GOL	B	226	-	-	0/4/4/4	0/0/0/0
5	NAG	B	300	2	-	0/6/23/26	0/1/1/1
5	NAG	B	304	2	-	0/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	304	NAG	C1-C2	2.96	1.56	1.52

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	304	NAG	O5-C5-C6	2.24	112.20	107.35
5	B	304	NAG	C1-O5-C5	5.57	119.32	112.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	183	GOL	2	0
3	B	225	GOL	2	0
3	B	226	GOL	1	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	143/196 (72%)	0.44	19 (13%) 4 6	28, 41, 93, 148	0
2	B	200/226 (88%)	0.31	28 (14%) 4 5	31, 58, 134, 203	0
All	All	343/422 (81%)	0.36	47 (13%) 4 5	28, 52, 128, 203	0

All (47) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	170	LEU	8.7
2	B	86	SER	8.1
2	B	87	SER	6.7
2	B	27	LEU	6.6
2	B	32	ASP	6.1
2	B	85	PRO	5.5
2	B	203	PRO	5.4
2	B	84	SER	5.3
2	B	33	VAL	5.0
1	A	115	PRO	4.7
1	A	168	GLY	4.6
2	B	28	GLY	4.4
1	A	44	LEU	4.2
1	A	111	ALA	4.1
2	B	202	VAL	4.1
1	A	169	ASN	3.9
1	A	114	GLN	3.7
2	B	56	GLY	3.6
2	B	4	LEU	3.6
2	B	168	PRO	3.6
1	A	154	LEU	3.6
2	B	166	LEU	3.4
1	A	167	ASP	3.4
2	B	83	VAL	3.3

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Mol	Chain	Res	Type	RSRZ
2	B	26	GLY	3.2
1	A	113	ILE	3.2
2	B	117	GLU	3.2
2	B	115	THR	3.1
2	B	31	GLN	3.1
1	A	158	LEU	3.1
2	B	148	GLY	3.1
1	A	46	ASN	3.0
2	B	29	ASN	3.0
1	A	157	LEU	2.9
2	B	169	ALA	2.9
2	B	5	ALA	2.9
2	B	88	LYS	2.8
1	A	66	ARG	2.8
1	A	112	CYS	2.7
2	B	58	LYS	2.7
1	A	31	PHE	2.7
2	B	116	GLU	2.6
1	A	155	PHE	2.4
1	A	65	VAL	2.4
2	B	24	LEU	2.3
1	A	102	LEU	2.2
2	B	34	THR	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](#)

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	NAG	B	301	14/15	0.93	0.17	13.41	63,71,75,77	0
4	BMA	B	303	11/12	0.69	0.42	-	125,139,144,145	0
4	NAG	B	302	14/15	0.80	0.28	-	57,90,101,114	0

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
3	GOL	B	226	6/6	0.84	0.20	7.63	81,87,95,99	0
3	GOL	B	225	6/6	0.84	0.27	3.20	51,82,86,90	0
3	GOL	A	183	6/6	0.90	0.20	2.73	90,95,102,106	0
5	NAG	B	304	14/15	0.81	0.20	-	82,96,103,111	0
5	NAG	B	300	14/15	0.76	0.37	-	116,133,144,148	0
3	GOL	A	182	6/6	0.61	0.44	-	97,106,115,118	0

6.5 Other polymers [i](#)

There are no such residues in this entry.