



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 06:55 pm GMT

PDB ID : 4R08
Title : Crystal structure of human TLR8 in complex with ssRNA40
Authors : Tanji, H.; Ohto, U.; Shimizu, T.
Deposited on : 2014-07-30
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

<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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : recalc28986

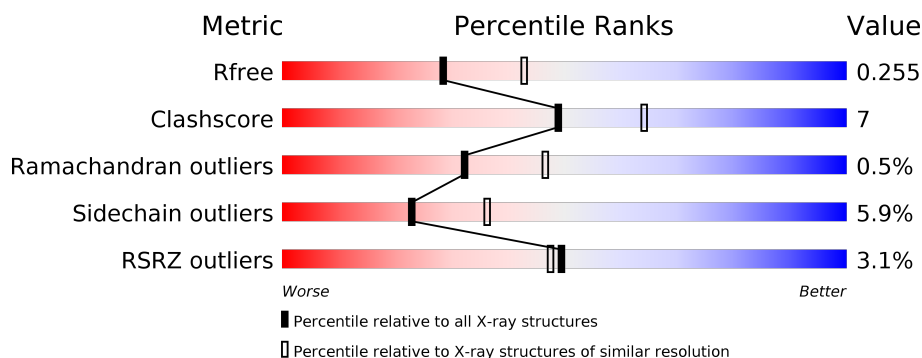
1 Overall quality at a glance

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}	100719	3166 (2.40-2.40)
Clashscore	112137	3674 (2.40-2.40)
Ramachandran outliers	110173	3616 (2.40-2.40)
Sidechain outliers	110143	3617 (2.40-2.40)
RSRZ outliers	101464	3195 (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. 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	811	<div> <div>77%</div> <div>14%</div> <div>8%</div> </div>
1	B	811	<div> <div>75%</div> <div>15%</div> <div>9%</div> </div>
1	C	811	<div> <div>66%</div> <div>20%</div> <div>10%</div> </div>
1	D	811	<div> <div>72%</div> <div>18%</div> <div>8%</div> </div>

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	NAG	A	911	-	-	-	X
5	NAG	D	913	-	-	-	X

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 25676 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 Toll-like receptor 8.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	743	Total	C	N	O	S	4	0	0
			5992	3835	1017	1121	19			
1	B	737	Total	C	N	O	S	4	0	0
			5941	3806	1009	1107	19			
1	C	731	Total	C	N	O	S	4	0	0
			5898	3779	1000	1100	19			
1	D	743	Total	C	N	O	S	4	0	0
			5989	3834	1016	1120	19			

There are 40 discrepancies between the modelled and reference sequences:

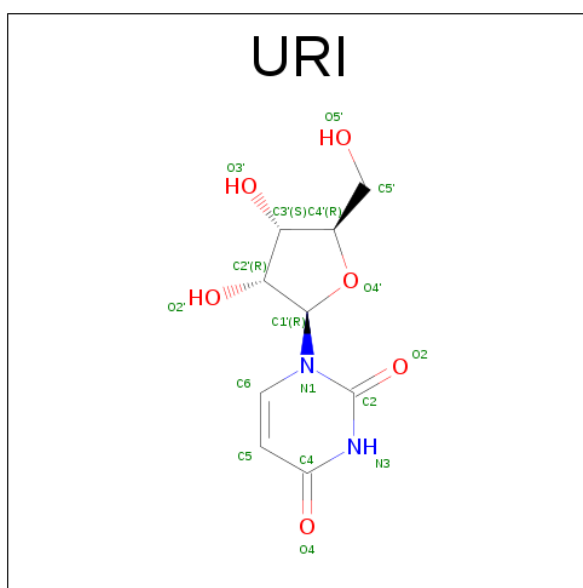
Chain	Residue	Modelled	Actual	Comment	Reference
A	22	ARG	-	EXPRESSION TAG	UNP Q9NR97
A	23	SER	-	EXPRESSION TAG	UNP Q9NR97
A	24	PRO	-	EXPRESSION TAG	UNP Q9NR97
A	25	TRP	-	EXPRESSION TAG	UNP Q9NR97
A	828	GLU	-	EXPRESSION TAG	UNP Q9NR97
A	829	PHE	-	EXPRESSION TAG	UNP Q9NR97
A	830	LEU	-	EXPRESSION TAG	UNP Q9NR97
A	831	VAL	-	EXPRESSION TAG	UNP Q9NR97
A	832	PRO	-	EXPRESSION TAG	UNP Q9NR97
A	833	ARG	-	EXPRESSION TAG	UNP Q9NR97
B	22	ARG	-	EXPRESSION TAG	UNP Q9NR97
B	23	SER	-	EXPRESSION TAG	UNP Q9NR97
B	24	PRO	-	EXPRESSION TAG	UNP Q9NR97
B	25	TRP	-	EXPRESSION TAG	UNP Q9NR97
B	828	GLU	-	EXPRESSION TAG	UNP Q9NR97
B	829	PHE	-	EXPRESSION TAG	UNP Q9NR97
B	830	LEU	-	EXPRESSION TAG	UNP Q9NR97
B	831	VAL	-	EXPRESSION TAG	UNP Q9NR97
B	832	PRO	-	EXPRESSION TAG	UNP Q9NR97
B	833	ARG	-	EXPRESSION TAG	UNP Q9NR97
C	22	ARG	-	EXPRESSION TAG	UNP Q9NR97

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Chain	Residue	Modelled	Actual	Comment	Reference
C	23	SER	-	EXPRESSION TAG	UNP Q9NR97
C	24	PRO	-	EXPRESSION TAG	UNP Q9NR97
C	25	TRP	-	EXPRESSION TAG	UNP Q9NR97
C	828	GLU	-	EXPRESSION TAG	UNP Q9NR97
C	829	PHE	-	EXPRESSION TAG	UNP Q9NR97
C	830	LEU	-	EXPRESSION TAG	UNP Q9NR97
C	831	VAL	-	EXPRESSION TAG	UNP Q9NR97
C	832	PRO	-	EXPRESSION TAG	UNP Q9NR97
C	833	ARG	-	EXPRESSION TAG	UNP Q9NR97
D	22	ARG	-	EXPRESSION TAG	UNP Q9NR97
D	23	SER	-	EXPRESSION TAG	UNP Q9NR97
D	24	PRO	-	EXPRESSION TAG	UNP Q9NR97
D	25	TRP	-	EXPRESSION TAG	UNP Q9NR97
D	828	GLU	-	EXPRESSION TAG	UNP Q9NR97
D	829	PHE	-	EXPRESSION TAG	UNP Q9NR97
D	830	LEU	-	EXPRESSION TAG	UNP Q9NR97
D	831	VAL	-	EXPRESSION TAG	UNP Q9NR97
D	832	PRO	-	EXPRESSION TAG	UNP Q9NR97
D	833	ARG	-	EXPRESSION TAG	UNP Q9NR97

- Molecule 2 is URIDINE (three-letter code: URI) (formula: C₉H₁₂N₂O₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			17	9	2	6		
2	B	1	Total	C	N	O	0	0
			17	9	2	6		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	C	1	Total	C	N	O	0	0
			17	9	2	6		
2	D	1	Total	C	N	O	0	0
			17	9	2	6		

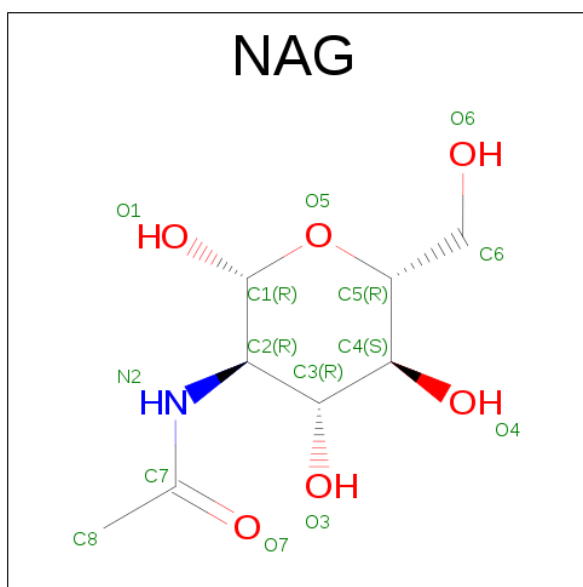
- Molecule 3 is a polymer of unknown type called SUGAR (3-MER).

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

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

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

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



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

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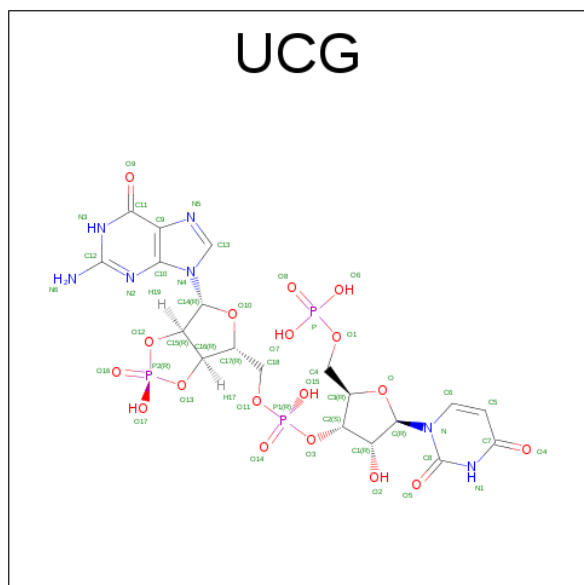
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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	D	1	Total	C	N	O	0	0
			14	8	1	5		
5	D	1	Total	C	N	O	0	0
			14	8	1	5		
5	D	1	Total	C	N	O	0	0
			14	8	1	5		
5	D	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 6 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	2	Total	C	N	O	0	0
			28	16	2	10		
6	B	2	Total	C	N	O	0	0
			28	16	2	10		
6	C	2	Total	C	N	O	0	0
			28	16	2	10		
6	D	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 7 is 3'-O-[(R)-{[(2R,3AR,4R,6R,6AR)-6-(2-AMINO-6-OXO-1,6-DIHYDRO-9H-PURIN-9-YL)-2-HYDROXY-2-OXIDOTETRAHYDROFURO[3,4-D][1,3,2]DIOXAPHOSPHOL-4-YL]METHOXY}(HYDROXY)PHOSPHORYL]URIDINE 5'-(DIHYDROGEN PHOSPHATE) (three-letter code: UCG) (formula: C₁₉H₂₄N₇O₁₈P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	A	1	Total 47	C 19	N 7	O 18	P 3	0	0
7	B	1	Total 47	C 19	N 7	O 18	P 3	0	0
7	C	1	Total 47	C 19	N 7	O 18	P 3	0	0
7	D	1	Total 47	C 19	N 7	O 18	P 3	0	0

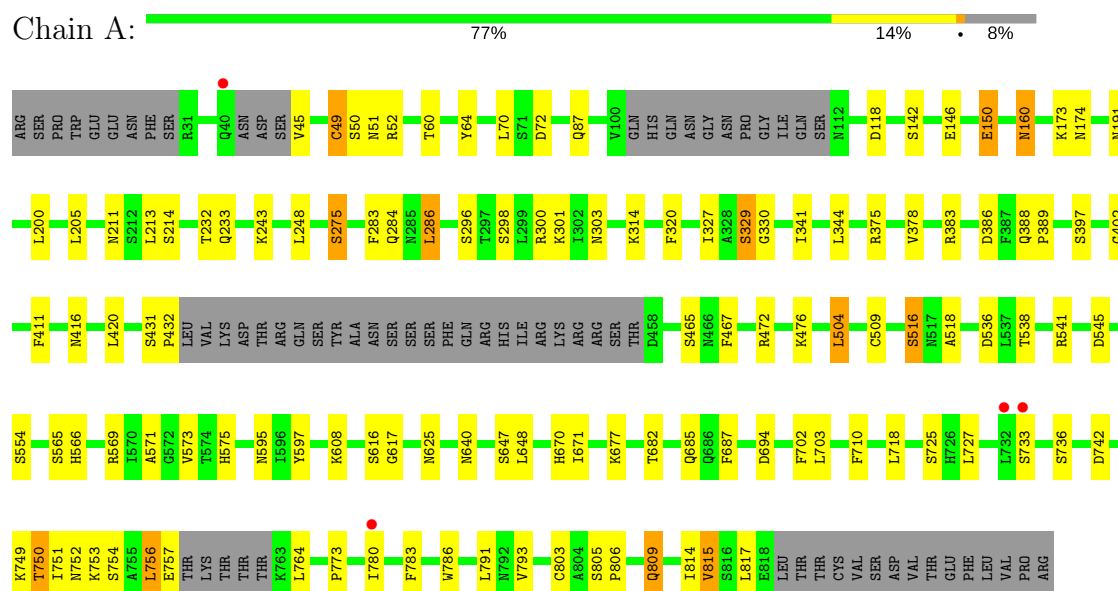
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	300	Total 300	O 300	0	0
8	B	250	Total 250	O 250	0	0
8	C	192	Total 192	O 192	0	0
8	D	182	Total 182	O 182	0	0

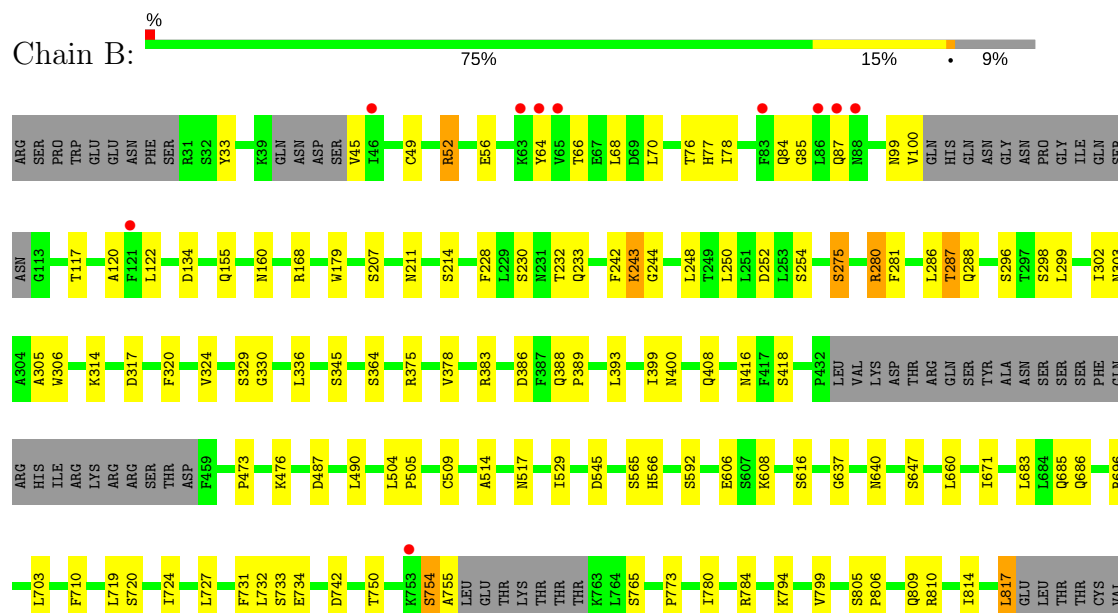
3 Residue-property plots

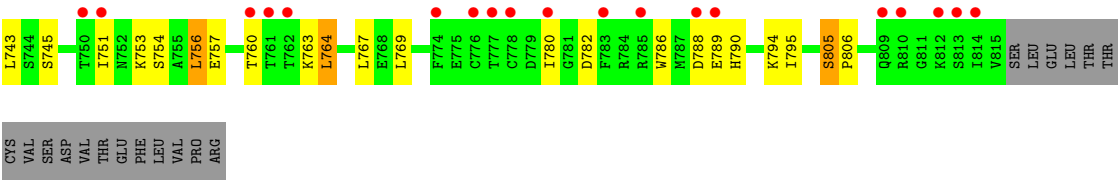
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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: Toll-like receptor 8



• Molecule 1: Toll-like receptor 8





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	86.32Å 139.69Å 167.57Å 90.00° 91.44° 90.00°	Depositor
Resolution (Å)	40.98 – 2.40 40.98 – 2.40	Depositor EDS
% Data completeness (in resolution range)	92.8 (40.98-2.40) 92.9 (40.98-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.83 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.8.0073	Depositor
R, R_{free}	0.190 , 0.258 0.194 , 0.255	Depositor DCC
R_{free} test set	7227 reflections (5.29%)	DCC
Wilson B-factor (Å ²)	34.6	Xtriage
Anisotropy	0.031	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 49.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.037 for h,-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	25676	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.15% 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: UCG, BMA, NAG, URI

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	0/6115	0.64	0/8289
1	B	0.48	1/6064 (0.0%)	0.60	0/8220
1	C	0.45	0/6021	0.58	0/8164
1	D	0.48	1/6113 (0.0%)	0.61	2/8289 (0.0%)
All	All	0.48	2/24313 (0.0%)	0.61	2/32962 (0.0%)

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	C	0	2
1	D	0	1
All	All	0	3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	476	LYS	CB-CG	-11.87	1.20	1.52
1	B	476	LYS	CB-CG	-6.11	1.36	1.52

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	805	SER	C-N-CD	-6.52	106.25	120.60
1	D	476	LYS	CA-CB-CG	6.21	127.06	113.40

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	218	PRO	Peptide
1	C	735	VAL	Peptide
1	D	169	LEU	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	5992	0	5972	57	0
1	B	5941	0	5931	67	0
1	C	5898	0	5878	126	0
1	D	5989	0	5978	84	0
2	A	17	0	12	0	0
2	B	17	0	12	0	0
2	C	17	0	12	0	0
2	D	17	0	12	0	0
3	A	39	0	34	2	0
3	B	78	0	68	0	0
3	C	78	0	68	0	0
3	D	78	0	68	2	0
4	A	39	0	34	0	0
5	A	70	0	65	1	0
5	B	70	0	65	2	0
5	C	56	0	52	0	0
5	D	56	0	52	1	0
6	A	28	0	25	0	0
6	B	28	0	25	0	0
6	C	28	0	25	0	0
6	D	28	0	25	0	0
7	A	47	0	20	2	0
7	B	47	0	20	1	0
7	C	47	0	20	3	0
7	D	47	0	20	1	0
8	A	300	0	0	6	0
8	B	250	0	0	5	0
8	C	192	0	0	7	0
8	D	182	0	0	4	0
All	All	25676	0	24493	336	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 7.

All (336) 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:C:219:LYS:HA	1:C:220:LEU:HB2	1.26	1.09
1:B:280:ARG:O	1:B:280:ARG:HD3	1.66	0.94
1:C:219:LYS:HA	1:C:220:LEU:CB	2.03	0.88
1:C:219:LYS:CA	1:C:220:LEU:HB2	2.08	0.81
1:C:708:SER:HB3	1:C:734:GLU:HG3	1.62	0.81
1:A:536:ASP:OD2	8:A:1049:HOH:O	1.97	0.80
1:C:99:ASN:O	1:C:100:VAL:HG13	1.80	0.80
1:C:303:ASN:OD1	8:C:1188:HOH:O	2.00	0.79
1:B:731:PHE:HA	1:B:734:GLU:HG3	1.66	0.76
1:B:280:ARG:HD3	1:B:280:ARG:C	2.05	0.76
1:C:230:SER:OG	1:C:254:SER:OG	1.76	0.75
1:B:242:PHE:O	1:B:244:GLY:N	2.21	0.73
1:D:150:GLU:HG2	1:D:174:ASN:HB2	1.70	0.73
1:D:751:ILE:HG22	1:D:756:LEU:HD21	1.69	0.73
1:D:239:GLU:HG3	1:D:284:GLN:HE21	1.55	0.71
1:D:730:GLY:N	1:D:754:SER:O	2.24	0.70
1:B:732:LEU:HB2	1:B:755:ALA:O	1.91	0.70
1:C:219:LYS:NZ	1:C:240:GLU:O	2.25	0.70
1:B:487:ASP:OD2	8:B:1157:HOH:O	2.10	0.69
1:B:660:LEU:CD2	1:B:686:GLN:HG3	2.24	0.67
1:C:192:ILE:HG21	1:C:217:PRO:HB3	1.77	0.66
1:C:715:ARG:NH2	1:C:715:ARG:HG3	2.10	0.66
1:C:50:SER:O	1:C:52:ARG:HG3	1.96	0.65
1:D:622:ILE:O	8:D:1158:HOH:O	2.14	0.65
1:D:693:LEU:HD21	1:D:695:LEU:HD11	1.79	0.65
1:C:150:GLU:CG	1:C:174:ASN:HD22	2.11	0.64
1:D:205:LEU:HD23	1:D:205:LEU:C	2.18	0.64
1:C:732:LEU:C	1:C:734:GLU:H	2.02	0.63
1:C:715:ARG:HH21	1:C:715:ARG:HG3	1.62	0.63
1:B:383:ARG:HB2	1:B:386:ASP:OD2	1.99	0.63
1:D:416:ASN:ND2	8:D:1023:HOH:O	2.31	0.63
1:C:728:PRO:O	1:C:731:PHE:HB3	1.98	0.62
1:B:720:SER:HB2	1:B:742:ASP:OD2	1.99	0.62
1:C:300:ARG:NE	8:C:1085:HOH:O	2.19	0.61
1:D:736:SER:OG	1:D:760:THR:HG21	2.01	0.61
1:C:270:CYS:O	1:C:273:GLY:N	2.32	0.60
1:D:751:ILE:CG2	1:D:756:LEU:HD21	2.32	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:49:CYS:HB3	1:A:70:LEU:HD23	1.83	0.60
1:D:467:PHE:HB3	3:D:903:NAG:H81	1.84	0.60
1:A:383:ARG:HB2	1:A:386:ASP:OD2	2.02	0.59
1:A:411:PHE:HB3	1:A:504:LEU:HD13	1.84	0.59
1:A:160:ASN:ND2	1:A:191:ASN:OD1	2.36	0.59
1:B:299:LEU:HD13	1:B:302:ILE:CD1	2.32	0.59
1:D:275:SER:HA	1:D:298:SER:HB2	1.83	0.59
1:D:536:ASP:OD1	1:D:538:THR:HG23	2.03	0.58
1:A:205:LEU:C	1:A:205:LEU:HD23	2.24	0.58
1:C:81:GLU:O	1:C:84:GLN:HB3	2.03	0.58
1:D:411:PHE:HB3	1:D:504:LEU:HD13	1.86	0.58
1:B:84:GLN:HG2	1:B:85:GLY:N	2.19	0.57
1:B:99:ASN:O	1:B:100:VAL:HB	2.04	0.57
1:C:162:THR:O	1:C:196:VAL:HG21	2.04	0.57
1:C:595:ASN:HB3	8:C:1042:HOH:O	2.05	0.57
1:C:326:GLU:OE1	1:C:330:GLY:HA2	2.04	0.57
1:D:211:ASN:O	1:D:232:THR:HA	2.05	0.57
1:C:341:ILE:HD12	7:C:901:UCG:C8	2.36	0.56
1:C:280:ARG:O	1:C:280:ARG:HD3	2.06	0.56
1:C:219:LYS:HG3	1:C:220:LEU:O	2.05	0.56
1:B:287:THR:HG22	1:B:288:GLN:HE21	1.71	0.56
1:A:687:PHE:O	8:A:1132:HOH:O	2.18	0.56
1:C:79:THR:C	1:C:120:ALA:HB1	2.27	0.56
1:C:317:ASP:OD1	1:C:319:GLU:OE1	2.24	0.55
1:D:207:SER:HA	1:D:228:PHE:HB2	1.88	0.55
1:A:211:ASN:O	1:A:232:THR:HA	2.06	0.55
1:A:518:ALA:HA	1:A:541:ARG:O	2.05	0.55
1:C:249:THR:HA	1:C:288:GLN:O	2.07	0.55
1:C:731:PHE:O	1:C:734:GLU:HB2	2.07	0.55
1:A:752:ASN:HB2	1:A:754:SER:OG	2.05	0.54
1:B:211:ASN:O	1:B:232:THR:HA	2.07	0.54
1:C:66:THR:O	1:C:90:THR:HG22	2.07	0.54
1:B:345:SER:HB3	1:B:375:ARG:HB2	1.89	0.54
1:D:741:LEU:HD21	1:D:743:LEU:HD11	1.88	0.54
1:B:280:ARG:CD	1:B:280:ARG:C	2.76	0.54
1:C:45:VAL:CG1	1:C:65:VAL:HA	2.38	0.53
1:C:731:PHE:C	1:C:731:PHE:CD2	2.81	0.53
1:C:303:ASN:HB3	1:C:306:TRP:CE2	2.44	0.53
1:A:805:SER:HB2	1:A:806:PRO:HA	1.91	0.53
1:B:52:ARG:HG2	1:B:799:VAL:HG21	1.91	0.53
1:B:275:SER:HA	1:B:298:SER:HB2	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:660:LEU:HD22	1:B:686:GLN:HG3	1.90	0.52
1:D:767:LEU:O	1:D:795:ILE:HD13	2.10	0.52
1:A:467:PHE:HB3	3:A:902:NAG:H81	1.90	0.52
1:C:61:VAL:HG12	1:C:62:GLY:O	2.10	0.52
1:C:219:LYS:HG3	1:C:220:LEU:HB2	1.90	0.52
1:D:296:SER:HA	1:D:320:PHE:O	2.10	0.52
1:A:608:LYS:O	1:A:640:ASN:HB2	2.09	0.52
1:C:218:PRO:O	1:C:219:LYS:O	2.28	0.52
1:C:377:TYR:O	1:C:378:VAL:HB	2.10	0.52
1:B:324:VAL:HG23	8:B:1009:HOH:O	2.10	0.51
1:C:59:GLN:HG3	1:C:84:GLN:NE2	2.24	0.51
1:C:411:PHE:O	8:C:1068:HOH:O	2.19	0.51
1:D:604:ASN:HB3	8:D:1116:HOH:O	2.09	0.51
1:A:341:ILE:HD11	7:A:915:UCG:C5	2.40	0.51
1:B:207:SER:HA	1:B:228:PHE:HB2	1.92	0.51
1:A:750:THR:OG1	1:A:751:ILE:N	2.36	0.51
1:A:411:PHE:HB3	1:A:504:LEU:CD1	2.41	0.51
1:C:118:ASP:OD1	1:C:118:ASP:N	2.44	0.51
1:A:803:CYS:HB2	1:A:809:GLN:O	2.12	0.50
1:D:341:ILE:HD12	7:D:902:UCG:H5	1.92	0.50
1:C:341:ILE:HD13	7:C:901:UCG:H5	1.93	0.50
1:D:549:ALA:O	1:D:550:LEU:HB2	2.12	0.50
1:D:741:LEU:HD23	1:D:767:LEU:CD1	2.41	0.50
1:A:670:HIS:HA	1:A:694:ASP:HB3	1.94	0.50
1:C:343:ASP:HA	1:C:373:HIS:HB2	1.94	0.50
1:C:732:LEU:HD22	1:C:732:LEU:C	2.31	0.50
1:D:571:ALA:HB2	1:D:597:TYR:OH	2.12	0.50
5:B:911:NAG:H81	8:B:1181:HOH:O	2.12	0.50
1:C:159:TYR:CE1	1:C:187:CYS:HB2	2.47	0.50
1:C:473:PRO:HA	7:C:901:UCG:O17	2.12	0.50
1:D:682:THR:HA	1:D:710:PHE:CD1	2.47	0.50
1:A:571:ALA:HB2	1:A:597:TYR:OH	2.12	0.49
1:C:225:ARG:HA	1:C:247:ASN:O	2.11	0.49
1:B:68:LEU:HD21	1:B:70:LEU:HD11	1.95	0.49
1:B:703:LEU:CD2	1:B:724:ILE:HG21	2.43	0.49
1:C:314:LYS:HD3	1:C:340:GLU:OE1	2.12	0.49
1:D:728:PRO:O	1:D:731:PHE:HB2	2.12	0.49
1:D:278:ILE:HB	1:D:306:TRP:CZ2	2.48	0.49
1:C:200:LEU:O	1:C:221:PRO:HG3	2.12	0.49
1:D:688:PRO:O	1:D:713:SER:CB	2.61	0.49
1:B:134:ASP:HA	1:B:155:GLN:O	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:303:ASN:HB3	1:C:306:TRP:CD2	2.48	0.48
1:C:732:LEU:HD13	1:C:733:SER:N	2.27	0.48
1:D:155:GLN:HA	1:D:179:TRP:O	2.12	0.48
1:A:329:SER:OG	1:A:330:GLY:N	2.44	0.48
1:D:677:LYS:O	1:D:700:LEU:HD23	2.14	0.48
1:C:514:ALA:HA	1:C:539:ASN:O	2.14	0.48
1:D:51:ASN:HA	1:D:72:ASP:O	2.13	0.48
1:B:250:LEU:HD23	1:B:250:LEU:C	2.34	0.48
1:C:45:VAL:HG11	1:C:65:VAL:HA	1.95	0.48
1:D:718:LEU:HA	1:D:742:ASP:HB3	1.96	0.48
1:B:45:VAL:HG11	1:B:64:TYR:O	2.14	0.47
1:D:705:ASP:HA	1:D:728:PRO:HB2	1.96	0.47
1:A:146:GLU:HG2	1:A:146:GLU:O	2.13	0.47
1:B:616:SER:HA	1:B:647:SER:O	2.13	0.47
1:C:813:SER:O	1:C:815:VAL:N	2.46	0.47
1:D:512:LEU:HB2	1:D:537:LEU:HD23	1.97	0.47
1:A:303:ASN:ND2	8:A:1278:HOH:O	2.44	0.47
1:C:49:CYS:HA	1:C:52:ARG:NE	2.29	0.47
1:C:152:SER:HA	1:C:176:TYR:HB2	1.96	0.47
1:C:206:LEU:HD21	1:C:208:LEU:HD11	1.96	0.47
1:C:359:ILE:HG23	1:C:363:PHE:CD1	2.49	0.47
1:A:51:ASN:HA	1:A:72:ASP:O	2.14	0.47
1:B:77:HIS:O	1:B:78:ILE:HD13	2.14	0.47
1:C:49:CYS:HB3	1:C:69:ASP:O	2.15	0.47
1:D:786:TRP:O	1:D:790:HIS:ND1	2.48	0.47
5:A:914:NAG:H81	8:A:1167:HOH:O	2.15	0.47
1:C:129:GLU:HA	1:C:150:GLU:O	2.14	0.47
1:C:33:TYR:HE1	1:C:813:SER:OG	1.98	0.47
1:C:175:LEU:HB2	1:C:203:LEU:HD11	1.97	0.47
1:C:155:GLN:HG3	1:C:459:PHE:CZ	2.50	0.47
1:D:597:TYR:HB3	1:D:619:ARG:HB2	1.96	0.47
1:D:606:GLU:HG2	1:D:637:GLY:HA3	1.97	0.46
1:B:683:LEU:HD21	5:B:914:NAG:O5	2.15	0.46
1:C:150:GLU:HG2	1:C:174:ASN:HD22	1.80	0.46
1:A:45:VAL:HG11	1:A:64:TYR:HD2	1.80	0.46
1:B:608:LYS:O	1:B:640:ASN:HB2	2.16	0.46
1:D:77:HIS:O	1:D:78:ILE:HD13	2.15	0.46
1:C:49:CYS:HA	1:C:52:ARG:HE	1.80	0.46
1:B:230:SER:HA	1:B:254:SER:O	2.15	0.46
1:C:720:SER:HA	1:C:744:SER:O	2.15	0.46
1:B:809:GLN:O	1:B:810:ARG:C	2.54	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:750:THR:HG22	1:C:751:ILE:N	2.30	0.46
1:D:586:LEU:O	1:D:610:LEU:HA	2.16	0.46
1:D:592:SER:HA	1:D:616:SER:O	2.16	0.46
1:C:336:LEU:N	1:C:337:PRO:CD	2.79	0.46
1:C:539:ASN:HA	1:C:563:TYR:O	2.16	0.46
1:C:814:ILE:HG13	1:C:815:VAL:N	2.31	0.46
1:C:806:PRO:O	1:C:809:GLN:N	2.43	0.46
1:C:366:LEU:O	1:C:393:LEU:HD22	2.16	0.46
1:D:656:ASN:HD21	1:D:680:ASN:HD22	1.63	0.46
1:D:763:LYS:HA	1:D:763:LYS:HE3	1.97	0.46
1:B:242:PHE:O	1:B:243:LYS:C	2.54	0.45
1:C:113:GLY:HA2	1:C:136:GLN:HB2	1.98	0.45
1:D:767:LEU:HD21	1:D:769:LEU:HD11	1.98	0.45
1:C:775:GLU:HA	1:C:805:SER:O	2.16	0.45
1:A:809:GLN:HE21	1:A:809:GLN:HA	1.81	0.45
1:B:490:LEU:HD23	1:B:514:ALA:HB1	1.99	0.45
1:B:685:GLN:HE21	1:B:710:PHE:HD1	1.65	0.45
1:C:150:GLU:HG2	1:C:174:ASN:HB2	1.99	0.45
1:D:705:ASP:N	1:D:705:ASP:OD1	2.46	0.45
1:C:632:ILE:HD13	1:C:632:ILE:HA	1.72	0.45
1:B:388:GLN:CB	1:B:389:PRO:HD3	2.47	0.45
1:D:670:HIS:HA	1:D:694:ASP:HB3	1.99	0.45
1:D:94:LEU:HB2	1:D:132:LEU:HD23	1.98	0.45
1:C:720:SER:OG	1:C:721:HIS:HD2	2.00	0.45
1:C:732:LEU:C	1:C:734:GLU:N	2.68	0.45
1:D:432:PRO:HG3	1:D:495:PHE:O	2.17	0.45
1:C:327:ILE:HG12	1:C:344:LEU:HD13	1.99	0.45
1:B:784:ARG:HD3	1:B:814:ILE:O	2.17	0.45
1:D:786:TRP:O	1:D:790:HIS:HB2	2.16	0.45
1:A:516:SER:O	1:A:516:SER:OG	2.29	0.44
1:A:727:LEU:HD12	1:A:751:ILE:HG12	1.99	0.44
1:B:696:ARG:HG2	1:B:720:SER:HB3	1.99	0.44
1:C:140:ILE:HG21	1:C:166:ILE:HD11	1.98	0.44
1:D:756:LEU:HD23	1:D:756:LEU:N	2.32	0.44
1:D:789:GLU:O	1:D:790:HIS:ND1	2.49	0.44
1:B:314:LYS:HG3	8:B:1109:HOH:O	2.17	0.44
1:D:205:LEU:CD2	1:D:205:LEU:C	2.86	0.44
1:A:300:ARG:NE	8:A:1172:HOH:O	2.15	0.44
1:C:145:PRO:O	1:C:169:LEU:HD22	2.17	0.44
1:C:808:ASP:O	1:C:812:LYS:NZ	2.49	0.44
1:D:197:PHE:HA	1:D:200:LEU:HD22	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:388:GLN:N	1:D:389:PRO:CD	2.80	0.44
1:A:214:SER:HA	1:A:233:GLN:O	2.18	0.44
1:A:671:ILE:O	1:A:671:ILE:HG22	2.18	0.44
1:D:754:SER:O	1:D:757:GLU:HB2	2.18	0.44
1:B:280:ARG:HD2	1:B:281:PHE:CE2	2.53	0.44
1:B:296:SER:HA	1:B:320:PHE:O	2.17	0.44
1:A:616:SER:HA	1:A:647:SER:O	2.18	0.44
1:C:778:CYS:O	1:C:779:ASP:C	2.56	0.43
1:A:283:PHE:HA	1:A:286:LEU:HD22	1.98	0.43
1:C:731:PHE:C	1:C:734:GLU:HB2	2.39	0.43
1:D:239:GLU:HG3	1:D:284:GLN:NE2	2.30	0.43
1:D:513:SER:HA	1:D:538:THR:O	2.18	0.43
1:A:749:LYS:HA	1:A:773:PRO:O	2.18	0.43
1:B:155:GLN:HA	1:B:179:TRP:O	2.17	0.43
1:B:280:ARG:NH2	1:B:305:ALA:HB1	2.33	0.43
1:C:51:ASN:HA	1:C:72:ASP:O	2.18	0.43
1:C:731:PHE:HA	1:C:734:GLU:OE2	2.19	0.43
1:D:745:SER:C	8:D:1096:HOH:O	2.56	0.43
1:D:467:PHE:CB	3:D:903:NAG:H81	2.49	0.43
1:A:718:LEU:HA	1:A:742:ASP:HB3	2.00	0.43
1:C:290:ARG:HB2	1:C:291:TYR:CD1	2.53	0.43
1:C:496:ILE:HG23	1:C:500:GLN:HB3	2.01	0.43
1:C:612:GLU:OE2	1:C:643:ARG:NH1	2.52	0.43
1:B:303:ASN:HB3	1:B:306:TRP:CE2	2.54	0.43
1:C:121:PHE:CE2	1:C:130:LEU:HD21	2.53	0.43
1:C:156:ASN:O	1:C:180:ASN:OD1	2.37	0.43
1:D:611:VAL:O	1:D:641:LEU:HD12	2.19	0.43
1:B:364:SER:HA	1:B:393:LEU:HD21	2.00	0.43
1:C:161:ILE:HG22	1:C:196:VAL:HG11	2.00	0.43
1:C:222:SER:HA	1:C:245:LEU:HD23	2.00	0.43
1:C:411:PHE:HB3	1:C:504:LEU:HD13	2.00	0.43
1:B:399:ILE:O	1:B:399:ILE:HG23	2.18	0.43
1:B:400:ASN:C	1:B:400:ASN:OD1	2.57	0.43
1:C:740:HIS:HD2	1:C:741:LEU:N	2.15	0.43
1:C:75:ILE:HG21	1:C:78:ILE:HD11	2.01	0.43
1:D:314:LYS:HG2	1:D:315:VAL:HG23	2.01	0.43
1:A:213:LEU:O	1:A:214:SER:HB2	2.19	0.43
1:A:388:GLN:N	1:A:389:PRO:CD	2.82	0.43
1:C:194:ASP:OD2	1:C:194:ASP:C	2.57	0.43
1:C:233:GLN:HA	8:C:1094:HOH:O	2.18	0.43
1:C:539:ASN:HB3	8:C:1141:HOH:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:612:GLU:HA	1:C:643:ARG:O	2.19	0.43
1:D:370:ARG:NH2	5:D:909:NAG:H81	2.34	0.43
1:A:536:ASP:OD1	1:A:538:THR:HG23	2.18	0.42
1:A:467:PHE:CB	3:A:902:NAG:H81	2.49	0.42
1:B:214:SER:HA	1:B:233:GLN:O	2.19	0.42
1:C:289:LEU:HD23	1:C:310:MET:SD	2.58	0.42
1:C:545:ASP:O	1:C:545:ASP:CG	2.56	0.42
1:C:731:PHE:O	1:C:734:GLU:CB	2.67	0.42
1:A:150:GLU:HG2	1:A:174:ASN:HB2	2.00	0.42
1:C:534:TYR:CE2	1:C:558:VAL:HG11	2.54	0.42
1:C:784:ARG:HE	1:C:784:ARG:HA	1.84	0.42
1:D:565:SER:HB3	1:D:569:ARG:NH1	2.35	0.42
1:D:616:SER:HA	1:D:647:SER:O	2.19	0.42
1:A:388:GLN:HB2	1:A:389:PRO:HD3	2.02	0.42
1:A:420:LEU:HA	1:A:420:LEU:HD23	1.85	0.42
1:B:336:LEU:HD23	8:B:1046:HOH:O	2.18	0.42
1:C:275:SER:HA	1:C:298:SER:HB2	2.02	0.42
1:A:275:SER:HA	1:A:298:SER:HB2	2.01	0.42
1:C:149:THR:HA	1:C:171:ASN:O	2.20	0.42
1:C:357:ILE:HG13	1:C:377:TYR:CZ	2.55	0.42
1:D:251:LEU:HD21	1:D:253:LEU:HD11	2.00	0.42
1:D:38:LYS:O	1:D:45:VAL:HA	2.19	0.42
1:A:565:SER:O	1:A:566:HIS:C	2.58	0.42
1:C:225:ARG:O	1:C:248:LEU:HD22	2.19	0.42
1:D:280:ARG:HG3	1:D:281:PHE:CD1	2.54	0.42
1:A:545:ASP:O	1:A:545:ASP:CG	2.57	0.42
1:A:764:LEU:O	1:A:793:VAL:HG22	2.19	0.42
1:B:545:ASP:O	1:B:545:ASP:CG	2.58	0.42
1:B:784:ARG:HH12	1:B:817:LEU:HB2	1.84	0.42
1:D:213:LEU:O	1:D:214:SER:HB2	2.19	0.42
1:A:296:SER:HA	1:A:320:PHE:O	2.20	0.42
1:A:431:SER:HB2	1:A:432:PRO:CD	2.49	0.42
1:A:783:PHE:O	1:A:786:TRP:HB3	2.20	0.42
1:C:211:ASN:O	1:C:232:THR:HA	2.19	0.42
1:D:611:VAL:HG13	1:D:642:THR:OG1	2.20	0.42
1:C:135:ASN:O	1:C:136:GLN:HB2	2.20	0.41
1:D:716:THR:HG23	1:D:740:HIS:HB3	2.01	0.41
1:D:71:SER:O	1:D:72:ASP:HB2	2.19	0.41
1:D:491:ASN:O	1:D:515:ASN:HA	2.20	0.41
1:D:598:THR:O	1:D:598:THR:HG23	2.19	0.41
1:C:283:PHE:HA	1:C:286:LEU:HD22	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:318:LEU:HB2	1:C:344:LEU:HD23	2.01	0.41
1:D:234:ILE:O	1:D:256:ASN:HB3	2.20	0.41
1:D:575:HIS:CE1	1:D:598:THR:HG23	2.55	0.41
1:D:669:LEU:HD21	1:D:671:ILE:HD11	2.02	0.41
1:A:375:ARG:HA	1:A:402:GLY:O	2.21	0.41
1:A:685:GLN:HG3	1:A:710:PHE:HA	2.03	0.41
1:A:814:ILE:HG13	1:A:815:VAL:N	2.35	0.41
1:B:117:THR:OG1	1:B:120:ALA:HB2	2.20	0.41
1:C:219:LYS:CG	1:C:220:LEU:HB2	2.50	0.41
1:C:397:SER:HA	1:C:419:ASN:O	2.21	0.41
1:D:743:LEU:HB2	1:D:769:LEU:HD23	2.03	0.41
1:A:569:ARG:NH1	8:A:1096:HOH:O	2.37	0.41
1:B:671:ILE:HG22	1:B:671:ILE:O	2.21	0.41
1:B:703:LEU:HD11	1:B:719:LEU:HD13	2.01	0.41
1:B:45:VAL:HG13	1:B:45:VAL:O	2.21	0.41
1:B:592:SER:HA	1:B:616:SER:O	2.21	0.41
1:B:606:GLU:HG2	1:B:637:GLY:HA3	2.03	0.41
1:D:340:GLU:C	1:D:341:ILE:HG12	2.40	0.41
1:B:45:VAL:HG21	1:B:64:TYR:HD2	1.86	0.41
1:D:780:ILE:O	1:D:782:ASP:N	2.53	0.41
1:B:280:ARG:NH2	1:B:305:ALA:CB	2.84	0.41
1:B:329:SER:OG	1:B:330:GLY:N	2.54	0.41
1:B:517:ASN:OD1	1:B:517:ASN:C	2.59	0.41
1:B:529:ILE:HG13	1:B:529:ILE:O	2.21	0.41
1:C:302:ILE:HG21	1:C:332:PHE:CD2	2.56	0.41
1:A:327:ILE:HD11	1:A:344:LEU:HD22	2.02	0.41
1:C:413:LEU:C	1:C:413:LEU:HD12	2.41	0.41
1:C:719:LEU:HB2	1:C:742:ASP:O	2.21	0.41
1:D:644:LEU:HD21	1:D:646:LEU:HD21	2.02	0.41
1:A:573:VAL:O	1:A:575:HIS:CE1	2.74	0.41
1:A:756:LEU:O	1:A:757:GLU:C	2.60	0.41
1:C:173:LYS:O	1:C:204:GLU:N	2.40	0.41
1:C:300:ARG:CD	8:C:1085:HOH:O	2.63	0.41
1:B:33:TYR:CD2	1:B:33:TYR:C	2.95	0.40
1:C:425:LEU:HB2	1:C:488:LEU:HD23	2.03	0.40
1:C:720:SER:HB3	1:C:742:ASP:OD2	2.20	0.40
1:C:96:HIS:HB3	1:C:99:ASN:OD1	2.21	0.40
1:A:617:GLY:HA2	1:A:648:LEU:O	2.21	0.40
1:B:228:PHE:HA	1:B:252:ASP:HB3	2.03	0.40
1:B:565:SER:O	1:B:566:HIS:C	2.59	0.40
1:D:738:LEU:O	1:D:764:LEU:HD12	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:740:HIS:C	1:C:740:HIS:CD2	2.94	0.40
1:A:341:ILE:HD11	7:A:915:UCG:C6	2.51	0.40
1:B:473:PRO:HA	7:B:915:UCG:O17	2.21	0.40
1:C:147:SER:HA	1:C:171:ASN:HD22	1.86	0.40
1:C:313:LEU:HD23	1:C:336:LEU:CD2	2.52	0.40
1:C:732:LEU:O	1:C:734:GLU:N	2.48	0.40
1:D:671:ILE:HB	1:D:695:LEU:HD23	2.03	0.40
1:D:688:PRO:O	1:D:713:SER:OG	2.37	0.40
1:B:805:SER:HB2	1:B:806:PRO:HA	2.02	0.40
1:C:813:SER:O	1:C:814:ILE:C	2.60	0.40
1:D:476:LYS:HA	1:D:477:PRO:HD2	1.84	0.40
1:D:494:PHE:O	1:D:517:ASN:HA	2.22	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	733/811 (90%)	672 (92%)	59 (8%)	2 (0%)	44	60
1	B	727/811 (90%)	669 (92%)	53 (7%)	5 (1%)	25	37
1	C	721/811 (89%)	643 (89%)	72 (10%)	6 (1%)	22	33
1	D	735/811 (91%)	671 (91%)	61 (8%)	3 (0%)	38	54
All	All	2916/3244 (90%)	2655 (91%)	245 (8%)	16 (0%)	32	46

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	243	LYS
1	C	220	LEU
1	C	219	LYS

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Mol	Chain	Res	Type
1	C	733	SER
1	D	243	LYS
1	B	754	SER
1	A	284	GLN
1	B	505	PRO
1	C	378	VAL
1	A	378	VAL
1	B	378	VAL
1	B	773	PRO
1	D	378	VAL
1	C	735	VAL
1	C	814	ILE
1	D	806	PRO

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	690/755 (91%)	646 (94%)	44 (6%)	20	32
1	B	684/755 (91%)	656 (96%)	28 (4%)	35	54
1	C	679/755 (90%)	627 (92%)	52 (8%)	15	23
1	D	690/755 (91%)	652 (94%)	38 (6%)	25	40
All	All	2743/3020 (91%)	2581 (94%)	162 (6%)	23	36

All (162) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	49	CYS
1	A	50	SER
1	A	52	ARG
1	A	60	THR
1	A	87	GLN
1	A	118	ASP
1	A	142	SER
1	A	150	GLU

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Mol	Chain	Res	Type
1	A	160	ASN
1	A	173	LYS
1	A	200	LEU
1	A	243	LYS
1	A	248	LEU
1	A	275	SER
1	A	286	LEU
1	A	301	LYS
1	A	314	LYS
1	A	329	SER
1	A	397	SER
1	A	416	ASN
1	A	465	SER
1	A	472	ARG
1	A	476	LYS
1	A	504	LEU
1	A	509	CYS
1	A	516	SER
1	A	554	SER
1	A	595	ASN
1	A	625	ASN
1	A	677	LYS
1	A	682	THR
1	A	702	PHE
1	A	703	LEU
1	A	725	SER
1	A	733	SER
1	A	736	SER
1	A	750	THR
1	A	753	LYS
1	A	756	LEU
1	A	780	ILE
1	A	791	LEU
1	A	809	GLN
1	A	815	VAL
1	A	817	LEU
1	B	49	CYS
1	B	52	ARG
1	B	56	GLU
1	B	66	THR
1	B	76	THR
1	B	87	GLN

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Mol	Chain	Res	Type
1	B	122	LEU
1	B	160	ASN
1	B	168	ARG
1	B	248	LEU
1	B	275	SER
1	B	280	ARG
1	B	286	LEU
1	B	287	THR
1	B	317	ASP
1	B	408	GLN
1	B	416	ASN
1	B	418	SER
1	B	504	LEU
1	B	509	CYS
1	B	727	LEU
1	B	733	SER
1	B	750	THR
1	B	754	SER
1	B	765	SER
1	B	780	ILE
1	B	794	LYS
1	B	817	LEU
1	C	33	TYR
1	C	76	THR
1	C	81	GLU
1	C	86	LEU
1	C	100	VAL
1	C	116	ILE
1	C	118	ASP
1	C	147	SER
1	C	150	GLU
1	C	164	GLU
1	C	167	SER
1	C	173	LYS
1	C	188	GLU
1	C	191	ASN
1	C	200	LEU
1	C	212	SER
1	C	214	SER
1	C	219	LYS
1	C	245	LEU
1	C	246	ILE

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Mol	Chain	Res	Type
1	C	248	LEU
1	C	259	ARG
1	C	275	SER
1	C	278	ILE
1	C	279	ASP
1	C	280	ARG
1	C	285	ASN
1	C	314	LYS
1	C	317	ASP
1	C	365	LYS
1	C	387	PHE
1	C	397	SER
1	C	412	LYS
1	C	416	ASN
1	C	418	SER
1	C	504	LEU
1	C	569	ARG
1	C	632	ILE
1	C	681	TRP
1	C	682	THR
1	C	706	SER
1	C	712	SER
1	C	715	ARG
1	C	720	SER
1	C	725	SER
1	C	729	SER
1	C	731	PHE
1	C	732	LEU
1	C	735	VAL
1	C	780	ILE
1	C	784	ARG
1	C	810	ARG
1	D	36	ASP
1	D	49	CYS
1	D	52	ARG
1	D	64	TYR
1	D	88	ASN
1	D	117	THR
1	D	122	LEU
1	D	150	GLU
1	D	170	ILE
1	D	200	LEU

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Mol	Chain	Res	Type
1	D	235	LYS
1	D	243	LYS
1	D	248	LEU
1	D	300	ARG
1	D	308	LYS
1	D	418	SER
1	D	419	ASN
1	D	472	ARG
1	D	476	LYS
1	D	509	CYS
1	D	582	ASN
1	D	609	SER
1	D	653	HIS
1	D	665	SER
1	D	686	GLN
1	D	702	PHE
1	D	705	ASP
1	D	712	SER
1	D	715	ARG
1	D	733	SER
1	D	735	VAL
1	D	736	SER
1	D	753	LYS
1	D	756	LEU
1	D	764	LEU
1	D	788	ASP
1	D	794	LYS
1	D	805	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	55	GLN
1	A	87	GLN
1	A	160	ASN
1	A	191	ASN
1	A	288	GLN
1	A	721	HIS
1	A	809	GLN
1	B	288	GLN
1	B	312	HIS
1	B	503	ASN

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Mol	Chain	Res	Type
1	C	84	GLN
1	C	174	ASN
1	C	202	ASN
1	C	284	GLN
1	C	285	ASN
1	C	721	HIS
1	D	123	ASN
1	D	247	ASN
1	D	284	GLN
1	D	416	ASN
1	D	581	GLN
1	D	629	ASN
1	D	656	ASN
1	D	661	ASN
1	D	686	GLN

5.3.3 RNA ⓘ

There are no RNA molecules 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 ⓘ

32 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	$\# Z > 2$
3	NAG	A	902	1,3	14,14,15	0.97	1 (7%)	15,19,21	2.03	3 (20%)
3	NAG	A	903	3	14,14,15	0.69	0	15,19,21	1.38	2 (13%)
3	BMA	A	904	3	11,11,12	1.06	1 (9%)	13,15,17	1.96	7 (53%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	BMA	A	905	4	11,11,12	0.55	0	13,15,17	1.43	2 (15%)
4	NAG	A	906	1,4	14,14,15	0.74	0	15,19,21	1.51	2 (13%)
4	NAG	A	907	4	14,14,15	0.69	0	15,19,21	1.02	1 (6%)
6	NAG	A	909	1,6	14,14,15	0.64	0	15,19,21	1.11	1 (6%)
6	NAG	A	910	6	14,14,15	0.73	1 (7%)	15,19,21	1.55	5 (33%)
3	NAG	B	902	1,3	14,14,15	0.77	0	15,19,21	1.43	4 (26%)
3	NAG	B	903	3	14,14,15	0.61	0	15,19,21	1.35	2 (13%)
3	BMA	B	904	3	11,11,12	0.25	0	13,15,17	0.54	0
3	NAG	B	905	1,3	14,14,15	0.74	0	15,19,21	1.56	5 (33%)
3	NAG	B	906	3	14,14,15	0.97	0	15,19,21	1.45	3 (20%)
3	BMA	B	907	3	11,11,12	0.44	0	13,15,17	1.12	1 (7%)
6	NAG	B	909	1,6	14,14,15	0.95	1 (7%)	15,19,21	1.62	2 (13%)
6	NAG	B	910	6	14,14,15	0.66	0	15,19,21	1.68	4 (26%)
3	NAG	C	903	1,3	14,14,15	0.57	0	15,19,21	1.54	4 (26%)
3	NAG	C	904	3	14,14,15	0.67	0	15,19,21	1.34	1 (6%)
3	BMA	C	905	3	11,11,12	0.70	0	13,15,17	1.42	2 (15%)
3	NAG	C	906	1,3	14,14,15	0.49	0	15,19,21	1.56	3 (20%)
3	NAG	C	907	3	14,14,15	0.77	0	15,19,21	0.92	1 (6%)
3	BMA	C	908	3	11,11,12	0.53	0	13,15,17	0.96	1 (7%)
6	NAG	C	910	1,6	14,14,15	0.84	0	15,19,21	1.75	4 (26%)
6	NAG	C	911	6	14,14,15	0.56	0	15,19,21	1.51	3 (20%)
3	NAG	D	903	1,3	14,14,15	0.86	1 (7%)	15,19,21	2.10	4 (26%)
3	NAG	D	904	3	14,14,15	0.77	0	15,19,21	1.68	4 (26%)
3	BMA	D	905	3	11,11,12	0.96	1 (9%)	13,15,17	2.62	5 (38%)
3	NAG	D	906	1,3	14,14,15	0.71	0	15,19,21	1.06	1 (6%)
3	NAG	D	907	3	14,14,15	0.58	0	15,19,21	0.97	0
3	BMA	D	908	3	11,11,12	0.44	0	13,15,17	1.18	1 (7%)
6	NAG	D	910	1,6	14,14,15	0.69	0	15,19,21	1.37	2 (13%)
6	NAG	D	911	6	14,14,15	0.65	0	15,19,21	1.15	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	NAG	A	902	1,3	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	A	903	3	-	0/6/23/26	0/1/1/1
3	BMA	A	904	3	-	0/2/19/22	0/1/1/1
4	BMA	A	905	4	-	0/2/19/22	0/1/1/1
4	NAG	A	906	1,4	-	0/6/23/26	0/1/1/1
4	NAG	A	907	4	-	0/6/23/26	0/1/1/1
6	NAG	A	909	1,6	-	0/6/23/26	0/1/1/1
6	NAG	A	910	6	-	0/6/23/26	0/1/1/1
3	NAG	B	902	1,3	-	0/6/23/26	0/1/1/1
3	NAG	B	903	3	-	0/6/23/26	0/1/1/1
3	BMA	B	904	3	-	0/2/19/22	0/1/1/1
3	NAG	B	905	1,3	-	0/6/23/26	0/1/1/1
3	NAG	B	906	3	-	0/6/23/26	0/1/1/1
3	BMA	B	907	3	-	0/2/19/22	0/1/1/1
6	NAG	B	909	1,6	-	0/6/23/26	0/1/1/1
6	NAG	B	910	6	-	0/6/23/26	0/1/1/1
3	NAG	C	903	1,3	-	0/6/23/26	0/1/1/1
3	NAG	C	904	3	-	0/6/23/26	0/1/1/1
3	BMA	C	905	3	-	0/2/19/22	0/1/1/1
3	NAG	C	906	1,3	-	0/6/23/26	0/1/1/1
3	NAG	C	907	3	-	0/6/23/26	0/1/1/1
3	BMA	C	908	3	-	0/2/19/22	0/1/1/1
6	NAG	C	910	1,6	-	0/6/23/26	0/1/1/1
6	NAG	C	911	6	-	0/6/23/26	0/1/1/1
3	NAG	D	903	1,3	-	0/6/23/26	0/1/1/1
3	NAG	D	904	3	-	0/6/23/26	0/1/1/1
3	BMA	D	905	3	-	0/2/19/22	0/1/1/1
3	NAG	D	906	1,3	-	0/6/23/26	0/1/1/1
3	NAG	D	907	3	-	0/6/23/26	0/1/1/1
3	BMA	D	908	3	-	0/2/19/22	0/1/1/1
6	NAG	D	910	1,6	-	0/6/23/26	0/1/1/1
6	NAG	D	911	6	-	0/6/23/26	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	902	NAG	O5-C1	-2.89	1.39	1.43
3	A	904	BMA	O5-C1	-2.75	1.39	1.43
3	D	903	NAG	O5-C1	-2.63	1.39	1.43
6	B	909	NAG	O5-C1	-2.42	1.39	1.43
3	D	905	BMA	O5-C1	-2.37	1.39	1.43
6	A	910	NAG	O5-C1	-2.30	1.40	1.43

All (80) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	905	BMA	O5-C1-C2	-5.59	102.04	110.79
3	D	903	NAG	O5-C1-C2	-5.25	104.17	111.47
3	D	905	BMA	O2-C2-C1	-4.24	100.55	109.18
6	C	910	NAG	O6-C6-C5	-4.17	97.30	111.34
6	B	910	NAG	O5-C1-C2	-3.58	106.50	111.47
6	B	910	NAG	C3-C4-C5	-3.53	103.99	110.22
3	D	904	NAG	O5-C1-C2	-3.39	106.76	111.47
3	A	904	BMA	O4-C4-C3	-3.35	103.08	110.36
6	A	910	NAG	O5-C1-C2	-3.25	106.95	111.47
6	D	910	NAG	C1-C2-N2	-3.21	105.01	110.49
3	B	906	NAG	O5-C1-C2	-3.21	107.01	111.47
4	A	905	BMA	C1-O5-C5	-3.11	107.89	112.17
3	D	903	NAG	C6-C5-C4	-3.09	105.78	113.00
3	B	905	NAG	O5-C1-C2	-3.01	107.28	111.47
3	D	904	NAG	O6-C6-C5	-3.01	101.20	111.34
6	C	910	NAG	C1-C2-N2	-2.96	105.43	110.49
3	D	906	NAG	O5-C1-C2	-2.93	107.40	111.47
6	D	910	NAG	O5-C1-C2	-2.92	107.41	111.47
3	A	904	BMA	O5-C1-C2	-2.91	106.23	110.79
3	C	904	NAG	O5-C1-C2	-2.87	107.48	111.47
3	C	903	NAG	O5-C1-C2	-2.77	107.62	111.47
3	B	905	NAG	O4-C4-C3	-2.73	104.41	110.36
6	C	911	NAG	C1-C2-N2	-2.73	105.83	110.49
3	B	902	NAG	O5-C1-C2	-2.70	107.71	111.47
3	A	903	NAG	O7-C7-C8	-2.68	117.17	122.06
3	B	902	NAG	O6-C6-C5	-2.65	102.44	111.34
4	A	906	NAG	O5-C1-C2	-2.62	107.83	111.47
6	B	909	NAG	C1-C2-N2	-2.59	106.06	110.49
3	A	903	NAG	O5-C1-C2	-2.49	108.01	111.47
3	D	904	NAG	C3-C4-C5	-2.47	105.87	110.22
3	C	903	NAG	O6-C6-C5	-2.44	103.13	111.34
3	B	907	BMA	O2-C2-C1	-2.42	104.25	109.18
4	A	906	NAG	O4-C4-C3	-2.42	105.09	110.36
3	C	906	NAG	O5-C1-C2	-2.41	108.11	111.47
3	C	908	BMA	C2-C3-C4	-2.36	106.77	110.88
3	B	902	NAG	C6-C5-C4	-2.35	107.51	113.00
3	A	902	NAG	C6-C5-C4	-2.34	107.52	113.00
3	B	903	NAG	O5-C1-C2	-2.27	108.31	111.47
3	D	904	NAG	C6-C5-C4	-2.26	107.72	113.00
6	A	910	NAG	O7-C7-C8	-2.26	117.95	122.06
3	C	907	NAG	O5-C1-C2	-2.23	108.37	111.47
6	C	910	NAG	C6-C5-C4	-2.16	107.96	113.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	909	NAG	C1-O5-C5	-2.15	109.20	112.17
3	A	902	NAG	O6-C6-C5	-2.14	104.15	111.34
3	C	903	NAG	C6-C5-C4	-2.10	108.08	113.00
3	D	903	NAG	O6-C6-C5	-2.08	104.34	111.34
6	B	910	NAG	O7-C7-C8	-2.08	118.27	122.06
3	C	906	NAG	O7-C7-C8	-2.07	118.29	122.06
4	A	907	NAG	O4-C4-C3	-2.04	105.92	110.36
3	B	906	NAG	O3-C3-C4	-2.03	105.93	110.36
6	A	910	NAG	O3-C3-C4	-2.01	105.98	110.36
6	B	910	NAG	O4-C4-C5	2.01	114.35	109.28
3	D	908	BMA	C1-C2-C3	2.02	112.21	109.65
3	A	904	BMA	O3-C3-C4	2.05	114.82	110.36
3	B	905	NAG	C2-N2-C7	2.05	125.93	122.94
6	A	910	NAG	O4-C4-C5	2.10	114.56	109.28
3	B	905	NAG	O7-C7-N2	2.11	125.98	121.92
4	A	905	BMA	O3-C3-C4	2.17	115.07	110.36
6	C	911	NAG	O4-C4-C5	2.20	114.82	109.28
3	B	905	NAG	C1-O5-C5	2.22	115.23	112.17
3	A	904	BMA	O6-C6-C5	2.23	118.83	111.34
3	C	905	BMA	O3-C3-C4	2.23	115.21	110.36
3	A	904	BMA	C6-C5-C4	2.23	118.22	113.00
3	B	903	NAG	C1-C2-N2	2.23	114.30	110.49
3	D	905	BMA	O3-C3-C4	2.25	115.24	110.36
3	A	904	BMA	O3-C3-C2	2.30	114.20	110.02
3	D	905	BMA	O6-C6-C5	2.30	119.07	111.34
3	C	903	NAG	C1-O5-C5	2.31	115.36	112.17
3	B	906	NAG	O7-C7-N2	2.34	126.43	121.92
6	A	910	NAG	C1-O5-C5	2.35	115.40	112.17
3	A	904	BMA	C1-O5-C5	2.48	115.58	112.17
6	C	911	NAG	C1-O5-C5	2.52	115.63	112.17
3	B	902	NAG	C1-O5-C5	2.58	115.73	112.17
3	C	905	BMA	C3-C4-C5	2.71	115.00	110.22
6	C	910	NAG	C1-O5-C5	2.74	115.95	112.17
3	C	906	NAG	C1-O5-C5	3.13	116.48	112.17
3	D	903	NAG	C1-O5-C5	3.77	117.36	112.17
3	D	905	BMA	C1-O5-C5	4.45	118.30	112.17
6	B	909	NAG	C1-O5-C5	4.75	118.72	112.17
3	A	902	NAG	C1-O5-C5	6.67	121.36	112.17

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	902	NAG	2	0
3	D	903	NAG	2	0

5.6 Ligand geometry

26 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	URI	A	901	-	14,18,18	0.73	0	15,26,26	2.28	1 (6%)
5	NAG	A	908	1	14,14,15	0.28	0	15,19,21	0.57	0
5	NAG	A	911	1	14,14,15	0.28	0	15,19,21	0.57	0
5	NAG	A	912	1	14,14,15	0.28	0	15,19,21	0.56	0
5	NAG	A	913	1	14,14,15	0.58	0	15,19,21	1.29	2 (13%)
5	NAG	A	914	1	14,14,15	0.70	0	15,19,21	1.38	3 (20%)
7	UCG	A	915	-	41,52,52	1.15	4 (9%)	50,82,82	2.79	19 (38%)
2	URI	B	901	-	14,18,18	0.83	0	15,26,26	2.54	1 (6%)
5	NAG	B	908	1	14,14,15	0.63	0	15,19,21	0.42	0
5	NAG	B	911	1	14,14,15	0.29	0	15,19,21	0.57	0
5	NAG	B	912	1	14,14,15	0.53	0	15,19,21	1.50	3 (20%)
5	NAG	B	913	1	14,14,15	0.57	0	15,19,21	1.31	1 (6%)
5	NAG	B	914	1	14,14,15	0.29	0	15,19,21	0.57	0
7	UCG	B	915	-	41,52,52	1.08	2 (4%)	50,82,82	2.83	16 (32%)
7	UCG	C	901	-	41,52,52	1.02	2 (4%)	50,82,82	2.88	11 (22%)
2	URI	C	902	-	14,18,18	0.93	1 (7%)	15,26,26	2.48	2 (13%)
5	NAG	C	909	1	14,14,15	0.53	0	15,19,21	1.25	2 (13%)
5	NAG	C	912	1	14,14,15	0.82	1 (7%)	15,19,21	1.25	1 (6%)
5	NAG	C	913	1	14,14,15	0.56	0	15,19,21	1.50	3 (20%)
5	NAG	C	914	1	14,14,15	0.50	0	15,19,21	1.70	3 (20%)
2	URI	D	901	-	14,18,18	0.96	0	15,26,26	2.16	3 (20%)
7	UCG	D	902	-	41,52,52	1.09	2 (4%)	50,82,82	2.78	13 (26%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	NAG	D	909	1	14,14,15	0.79	0	15,19,21	2.61	5 (33%)
5	NAG	D	912	1	14,14,15	0.63	0	15,19,21	1.44	1 (6%)
5	NAG	D	913	1	14,14,15	0.28	0	15,19,21	0.57	0
5	NAG	D	914	1	14,14,15	0.53	0	15,19,21	2.29	3 (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
2	URI	A	901	-	-	0/2/22/22	0/2/2/2
5	NAG	A	908	1	-	0/6/23/26	0/1/1/1
5	NAG	A	911	1	-	0/6/23/26	0/1/1/1
5	NAG	A	912	1	-	0/6/23/26	0/1/1/1
5	NAG	A	913	1	-	0/6/23/26	0/1/1/1
5	NAG	A	914	1	-	0/6/23/26	0/1/1/1
7	UCG	A	915	-	-	0/17/67/67	0/6/6/6
2	URI	B	901	-	-	0/2/22/22	0/2/2/2
5	NAG	B	908	1	-	0/6/23/26	0/1/1/1
5	NAG	B	911	1	-	0/6/23/26	0/1/1/1
5	NAG	B	912	1	-	0/6/23/26	0/1/1/1
5	NAG	B	913	1	-	0/6/23/26	0/1/1/1
5	NAG	B	914	1	-	0/6/23/26	0/1/1/1
7	UCG	B	915	-	-	0/17/67/67	0/6/6/6
7	UCG	C	901	-	-	0/17/67/67	0/6/6/6
2	URI	C	902	-	-	0/2/22/22	0/2/2/2
5	NAG	C	909	1	-	0/6/23/26	0/1/1/1
5	NAG	C	912	1	-	0/6/23/26	0/1/1/1
5	NAG	C	913	1	-	0/6/23/26	0/1/1/1
5	NAG	C	914	1	-	0/6/23/26	0/1/1/1
2	URI	D	901	-	-	0/2/22/22	0/2/2/2
7	UCG	D	902	-	-	0/17/67/67	0/6/6/6
5	NAG	D	909	1	-	0/6/23/26	0/1/1/1
5	NAG	D	912	1	-	0/6/23/26	0/1/1/1
5	NAG	D	913	1	-	0/6/23/26	0/1/1/1
5	NAG	D	914	1	-	0/6/23/26	0/1/1/1

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	902	URI	C2-N3	-2.16	1.33	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	C	912	NAG	O5-C1	-2.02	1.40	1.43
7	A	915	UCG	C16-C15	2.08	1.57	1.53
7	A	915	UCG	O-C	2.11	1.44	1.41
7	B	915	UCG	P2-O17	2.13	1.65	1.55
7	A	915	UCG	C7-N1	2.15	1.36	1.33
7	C	901	UCG	P2-O17	2.25	1.66	1.55
7	D	902	UCG	P2-O17	2.42	1.67	1.55
7	D	902	UCG	P2-O16	3.95	1.65	1.50
7	C	901	UCG	P2-O16	4.00	1.65	1.50
7	B	915	UCG	P2-O16	4.12	1.66	1.50
7	A	915	UCG	P2-O16	4.23	1.66	1.50

All (93) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	D	902	UCG	O12-P2-O16	-7.90	94.91	115.76
7	C	901	UCG	O12-P2-O16	-7.66	95.53	115.76
7	B	915	UCG	O12-P2-O16	-6.00	99.91	115.76
7	A	915	UCG	O12-P2-O16	-5.73	100.62	115.76
7	B	915	UCG	N2-C12-N3	-5.32	119.69	127.46
7	C	901	UCG	O13-P2-O16	-5.29	101.78	115.76
7	A	915	UCG	C17-O10-C14	-4.86	104.60	109.77
7	D	902	UCG	N2-C12-N3	-4.40	121.03	127.46
5	D	912	NAG	O5-C1-C2	-4.15	105.70	111.47
7	B	915	UCG	O13-P2-O16	-3.92	105.41	115.76
7	C	901	UCG	N2-C12-N3	-3.86	121.83	127.46
7	A	915	UCG	N2-C12-N3	-3.69	122.07	127.46
7	A	915	UCG	O17-P2-O16	-3.44	98.78	109.89
5	D	914	NAG	C4-C3-C2	-3.42	106.00	111.02
7	D	902	UCG	O13-P2-O16	-3.35	106.93	115.76
7	C	901	UCG	C9-C11-N3	-3.24	118.87	123.48
5	C	914	NAG	O5-C1-C2	-3.16	107.08	111.47
7	D	902	UCG	O17-P2-O16	-3.02	100.15	109.89
7	A	915	UCG	C9-C11-N3	-2.94	119.30	123.48
7	B	915	UCG	C9-C11-N3	-2.84	119.44	123.48
7	A	915	UCG	C15-C16-C17	-2.82	96.75	103.68
7	A	915	UCG	O10-C14-C15	-2.78	101.72	106.59
5	C	913	NAG	O5-C1-C2	-2.78	107.61	111.47
5	D	909	NAG	O3-C3-C4	-2.63	104.64	110.36
7	D	902	UCG	C9-C11-N3	-2.58	119.81	123.48
5	C	913	NAG	C2-N2-C7	-2.56	119.21	122.94
7	B	915	UCG	O17-P2-O16	-2.48	101.88	109.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	901	URI	O3'-C3'-C2'	-2.48	103.88	111.83
5	D	914	NAG	C3-C4-C5	-2.40	105.99	110.22
7	B	915	UCG	C10-C9-N5	-2.40	107.09	109.41
7	B	915	UCG	C17-O10-C14	-2.37	107.25	109.77
5	B	912	NAG	O5-C1-C2	-2.37	108.18	111.47
7	B	915	UCG	C15-C16-C17	-2.30	98.05	103.68
5	C	909	NAG	C2-N2-C7	-2.29	119.61	122.94
5	C	909	NAG	O5-C1-C2	-2.23	108.37	111.47
2	D	901	URI	O2'-C2'-C3'	-2.20	104.80	111.83
7	A	915	UCG	O13-P2-O16	-2.13	110.14	115.76
7	D	902	UCG	O10-C17-C18	-2.13	102.22	109.40
2	C	902	URI	O2'-C2'-C3'	-2.13	105.02	111.83
5	D	909	NAG	C6-C5-C4	-2.11	108.06	113.00
7	A	915	UCG	C6-N-C8	-2.07	117.92	121.28
5	C	914	NAG	C6-C5-C4	-2.03	108.25	113.00
5	A	914	NAG	C6-C5-C4	-2.02	108.28	113.00
7	A	915	UCG	O7-P-O6	2.01	115.73	107.61
5	A	913	NAG	C3-C4-C5	2.04	113.81	110.22
5	C	913	NAG	C4-C3-C2	2.06	114.04	111.02
7	A	915	UCG	N6-C12-N3	2.12	120.63	117.24
7	C	901	UCG	P-O1-C4	2.13	124.16	118.30
7	A	915	UCG	C2-C1-C	2.17	104.82	99.95
7	D	902	UCG	C2-C1-C	2.25	105.01	99.95
7	A	915	UCG	P-O1-C4	2.33	124.71	118.30
5	A	914	NAG	C1-O5-C5	2.38	115.45	112.17
5	A	913	NAG	C1-O5-C5	2.39	115.46	112.17
7	B	915	UCG	O-C-N	2.40	112.89	108.08
5	B	912	NAG	C1-C2-N2	2.42	114.62	110.49
5	B	912	NAG	O7-C7-N2	2.44	126.61	121.92
7	D	902	UCG	C11-N3-C12	2.47	119.61	116.06
5	A	914	NAG	C3-C4-C5	2.51	114.64	110.22
7	A	915	UCG	C11-N3-C12	2.56	119.75	116.06
5	D	909	NAG	O5-C1-C2	2.62	115.11	111.47
7	B	915	UCG	C2-C1-C	2.64	105.87	99.95
7	C	901	UCG	O10-C17-C16	2.80	110.99	104.81
5	D	909	NAG	C3-C4-C5	2.83	115.21	110.22
5	C	912	NAG	C1-O5-C5	3.12	116.46	112.17
7	C	901	UCG	C11-N3-C12	3.25	120.73	116.06
7	A	915	UCG	O10-C17-C16	3.35	112.20	104.81
5	B	913	NAG	C3-C4-C5	3.39	116.19	110.22
7	B	915	UCG	C11-N3-C12	3.49	121.07	116.06
7	D	902	UCG	O10-C17-C16	3.70	112.97	104.81

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	C	901	UCG	C12-N2-C10	3.93	119.75	115.16
7	B	915	UCG	O10-C17-C16	3.98	113.60	104.81
5	C	914	NAG	C1-O5-C5	4.04	117.73	112.17
7	A	915	UCG	C12-N2-C10	4.38	120.27	115.16
7	D	902	UCG	C12-N2-C10	4.76	120.72	115.16
7	B	915	UCG	C12-N2-C10	5.24	121.28	115.16
7	D	902	UCG	C7-N1-C8	5.89	119.19	114.13
5	D	914	NAG	C1-O5-C5	6.55	121.19	112.17
7	A	915	UCG	O13-C16-C15	6.56	117.49	105.13
7	A	915	UCG	C7-N1-C8	6.69	119.88	114.13
2	D	901	URI	C4-N3-C2	7.08	120.21	114.13
7	B	915	UCG	C7-N1-C8	7.23	120.34	114.13
7	C	901	UCG	C7-N1-C8	7.25	120.36	114.13
7	B	915	UCG	O13-C16-C15	7.39	119.05	105.13
5	D	909	NAG	C1-O5-C5	7.89	123.04	112.17
7	D	902	UCG	O13-C16-C15	8.28	120.73	105.13
2	C	902	URI	C4-N3-C2	8.45	121.39	114.13
2	A	901	URI	C4-N3-C2	8.46	121.40	114.13
7	B	915	UCG	O12-C15-C16	8.82	121.74	105.13
7	D	902	UCG	O12-C15-C16	8.82	121.76	105.13
7	C	901	UCG	O12-C15-C16	8.99	122.07	105.13
7	C	901	UCG	O13-C16-C15	9.11	122.29	105.13
2	B	901	URI	C4-N3-C2	9.30	122.12	114.13
7	A	915	UCG	O12-C15-C16	9.80	123.61	105.13

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

8 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	914	NAG	1	0
7	A	915	UCG	2	0
5	B	911	NAG	1	0
5	B	914	NAG	1	0
7	B	915	UCG	1	0
7	C	901	UCG	3	0
7	D	902	UCG	1	0
5	D	909	NAG	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	743/811 (91%)	-0.47	4 (0%) 90 89	17, 35, 71, 113	1 (0%)
1	B	737/811 (90%)	-0.44	10 (1%) 75 74	17, 41, 77, 108	1 (0%)
1	C	731/811 (90%)	0.05	50 (6%) 18 16	18, 54, 110, 157	1 (0%)
1	D	743/811 (91%)	-0.24	29 (3%) 40 39	18, 44, 97, 146	1 (0%)
All	All	2954/3244 (91%)	-0.28	93 (3%) 49 47	17, 42, 96, 157	4 (0%)

All (93) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	64	TYR	8.8
1	D	751	ILE	8.1
1	C	733	SER	6.3
1	B	65	VAL	6.0
1	C	81	GLU	5.6
1	C	458	ASP	5.4
1	D	774	PHE	5.3
1	C	123	ASN	5.3
1	C	84	GLN	5.1
1	C	807	GLY	4.9
1	C	85	GLY	4.7
1	C	806	PRO	4.7
1	C	188	GLU	4.7
1	C	63	LYS	4.6
1	B	86	LEU	4.5
1	D	783	PHE	4.4
1	C	86	LEU	4.4
1	B	64	TYR	4.3
1	C	780	ILE	4.3
1	C	46	ILE	4.3
1	D	809	GLN	4.2

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Mol	Chain	Res	Type	RSRZ
1	D	39	LYS	4.1
1	C	65	VAL	4.1
1	C	87	GLN	4.0
1	C	78	ILE	4.0
1	D	761	THR	3.9
1	D	750	THR	3.8
1	A	40	GLN	3.7
1	D	736	SER	3.7
1	D	64	TYR	3.7
1	D	678	PHE	3.6
1	B	46	ILE	3.6
1	D	785	ARG	3.6
1	D	776	CYS	3.6
1	D	777	THR	3.5
1	C	201	THR	3.4
1	C	125	LYS	3.3
1	C	121	PHE	3.3
1	C	186	VAL	3.3
1	D	778	CYS	3.3
1	C	785	ARG	3.3
1	D	760	THR	3.2
1	C	89	LEU	3.1
1	C	808	ASP	3.1
1	D	780	ILE	3.1
1	D	789	GLU	3.1
1	C	732	LEU	3.1
1	B	83	PHE	3.0
1	C	83	PHE	3.0
1	C	786	TRP	3.0
1	D	812	LYS	3.0
1	C	37	GLU	2.9
1	C	82	SER	2.9
1	C	114	LEU	2.9
1	C	122	LEU	2.9
1	D	38	LYS	2.9
1	D	46	ILE	2.8
1	C	781	GLY	2.8
1	C	245	LEU	2.7
1	D	810	ARG	2.7
1	C	38	LYS	2.7
1	D	656	ASN	2.7
1	C	225	ARG	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	733	SER	2.6
1	C	70	LEU	2.5
1	D	170	ILE	2.5
1	C	243	LYS	2.5
1	D	703	LEU	2.5
1	B	88	ASN	2.5
1	D	762	THR	2.5
1	D	788	ASP	2.5
1	B	87	GLN	2.5
1	C	166	ILE	2.5
1	C	58	PRO	2.5
1	C	459	PHE	2.4
1	B	63	LYS	2.4
1	B	753	LYS	2.4
1	D	813	SER	2.4
1	C	168	ARG	2.3
1	C	734	GLU	2.3
1	C	805	SER	2.3
1	C	731	PHE	2.3
1	C	77	HIS	2.2
1	C	54	LEU	2.2
1	B	121	PHE	2.2
1	D	100	VAL	2.2
1	C	76	THR	2.1
1	D	814	ILE	2.1
1	C	45	VAL	2.1
1	C	199	THR	2.1
1	A	780	ILE	2.1
1	A	732	LEU	2.1
1	C	68	LEU	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
3	NAG	C	903	14/15	0.94	0.19	1.29	39,44,49,50	0
3	NAG	B	902	14/15	0.97	0.14	0.18	23,26,31,45	0
6	NAG	B	909	14/15	0.97	0.12	0.06	19,23,28,35	0
3	NAG	D	903	14/15	0.97	0.12	-0.12	17,23,29,38	0
6	NAG	D	910	14/15	0.97	0.13	-0.13	24,31,38,38	0
6	NAG	C	910	14/15	0.97	0.09	-0.26	22,28,36,42	0
6	NAG	A	909	14/15	0.98	0.11	-0.53	17,20,25,25	0
3	NAG	C	906	14/15	0.97	0.08	-0.55	19,23,25,26	0
3	NAG	A	902	14/15	0.98	0.12	-0.56	22,24,28,31	0
3	NAG	C	904	14/15	0.97	0.12	-0.62	46,49,52,53	0
3	NAG	D	904	14/15	0.97	0.10	-0.75	27,32,33,41	0
3	NAG	B	903	14/15	0.98	0.10	-0.86	29,35,42,48	0
3	NAG	A	903	14/15	0.97	0.10	-0.97	23,27,32,36	0
4	NAG	A	907	14/15	0.96	0.10	-1.22	23,27,35,36	0
4	NAG	A	906	14/15	0.96	0.09	-1.50	18,25,27,28	0
3	NAG	B	905	14/15	0.98	0.07	-1.57	23,25,27,27	0
3	NAG	D	906	14/15	0.96	0.08	-1.80	29,33,35,36	0
6	NAG	C	911	14/15	0.92	0.13	-	33,40,46,51	0
3	NAG	B	906	14/15	0.98	0.13	-	22,27,31,34	0
4	BMA	A	905	11/12	0.92	0.17	-	41,43,50,51	0
3	BMA	D	908	11/12	0.91	0.12	-	50,57,63,68	0
6	NAG	B	910	14/15	0.96	0.11	-	32,37,41,44	0
3	BMA	D	905	11/12	0.83	0.14	-	40,48,51,56	0
6	NAG	D	911	14/15	0.94	0.13	-	36,46,52,52	0
3	BMA	C	908	11/12	0.92	0.19	-	43,50,57,59	0
3	BMA	B	904	11/12	0.82	0.16	-	44,50,54,55	0
6	NAG	A	910	14/15	0.95	0.10	-	27,33,40,43	0
3	NAG	C	907	14/15	0.96	0.13	-	28,32,38,39	0
3	BMA	C	905	11/12	0.90	0.10	-	52,57,59,61	0
3	BMA	A	904	11/12	0.85	0.10	-	34,37,43,45	0
3	BMA	B	907	11/12	0.94	0.17	-	40,45,54,54	0
3	NAG	D	907	14/15	0.97	0.10	-	39,42,48,50	0

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	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
5	NAG	A	911	14/15	0.88	0.17	3.42	46,52,56,60	0
5	NAG	D	913	14/15	0.84	0.24	2.32	62,72,76,80	0
5	NAG	C	914	14/15	0.92	0.12	0.73	52,60,68,68	0
5	NAG	D	909	14/15	0.95	0.14	0.63	37,41,47,51	0
5	NAG	B	913	14/15	0.92	0.12	0.56	46,53,59,61	0
5	NAG	C	913	14/15	0.95	0.12	0.38	38,57,68,76	0
5	NAG	C	909	14/15	0.86	0.14	0.14	62,71,82,85	0
7	UCG	D	902	47/47	0.95	0.12	0.03	21,38,65,81	0
2	URI	C	902	17/17	0.96	0.12	0.01	26,35,45,48	0
2	URI	D	901	17/17	0.98	0.10	-0.22	17,23,29,34	0
7	UCG	B	915	47/47	0.94	0.12	-0.42	36,49,83,96	0
7	UCG	A	915	47/47	0.96	0.12	-0.54	24,31,70,79	0
2	URI	B	901	17/17	0.98	0.11	-0.59	23,25,30,35	0
5	NAG	B	912	14/15	0.94	0.11	-0.59	40,45,51,56	0
7	UCG	C	901	47/47	0.95	0.11	-1.04	40,48,83,106	0
5	NAG	A	908	14/15	0.94	0.14	-1.38	32,38,46,52	0
2	URI	A	901	17/17	0.98	0.10	-2.33	21,28,34,36	0
5	NAG	D	914	14/15	0.94	0.13	-	59,69,79,79	0
5	NAG	B	911	14/15	0.84	0.25	-	50,61,76,77	0
5	NAG	A	913	14/15	0.85	0.19	-	54,65,69,69	0
5	NAG	C	912	14/15	0.94	0.12	-	44,50,78,87	0
5	NAG	B	908	14/15	0.89	0.33	-	71,83,94,111	0
5	NAG	A	914	14/15	0.95	0.15	-	38,49,54,57	0
5	NAG	A	912	14/15	0.83	0.22	-	64,73,80,83	0
5	NAG	B	914	14/15	0.91	0.23	-	53,60,69,71	0
5	NAG	D	912	14/15	0.94	0.12	-	54,62,64,66	0

6.5 Other polymers ⓘ

There are no such residues in this entry.