Full wwPDB X-ray Structure Validation Report (i)

Dec 15, 2020 – 02:21 PM JST

PDB ID	:	6AH4
Title	:	Structure of human P2X3 receptor in complex with ATP and Ca2+ ion
Authors	:	Hattori, M.
Deposited on	:	2018-08-16
Resolution	:	3.30 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

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

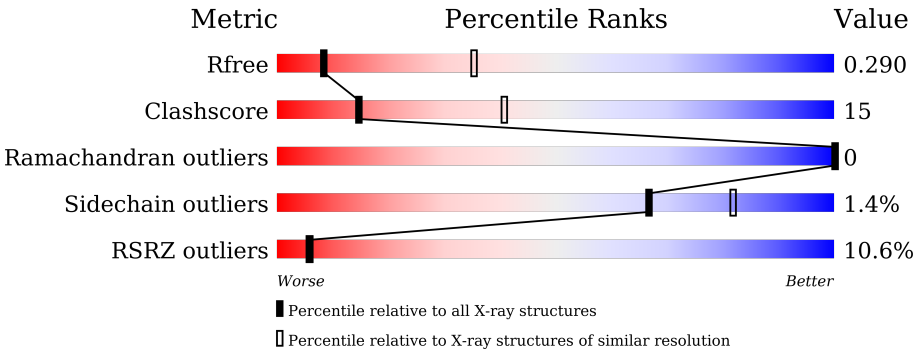
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.15.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.15.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.30 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1149 (3.34-3.26)
Clashscore	141614	1205 (3.34-3.26)
Ramachandran outliers	138981	1183 (3.34-3.26)
Sidechain outliers	138945	1182 (3.34-3.26)
RSRZ outliers	127900	1115 (3.34-3.26)

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 of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	362	4% 66% 31%
1	B	362	14% 64% 34%
1	C	362	13% 69% 29%
2	D	2	50% 50%
2	E	2	100%
2	F	2	100%

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
2	G	2	100%
2	H	2	50% 50%
2	I	2	100%

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
5	EDO	A	407	-	-	-	X
5	EDO	C	406	-	-	-	X

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 8668 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 P2X purinoceptor 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	357	Total	C	N	O	S	0	0	0
			2770	1788	456	508	18			
1	B	357	Total	C	N	O	S	0	0	0
			2774	1790	456	510	18			
1	C	357	Total	C	N	O	S	0	0	0
			2766	1786	456	506	18			

There are 45 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	2	SER	-	expression tag	UNP P56373
A	3	ARG	-	expression tag	UNP P56373
A	4	GLU	-	expression tag	UNP P56373
A	5	PHE	-	expression tag	UNP P56373
A	6	ASP	-	expression tag	UNP P56373
A	7	PHE	-	expression tag	UNP P56373
A	8	PHE	-	expression tag	UNP P56373
A	9	THR	-	expression tag	UNP P56373
A	10	TYR	-	expression tag	UNP P56373
A	11	GLU	-	expression tag	UNP P56373
A	12	THR	-	expression tag	UNP P56373
A	13	PRO	-	expression tag	UNP P56373
A	14	LYS	-	expression tag	UNP P56373
A	15	VAL	-	expression tag	UNP P56373
A	16	ILE	-	expression tag	UNP P56373
B	2	SER	-	expression tag	UNP P56373
B	3	ARG	-	expression tag	UNP P56373
B	4	GLU	-	expression tag	UNP P56373
B	5	PHE	-	expression tag	UNP P56373
B	6	ASP	-	expression tag	UNP P56373
B	7	PHE	-	expression tag	UNP P56373
B	8	PHE	-	expression tag	UNP P56373
B	9	THR	-	expression tag	UNP P56373

Continued on next page...

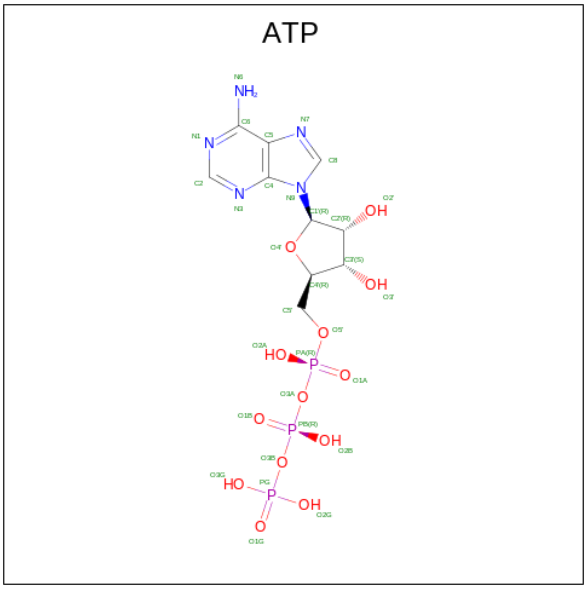
Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
B	10	TYR	-	expression tag	UNP P56373
B	11	GLU	-	expression tag	UNP P56373
B	12	THR	-	expression tag	UNP P56373
B	13	PRO	-	expression tag	UNP P56373
B	14	LYS	-	expression tag	UNP P56373
B	15	VAL	-	expression tag	UNP P56373
B	16	ILE	-	expression tag	UNP P56373
C	2	SER	-	expression tag	UNP P56373
C	3	ARG	-	expression tag	UNP P56373
C	4	GLU	-	expression tag	UNP P56373
C	5	PHE	-	expression tag	UNP P56373
C	6	ASP	-	expression tag	UNP P56373
C	7	PHE	-	expression tag	UNP P56373
C	8	PHE	-	expression tag	UNP P56373
C	9	THR	-	expression tag	UNP P56373
C	10	TYR	-	expression tag	UNP P56373
C	11	GLU	-	expression tag	UNP P56373
C	12	THR	-	expression tag	UNP P56373
C	13	PRO	-	expression tag	UNP P56373
C	14	LYS	-	expression tag	UNP P56373
C	15	VAL	-	expression tag	UNP P56373
C	16	ILE	-	expression tag	UNP P56373

- Molecule 2 is an oligosaccharide called alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose.

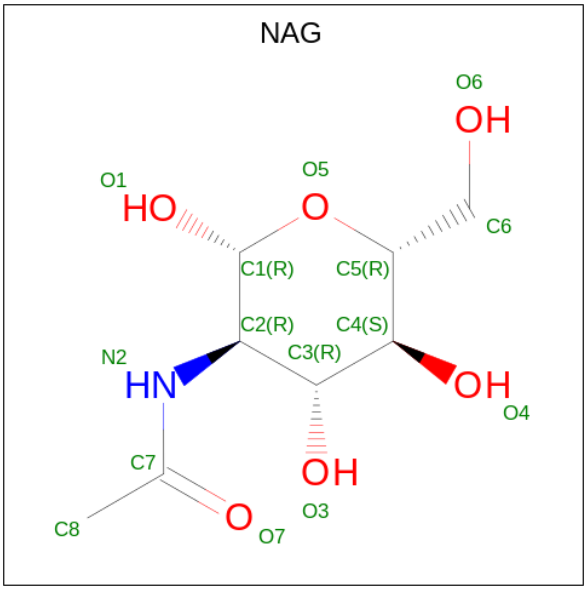
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	D	2	Total	C	O	0	0	0
			23	12	11			
2	E	2	Total	C	O	0	0	0
			23	12	11			
2	F	2	Total	C	O	0	0	0
			23	12	11			
2	G	2	Total	C	O	0	0	0
			23	12	11			
2	H	2	Total	C	O	0	0	0
			23	12	11			
2	I	2	Total	C	O	0	0	0
			23	12	11			

- Molecule 3 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$).



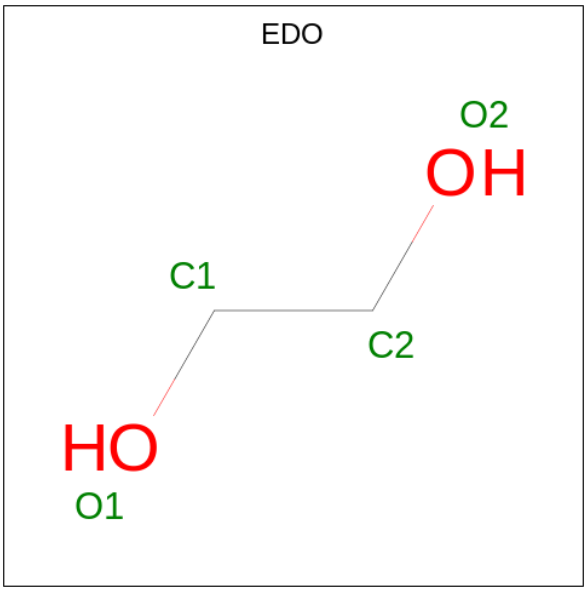
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
3	A	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
3	B	1	Total	C	N	O	P	0	0
			31	10	5	13	3		

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



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

- Molecule 5 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total C O 4 2 2	0	0
5	B	1	Total C O 4 2 2	0	0
5	C	1	Total C O 4 2 2	0	0

- Molecule 6 is CALCIUM ION (three-letter code: CA) (formula: Ca).

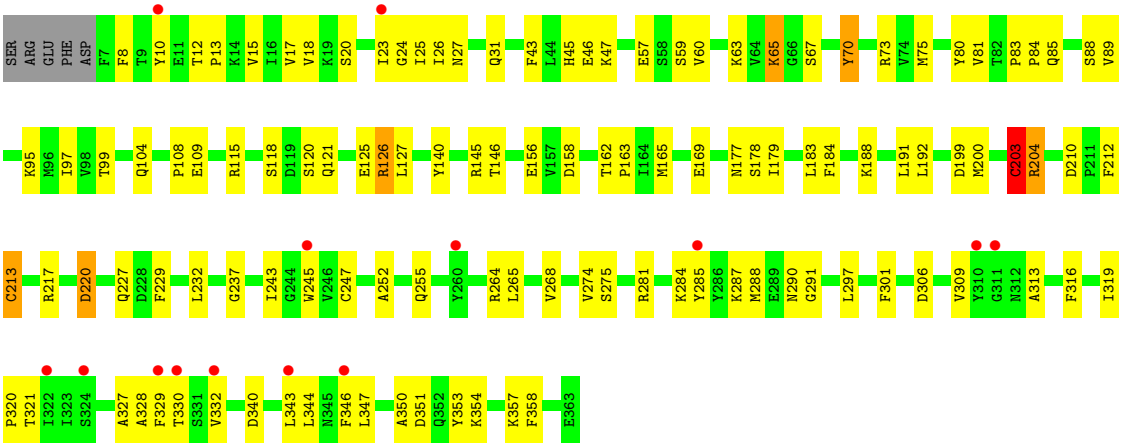
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	B	1	Total Ca 1 1	0	0
6	A	1	Total Ca 1 1	0	0
6	C	1	Total Ca 1 1	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.

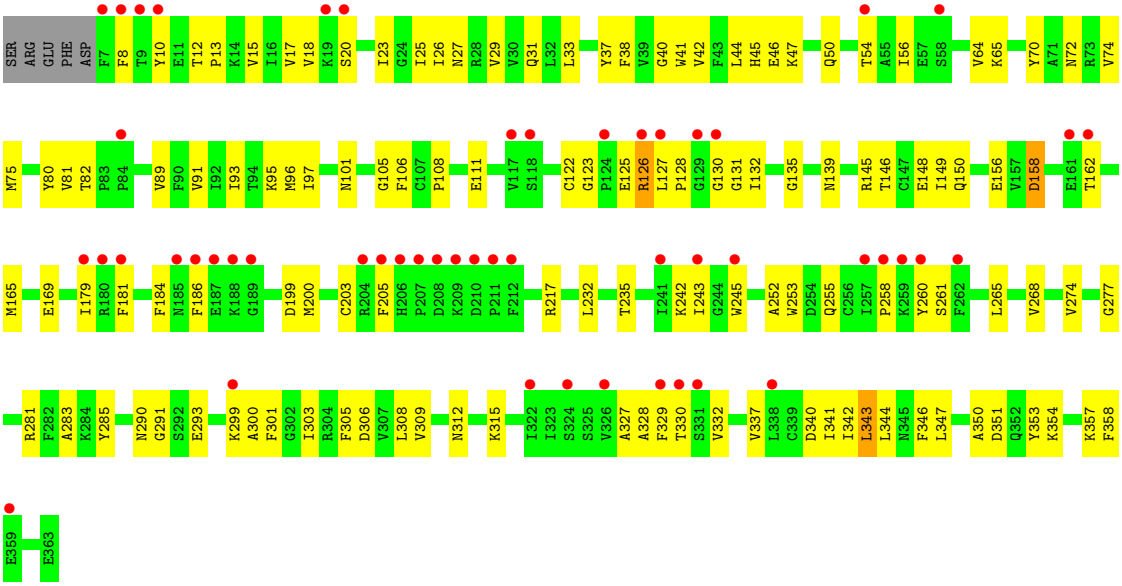
• Molecule 1: P2X purinoceptor 3

Chain A:



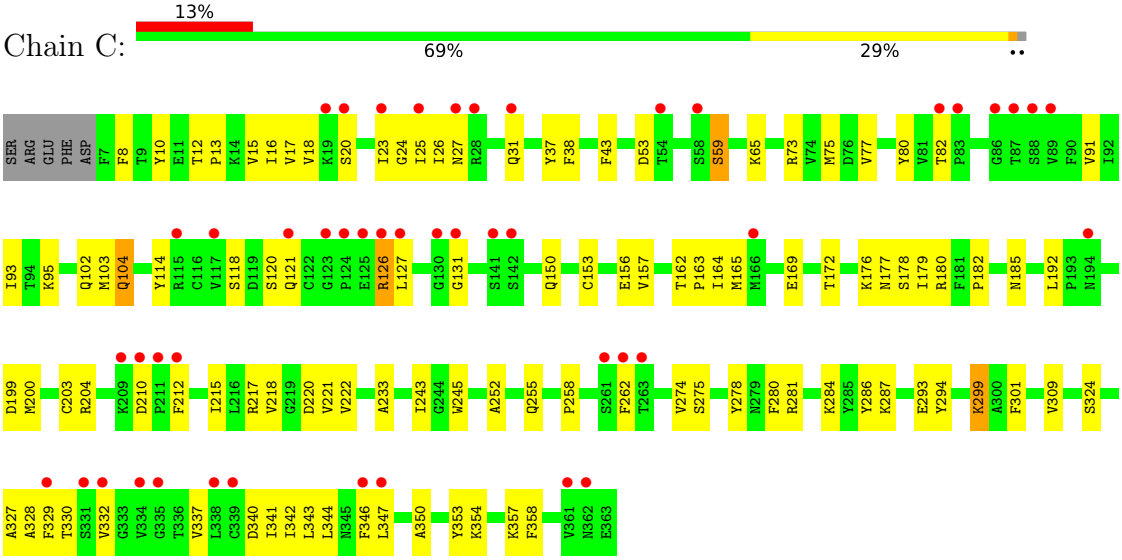
• Molecule 1: P2X purinoceptor 3

Chain B:



• Molecule 1: P2X purinoceptor 3

Chain C:

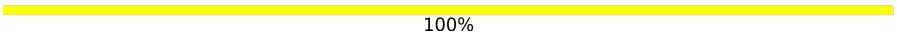


• Molecule 2: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain D:

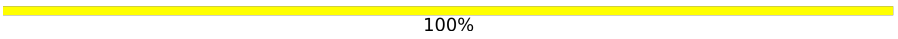
• Molecule 2: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain E:



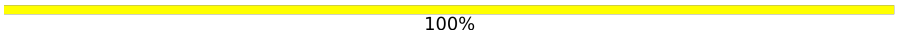
• Molecule 2: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain F:



• Molecule 2: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

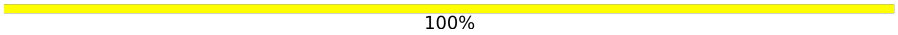
Chain G:



• Molecule 2: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain H:

• Molecule 2: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain I: 

4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α, β, γ	110.90Å 142.50Å 325.34Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.05 – 3.30 49.37 – 3.30	Depositor EDS
% Data completeness (in resolution range)	99.5 (48.05-3.30) 86.2 (49.37-3.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.19 (at 3.33Å)	Xtriage
Refinement program	PHENIX (1.12_2829: ???)	Depositor
R, R_{free}	0.240 , 0.290 0.239 , 0.290	Depositor DCC
R_{free} test set	2000 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å^2)	97.9	Xtriage
Anisotropy	0.997	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.26 , 113.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.84	EDS
Total number of atoms	8668	wwPDB-VP
Average B, all atoms (Å^2)	162.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.06% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: CA, GLC, NAG, ATP, EDO

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	0.57	2/2832 (0.1%)	0.78	4/3843 (0.1%)
1	B	0.51	0/2836	0.73	2/3848 (0.1%)
1	C	0.49	0/2828	0.72	1/3838 (0.0%)
All	All	0.52	2/8496 (0.0%)	0.74	7/11529 (0.1%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	B	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	A	204	ARG	CB-CG	7.20	1.72	1.52
1	A	213	CYS	CB-SG	-5.31	1.73	1.81

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	65	LYS	CD-CE-NZ	7.65	129.29	111.70
1	C	299	LYS	CD-CE-NZ	6.96	127.70	111.70
1	B	111	GLU	CA-CB-CG	6.54	127.80	113.40
1	A	191	LEU	CA-CB-CG	5.53	128.01	115.30
1	A	284	LYS	CD-CE-NZ	5.50	124.34	111.70
1	A	220	ASP	CB-CG-OD2	-5.17	113.65	118.30

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	343	LEU	CA-CB-CG	5.08	127.00	115.30

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	158	ASP	Peptide
1	A	203	CYS	Peptide
1	B	158	ASP	Peptide

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	2770	0	2735	107	0
1	B	2774	0	2742	106	0
1	C	2766	0	2732	90	0
2	D	23	0	21	1	0
2	E	23	0	21	0	0
2	F	23	0	21	0	0
2	G	23	0	21	0	0
2	H	23	0	21	0	0
2	I	23	0	21	0	0
3	A	62	0	24	4	0
3	B	31	0	12	1	0
4	A	28	0	26	0	0
4	B	42	0	39	2	0
4	C	42	0	39	1	0
5	A	4	0	6	0	0
5	B	4	0	6	1	0
5	C	4	0	6	0	0
6	A	1	0	0	0	0
6	B	1	0	0	0	0
6	C	1	0	0	0	0
All	All	8668	0	8493	263	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 15.

All (263) 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:252:ALA:HB3	1:A:255:GLN:HG2	1.60	0.82
1:A:13:PRO:HA	1:B:15:VAL:HA	1.63	0.80
1:A:192:LEU:HD11	1:A:212:PHE:CD1	2.19	0.78
1:B:12:THR:OG1	1:C:357:LYS:NZ	2.19	0.76
1:B:252:ALA:HB3	1:B:255:GLN:HG2	1.67	0.76
1:A:125:GLU:HG3	1:A:126:ARG:HD2	1.67	0.75
1:B:290:ASN:HD21	4:B:403:NAG:C7	2.02	0.72
1:B:10:TYR:OH	1:C:340:ASP:OD1	2.08	0.72
1:A:10:TYR:OH	1:B:340:ASP:OD1	2.08	0.71
1:A:73:ARG:NH1	1:B:293:GLU:OE2	2.24	0.71
1:C:118:SER:HG	1:C:120:SER:HG	1.34	0.71
1:C:252:ALA:HB3	1:C:255:GLN:HG2	1.73	0.69
1:C:192:LEU:HG	1:C:212:PHE:CE1	2.28	0.69
1:B:23:ILE:HA	1:B:26:ILE:HG22	1.73	0.68
1:A:125:GLU:HG3	1:A:126:ARG:CD	2.24	0.68
1:A:357:LYS:NZ	1:C:12:THR:OG1	2.27	0.68
1:A:13:PRO:O	1:B:357:LYS:NZ	2.25	0.68
1:B:351:ASP:HA	1:B:354:LYS:HD3	1.75	0.68
1:B:70:TYR:HE2	1:B:96:MET:HB2	1.59	0.68
1:A:125:GLU:CG	1:A:126:ARG:HD2	2.23	0.67
1:B:158:ASP:OD1	1:B:299:LYS:NZ	2.17	0.66
1:A:23:ILE:HA	1:A:26:ILE:HG22	1.76	0.66
1:B:156:GLU:OE2	1:B:281:ARG:NH1	2.29	0.66
1:A:274:VAL:HG21	1:C:200:MET:HG3	1.76	0.66
1:A:60:VAL:HB	1:A:88:SER:HB3	1.80	0.64
1:B:281:ARG:NH1	1:B:299:LYS:HD2	2.12	0.64
1:A:204:ARG:N	1:A:213:CYS:SG	2.71	0.64
1:B:27:ASN:O	1:B:31:GLN:HG3	1.98	0.63
1:C:169:GLU:HB3	4:C:401:NAG:H82	1.80	0.63
1:A:27:ASN:O	1:A:31:GLN:HG3	1.99	0.62
1:A:340:ASP:OD1	1:C:10:TYR:OH	2.17	0.62
1:B:169:GLU:HG2	1:B:217:ARG:NH2	2.15	0.62
1:A:287:LYS:HG2	1:C:286:TYR:CD1	2.35	0.62
1:A:350:ALA:HA	1:A:353:TYR:HD2	1.64	0.62
1:A:350:ALA:HA	1:A:353:TYR:CD2	2.35	0.62
1:B:108:PRO:HA	1:B:148:GLU:HA	1.81	0.62
1:C:23:ILE:HA	1:C:26:ILE:HG22	1.82	0.61
1:B:283:ALA:HB3	1:B:285:TYR:HE1	1.66	0.61

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:45:HIS:ND1	1:A:46:GLU:HG2	2.16	0.61
1:B:283:ALA:HB3	1:B:285:TYR:CE1	2.36	0.61
1:B:95:LYS:HB3	1:B:301:PHE:HB2	1.83	0.60
1:A:343:LEU:HD12	1:A:344:LEU:HD12	1.82	0.60
1:C:281:ARG:NH1	1:C:299:LYS:HD2	2.16	0.60
1:C:350:ALA:HA	1:C:353:TYR:HD2	1.66	0.60
1:C:329:PHE:O	1:C:332:VAL:HG22	2.02	0.60
1:B:127:LEU:HD23	1:B:127:LEU:H	1.66	0.59
1:A:12:THR:HG1	1:B:357:LYS:HZ3	1.47	0.59
1:A:127:LEU:HD23	1:A:127:LEU:H	1.67	0.59
1:C:350:ALA:HA	1:C:353:TYR:CD2	2.37	0.59
1:A:104:GLN:HE21	1:C:73:ARG:HB3	1.67	0.59
1:A:204:ARG:O	1:A:210:ASP:HB3	2.03	0.59
1:B:56:ILE:HD11	1:B:179:ILE:HB	1.85	0.59
1:B:328:ALA:O	1:B:332:VAL:HG13	2.03	0.58
1:C:127:LEU:H	1:C:127:LEU:HD23	1.68	0.58
1:A:70:TYR:HE1	1:A:163:PRO:HG2	1.68	0.58
1:B:8:PHE:CD2	1:C:25:ILE:HG22	2.37	0.58
1:A:109:GLU:O	1:A:146:THR:OG1	2.14	0.58
1:C:156:GLU:OE2	1:C:281:ARG:NH1	2.36	0.58
1:A:285:TYR:CD1	1:C:284:LYS:HE2	2.38	0.58
1:A:357:LYS:HD3	1:A:358:PHE:CE1	2.39	0.58
1:C:80:TYR:O	1:C:93:ILE:HG13	2.04	0.57
1:A:108:PRO:HG2	1:A:140:TYR:HD1	1.70	0.57
1:A:179:ILE:HG12	1:A:245:TRP:CE2	2.38	0.57
1:A:65:LYS:HG3	3:A:401:ATP:C6	2.39	0.57
1:A:316:PHE:HE2	1:A:321:THR:HG1	1.52	0.57
1:A:12:THR:OG1	1:B:357:LYS:NZ	2.19	0.57
1:A:23:ILE:HG13	1:A:343:LEU:HD23	1.87	0.57
1:C:164:ILE:HD12	1:C:233:ALA:HB3	1.85	0.57
1:A:59:SER:O	1:A:177:ASN:HA	2.04	0.56
1:B:343:LEU:HD12	1:B:344:LEU:HD12	1.85	0.56
1:A:67:SER:OG	1:B:130:GLY:O	2.16	0.56
1:B:89:VAL:HA	1:B:305:PHE:O	2.06	0.56
1:B:329:PHE:O	1:B:332:VAL:HG22	2.06	0.56
1:B:243:ILE:HG12	1:B:260:TYR:CE2	2.40	0.56
1:C:185:ASN:ND2	1:C:185:ASN:O	2.38	0.56
1:A:25:ILE:HG22	1:C:8:PHE:CD2	2.41	0.55
1:A:75:MET:HG3	1:A:165:MET:CE	2.37	0.55
1:A:104:GLN:NE2	1:C:73:ARG:HB3	2.22	0.55
1:B:350:ALA:HA	1:B:353:TYR:CD2	2.42	0.55

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:357:LYS:HD3	1:B:358:PHE:CE1	2.42	0.54
1:B:13:PRO:HA	1:C:15:VAL:HA	1.88	0.54
1:A:217:ARG:HB3	1:A:220:ASP:CG	2.27	0.54
1:B:199:ASP:O	1:B:203:CYS:N	2.39	0.54
1:C:327:ALA:O	1:C:330:THR:HG22	2.07	0.54
1:B:8:PHE:HD2	1:C:25:ILE:HG22	1.73	0.54
1:A:84:PRO:HD2	1:A:85:GLN:CD	2.27	0.54
1:B:327:ALA:HA	1:B:330:THR:HG22	1.90	0.54
1:C:131:GLY:HA3	1:C:150:GLN:O	2.09	0.53
1:C:102:GLN:HA	1:C:153:CYS:O	2.09	0.53
1:B:243:ILE:HG12	1:B:260:TYR:HE2	1.73	0.53
1:B:45:HIS:ND1	1:B:46:GLU:HG2	2.24	0.53
1:A:75:MET:HG3	1:A:165:MET:HE1	1.90	0.52
1:C:27:ASN:O	1:C:31:GLN:HG3	2.09	0.52
1:C:343:LEU:HD12	1:C:344:LEU:HD12	1.91	0.52
1:A:275:SER:OG	1:C:176:LYS:NZ	2.20	0.52
1:A:183:LEU:HD23	1:A:184:PHE:CZ	2.44	0.51
1:A:199:ASP:O	1:A:203:CYS:N	2.40	0.51
1:A:118:SER:N	1:A:121:GLN:OE1	2.37	0.51
1:B:337:VAL:O	1:B:341:ILE:HG12	2.11	0.51
1:C:337:VAL:O	1:C:341:ILE:HG12	2.10	0.51
1:C:95:LYS:HB3	1:C:301:PHE:HB2	1.92	0.51
1:A:179:ILE:HG21	1:A:245:TRP:CD1	2.45	0.51
1:A:351:ASP:HA	1:A:354:LYS:HD3	1.92	0.51
1:B:186:PHE:HE1	1:B:253:TRP:NE1	2.08	0.51
1:A:118:SER:OG	1:A:120:SER:OG	2.05	0.51
1:A:63:LYS:NZ	3:A:401:ATP:O2B	2.43	0.51
1:B:235:THR:O	1:B:277:GLY:HA3	2.11	0.51
1:A:243:ILE:HB	1:A:309:VAL:HG22	1.93	0.50
1:C:221:VAL:HG22	1:C:262:PHE:CE2	2.46	0.50
1:A:169:GLU:HB2	1:A:229:PHE:CE1	2.47	0.50
1:C:26:ILE:HD13	1:C:342:ILE:HD13	1.92	0.50
1:A:178:SER:OG	1:B:268:VAL:HG12	2.11	0.50
1:B:23:ILE:HG13	1:B:343:LEU:HD23	1.94	0.50
1:B:70:TYR:CE2	1:B:96:MET:HB2	2.45	0.50
1:C:192:LEU:HG	1:C:212:PHE:CZ	2.47	0.50
1:A:83:PRO:HB2	1:A:85:GLN:OE1	2.11	0.50
1:C:118:SER:N	1:C:121:GLN:OE1	2.26	0.50
1:B:131:GLY:HA3	1:B:150:GLN:O	2.12	0.49
1:C:217:ARG:HD3	1:C:220:ASP:OD1	2.12	0.49
1:B:64:VAL:O	1:B:65:LYS:HG2	2.12	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:156:GLU:OE2	1:A:281:ARG:NH1	2.45	0.49
1:A:57:GLU:O	1:A:179:ILE:HA	2.12	0.49
1:A:15:VAL:HA	1:C:13:PRO:HA	1.93	0.49
1:C:91:VAL:HG21	1:C:278:TYR:OH	2.12	0.49
1:B:8:PHE:O	1:C:24:GLY:HA3	2.11	0.49
1:B:350:ALA:HA	1:B:353:TYR:HD2	1.78	0.49
1:C:287:LYS:HE2	1:C:293:GLU:HG2	1.95	0.49
1:B:179:ILE:HG12	1:B:245:TRP:CE2	2.48	0.48
1:C:343:LEU:CD1	1:C:344:LEU:HD12	2.42	0.48
1:C:162:THR:HB	1:C:163:PRO:HD2	1.95	0.48
1:A:287:LYS:HD2	1:C:294:TYR:CE2	2.48	0.48
1:C:59:SER:O	1:C:177:ASN:HA	2.13	0.48
1:A:329:PHE:O	1:A:332:VAL:HG22	2.13	0.48
1:A:204:ARG:HG3	1:A:210:ASP:CB	2.43	0.48
1:A:343:LEU:CD1	1:A:344:LEU:HD12	2.43	0.48
1:B:242:LYS:HG2	1:B:308:LEU:HB2	1.96	0.48
1:A:95:LYS:HB3	1:A:301:PHE:HB2	1.94	0.48
1:B:290:ASN:OD1	1:B:291:GLY:N	2.47	0.48
1:B:81:VAL:HA	1:B:91:VAL:O	2.14	0.48
1:B:12:THR:HG23	1:C:16:ILE:HB	1.95	0.48
1:B:205:PHE:CE1	1:B:260:TYR:HD1	2.30	0.47
1:C:204:ARG:O	1:C:210:ASP:HB3	2.14	0.47
1:B:75:MET:HG3	1:B:165:MET:HE2	1.95	0.47
1:A:70:TYR:OH	1:A:163:PRO:O	2.22	0.47
1:B:343:LEU:CD1	1:B:344:LEU:HD12	2.43	0.47
1:C:179:ILE:HG12	1:C:245:TRP:CE2	2.49	0.47
1:B:108:PRO:HB2	1:B:146:THR:OG1	2.15	0.47
1:A:344:LEU:HD21	1:A:358:PHE:HE1	1.80	0.47
1:B:123:GLY:C	1:B:125:GLU:H	2.18	0.47
1:B:281:ARG:HG2	1:B:299:LYS:HA	1.96	0.47
1:A:340:ASP:O	1:A:344:LEU:HB2	2.15	0.47
1:B:50:GLN:HB2	1:B:315:LYS:C	2.35	0.47
1:B:342:ILE:HA	1:B:346:PHE:HD2	1.79	0.47
1:B:242:LYS:HB2	1:B:261:SER:OG	2.15	0.47
1:A:84:PRO:HD2	1:A:85:GLN:OE1	2.15	0.47
1:B:245:TRP:CZ2	1:B:258:PRO:HB3	2.51	0.46
1:B:346:PHE:O	1:B:347:LEU:HD12	2.15	0.46
1:A:287:LYS:HD2	1:C:294:TYR:HE2	1.79	0.46
1:B:54:THR:HG22	1:B:312:ASN:HB3	1.97	0.46
1:A:43:PHE:O	1:A:47:LYS:HA	2.15	0.46
1:B:290:ASN:ND2	4:B:403:NAG:O7	2.41	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:106:PHE:CE1	1:B:150:GLN:HB2	2.50	0.46
1:C:103:MET:HG2	1:C:294:TYR:HB3	1.98	0.46
1:A:288:MET:O	1:A:290:ASN:O	2.34	0.46
1:C:200:MET:CE	1:C:215:ILE:HD11	2.45	0.46
1:A:99:THR:HB	1:A:297:LEU:HB3	1.97	0.45
1:C:180:ARG:HH12	1:C:182:PRO:HB3	1.80	0.45
1:C:346:PHE:O	1:C:347:LEU:HD12	2.16	0.45
1:B:64:VAL:HB	1:B:81:VAL:HG21	1.99	0.45
1:C:82:THR:HG21	1:C:280:PHE:HZ	1.82	0.45
1:A:287:LYS:HE3	1:C:286:TYR:CZ	2.51	0.45
1:A:329:PHE:HA	1:A:332:VAL:HG13	1.99	0.45
1:B:200:MET:HG3	1:C:274:VAL:HG21	1.99	0.45
2:D:1:GLC:O3	2:D:2:GLC:O2	2.18	0.45
1:B:74:VAL:N	1:C:104:GLN:OE1	2.41	0.45
1:C:212:PHE:CE2	1:C:258:PRO:HD2	2.52	0.45
1:A:217:ARG:HD3	1:A:220:ASP:OD1	2.16	0.45
3:A:402:ATP:N6	1:C:172:THR:O	2.48	0.45
1:C:43:PHE:CE1	1:C:324:SER:HB3	2.52	0.45
1:A:247:CYS:O	1:A:313:ALA:HA	2.17	0.45
1:A:25:ILE:HG22	1:C:8:PHE:HD2	1.82	0.44
1:A:346:PHE:O	1:A:347:LEU:HD12	2.17	0.44
1:C:53:ASP:OD2	1:C:182:PRO:HD2	2.17	0.44
1:C:18:VAL:HG12	1:C:20:SER:H	1.81	0.44
1:C:218:VAL:O	1:C:222:VAL:HG23	2.18	0.44
1:A:8:PHE:CD2	1:B:25:ILE:HG22	2.52	0.44
1:B:344:LEU:HD21	1:B:358:PHE:CE1	2.53	0.44
1:B:50:GLN:HB2	1:B:315:LYS:O	2.18	0.44
1:C:200:MET:HE3	1:C:215:ILE:HD11	1.99	0.44
1:A:80:TYR:CD1	1:A:81:VAL:HG23	2.52	0.44
1:B:17:VAL:HG21	1:C:358:PHE:HB3	1.99	0.44
1:C:199:ASP:O	1:C:203:CYS:N	2.44	0.44
1:A:290:ASN:OD1	1:A:291:GLY:N	2.50	0.44
1:B:139:ASN:HA	1:B:145:ARG:HG2	1.99	0.44
1:B:101:ASN:HB3	5:B:406:EDO:H21	2.00	0.43
1:A:89:VAL:HG22	1:A:306:ASP:HB3	2.00	0.43
1:A:265:LEU:HD23	1:A:265:LEU:N	2.33	0.43
1:C:243:ILE:HB	1:C:309:VAL:HG22	1.99	0.43
1:B:97:ILE:HG12	1:B:162:THR:HG21	2.01	0.43
1:A:24:GLY:HA3	1:C:8:PHE:O	2.19	0.43
1:A:327:ALA:O	1:A:330:THR:HG22	2.18	0.43
1:C:114:TYR:HA	1:C:126:ARG:HH12	1.84	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:127:LEU:HG	1:B:128:PRO:HD3	1.99	0.43
1:B:232:LEU:HG	1:B:303:ILE:HG12	1.99	0.43
1:B:70:TYR:HB2	1:B:165:MET:HE2	2.01	0.43
1:B:243:ILE:HB	1:B:309:VAL:HG22	2.00	0.43
1:B:132:ILE:O	1:B:149:ILE:HA	2.19	0.43
1:B:265:LEU:HD23	1:B:265:LEU:N	2.34	0.43
1:A:125:GLU:HG2	1:A:126:ARG:HD2	2.00	0.42
1:B:242:LYS:HA	1:B:308:LEU:O	2.18	0.42
1:B:41:TRP:HD1	1:B:42:VAL:HG23	1.83	0.42
1:B:181:PHE:HD2	1:B:186:PHE:HD2	1.66	0.42
1:A:327:ALA:HA	1:A:330:THR:HG22	2.00	0.42
1:A:97:ILE:HG12	1:A:162:THR:CG2	2.49	0.42
1:A:97:ILE:HG12	1:A:162:THR:HG21	2.01	0.42
1:A:17:VAL:HG21	1:B:358:PHE:HD2	1.84	0.42
1:A:179:ILE:HG12	1:A:245:TRP:NE1	2.34	0.42
1:B:18:VAL:HG12	1:B:20:SER:H	1.83	0.42
1:C:82:THR:HG23	1:C:91:VAL:HB	2.02	0.42
1:A:200:MET:HG3	1:B:274:VAL:HG23	2.02	0.42
1:C:342:ILE:HA	1:C:346:PHE:HD2	1.84	0.42
1:A:319:ILE:N	1:A:320:PRO:HD2	2.35	0.42
1:C:118:SER:OG	1:C:120:SER:OG	2.18	0.42
1:C:344:LEU:HD21	1:C:358:PHE:HE1	1.84	0.42
1:C:77:VAL:HA	1:C:80:TYR:CZ	2.54	0.42
1:A:18:VAL:HG12	1:A:20:SER:H	1.85	0.42
1:A:328:ALA:O	1:A:332:VAL:HG13	2.19	0.42
1:B:106:PHE:HE1	1:B:150:GLN:HB2	1.84	0.42
1:A:268:VAL:HG12	1:C:178:SER:OG	2.19	0.42
1:A:179:ILE:HG13	1:A:188:LYS:O	2.20	0.41
1:A:358:PHE:HD2	1:C:17:VAL:HG21	1.85	0.41
1:B:40:GLY:O	1:B:44:LEU:HB2	2.20	0.41
1:C:157:VAL:HG23	1:C:157:VAL:O	2.20	0.41
1:A:84:PRO:HD2	1:A:85:GLN:NE2	2.35	0.41
1:C:118:SER:HG	1:C:120:SER:H	1.67	0.41
1:A:200:MET:HG3	1:B:274:VAL:CG2	2.50	0.41
1:A:169:GLU:HG2	1:A:217:ARG:NH2	2.36	0.41
1:B:82:THR:HG23	1:B:91:VAL:HB	2.02	0.41
1:A:232:LEU:HD11	1:A:237:GLY:HA3	2.02	0.41
1:B:29:VAL:O	1:B:33:LEU:HG	2.21	0.41
1:A:227:GLN:OE1	1:A:264:ARG:NH1	2.51	0.41
1:A:179:ILE:HG12	1:A:245:TRP:CD1	2.56	0.41
1:B:122:CYS:O	1:B:135:GLY:HA2	2.20	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:115:ARG:HA	1:A:145:ARG:O	2.20	0.41
1:A:344:LEU:HD21	1:A:358:PHE:CE1	2.55	0.41
1:B:126:ARG:HD2	1:B:126:ARG:H	1.85	0.41
1:B:93:ILE:HD13	1:B:300:ALA:HB1	2.03	0.41
1:B:80:TYR:CE1	1:B:81:VAL:HG23	2.56	0.41
1:C:328:ALA:O	1:C:332:VAL:HG13	2.21	0.41
3:A:402:ATP:C6	1:C:65:LYS:HG3	2.56	0.41
1:B:184:PHE:CD1	1:B:184:PHE:N	2.87	0.41
3:B:407:ATP:H5'1	1:C:275:SER:HB2	2.02	0.41
1:B:80:TYR:CD1	1:B:81:VAL:HG23	2.56	0.41
1:C:75:MET:HG3	1:C:165:MET:CE	2.51	0.41
1:C:354:LYS:O	1:C:358:PHE:HD1	2.04	0.40
1:A:204:ARG:HG3	1:A:210:ASP:HB3	2.02	0.40
1:B:47:LYS:O	1:B:50:GLN:HG2	2.22	0.40
1:A:274:VAL:CG2	1:C:200:MET:HG3	2.49	0.40
1:C:43:PHE:HE1	1:C:324:SER:HB3	1.86	0.40
1:B:105:GLY:O	1:B:150:GLN:HA	2.21	0.40
1:B:252:ALA:O	1:B:255:GLN:HB2	2.21	0.40
1:B:281:ARG:HH11	1:B:281:ARG:HD3	1.72	0.40
1:B:12:THR:O	1:C:16:ILE:N	2.55	0.40

There are no symmetry-related clashes.

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	355/362 (98%)	333 (94%)	22 (6%)	0	100	100
1	B	355/362 (98%)	335 (94%)	20 (6%)	0	100	100
1	C	355/362 (98%)	337 (95%)	18 (5%)	0	100	100
All	All	1065/1086 (98%)	1005 (94%)	60 (6%)	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	303/319 (95%)	300 (99%)	3 (1%)	76 86
1	B	304/319 (95%)	299 (98%)	5 (2%)	62 79
1	C	302/319 (95%)	297 (98%)	5 (2%)	60 78
All	All	909/957 (95%)	896 (99%)	13 (1%)	67 82

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	70	TYR
1	A	126	ARG
1	A	203	CYS
1	B	37	TYR
1	B	38	PHE
1	B	72	ASN
1	B	126	ARG
1	B	306	ASP
1	C	37	TYR
1	C	38	PHE
1	C	59	SER
1	C	104	GLN
1	C	126	ARG

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

5.3.3 RNA

There are no RNA 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)

12 monosaccharides 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	GLC	D	1	2	12,12,12	0.48	0	17,17,17	0.70	0
2	GLC	D	2	2	11,11,12	0.54	0	15,15,17	1.60	3 (20%)
2	GLC	E	1	2	12,12,12	0.46	0	17,17,17	1.24	2 (11%)
2	GLC	E	2	2	11,11,12	0.59	0	15,15,17	1.17	3 (20%)
2	GLC	F	1	2	12,12,12	0.62	0	17,17,17	0.84	1 (5%)
2	GLC	F	2	2	11,11,12	0.60	0	15,15,17	1.22	1 (6%)
2	GLC	G	1	2	12,12,12	0.53	0	17,17,17	0.91	1 (5%)
2	GLC	G	2	2	11,11,12	0.60	0	15,15,17	1.99	5 (33%)
2	GLC	H	1	2	12,12,12	0.48	0	17,17,17	0.71	0
2	GLC	H	2	2	11,11,12	0.56	0	15,15,17	1.32	1 (6%)
2	GLC	I	1	2	12,12,12	0.61	0	17,17,17	1.17	3 (17%)
2	GLC	I	2	2	11,11,12	0.67	0	15,15,17	1.43	1 (6%)

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	GLC	D	1	2	-	2/2/22/22	0/1/1/1
2	GLC	D	2	2	-	1/2/19/22	0/1/1/1
2	GLC	E	1	2	-	1/2/22/22	0/1/1/1
2	GLC	E	2	2	-	2/2/19/22	0/1/1/1
2	GLC	F	1	2	-	2/2/22/22	0/1/1/1

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GLC	F	2	2	-	2/2/19/22	0/1/1/1
2	GLC	G	1	2	-	2/2/22/22	0/1/1/1
2	GLC	G	2	2	-	0/2/19/22	0/1/1/1
2	GLC	H	1	2	-	2/2/22/22	0/1/1/1
2	GLC	H	2	2	-	2/2/19/22	0/1/1/1
2	GLC	I	1	2	-	0/2/22/22	0/1/1/1
2	GLC	I	2	2	-	1/2/19/22	0/1/1/1

There are no bond length outliers.

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	I	2	GLC	C1-C2-C3	4.31	114.96	109.67
2	G	2	GLC	C1-C2-C3	4.10	114.71	109.67
2	H	2	GLC	C1-O5-C5	3.97	117.57	112.19
2	G	2	GLC	C1-O5-C5	3.68	117.18	112.19
2	D	2	GLC	C1-C2-C3	3.49	113.95	109.67
2	G	2	GLC	O5-C1-C2	3.16	115.64	110.77
2	D	2	GLC	C2-C3-C4	2.91	115.94	110.89
2	E	1	GLC	C1-O5-C5	2.76	118.87	113.66
2	G	2	GLC	C2-C3-C4	2.44	115.11	110.89
2	E	2	GLC	C1-O5-C5	2.36	115.39	112.19
2	F	2	GLC	O5-C5-C6	2.35	110.89	107.20
2	I	1	GLC	C4-C3-C2	2.35	114.92	110.82
2	D	2	GLC	C3-C4-C5	2.29	114.32	110.24
2	G	1	GLC	O5-C5-C6	2.29	112.12	106.44
2	I	1	GLC	O5-C5-C6	2.28	112.09	106.44
2	E	2	GLC	O2-C2-C1	2.27	113.80	109.15
2	E	1	GLC	O5-C5-C4	2.23	113.75	109.69
2	G	2	GLC	C3-C4-C5	2.22	114.19	110.24
2	E	2	GLC	O5-C5-C6	2.18	110.62	107.20
2	F	1	GLC	O2-C2-C1	2.18	114.21	109.16
2	I	1	GLC	C3-C4-C5	2.02	113.84	110.24

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	F	1	GLC	C4-C5-C6-O6
2	D	1	GLC	O5-C5-C6-O6

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
2	H	2	GLC	C4-C5-C6-O6
2	H	2	GLC	O5-C5-C6-O6
2	F	1	GLC	O5-C5-C6-O6
2	H	1	GLC	O5-C5-C6-O6
2	F	2	GLC	C4-C5-C6-O6
2	D	1	GLC	C4-C5-C6-O6
2	H	1	GLC	C4-C5-C6-O6
2	G	1	GLC	C4-C5-C6-O6
2	E	2	GLC	C4-C5-C6-O6
2	F	2	GLC	O5-C5-C6-O6
2	E	1	GLC	C4-C5-C6-O6
2	I	2	GLC	O5-C5-C6-O6
2	D	2	GLC	O5-C5-C6-O6
2	E	2	GLC	O5-C5-C6-O6
2	G	1	GLC	O5-C5-C6-O6

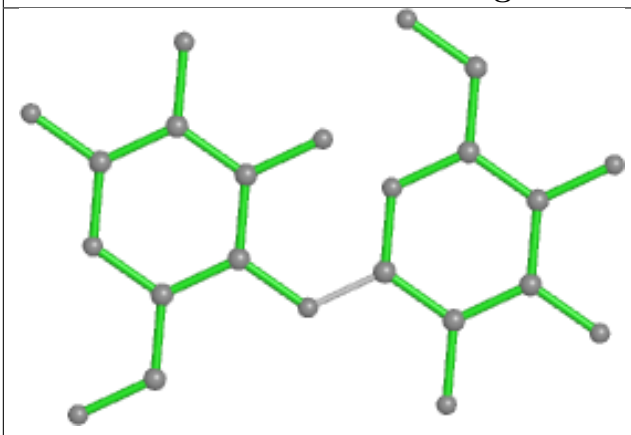
There are no ring outliers.

2 monomers are involved in 1 short contact:

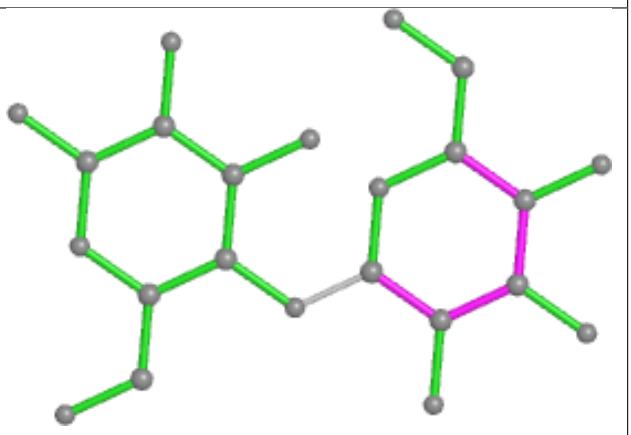
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	1	GLC	1	0
2	D	2	GLC	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

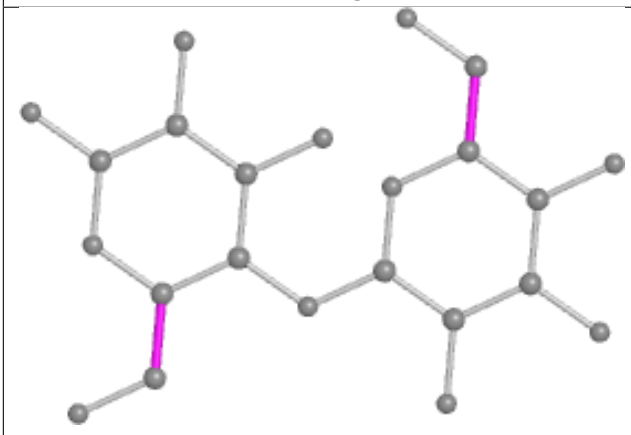
Oligosaccharide Chain D



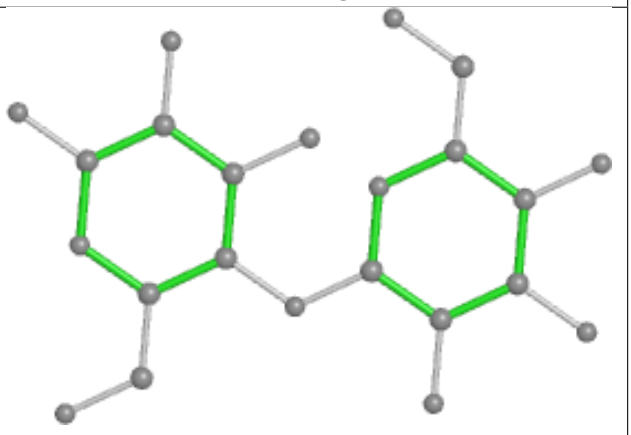
Bond lengths



Bond angles

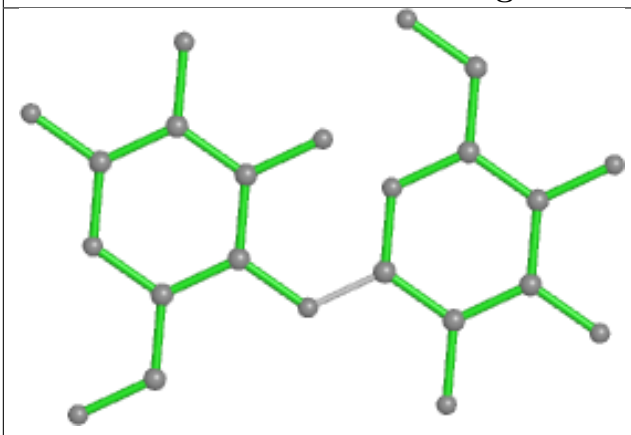


Torsions

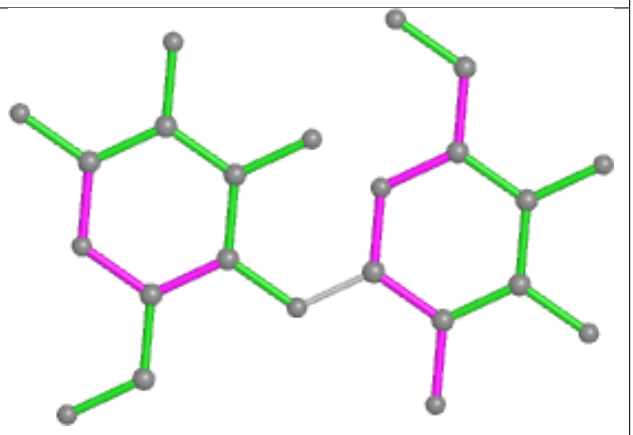


Rings

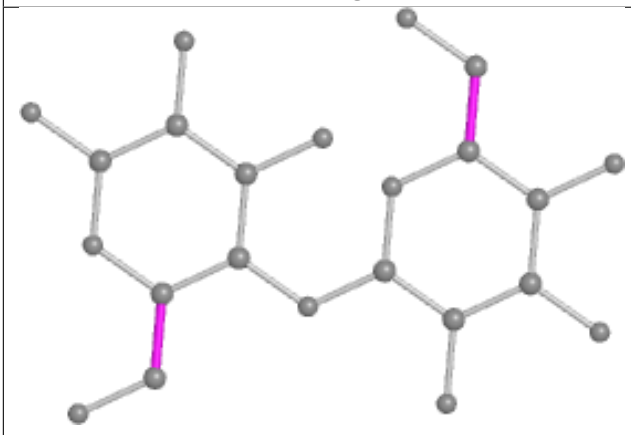
Oligosaccharide Chain E



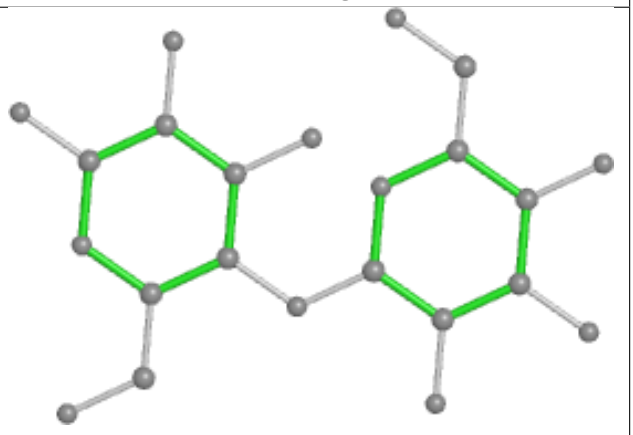
Bond lengths



Bond angles

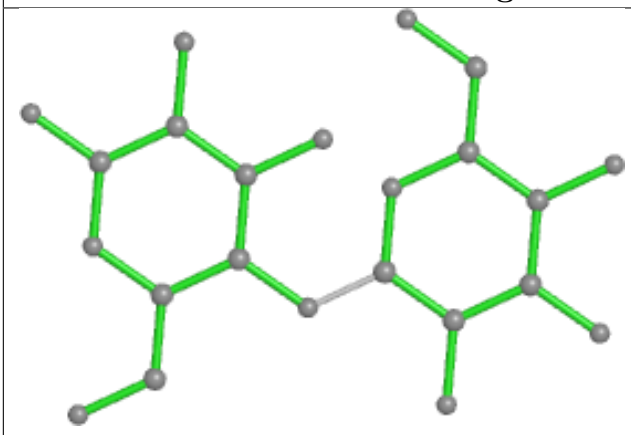


Torsions

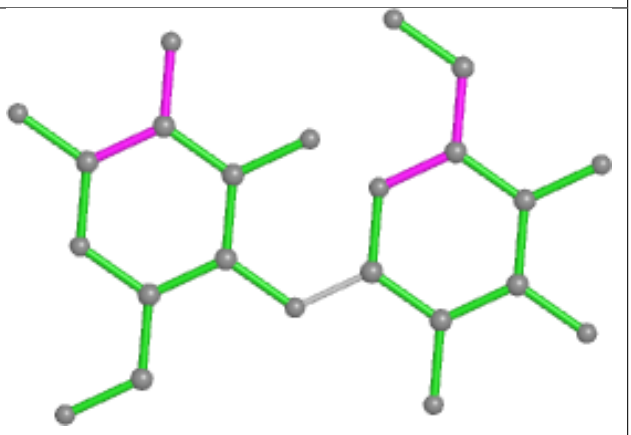


Rings

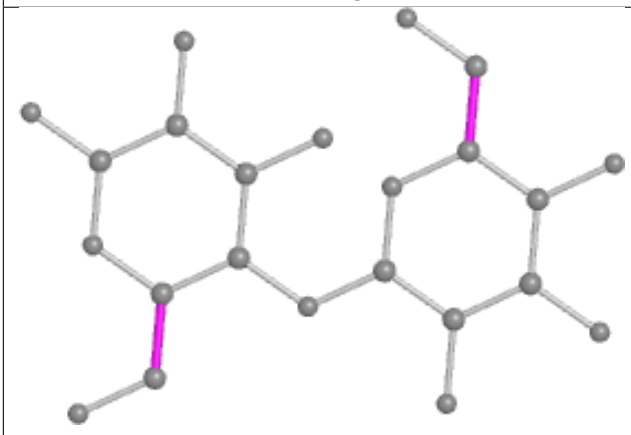
Oligosaccharide Chain F



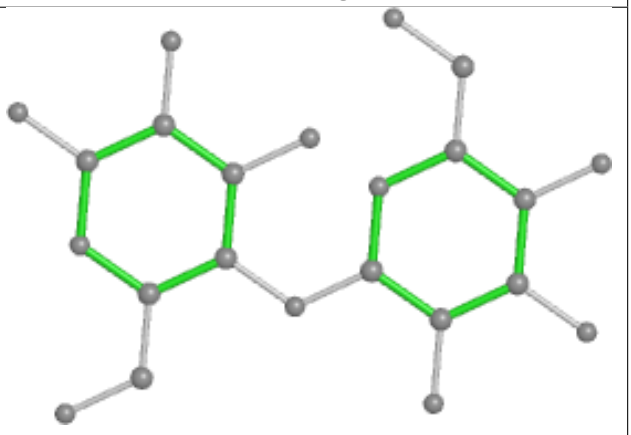
Bond lengths



Bond angles

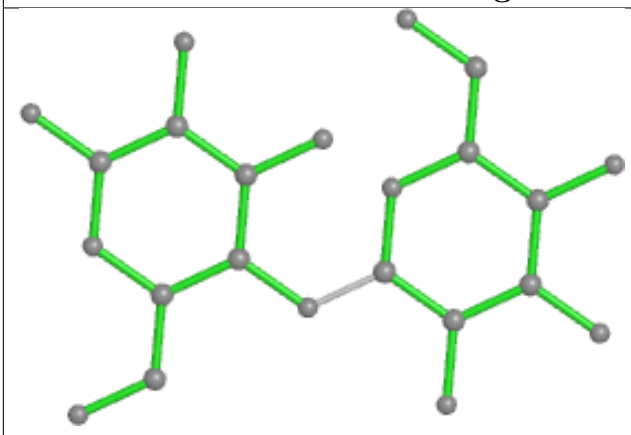


Torsions

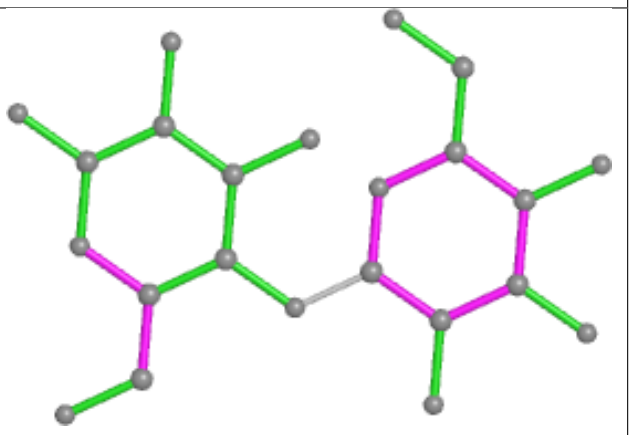


Rings

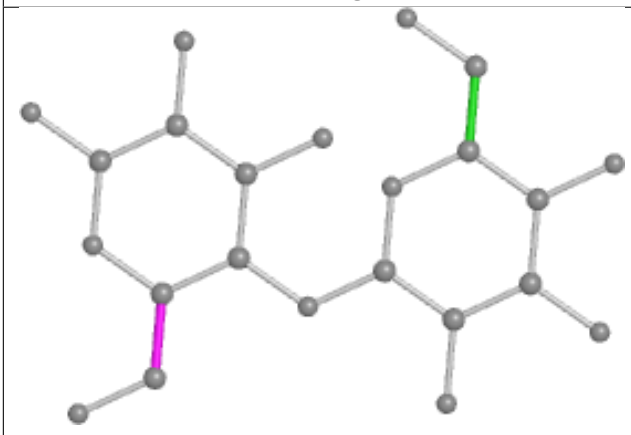
Oligosaccharide Chain G



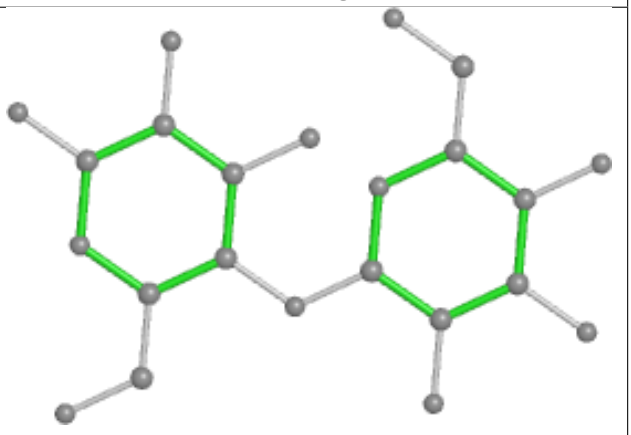
Bond lengths



Bond angles

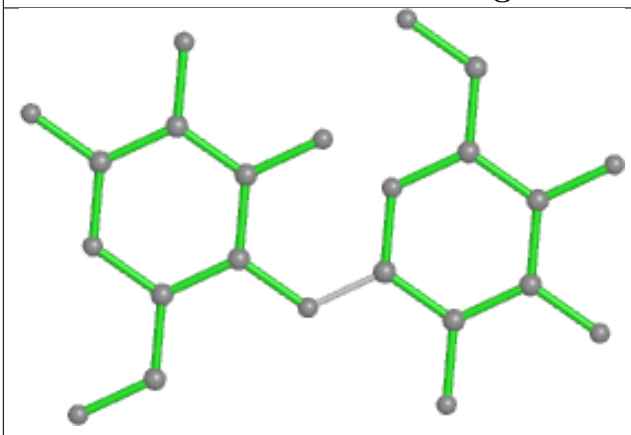


Torsions

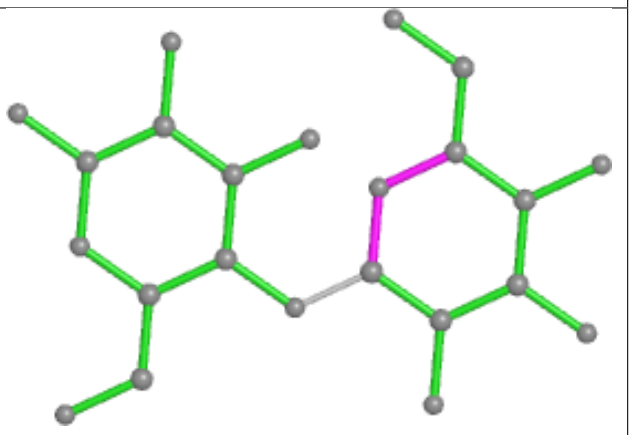


Rings

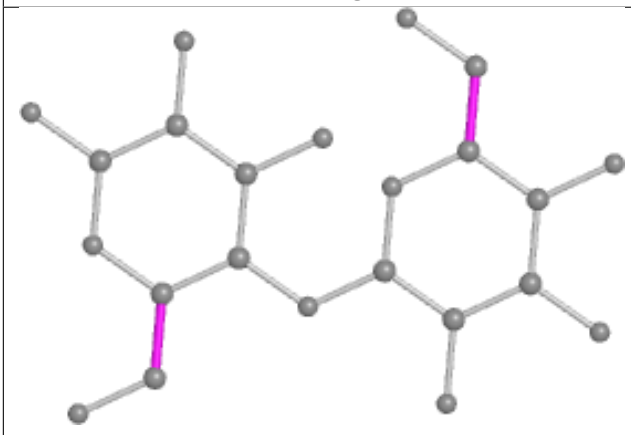
Oligosaccharide Chain H



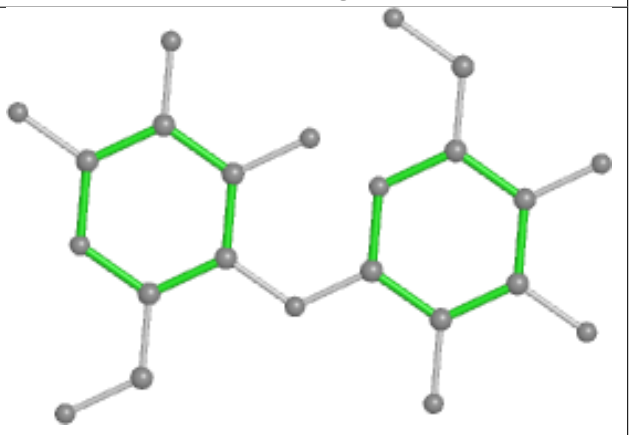
Bond lengths



Bond angles

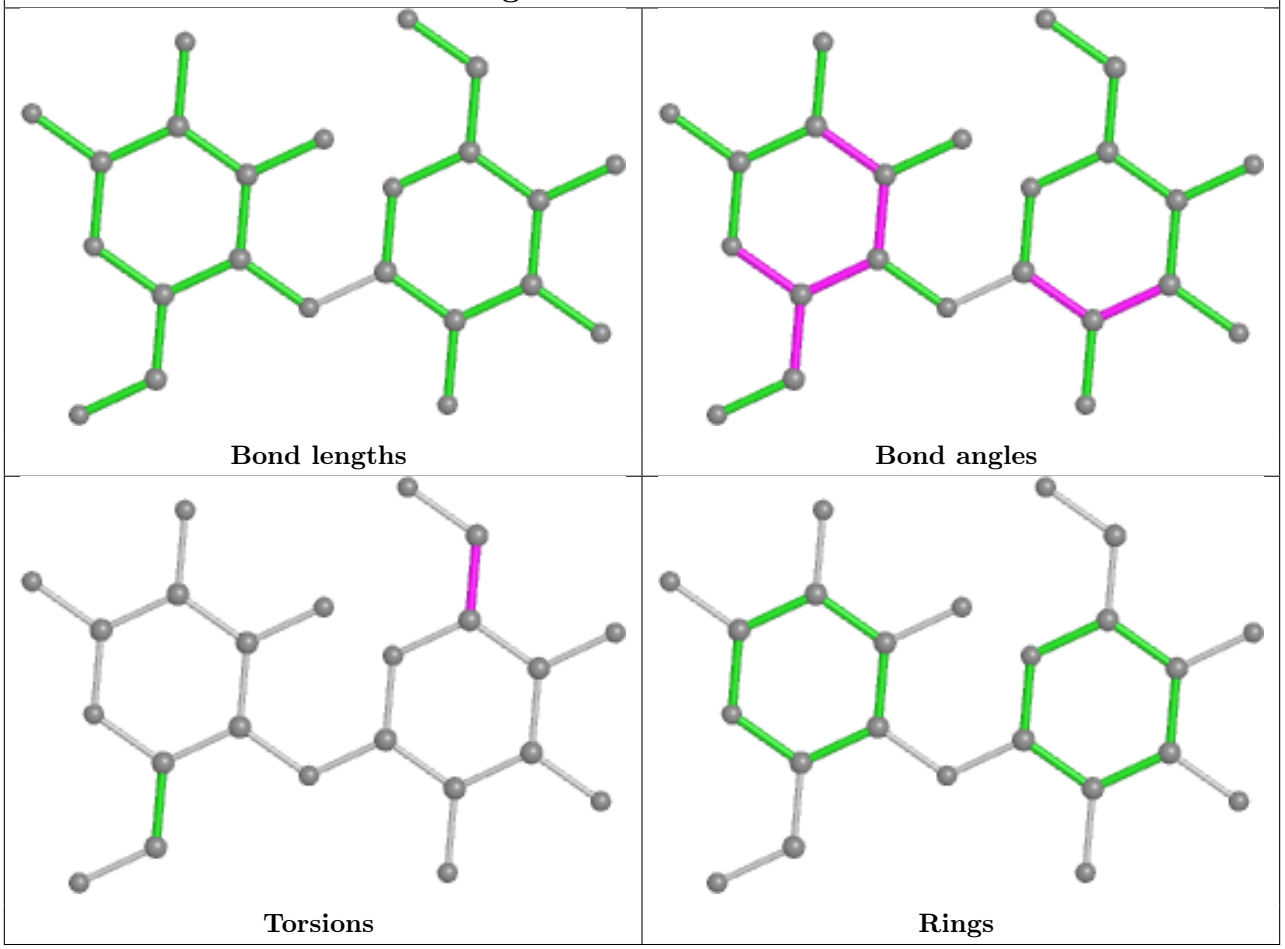


Torsions



Rings

Oligosaccharide Chain I



5.6 Ligand geometry (i)

Of 17 ligands modelled in this entry, 3 are monoatomic - leaving 14 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	NAG	C	402	1	14,14,15	0.54	0	17,19,21	0.68	0
4	NAG	B	401	1	14,14,15	0.43	0	17,19,21	0.38	0
4	NAG	C	401	1	14,14,15	0.34	0	17,19,21	0.70	1 (5%)
3	ATP	B	407	6	26,33,33	1.04	1 (3%)	31,52,52	1.27	3 (9%)
5	EDO	A	407	-	3,3,3	0.69	0	2,2,2	0.17	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
4	NAG	A	404	1	14,14,15	0.32	0	17,19,21	0.44	0
4	NAG	B	402	1	14,14,15	0.32	0	17,19,21	0.45	0
3	ATP	A	401	6	26,33,33	1.07	2 (7%)	31,52,52	1.55	9 (29%)
5	EDO	B	406	-	3,3,3	0.85	0	2,2,2	0.51	0
3	ATP	A	402	6	26,33,33	1.01	2 (7%)	31,52,52	1.49	5 (16%)
5	EDO	C	406	-	3,3,3	0.50	0	2,2,2	0.12	0
4	NAG	C	403	1	14,14,15	0.48	0	17,19,21	0.65	0
4	NAG	A	403	1	14,14,15	0.37	0	17,19,21	0.50	0
4	NAG	B	403	1	14,14,15	0.39	0	17,19,21	0.67	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	C	402	1	-	4/6/23/26	0/1/1/1
4	NAG	B	401	1	-	2/6/23/26	0/1/1/1
4	NAG	C	401	1	-	2/6/23/26	0/1/1/1
3	ATP	B	407	6	-	4/18/38/38	0/3/3/3
5	EDO	A	407	-	-	0/1/1/1	-
4	NAG	A	404	1	-	2/6/23/26	0/1/1/1
4	NAG	B	402	1	-	2/6/23/26	0/1/1/1
3	ATP	A	401	6	-	2/18/38/38	0/3/3/3
5	EDO	B	406	-	-	1/1/1/1	-
3	ATP	A	402	6	-	1/18/38/38	0/3/3/3
5	EDO	C	406	-	-	1/1/1/1	-
4	NAG	C	403	1	-	1/6/23/26	0/1/1/1
4	NAG	A	403	1	-	2/6/23/26	0/1/1/1
4	NAG	B	403	1	-	2/6/23/26	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	407	ATP	C5-C4	2.62	1.47	1.40
3	A	401	ATP	C5-C4	2.41	1.47	1.40
3	A	402	ATP	C5-C4	2.36	1.47	1.40
3	A	401	ATP	C2'-C1'	-2.27	1.50	1.53
3	A	402	ATP	C2'-C1'	-2.12	1.50	1.53

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	401	ATP	O3G-PG-O2G	3.40	120.62	107.64
3	A	402	ATP	PA-O3A-PB	-3.35	121.33	132.83
3	A	402	ATP	C4-C5-N7	-3.25	106.01	109.40
3	B	407	ATP	N3-C2-N1	-3.08	123.86	128.68
3	A	401	ATP	C4-C5-N7	-3.03	106.24	109.40
3	A	401	ATP	N3-C2-N1	-2.97	124.04	128.68
3	B	407	ATP	C4-C5-N7	-2.92	106.35	109.40
3	A	401	ATP	C3'-C2'-C1'	2.68	105.01	100.98
4	C	401	NAG	C1-O5-C5	2.38	115.41	112.19
3	A	402	ATP	N3-C2-N1	-2.32	125.06	128.68
3	B	407	ATP	PB-O3B-PG	-2.22	125.21	132.83
3	A	401	ATP	C1'-N9-C4	-2.16	122.85	126.64
3	A	401	ATP	C2'-C3'-C4'	2.09	106.70	102.64
3	A	401	ATP	O2A-PA-O1A	2.04	122.34	112.24
3	A	401	ATP	C2-N1-C6	2.02	122.21	118.75
3	A	402	ATP	O3B-PG-O1G	-2.02	100.00	111.19
3	A	401	ATP	PA-O3A-PB	-2.01	125.92	132.83
3	A	402	ATP	PB-O3B-PG	-2.01	125.93	132.83

There are no chirality outliers.

All (26) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	401	ATP	PB-O3B-PG-O3G
4	C	402	NAG	C4-C5-C6-O6
4	C	402	NAG	O5-C5-C6-O6
4	B	402	NAG	O5-C5-C6-O6
4	A	403	NAG	O5-C5-C6-O6
4	B	401	NAG	O5-C5-C6-O6
4	A	403	NAG	C4-C5-C6-O6
4	B	402	NAG	C4-C5-C6-O6
4	C	401	NAG	O5-C5-C6-O6
4	B	401	NAG	C4-C5-C6-O6
4	C	401	NAG	C4-C5-C6-O6
4	A	404	NAG	O5-C5-C6-O6
5	C	406	EDO	O1-C1-C2-O2
4	C	402	NAG	C1-C2-N2-C7
5	B	406	EDO	O1-C1-C2-O2
4	A	404	NAG	C4-C5-C6-O6
3	B	407	ATP	C3'-C4'-C5'-O5'
3	B	407	ATP	PA-O3A-PB-O3B

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
4	C	403	NAG	C4-C5-C6-O6
4	B	403	NAG	O5-C5-C6-O6
3	A	401	ATP	PB-O3B-PG-O2G
3	B	407	ATP	O4'-C4'-C5'-O5'
3	B	407	ATP	PA-O3A-PB-O1B
3	A	402	ATP	PB-O3A-PA-O2A
4	C	402	NAG	C3-C2-N2-C7
4	B	403	NAG	C3-C2-N2-C7

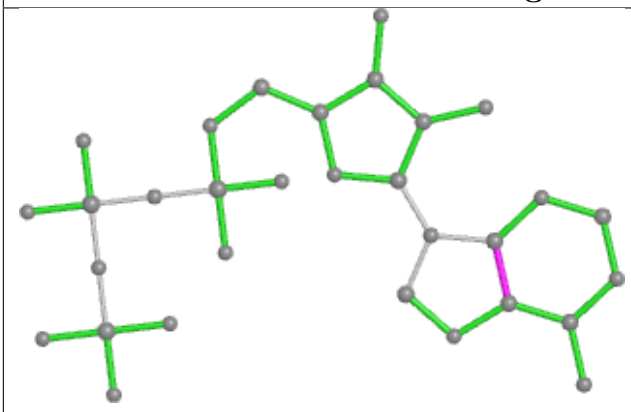
There are no ring outliers.

6 monomers are involved in 9 short contacts:

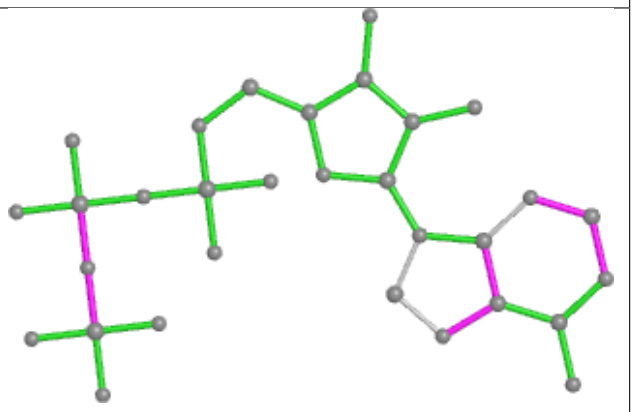
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	401	NAG	1	0
3	B	407	ATP	1	0
3	A	401	ATP	2	0
5	B	406	EDO	1	0
3	A	402	ATP	2	0
4	B	403	NAG	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

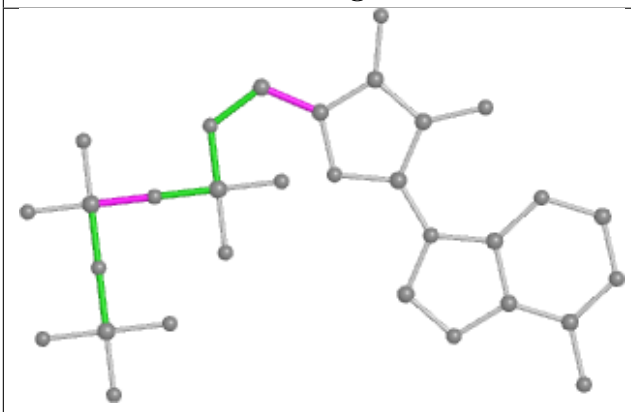
Ligand ATP B 407



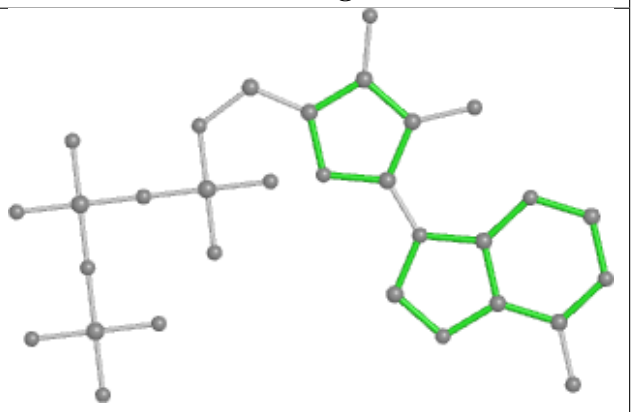
Bond lengths



Bond angles

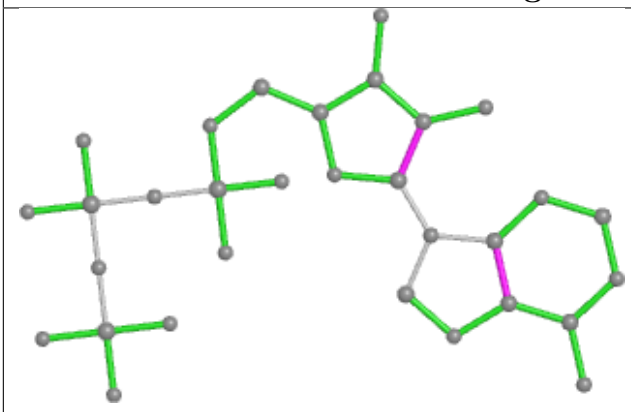


Torsions

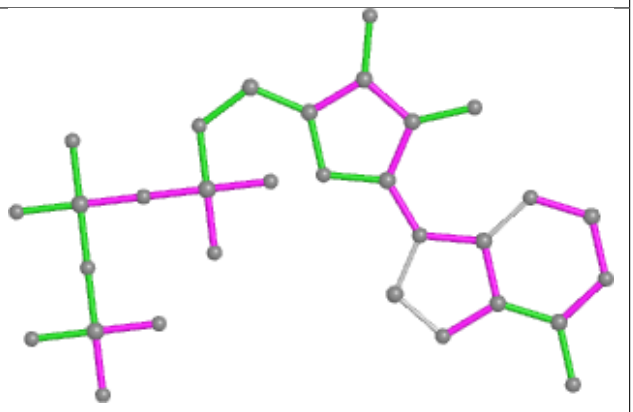


Rings

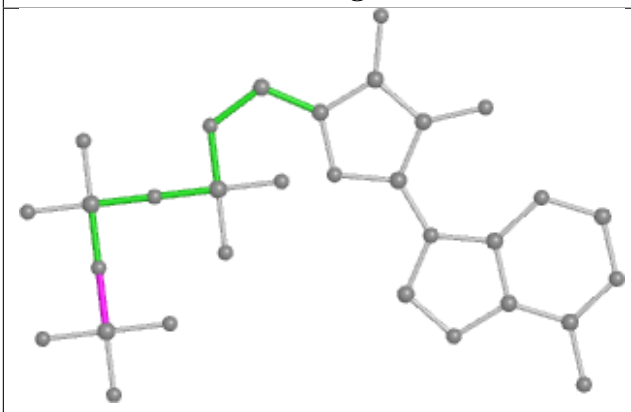
Ligand ATP A 401



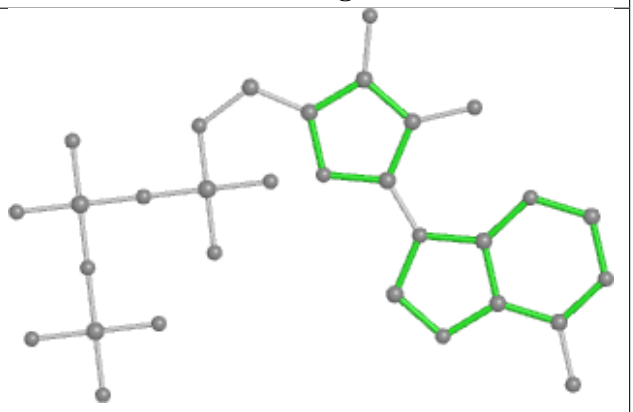
Bond lengths



Bond angles

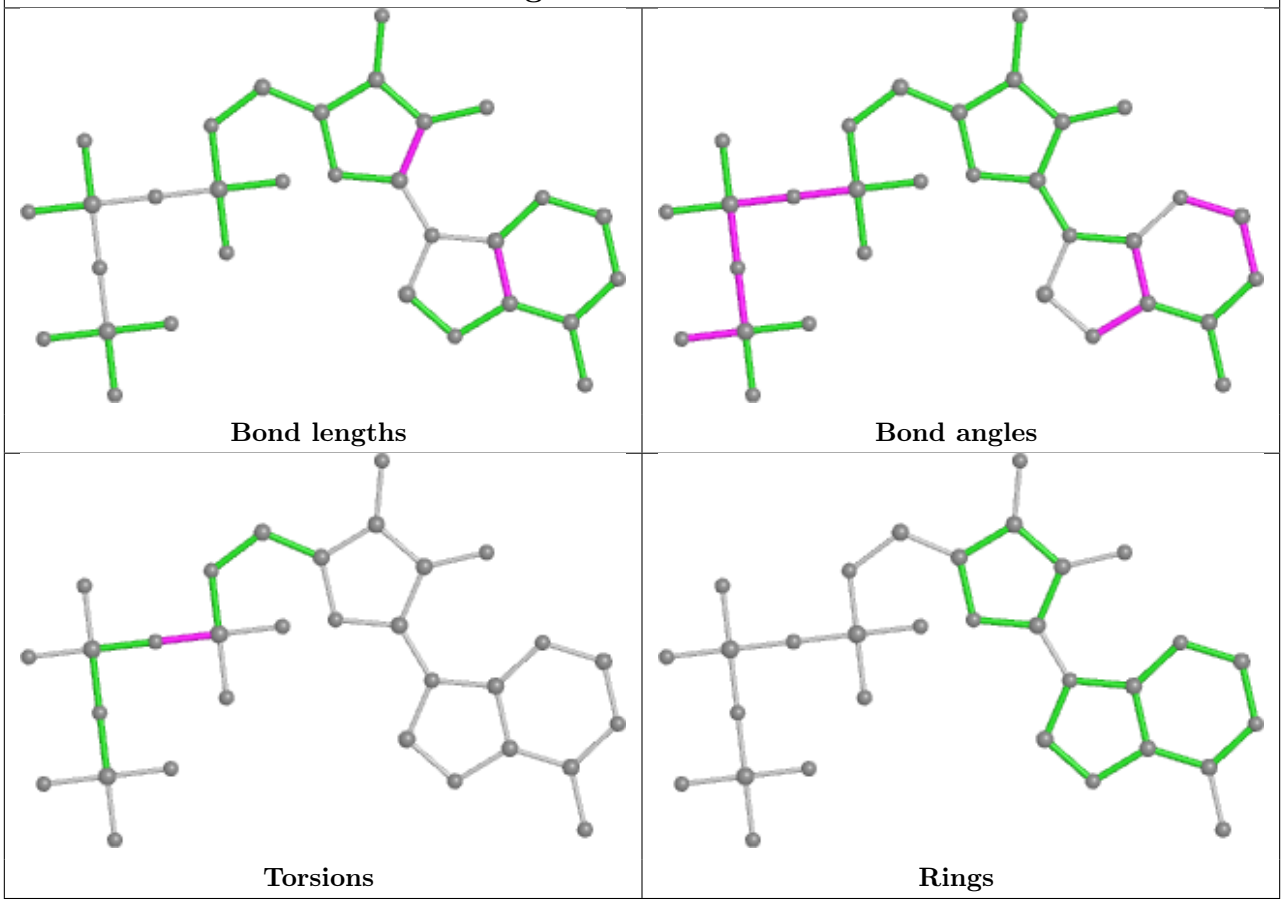


Torsions



Rings

Ligand ATP A 402



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, 95^{th} 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	357/362 (98%)	0.10	14 (3%) 39 37	101, 149, 203, 235	0
1	B	357/362 (98%)	0.45	52 (14%) 2 2	112, 159, 211, 241	0
1	C	357/362 (98%)	0.43	47 (13%) 3 3	105, 167, 222, 239	0
All	All	1071/1086 (98%)	0.33	113 (10%) 6 6	101, 161, 213, 241	0

All (113) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	260	TYR	5.7
1	C	331	SER	5.0
1	C	130	GLY	4.9
1	B	180	ARG	4.9
1	B	179	ILE	4.6
1	B	186	PHE	4.6
1	B	19	LYS	4.5
1	C	338	LEU	4.4
1	C	212	PHE	4.3
1	C	335	GLY	4.3
1	C	263	THR	4.2
1	C	126	ARG	4.2
1	B	126	ARG	4.2
1	B	209	LYS	4.2
1	C	209	LYS	4.1
1	C	339	CYS	4.1
1	B	330	THR	4.1
1	B	127	LEU	4.0
1	B	188	LYS	4.0
1	B	245	TRP	4.0
1	B	205	PHE	3.9
1	B	130	GLY	3.8
1	C	115	ARG	3.8

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	161	GLU	3.7
1	C	124	PRO	3.7
1	C	25	ILE	3.6
1	B	204	ARG	3.6
1	C	82	THR	3.6
1	C	87	THR	3.5
1	B	129	GLY	3.5
1	B	338	LEU	3.4
1	C	127	LEU	3.3
1	B	212	PHE	3.3
1	C	211	PRO	3.3
1	A	346	PHE	3.2
1	C	117	VAL	3.2
1	C	142	SER	3.2
1	B	7	PHE	3.1
1	B	331	SER	3.1
1	B	187	GLU	3.1
1	A	260	TYR	3.1
1	C	123	GLY	3.1
1	C	131	GLY	3.0
1	A	324	SER	3.0
1	B	257	ILE	3.0
1	B	189	GLY	3.0
1	C	88	SER	3.0
1	C	27	ASN	2.9
1	A	332	VAL	2.9
1	C	121	GLN	2.9
1	B	210	ASP	2.8
1	C	166	MET	2.7
1	B	118	SER	2.7
1	C	141	SER	2.7
1	B	211	PRO	2.7
1	B	117	VAL	2.7
1	A	330	THR	2.7
1	B	262	PHE	2.7
1	A	322	ILE	2.7
1	C	19	LYS	2.7
1	C	83	PRO	2.7
1	B	207	PRO	2.7
1	C	334	VAL	2.6
1	B	8	PHE	2.6
1	B	84	PRO	2.6

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	329	PHE	2.6
1	B	259	LYS	2.6
1	C	262	PHE	2.5
1	A	329	PHE	2.5
1	C	194	ASN	2.5
1	C	125	GLU	2.5
1	C	332	VAL	2.5
1	A	343	LEU	2.5
1	C	210	ASP	2.4
1	B	324	SER	2.4
1	C	54	THR	2.4
1	A	10	TYR	2.4
1	C	28	ARG	2.4
1	B	20	SER	2.4
1	B	10	TYR	2.4
1	B	9	THR	2.4
1	C	31	GLN	2.4
1	C	346	PHE	2.4
1	A	285	TYR	2.4
1	C	58	SER	2.3
1	B	162	THR	2.3
1	B	258	PRO	2.3
1	A	23	ILE	2.3
1	C	362	ASN	2.3
1	A	310	TYR	2.3
1	C	20	SER	2.2
1	B	241	ILE	2.2
1	B	124	PRO	2.2
1	C	89	VAL	2.2
1	B	208	ASP	2.2
1	B	54	THR	2.2
1	C	261	SER	2.2
1	C	329	PHE	2.2
1	B	185	ASN	2.2
1	C	361	VAL	2.1
1	A	245	TRP	2.1
1	B	181	PHE	2.1
1	B	359	GLU	2.1
1	C	86	GLY	2.1
1	C	347	LEU	2.1
1	B	206	HIS	2.1
1	B	58	SER	2.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	322	ILE	2.1
1	B	326	VAL	2.1
1	C	23	ILE	2.0
1	A	311	GLY	2.0
1	B	299	LYS	2.0
1	B	243	ILE	2.0

6.2 Non-standard residues in protein, DNA, RNA chains (i)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates (i)

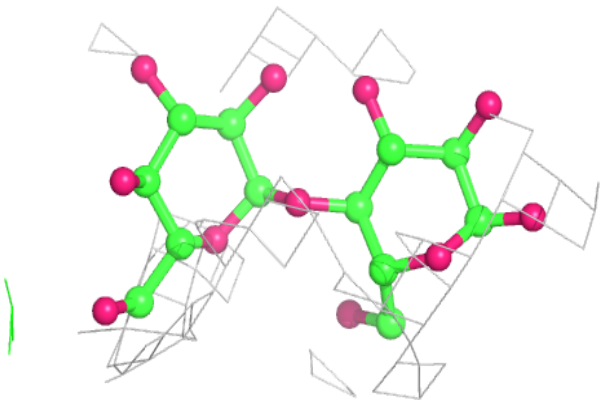
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

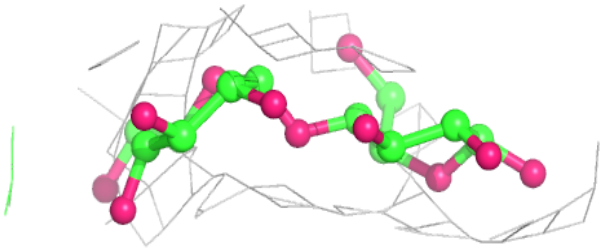
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	GLC	H	2	11/12	0.58	0.26	188,216,224,231	0
2	GLC	D	2	11/12	0.63	0.20	194,211,223,224	0
2	GLC	E	2	11/12	0.69	0.12	188,225,232,232	0
2	GLC	F	1	12/12	0.69	0.36	203,219,235,237	0
2	GLC	F	2	11/12	0.72	0.22	191,210,230,230	0
2	GLC	E	1	12/12	0.73	0.16	182,204,217,241	0
2	GLC	I	1	12/12	0.76	0.13	199,219,225,227	0
2	GLC	G	1	12/12	0.77	0.11	191,204,217,219	0
2	GLC	I	2	11/12	0.79	0.12	193,211,218,220	0
2	GLC	H	1	12/12	0.80	0.26	188,198,212,215	0
2	GLC	D	1	12/12	0.87	0.11	193,207,217,222	0
2	GLC	G	2	11/12	0.96	0.08	188,197,212,212	0

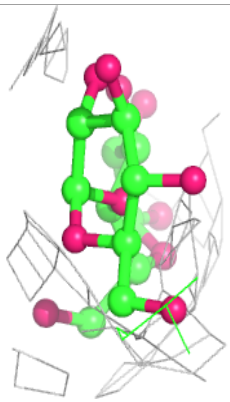
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain D:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

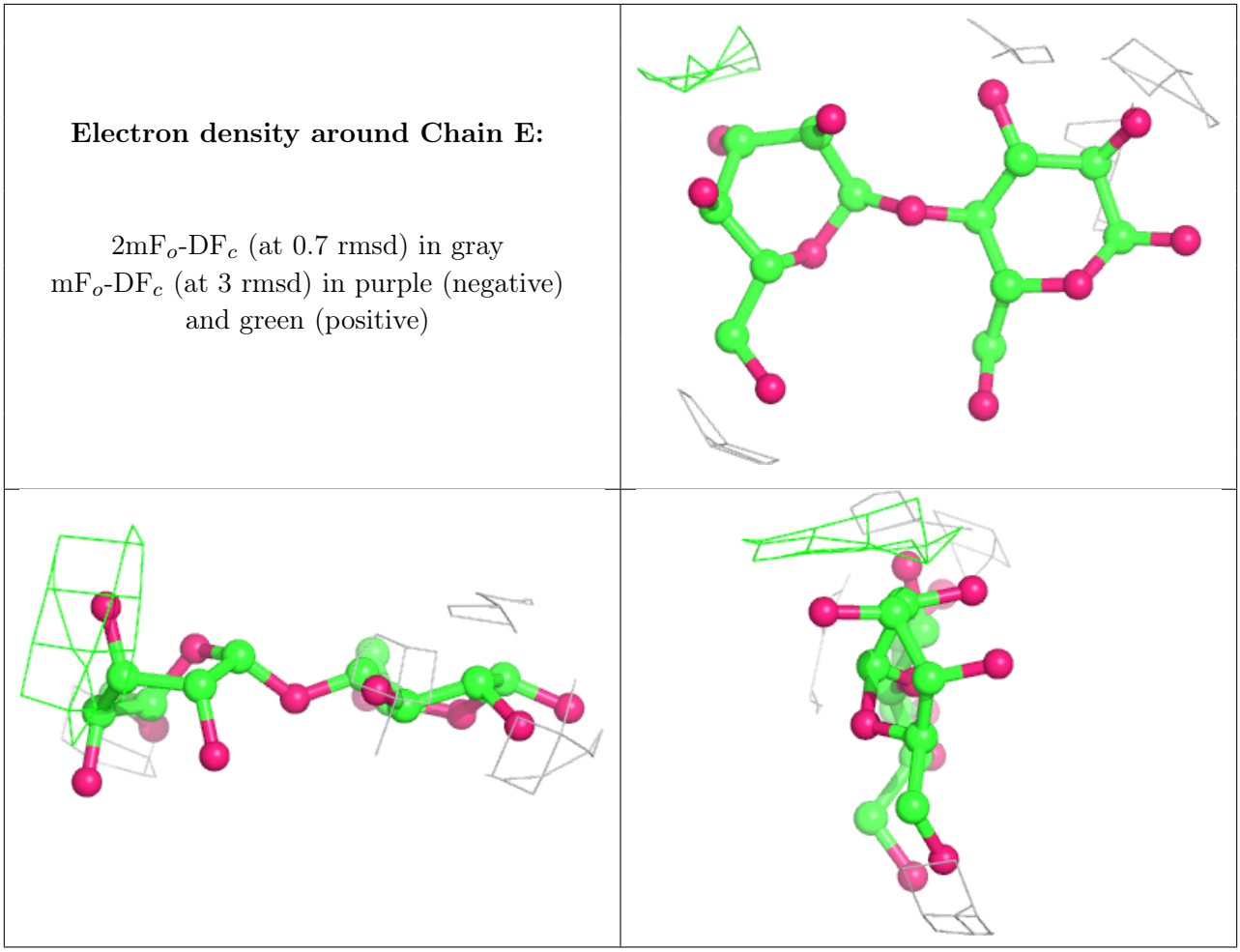






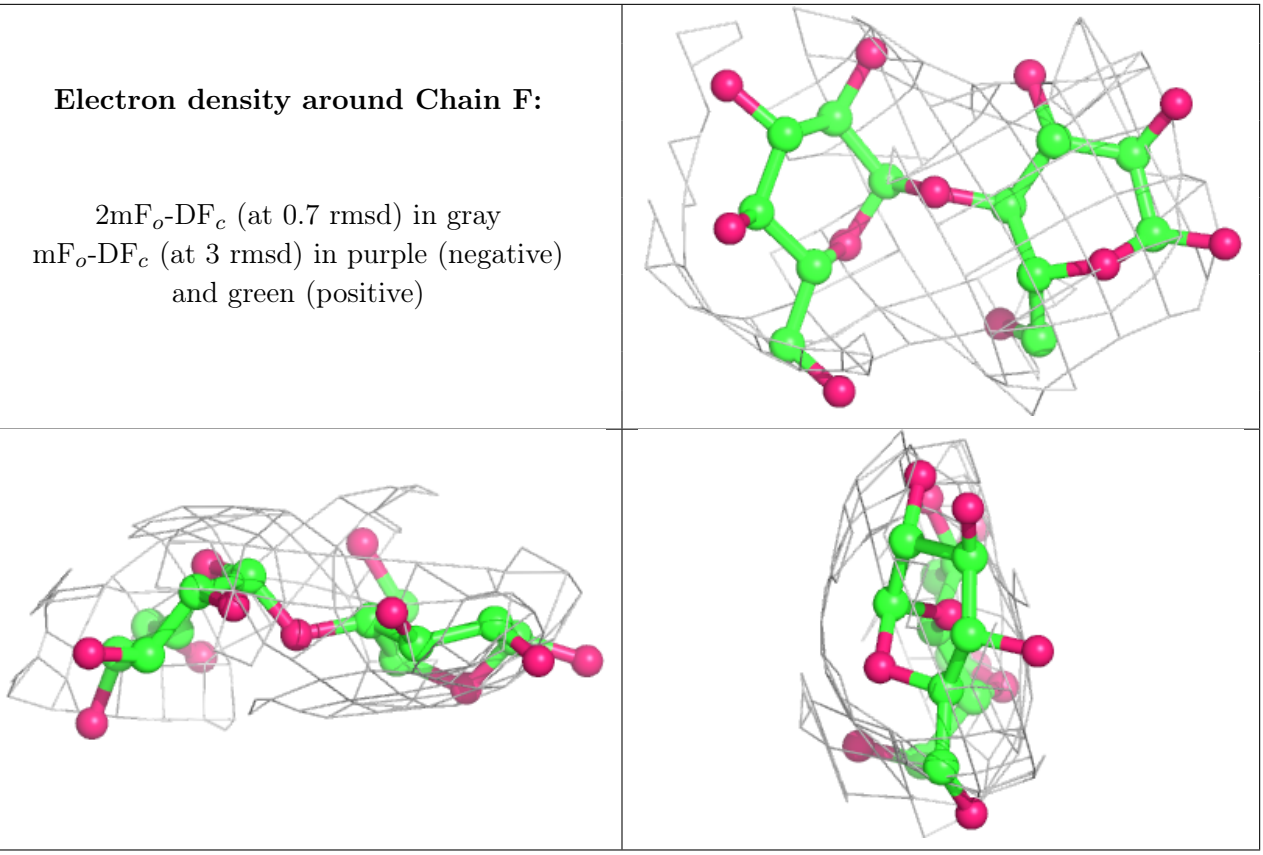
Electron density around Chain E:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



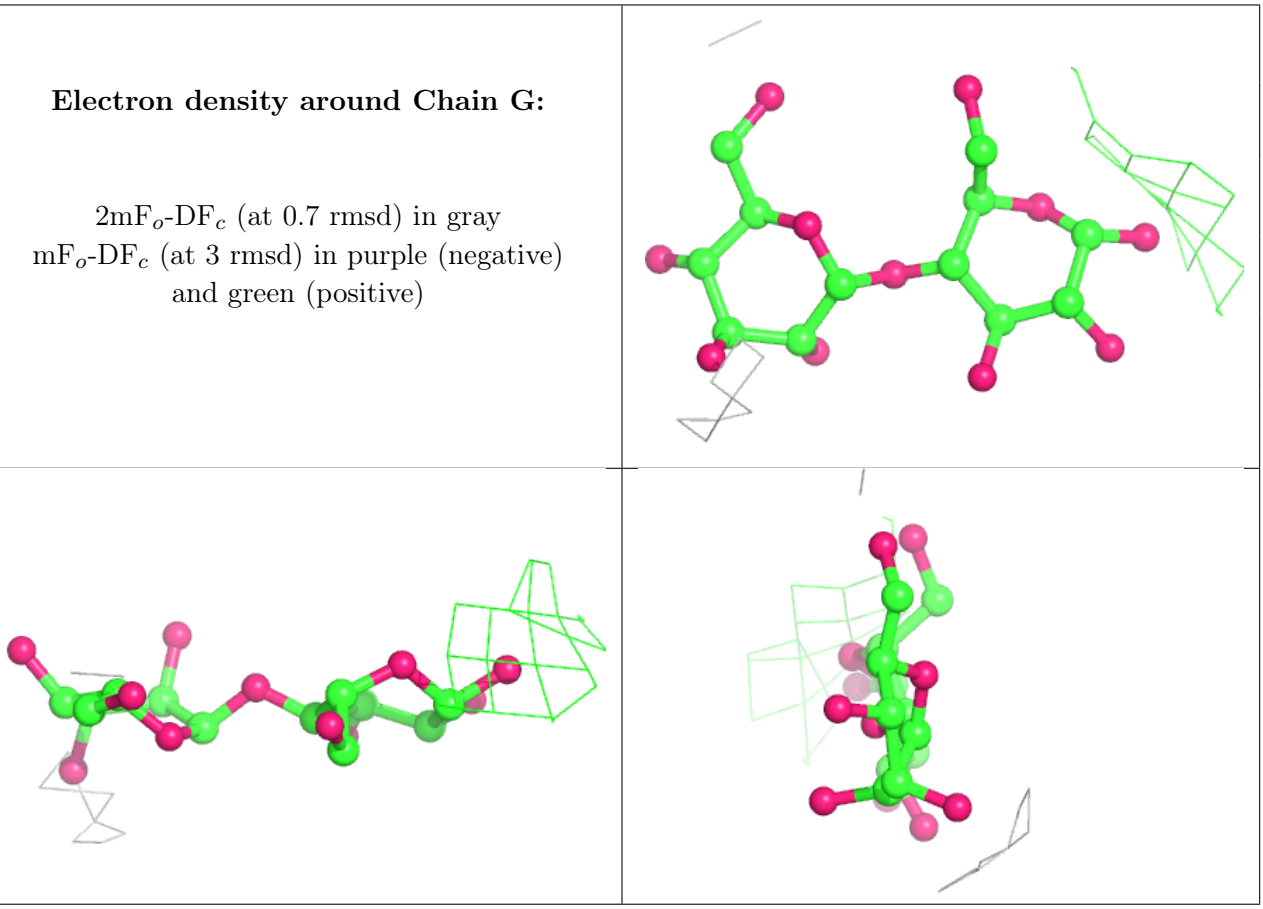
Electron density around Chain F:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



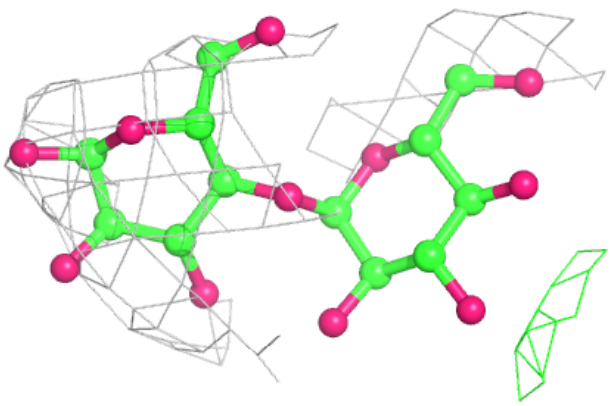
Electron density around Chain G:

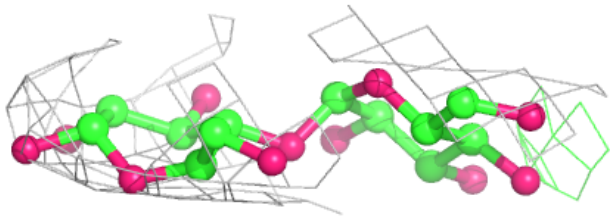
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

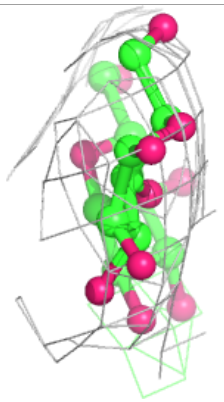


Electron density around Chain H:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

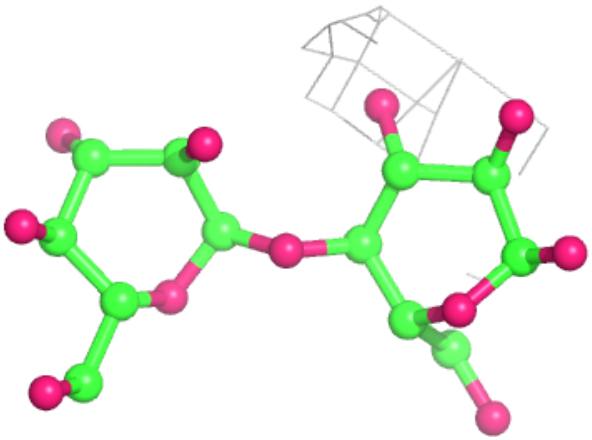


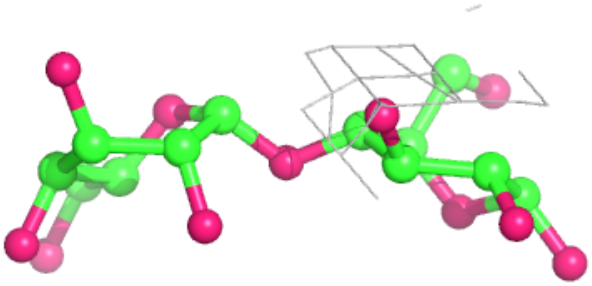


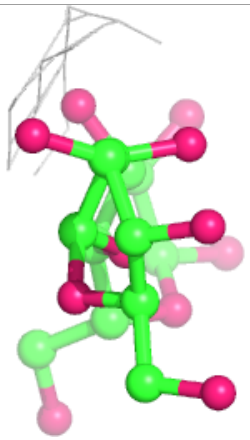


Electron density around Chain I:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)







6.4 Ligands (i)

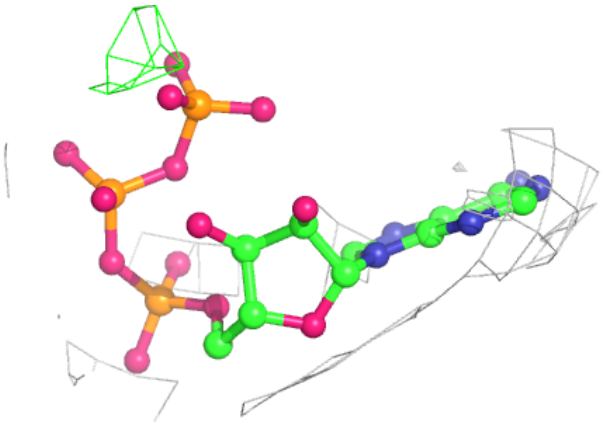
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

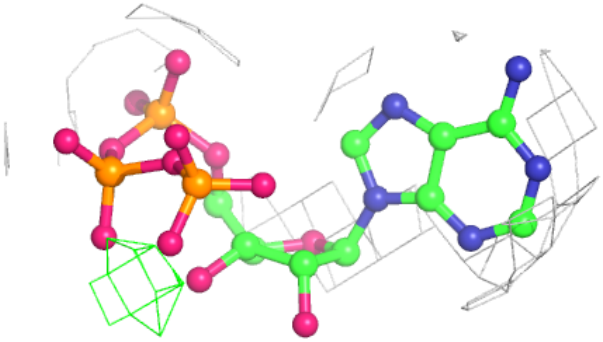
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
5	EDO	C	406	4/4	0.40	0.45	114,118,125,131	0
5	EDO	A	407	4/4	0.55	0.49	103,130,130,134	0
5	EDO	B	406	4/4	0.60	0.39	119,137,147,150	0
4	NAG	C	401	14/15	0.87	0.38	159,194,202,206	0
4	NAG	C	402	14/15	0.89	0.96	182,207,217,223	0
4	NAG	B	402	14/15	0.89	0.24	176,213,221,222	0
4	NAG	C	403	14/15	0.89	0.15	147,169,188,209	0
4	NAG	B	401	14/15	0.90	0.32	169,196,214,215	0
3	ATP	B	407	31/31	0.91	0.24	138,148,182,187	0
4	NAG	A	403	14/15	0.93	0.29	150,174,189,196	0
6	CA	C	407	1/1	0.94	0.46	192,192,192,192	0
4	NAG	A	404	14/15	0.94	0.29	112,125,150,164	0
3	ATP	A	402	31/31	0.94	0.23	111,124,143,146	0
4	NAG	B	403	14/15	0.95	0.23	145,162,197,202	0
3	ATP	A	401	31/31	0.96	0.17	98,114,138,189	0
6	CA	B	408	1/1	0.97	0.24	124,124,124,124	0
6	CA	A	408	1/1	0.99	0.39	153,153,153,153	0

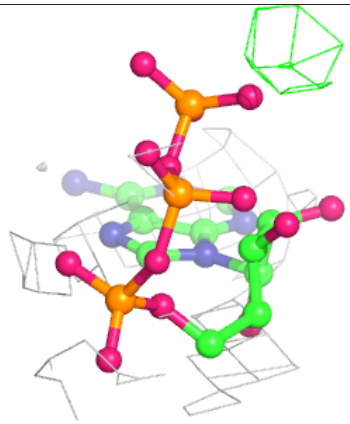
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around ATP B 407:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

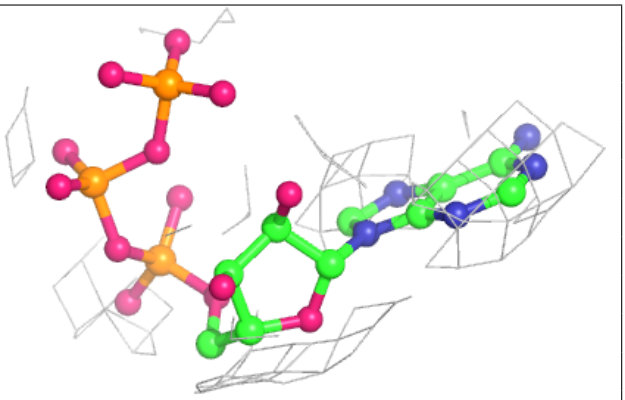


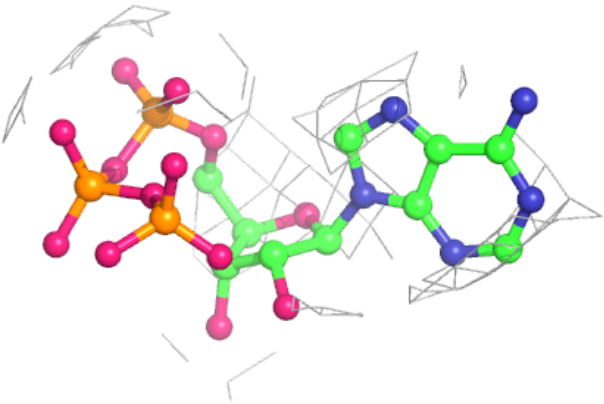


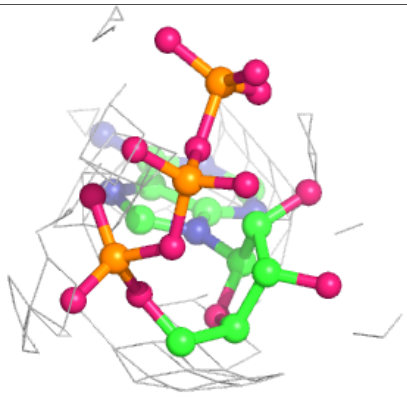


Electron density around ATP A 402:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

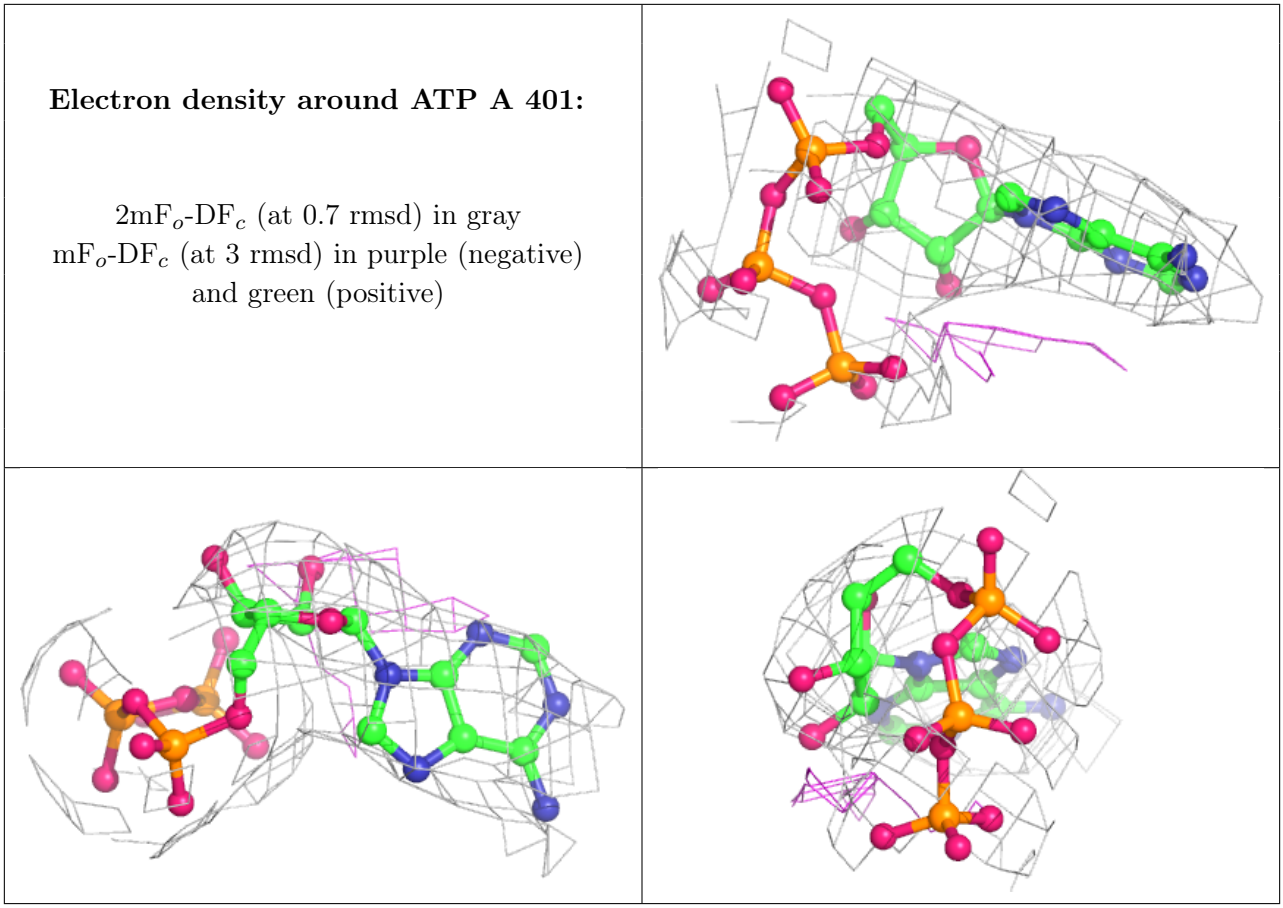






Electron density around ATP A 401:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



6.5 Other polymers

There are no such residues in this entry.