Full wwPDB X-ray Structure Validation Report (i)

Sep 23, 2023 – 08:34 PM EDT

PDB ID	:	5TGV
Title	:	Crystal structure of H10 hemagglutinin mutant (K158aA-D193T-Q226L-G228S) from Jiangxi-Donghu (2013) H10N8 influenza virus in complex with 3'-SLN
Authors	:	Tzarum, N.; Wilson, I.A.
Deposited on	:	2016-09-28
Resolution	:	2.97 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 types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

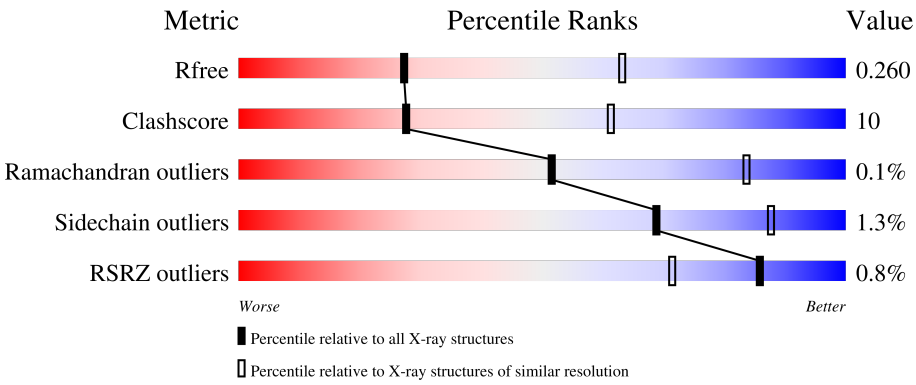
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

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 2.97 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	2754 (3.00-2.96)
Clashscore	141614	3103 (3.00-2.96)
Ramachandran outliers	138981	2993 (3.00-2.96)
Sidechain outliers	138945	2996 (3.00-2.96)
RSRZ outliers	127900	2644 (3.00-2.96)

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	323	78% 20% •
1	C	323	2% 73% 25% •
1	E	323	% 74% 24% • •
2	B	180	% 76% 18% • 5%
2	D	180	% 67% 25% • 7%

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Mol	Chain	Length	Quality of chain
2	F	180	% 76% 19% 5%
3	G	2	100%
4	H	3	100%
5	I	2	50% 50%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	NAG	G	2	-	-	-	X

2 Entry composition (i)

There are 7 unique types of molecules in this entry. The entry contains 11626 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 Hemagglutinin HA1 chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	318	2433	1505	447	464	17	0	0	0
1	C	319	2441	1511	448	465	17	0	0	0
1	E	318	2433	1505	447	464	17	0	0	0

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	158A	ALA	LYS	engineered mutation	UNP A0A0J9X252
A	193	THR	ASP	engineered mutation	UNP A0A0J9X252
A	226	LEU	GLN	engineered mutation	UNP A0A0J9X252
A	228	SER	GLY	engineered mutation	UNP A0A0J9X252
C	158A	ALA	LYS	engineered mutation	UNP A0A0J9X252
C	193	THR	ASP	engineered mutation	UNP A0A0J9X252
C	226	LEU	GLN	engineered mutation	UNP A0A0J9X252
C	228	SER	GLY	engineered mutation	UNP A0A0J9X252
E	158A	ALA	LYS	engineered mutation	UNP A0A0J9X252
E	193	THR	ASP	engineered mutation	UNP A0A0J9X252
E	226	LEU	GLN	engineered mutation	UNP A0A0J9X252
E	228	SER	GLY	engineered mutation	UNP A0A0J9X252

- Molecule 2 is a protein called Hemagglutinin HA2 chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	171	1382	854	239	281	8	0	0	0
2	D	168	1358	840	234	276	8	0	0	0
2	F	171	1382	854	239	281	8	0	0	0

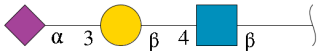
There are 21 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	175	SER	-	expression tag	UNP A0A0J9X253
B	176	GLY	-	expression tag	UNP A0A0J9X253
B	177	ARG	-	expression tag	UNP A0A0J9X253
B	178	LEU	-	expression tag	UNP A0A0J9X253
B	179	VAL	-	expression tag	UNP A0A0J9X253
B	180	PRO	-	expression tag	UNP A0A0J9X253
B	181	ARG	-	expression tag	UNP A0A0J9X253
D	175	SER	-	expression tag	UNP A0A0J9X253
D	176	GLY	-	expression tag	UNP A0A0J9X253
D	177	ARG	-	expression tag	UNP A0A0J9X253
D	178	LEU	-	expression tag	UNP A0A0J9X253
D	179	VAL	-	expression tag	UNP A0A0J9X253
D	180	PRO	-	expression tag	UNP A0A0J9X253
D	181	ARG	-	expression tag	UNP A0A0J9X253
F	175	SER	-	expression tag	UNP A0A0J9X253
F	176	GLY	-	expression tag	UNP A0A0J9X253
F	177	ARG	-	expression tag	UNP A0A0J9X253
F	178	LEU	-	expression tag	UNP A0A0J9X253
F	179	VAL	-	expression tag	UNP A0A0J9X253
F	180	PRO	-	expression tag	UNP A0A0J9X253
F	181	ARG	-	expression tag	UNP A0A0J9X253

- Molecule 3 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
3	G	2	28	16	2	10	0	0	0

- Molecule 4 is an oligosaccharide called N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.

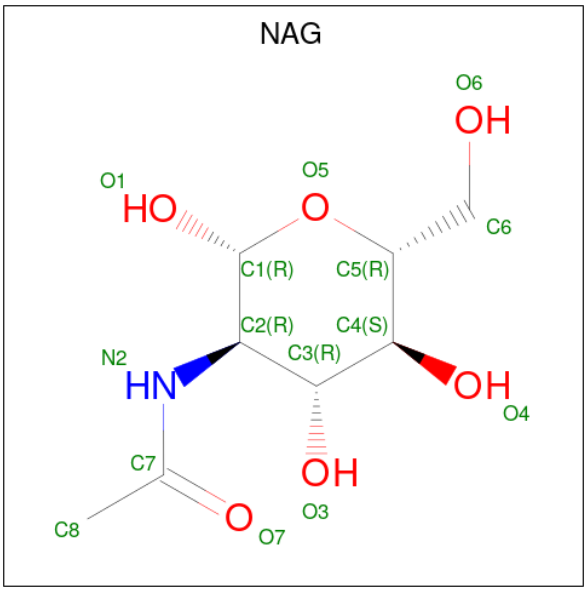


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
4	H	3	46	25	2	19	0	0	0

- Molecule 5 is an oligosaccharide called beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.

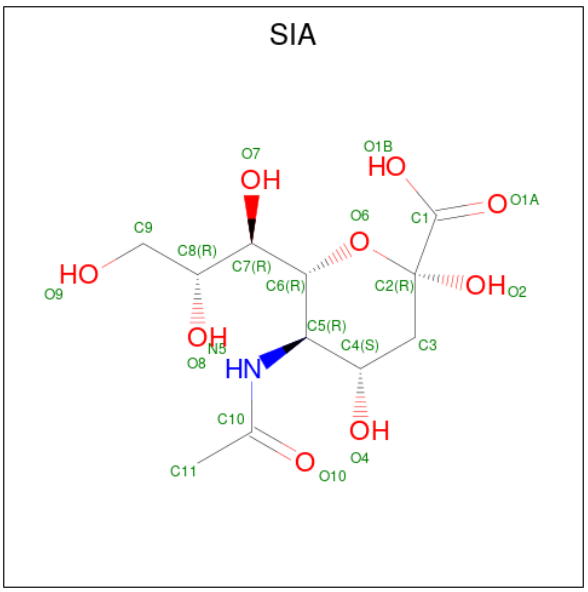
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
5	I	2	25	14	1	10	0	0	0

- Molecule 6 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



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

- Molecule 7 is N-acetyl-alpha-neuraminic acid (three-letter code: SIA) (formula: $C_{11}H_{19}NO_9$).

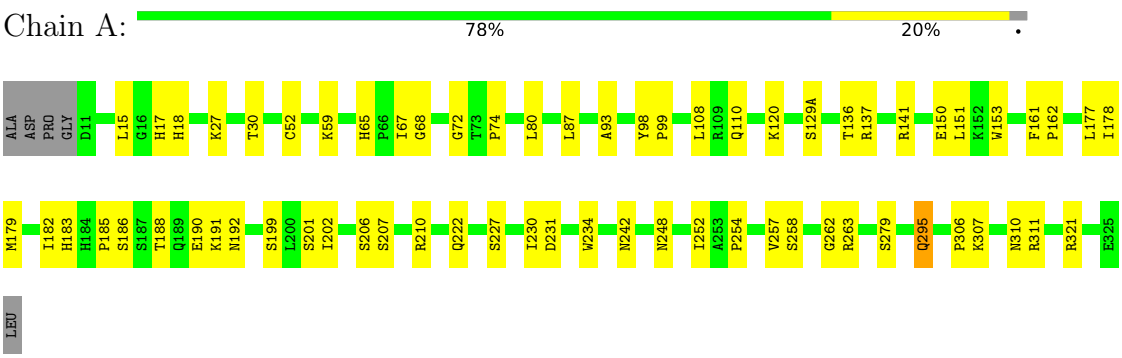


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
7	C	1	21	11	1	9	0	0
			Total	C	N	O		
7	E	1	21	11	1	9	0	0

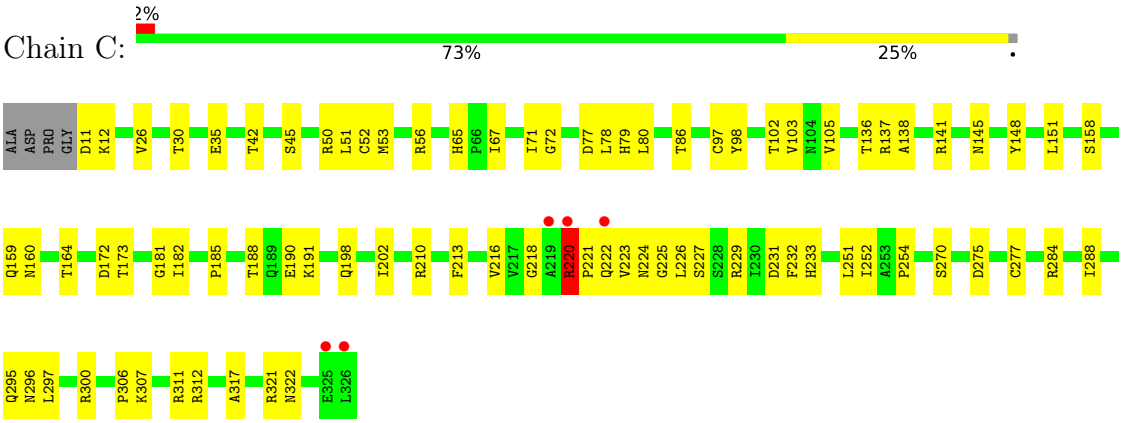
3 Residue-property plots (i)

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.

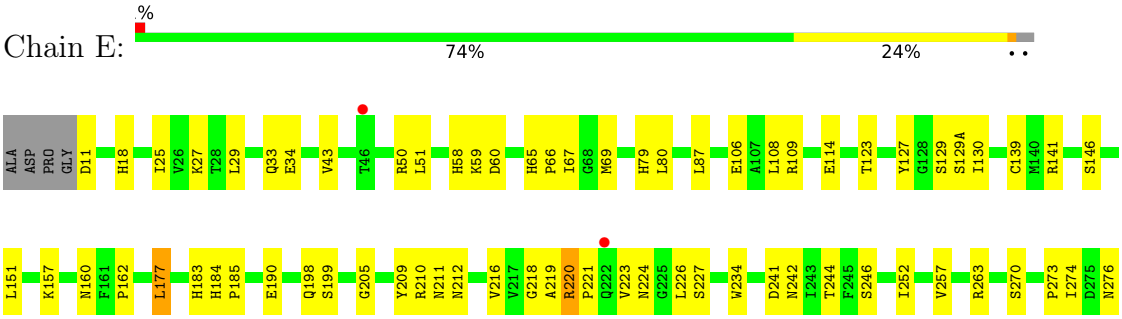
• Molecule 1: Hemagglutinin HA1 chain

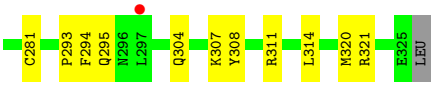


• Molecule 1: Hemagglutinin HA1 chain



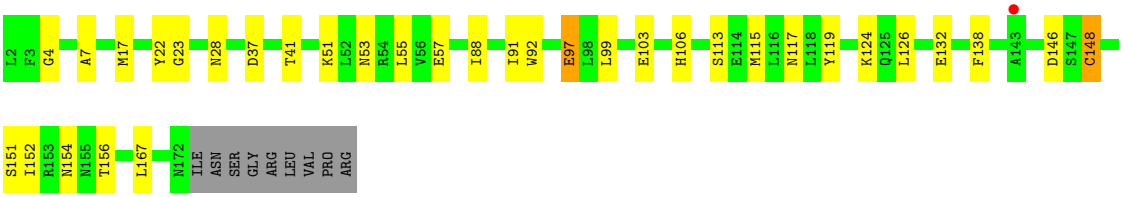
• Molecule 1: Hemagglutinin HA1 chain





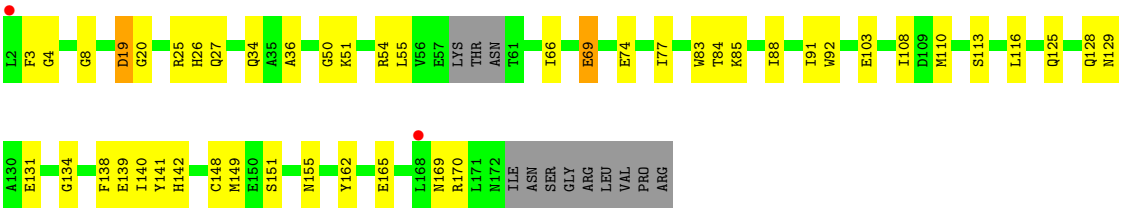
• Molecule 2: Hemagglutinin HA2 chain

Chain B:



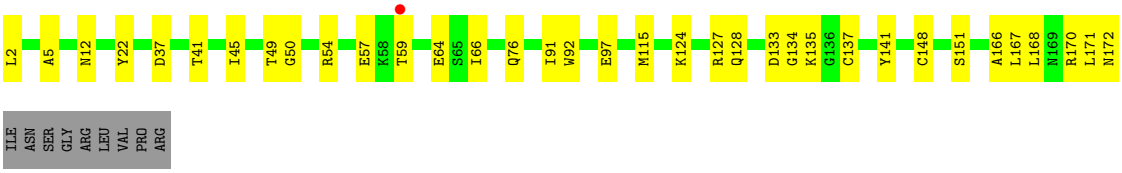
• Molecule 2: Hemagglutinin HA2 chain

Chain D:



• Molecule 2: Hemagglutinin HA2 chain

Chain F:

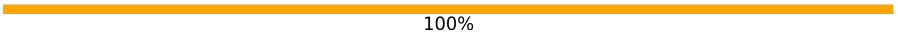


• Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain G:

• Molecule 4: N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain H:



• Molecule 5: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain I:

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	63.51Å 253.17Å 69.96Å 90.00° 111.97° 90.00°	Depositor
Resolution (Å)	35.72 – 2.97 48.30 – 2.97	Depositor EDS
% Data completeness (in resolution range)	92.2 (35.72-2.97) 86.0 (48.30-2.97)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.18	Depositor
$< I/\sigma(I) >$ [1]	1.42 (at 2.96Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155: ???)	Depositor
R, R_{free}	0.187 , 0.260 0.187 , 0.260	Depositor DCC
R_{free} test set	2008 reflections (5.19%)	wwPDB-VP
Wilson B-factor (Å^2)	46.2	Xtriage
Anisotropy	0.174	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.30 , 26.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.91	EDS
Total number of atoms	11626	wwPDB-VP
Average B, all atoms (Å^2)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.58% 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 (i)

5.1 Standard geometry (i)

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

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	0.62	0/2482	0.77	0/3365
1	C	0.58	0/2490	0.74	1/3376 (0.0%)
1	E	0.63	0/2482	0.77	2/3365 (0.1%)
2	B	0.59	0/1407	0.71	0/1899
2	D	0.59	0/1382	0.67	0/1864
2	F	0.62	0/1407	0.71	0/1899
All	All	0.61	0/11650	0.74	3/15768 (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	E	0	1
2	B	0	1
All	All	0	2

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	220	ARG	C-N-CD	6.47	141.98	128.40
1	E	177	LEU	CB-CG-CD1	-6.46	100.02	111.00
1	E	80	LEU	CA-CB-CG	-5.49	102.67	115.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	4	GLY	Peptide
1	E	11	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	2433	0	2385	52	0
1	C	2441	0	2396	59	0
1	E	2433	0	2386	54	1
2	B	1382	0	1279	30	0
2	D	1358	0	1252	37	0
2	F	1382	0	1279	27	0
3	G	28	0	25	4	0
4	H	46	0	40	8	1
5	I	25	0	22	0	0
6	A	14	0	13	0	0
6	C	28	0	25	0	0
6	E	14	0	13	1	0
7	C	21	0	18	2	0
7	E	21	0	18	4	0
All	All	11626	0	11151	227	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 10.

All (227) 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)
3:G:1:NAG:H62	3:G:2:NAG:HN2	0.99	1.07
1:C:220:ARG:HB3	1:C:221:PRO:HD3	1.49	0.94
3:G:1:NAG:H62	3:G:2:NAG:N2	1.85	0.92
4:H:2:GAL:H4	4:H:3:SIA:C1	1.99	0.92
1:C:190:GLU:OE1	7:C:405:SIA:O9	1.88	0.91
3:G:1:NAG:C6	3:G:2:NAG:HN2	1.83	0.91
1:A:190:GLU:OE2	4:H:3:SIA:O9	1.96	0.84
1:A:210:ARG:NH2	1:E:216:VAL:O	2.16	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:220:ARG:HB3	1:C:221:PRO:CD	2.12	0.79
2:B:53:ASN:O	2:B:57:GLU:HG2	1.84	0.78
2:D:25:ARG:NE	2:D:34:GLN:OE1	2.20	0.73
2:B:51:LYS:HZ1	2:B:106:HIS:HB3	1.55	0.72
1:A:129(A):SER:HB2	1:A:162:PRO:HG3	1.73	0.71
1:C:221:PRO:HB2	1:C:223:VAL:HG12	1.73	0.70
1:C:224:ASN:HB2	1:E:244:THR:OG1	1.92	0.70
1:C:284:ARG:HD2	2:D:69:GLU:HG2	1.75	0.69
2:F:22:TYR:H	2:F:41:THR:HG22	1.59	0.68
2:F:128:GLN:O	2:F:170:ARG:NH1	2.26	0.68
1:E:190:GLU:CD	7:E:602:SIA:HO9	1.97	0.68
1:E:50:ARG:HG2	1:E:273:PRO:HG2	1.77	0.67
1:E:190:GLU:CD	7:E:602:SIA:O9	2.33	0.67
1:C:137:ARG:HD2	1:C:145:ASN:OD1	1.95	0.66
1:A:186:SER:OG	1:A:190:GLU:OE1	2.11	0.66
1:E:25:ILE:HG23	1:E:33:GLN:HA	1.78	0.66
1:C:65:HIS:CE1	1:C:67:ILE:HG12	2.31	0.66
1:A:179:MET:HG2	1:A:234:TRP:HB3	1.78	0.65
1:A:310:ASN:O	1:A:311:ARG:HD3	1.97	0.65
2:B:55:LEU:HD22	2:B:99:LEU:HD21	1.77	0.65
2:D:125:GLN:OE1	2:D:155:ASN:HA	1.97	0.65
1:A:161:PHE:HB3	1:A:248:ASN:O	1.97	0.64
1:A:321:ARG:HD2	2:B:7:ALA:HB2	1.80	0.64
4:H:2:GAL:H4	4:H:3:SIA:O1A	1.98	0.63
1:E:123:THR:HG22	1:E:257:VAL:HG13	1.79	0.63
2:D:19:ASP:OD1	2:D:19:ASP:N	2.28	0.63
2:B:126:LEU:HD21	2:B:152:ILE:HD13	1.79	0.63
1:C:284:ARG:HD2	2:D:69:GLU:CG	2.28	0.63
2:D:54:ARG:NH2	2:D:103:GLU:OE2	2.31	0.63
1:A:108:LEU:HD11	1:A:177:LEU:HD11	1.82	0.62
2:D:110:MET:O	2:D:113:SER:HB3	1.99	0.62
2:D:134:GLY:HA2	2:F:124:LYS:HD3	1.82	0.62
1:C:307:LYS:HG3	2:D:92:TRP:CE2	2.36	0.61
2:B:91:ILE:HD13	2:D:91:ILE:HG21	1.83	0.60
2:B:51:LYS:NZ	2:B:106:HIS:HB3	2.16	0.60
1:C:42:THR:HG22	2:D:55:LEU:HD21	1.84	0.59
1:C:35:GLU:OE2	1:C:322:ASN:ND2	2.33	0.59
1:E:27:LYS:NZ	2:F:97:GLU:OE2	2.36	0.59
4:H:2:GAL:C4	4:H:3:SIA:C1	2.73	0.59
1:A:65:HIS:O	1:A:68:GLY:N	2.30	0.59
2:F:133:ASP:OD2	2:F:137:CYS:HB2	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:210:ARG:HD3	1:E:216:VAL:HB	1.85	0.59
1:C:151:LEU:HD23	1:C:254:PRO:HA	1.84	0.59
2:F:168:LEU:O	2:F:172:ASN:HB2	2.02	0.58
1:A:74:PRO:HG3	1:A:141:ARG:HE	1.67	0.58
4:H:1:NAG:H83	4:H:1:NAG:O1	2.03	0.58
1:C:300:ARG:NH1	2:D:69:GLU:OE2	2.37	0.58
1:C:190:GLU:CD	7:C:405:SIA:O9	2.42	0.57
2:B:51:LYS:HD3	2:B:103:GLU:HG3	1.86	0.57
1:E:108:LEU:HD11	1:E:177:LEU:HD11	1.85	0.57
2:D:66:ILE:HD11	2:D:85:LYS:HE3	1.87	0.57
1:E:106:GLU:OE2	1:E:109:ARG:NH2	2.32	0.57
2:B:103:GLU:OE2	1:E:29:LEU:HD22	2.05	0.56
2:D:51:LYS:HD3	2:D:103:GLU:HG3	1.87	0.55
2:D:83:TRP:CG	2:F:66:ILE:HD13	2.42	0.55
1:A:307:LYS:HG3	2:B:92:TRP:CE2	2.41	0.55
1:C:30:THR:O	2:F:50:GLY:HA3	2.06	0.55
1:A:27:LYS:NZ	2:B:97:GLU:OE2	2.38	0.55
1:A:98:TYR:HD2	1:A:136:THR:HG21	1.72	0.55
1:E:114:GLU:HB3	1:E:263:ARG:HH12	1.70	0.55
2:B:148:CYS:O	2:B:151:SER:HB3	2.07	0.55
1:C:225:GLY:HA3	1:E:205:GLY:O	2.06	0.54
1:A:242:ASN:HD22	3:G:1:NAG:H83	1.73	0.54
1:E:51:LEU:HD11	1:E:270:SER:HB3	1.90	0.54
1:C:218:GLY:HA3	1:C:226:LEU:HB3	1.89	0.54
2:F:148:CYS:O	2:F:151:SER:HB3	2.07	0.54
1:C:185:PRO:HG2	1:C:191:LYS:HB2	1.90	0.53
1:E:127:TYR:HB3	1:E:130:ILE:HD12	1.91	0.53
1:E:220:ARG:CB	1:E:221:PRO:HD3	2.38	0.53
1:A:59:LYS:HB3	1:A:87:LEU:HD12	1.89	0.53
2:D:131:GLU:HG2	2:D:141:TYR:HE1	1.72	0.53
2:B:132:GLU:HG2	2:B:138:PHE:CE1	2.43	0.52
2:B:22:TYR:H	2:B:41:THR:HG22	1.74	0.52
1:A:15:LEU:HD13	2:B:119:TYR:HA	1.92	0.52
1:E:190:GLU:OE2	7:E:602:SIA:O9	2.28	0.52
1:A:230:ILE:HD13	1:A:252:ILE:HG13	1.91	0.52
1:C:229:ARG:NH1	1:C:231:ASP:OD1	2.43	0.52
1:C:172:ASP:OD1	1:C:173:THR:N	2.37	0.51
1:E:221:PRO:HB2	1:E:223:VAL:HG12	1.92	0.51
1:A:201:SER:H	1:A:248:ASN:HB2	1.75	0.51
2:B:88:ILE:HG22	2:B:92:TRP:HD1	1.74	0.51
1:A:72:GLY:HA2	1:A:80:LEU:HD21	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:78:LEU:O	1:C:79:HIS:ND1	2.43	0.51
1:C:202:ILE:HG21	1:C:251:LEU:HD13	1.93	0.51
1:E:183:HIS:O	1:E:185:PRO:HD3	2.10	0.51
1:C:12:LYS:HB2	2:D:138:PHE:O	2.11	0.50
2:F:37:ASP:O	2:F:41:THR:HG23	2.12	0.50
2:F:45:ILE:O	2:F:49:THR:HG23	2.11	0.50
1:A:262:GLY:O	1:A:263:ARG:HD3	2.12	0.50
1:C:65:HIS:HE1	1:C:67:ILE:HG12	1.75	0.50
2:D:141:TYR:CD2	2:D:170:ARG:HG2	2.46	0.50
1:E:129:SER:O	1:E:157:LYS:HD3	2.12	0.50
1:A:17:HIS:HB3	2:B:115:MET:HE1	1.93	0.50
2:D:148:CYS:O	2:D:151:SER:HB3	2.11	0.50
2:B:28:ASN:ND2	2:B:146:ASP:OD1	2.35	0.49
1:A:295:GLN:HG2	1:A:306:PRO:HG2	1.93	0.49
1:E:220:ARG:HB2	1:E:221:PRO:HD3	1.94	0.49
1:E:65:HIS:CE1	1:E:67:ILE:HG12	2.47	0.49
1:A:199:SER:O	1:A:248:ASN:ND2	2.46	0.49
2:B:99:LEU:O	2:B:103:GLU:HB2	2.13	0.49
1:C:77:ASP:OD1	1:C:141:ARG:NH2	2.43	0.49
1:C:221:PRO:HD2	1:C:224:ASN:ND2	2.27	0.49
1:C:295:GLN:HG2	1:C:306:PRO:HG2	1.95	0.49
1:E:281:CYS:HB2	1:E:304:GLN:O	2.13	0.48
1:C:11:ASP:HB2	2:D:27:GLN:O	2.13	0.48
1:C:311:ARG:NH2	2:F:57:GLU:OE2	2.46	0.48
2:F:54:ARG:HB3	2:F:54:ARG:NH1	2.29	0.48
1:C:160:ASN:HD21	1:C:198:GLN:HB3	1.78	0.48
1:C:321:ARG:HB2	2:D:108:ILE:HG23	1.96	0.48
1:C:181:GLY:HA3	1:C:252:ILE:HB	1.95	0.48
1:C:102:THR:HB	1:C:105:VAL:HB	1.96	0.48
1:E:242:ASN:OD1	1:E:242:ASN:N	2.46	0.48
1:A:15:LEU:CD1	2:B:119:TYR:HA	2.43	0.48
2:D:165:GLU:O	2:D:169:ASN:ND2	2.45	0.48
1:A:65:HIS:HB2	1:A:93:ALA:HB1	1.95	0.47
2:D:3:PHE:HB3	2:D:116:LEU:HD23	1.96	0.47
2:D:141:TYR:CG	2:D:170:ARG:HG2	2.50	0.47
1:E:320:MET:HG3	1:E:321:ARG:O	2.13	0.47
1:C:50:ARG:NE	1:C:275:ASP:OD2	2.47	0.47
1:A:137:ARG:HE	4:H:2:GAL:H61	1.79	0.47
1:E:66:PRO:O	1:E:69:MET:HB2	2.15	0.47
1:E:293:PRO:HG2	1:E:294:PHE:CD2	2.49	0.47
1:A:52:CYS:HB2	1:A:279:SER:HB3	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:153:TRP:HZ2	1:A:183:HIS:CE1	2.33	0.47
1:C:71:ILE:O	1:C:148:TYR:HB3	2.15	0.47
1:C:296:ASN:HD22	1:C:312:ARG:HA	1.80	0.47
2:D:4:GLY:O	2:D:8:GLY:HA3	2.15	0.46
1:E:58:HIS:HE1	1:E:276:ASN:OD1	1.99	0.46
1:E:209:TYR:CD1	1:E:210:ARG:N	2.83	0.46
1:A:67:ILE:HD12	1:A:108:LEU:HD23	1.98	0.46
1:A:98:TYR:CD1	1:A:99:PRO:HD2	2.50	0.46
2:D:84:THR:O	2:D:88:ILE:HG12	2.14	0.46
2:D:131:GLU:OE1	2:F:127:ARG:NE	2.41	0.46
1:A:188:THR:HG22	1:A:192:ASN:ND2	2.30	0.46
2:B:167:LEU:HD23	2:B:167:LEU:HA	1.66	0.46
1:C:98:TYR:HD2	1:C:136:THR:HG21	1.79	0.46
2:D:26:HIS:CD2	2:D:149:MET:HG3	2.50	0.46
2:B:37:ASP:O	2:B:41:THR:HG23	2.16	0.45
2:B:55:LEU:CD2	2:B:99:LEU:HD21	2.43	0.45
1:C:72:GLY:HA2	1:C:80:LEU:HD21	1.97	0.45
1:E:69:MET:HE1	1:E:87:LEU:HD21	1.98	0.45
1:E:241:ASP:HA	6:E:601:NAG:H82	1.98	0.45
1:E:263:ARG:NH2	2:F:64:GLU:OE2	2.49	0.45
1:A:65:HIS:CE1	1:A:67:ILE:HG12	2.50	0.45
1:C:213:PHE:HE2	1:C:233:HIS:CD2	2.35	0.45
2:D:20:GLY:HA3	2:D:36:ALA:HB1	1.99	0.45
1:C:53:MET:HE1	1:C:56:ARG:HD3	1.97	0.45
2:D:129:ASN:HB3	2:D:142:HIS:CD2	2.52	0.45
4:H:3:SIA:H7	4:H:3:SIA:O10	2.16	0.45
1:E:129(A):SER:HB2	1:E:162:PRO:HG3	1.97	0.45
1:A:182:ILE:HD12	1:A:202:ILE:HD13	1.99	0.45
1:A:210:ARG:HH21	1:E:184:HIS:HB3	1.81	0.45
2:D:139:GLU:OE1	2:F:127:ARG:NH2	2.50	0.45
1:E:218:GLY:HA2	1:E:226:LEU:HB3	1.98	0.45
2:B:154:ASN:HB3	2:B:156:THR:HG22	1.98	0.44
2:D:142:HIS:CD2	2:D:162:TYR:HB3	2.53	0.44
1:E:190:GLU:OE1	7:E:602:SIA:O9	2.22	0.44
2:F:134:GLY:O	2:F:135:LYS:HD2	2.18	0.44
1:A:188:THR:O	1:A:192:ASN:ND2	2.50	0.44
1:C:225:GLY:O	1:C:227:SER:N	2.49	0.44
1:E:139:CYS:O	1:E:146:SER:HB3	2.18	0.44
2:B:17:MET:SD	2:B:23:GLY:HA3	2.58	0.44
2:B:117:ASN:HD21	2:F:2:LEU:HD23	1.82	0.43
2:B:124:LYS:HD3	2:F:134:GLY:HA2	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:288:ILE:HG12	1:C:297:LEU:HD12	2.00	0.43
1:E:60:ASP:HB2	1:E:274:ILE:HD11	1.99	0.43
1:E:151:LEU:HB3	1:E:252:ILE:HG22	2.00	0.43
1:A:185:PRO:HB3	1:A:190:GLU:HG2	1.99	0.43
2:B:91:ILE:HG21	2:F:91:ILE:HD13	2.01	0.43
2:F:167:LEU:O	2:F:171:LEU:HG	2.18	0.43
1:E:43:VAL:HG23	1:E:314:LEU:HB2	2.00	0.43
1:E:311:ARG:NE	2:F:97:GLU:OE1	2.43	0.43
1:A:27:LYS:HE3	1:A:27:LYS:HB2	1.76	0.43
1:A:120:LYS:HE2	1:A:258:SER:OG	2.18	0.43
1:C:103:VAL:HG21	1:C:233:HIS:CD2	2.53	0.43
1:C:182:ILE:HD12	1:C:202:ILE:HD13	2.00	0.43
2:F:141:TYR:O	2:F:166:ALA:HA	2.19	0.43
1:A:222:GLN:HG3	1:A:227:SER:OG	2.19	0.43
1:C:51:LEU:HD11	1:C:270:SER:HB3	2.00	0.43
1:E:307:LYS:HG3	2:F:92:TRP:CE2	2.54	0.43
2:F:5:ALA:HB1	2:F:115:MET:HG2	2.01	0.42
1:C:52:CYS:HB3	1:C:277:CYS:O	2.19	0.42
1:E:209:TYR:CE1	1:E:211:ASN:HB2	2.54	0.42
1:A:110:GLN:OE1	2:F:76:GLN:HG3	2.20	0.42
1:A:321:ARG:HH11	2:B:7:ALA:CB	2.32	0.42
1:C:53:MET:HE1	1:C:86:THR:HG21	2.02	0.42
1:E:34:GLU:OE2	1:E:321:ARG:NE	2.52	0.42
1:A:210:ARG:NH1	1:E:227:SER:O	2.52	0.42
2:D:26:HIS:CG	2:D:149:MET:HG3	2.55	0.42
1:E:65:HIS:HE1	1:E:67:ILE:HG12	1.85	0.42
1:C:97:CYS:HB2	1:C:138:ALA:O	2.20	0.41
1:E:59:LYS:HZ2	1:E:79:HIS:CE1	2.38	0.41
1:E:219:ALA:O	1:E:224:ASN:ND2	2.52	0.41
4:H:3:SIA:O1B	4:H:3:SIA:H6	2.19	0.41
1:A:151:LEU:HB3	1:A:252:ILE:HG22	2.01	0.41
1:C:102:THR:HG23	1:C:232:PHE:HB2	2.01	0.41
1:C:216:VAL:CG1	1:E:212:ASN:HB2	2.50	0.41
1:E:295:GLN:O	1:E:308:TYR:HA	2.19	0.41
1:C:158:SER:O	1:C:159:GLN:HB2	2.21	0.41
1:C:222:GLN:O	1:C:227:SER:OG	2.33	0.41
1:E:108:LEU:HD22	1:E:234:TRP:CD2	2.56	0.41
1:A:30:THR:O	2:D:50:GLY:HA3	2.21	0.41
1:A:183:HIS:HB2	1:A:252:ILE:HD11	2.02	0.41
1:A:185:PRO:HG2	1:A:191:LYS:HB2	2.02	0.41
1:A:231:ASP:OD2	1:C:210:ARG:NH2	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:160:ASN:ND2	1:C:198:GLN:HB3	2.35	0.41
1:C:213:PHE:CE2	1:C:233:HIS:CD2	3.09	0.41
2:D:74:GLU:O	2:D:77:ILE:HG22	2.20	0.41
2:D:140:ILE:HG22	2:D:142:HIS:H	1.86	0.41
1:A:179:MET:O	1:A:254:PRO:HB3	2.21	0.40
1:A:206:SER:OG	1:A:207:SER:N	2.53	0.40
1:E:160:ASN:HB3	1:E:198:GLN:NE2	2.36	0.40
1:A:178:ILE:HG13	1:A:257:VAL:HG12	2.03	0.40
1:C:26:VAL:HG21	1:C:317:ALA:HB2	2.02	0.40
2:B:117:ASN:ND2	2:F:2:LEU:HD23	2.37	0.40
1:C:45:SER:HA	1:C:296:ASN:OD1	2.22	0.40
1:C:296:ASN:ND2	1:C:312:ARG:HA	2.37	0.40
1:E:141:ARG:NH1	1:E:146:SER:OG	2.54	0.40
2:D:129:ASN:HB3	2:D:142:HIS:HD2	1.86	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:129:SER:OG	4:H:1:NAG:O7[1_655]	2.02	0.18

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	316/323 (98%)	304 (96%)	12 (4%)	0	100	100
1	C	317/323 (98%)	297 (94%)	19 (6%)	1 (0%)	41	74
1	E	316/323 (98%)	304 (96%)	11 (4%)	1 (0%)	41	74
2	B	169/180 (94%)	165 (98%)	4 (2%)	0	100	100
2	D	164/180 (91%)	157 (96%)	7 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	F	169/180 (94%)	161 (95%)	8 (5%)	0	100	100
All	All	1451/1509 (96%)	1388 (96%)	61 (4%)	2 (0%)	51	83

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	220	ARG
1	E	220	ARG

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	269/272 (99%)	266 (99%)	3 (1%)	73	90
1	C	270/272 (99%)	268 (99%)	2 (1%)	84	93
1	E	269/272 (99%)	266 (99%)	3 (1%)	73	90
2	B	146/154 (95%)	143 (98%)	3 (2%)	53	80
2	D	143/154 (93%)	140 (98%)	3 (2%)	53	80
2	F	146/154 (95%)	144 (99%)	2 (1%)	67	86
All	All	1243/1278 (97%)	1227 (99%)	16 (1%)	69	88

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	HIS
1	A	150	GLU
1	A	295	GLN
2	B	97	GLU
2	B	113	SER
2	B	148	CYS
1	C	164	THR
1	C	188	THR
2	D	19	ASP

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Mol	Chain	Res	Type
2	D	69	GLU
2	D	128	GLN
1	E	18	HIS
1	E	199	SER
1	E	246	SER
2	F	12	ASN
2	F	59	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	192	ASN
1	A	212	ASN
2	B	95	GLN
2	D	79	ASN
1	E	58	HIS
1	E	248	ASN
1	E	276	ASN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

7 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	$\#\|Z\| > 2$
3	NAG	G	1	1,3	14,14,15	0.91	2 (14%)	17,19,21	1.16	1 (5%)
3	NAG	G	2	3	14,14,15	0.39	0	17,19,21	0.97	1 (5%)
4	NAG	H	1	4	15,15,15	0.50	0	21,21,21	1.81	5 (23%)
4	GAL	H	2	4	11,11,12	1.62	2 (18%)	15,15,17	1.66	4 (26%)
4	SIA	H	3	4	20,20,21	1.57	3 (15%)	24,28,31	1.48	5 (20%)
5	NAG	I	1	5	14,14,15	0.22	0	17,19,21	0.53	0
5	BMA	I	2	5	11,11,12	1.47	2 (18%)	15,15,17	0.91	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
3	NAG	G	1	1,3	-	4/6/23/26	0/1/1/1
3	NAG	G	2	3	-	4/6/23/26	0/1/1/1
4	NAG	H	1	4	-	3/6/26/26	0/1/1/1
4	GAL	H	2	4	-	0/2/19/22	0/1/1/1
4	SIA	H	3	4	-	4/18/34/38	0/1/1/1
5	NAG	I	1	5	-	2/6/23/26	0/1/1/1
5	BMA	I	2	5	-	0/2/19/22	0/1/1/1

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	H	3	SIA	C2-C1	-5.51	1.47	1.52
4	H	2	GAL	O5-C1	-3.86	1.37	1.43
4	H	2	GAL	C1-C2	3.29	1.59	1.52
5	I	2	BMA	O5-C1	3.11	1.48	1.43
5	I	2	BMA	O5-C5	2.59	1.48	1.43
3	G	1	NAG	O5-C1	-2.38	1.39	1.43
4	H	3	SIA	O10-C10	-2.32	1.18	1.23
4	H	3	SIA	O1B-C1	-2.17	1.23	1.30
3	G	1	NAG	C2-N2	-2.10	1.42	1.46

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	H	3	SIA	C6-C5-N5	-4.42	103.57	110.91

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	H	1	NAG	O5-C1-C2	-4.06	105.44	109.52
4	H	1	NAG	O3-C3-C4	-3.99	101.12	110.35
4	H	1	NAG	O4-C4-C3	-3.41	102.47	110.35
3	G	1	NAG	C1-C2-N2	-3.26	104.91	110.49
4	H	2	GAL	C1-O5-C5	-3.18	107.88	112.19
4	H	2	GAL	O2-C2-C3	-3.17	103.78	110.14
3	G	2	NAG	C2-N2-C7	-2.91	118.75	122.90
4	H	2	GAL	O2-C2-C1	2.55	114.38	109.15
4	H	1	NAG	O1-C1-C2	2.51	114.44	109.22
4	H	2	GAL	O5-C1-C2	-2.42	107.03	110.77
4	H	3	SIA	O1A-C1-C2	-2.37	116.97	122.57
4	H	3	SIA	C6-O6-C2	2.30	116.26	111.34
5	I	2	BMA	C1-O5-C5	2.21	115.19	112.19
4	H	3	SIA	O1B-C1-C2	2.13	119.10	113.03
4	H	3	SIA	O7-C7-C8	-2.12	103.69	108.81
4	H	1	NAG	O5-C5-C6	2.03	111.47	106.44

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	G	1	NAG	C8-C7-N2-C2
3	G	1	NAG	O7-C7-N2-C2
4	H	1	NAG	C1-C2-N2-C7
4	H	1	NAG	C8-C7-N2-C2
4	H	1	NAG	O7-C7-N2-C2
4	H	3	SIA	C7-C8-C9-O9
4	H	3	SIA	O8-C8-C9-O9
3	G	2	NAG	O5-C5-C6-O6
5	I	1	NAG	O5-C5-C6-O6
3	G	2	NAG	C4-C5-C6-O6
3	G	1	NAG	O5-C5-C6-O6
5	I	1	NAG	C4-C5-C6-O6
4	H	3	SIA	C11-C10-N5-C5
3	G	2	NAG	C8-C7-N2-C2
3	G	1	NAG	C4-C5-C6-O6
4	H	3	SIA	O10-C10-N5-C5
3	G	2	NAG	O7-C7-N2-C2

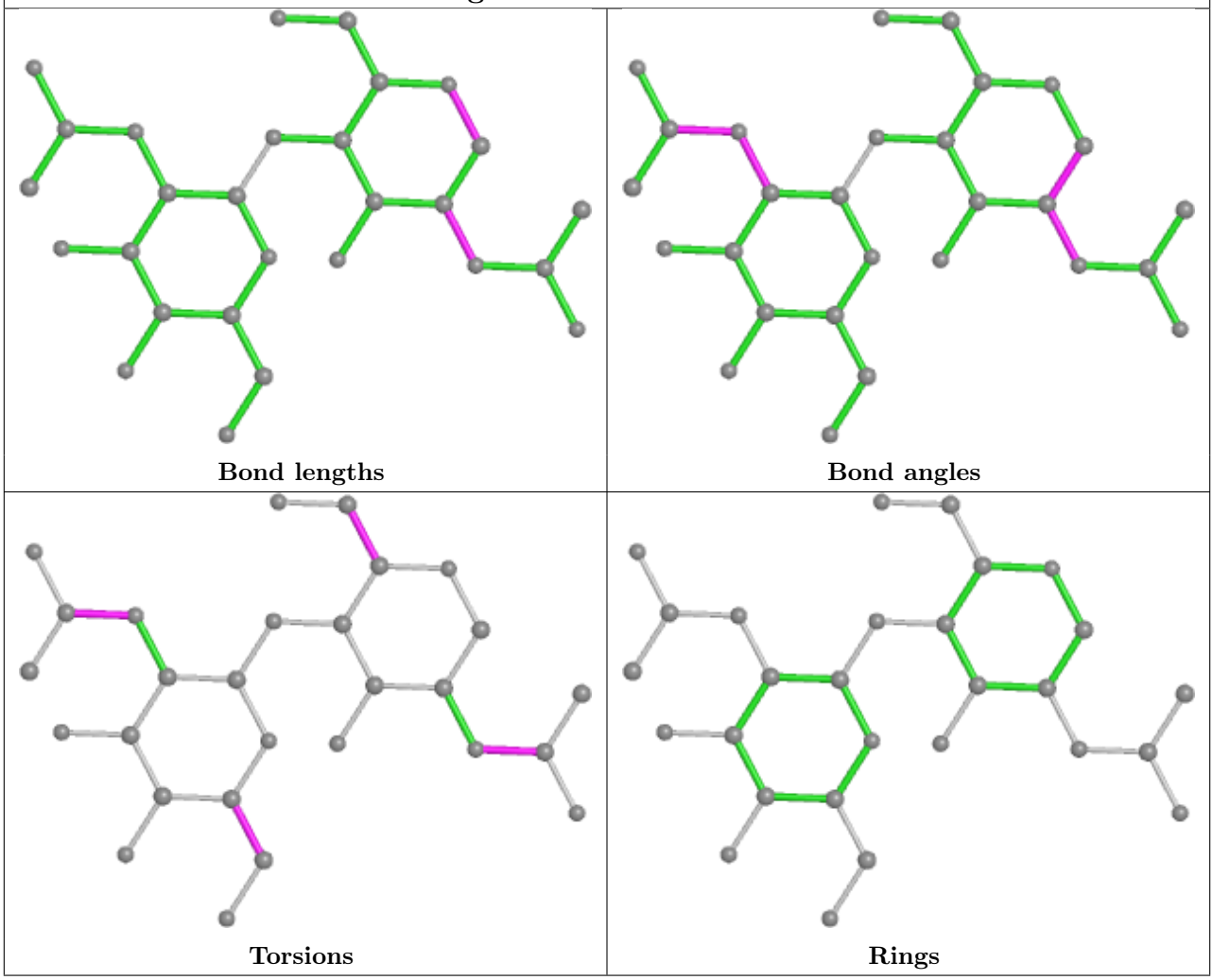
There are no ring outliers.

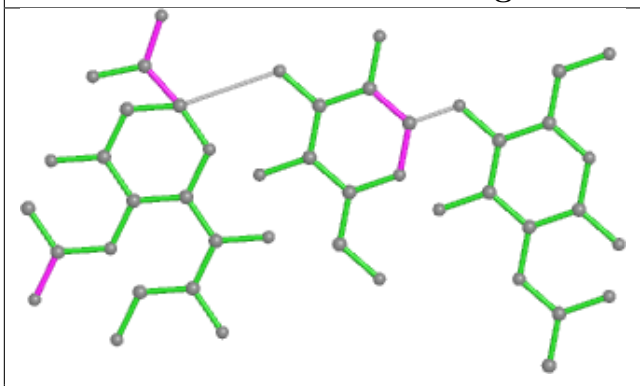
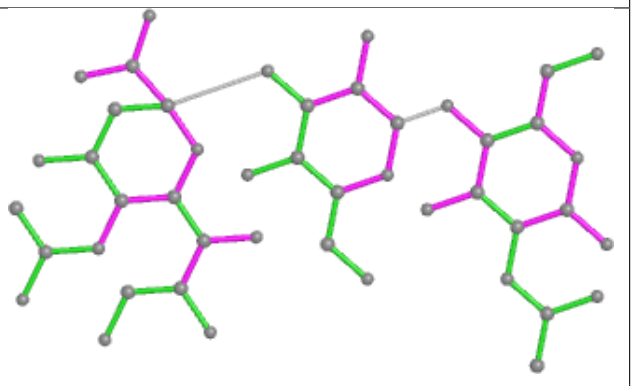
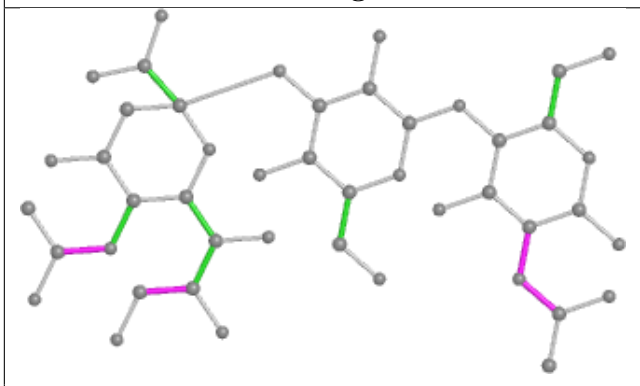
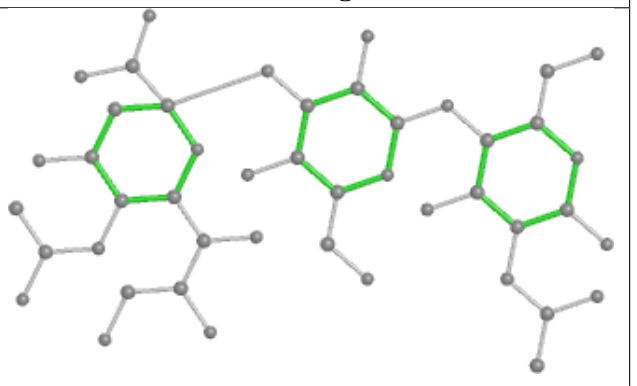
5 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	H	3	SIA	6	0
4	H	2	GAL	4	0
3	G	1	NAG	4	0
3	G	2	NAG	3	0
4	H	1	NAG	1	1

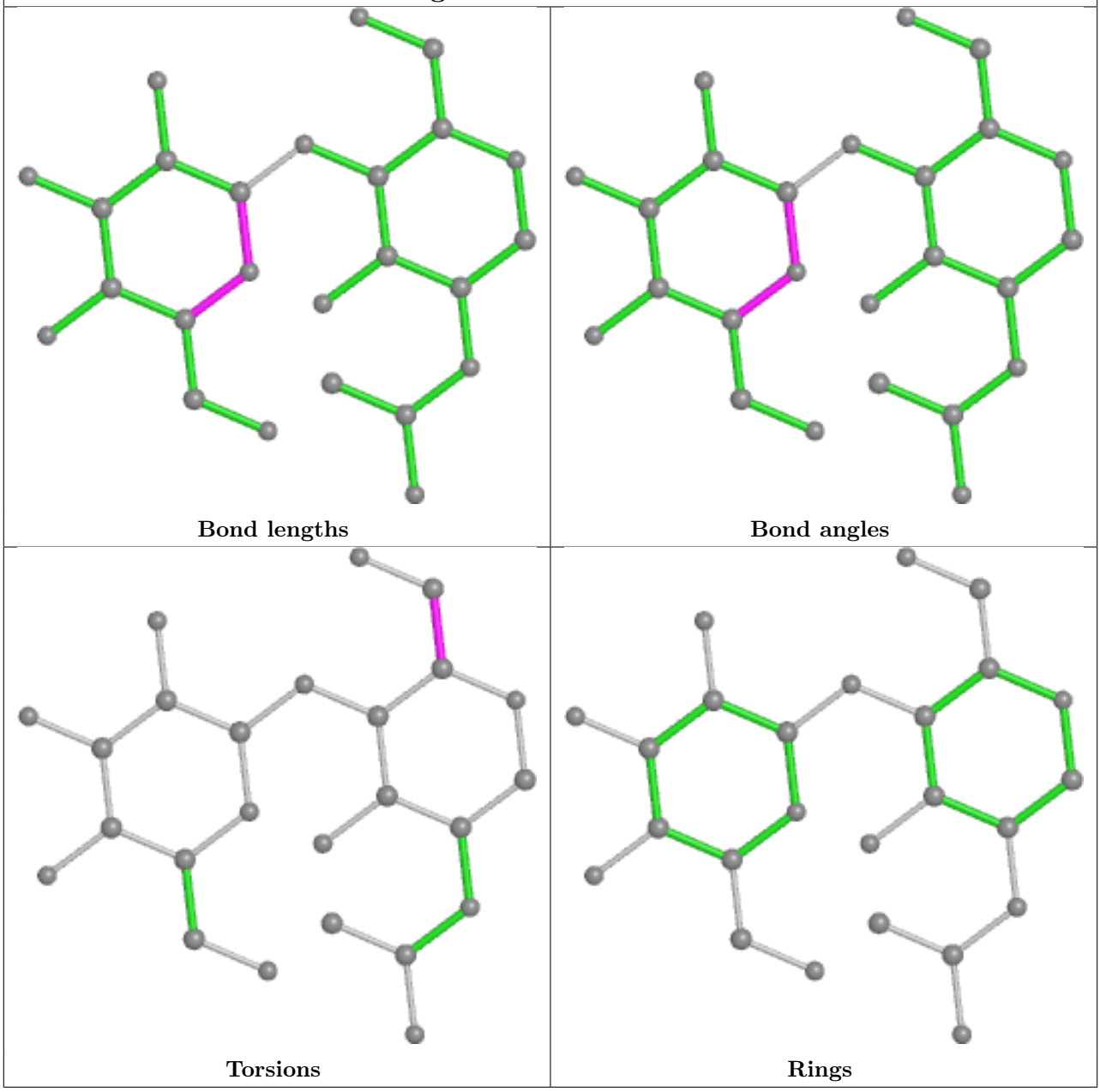
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

Oligosaccharide Chain G



Oligosaccharide Chain H	
 Bond lengths	 Bond angles
 Torsions	 Rings

Oligosaccharide Chain I



5.6 Ligand geometry (i)

6 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
6	NAG	E	601	1	14,14,15	1.16	2 (14%)	17,19,21	1.21	1 (5%)
6	NAG	C	402	1	14,14,15	0.76	1 (7%)	17,19,21	0.84	0
6	NAG	A	401	1	14,14,15	0.69	1 (7%)	17,19,21	0.79	0
6	NAG	C	401	1	14,14,15	0.86	1 (7%)	17,19,21	1.11	1 (5%)
7	SIA	C	405	-	21,21,21	1.68	3 (14%)	25,31,31	2.12	7 (28%)
7	SIA	E	602	-	21,21,21	1.78	4 (19%)	25,31,31	2.21	7 (28%)

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
6	NAG	E	601	1	-	2/6/23/26	0/1/1/1
6	NAG	C	402	1	-	0/6/23/26	0/1/1/1
6	NAG	A	401	1	-	4/6/23/26	0/1/1/1
6	NAG	C	401	1	-	5/6/23/26	0/1/1/1
7	SIA	C	405	-	-	11/20/38/38	0/1/1/1
7	SIA	E	602	-	-	6/20/38/38	0/1/1/1

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	C	405	SIA	C2-C1	-5.15	1.45	1.53
7	E	602	SIA	C2-C1	-4.57	1.46	1.53
7	E	602	SIA	O2-C2	3.86	1.44	1.39
7	C	405	SIA	O2-C2	3.84	1.44	1.39
7	E	602	SIA	C8-C7	-2.57	1.48	1.53
6	C	401	NAG	O5-C1	-2.45	1.39	1.43
6	C	402	NAG	O5-C1	-2.44	1.39	1.43
6	E	601	NAG	O5-C1	-2.42	1.39	1.43
7	E	602	SIA	C5-N5	-2.24	1.42	1.45
6	A	401	NAG	O5-C1	-2.20	1.40	1.43
6	E	601	NAG	C2-N2	-2.17	1.42	1.46
7	C	405	SIA	O10-C10	-2.06	1.18	1.23

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
7	E	602	SIA	O1A-C1-C2	-6.09	114.36	123.59
7	C	405	SIA	O1A-C1-C2	-5.22	115.68	123.59
7	C	405	SIA	C6-C5-N5	-5.04	102.54	110.91
7	E	602	SIA	O6-C6-C7	4.63	114.44	107.29
7	C	405	SIA	C4-C5-N5	-3.94	102.58	110.38
7	E	602	SIA	C6-C5-N5	-3.89	104.44	110.91
7	C	405	SIA	O2-C2-C1	-2.78	105.12	110.76
7	E	602	SIA	C3-C2-C1	-2.58	108.19	113.00
7	C	405	SIA	O7-C7-C6	-2.57	103.95	109.50
7	E	602	SIA	O6-C6-C5	2.54	112.25	109.78
6	C	401	NAG	C1-C2-N2	-2.53	106.16	110.49
7	C	405	SIA	O6-C6-C5	2.52	112.23	109.78
7	E	602	SIA	O4-C4-C5	2.43	115.35	109.77
7	E	602	SIA	C4-C5-N5	-2.42	105.59	110.38
7	C	405	SIA	O4-C4-C5	2.39	115.27	109.77
6	E	601	NAG	O5-C5-C6	-2.21	103.74	107.20

There are no chirality outliers.

All (28) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	401	NAG	C8-C7-N2-C2
6	A	401	NAG	O7-C7-N2-C2
6	C	401	NAG	C8-C7-N2-C2
6	C	401	NAG	O7-C7-N2-C2
7	C	405	SIA	O7-C7-C8-C9
7	C	405	SIA	O8-C8-C9-O9
7	E	602	SIA	O1A-C1-C2-O2
7	E	602	SIA	O1A-C1-C2-O6
7	E	602	SIA	O1B-C1-C2-O2
7	E	602	SIA	O1B-C1-C2-O6
6	E	601	NAG	O5-C5-C6-O6
7	E	602	SIA	O8-C8-C9-O9
6	C	401	NAG	O5-C5-C6-O6
7	C	405	SIA	C7-C8-C9-O9
7	E	602	SIA	C7-C8-C9-O9
6	E	601	NAG	C4-C5-C6-O6
6	C	401	NAG	C4-C5-C6-O6
6	A	401	NAG	O5-C5-C6-O6
7	C	405	SIA	O7-C7-C8-O8
6	A	401	NAG	C4-C5-C6-O6
7	C	405	SIA	C6-C7-C8-C9
7	C	405	SIA	C11-C10-N5-C5

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Mol	Chain	Res	Type	Atoms
7	C	405	SIA	C6-C7-C8-O8
7	C	405	SIA	O10-C10-N5-C5
6	C	401	NAG	C3-C2-N2-C7
7	C	405	SIA	O1A-C1-C2-O6
7	C	405	SIA	O1A-C1-C2-O2
7	C	405	SIA	O1B-C1-C2-O6

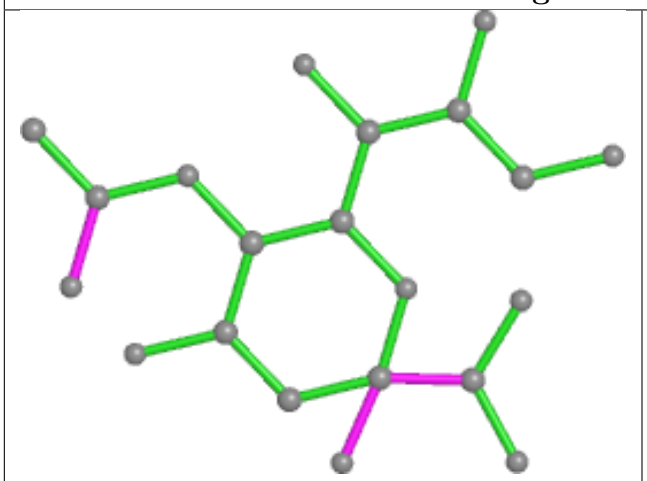
There are no ring outliers.

3 monomers are involved in 7 short contacts:

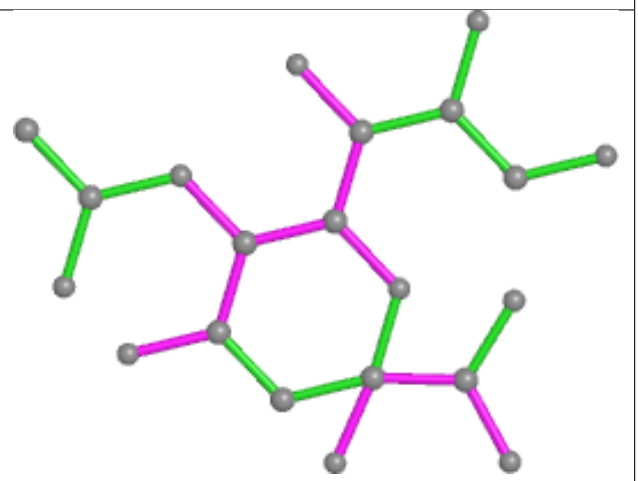
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	E	601	NAG	1	0
7	C	405	SIA	2	0
7	E	602	SIA	4	0

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.

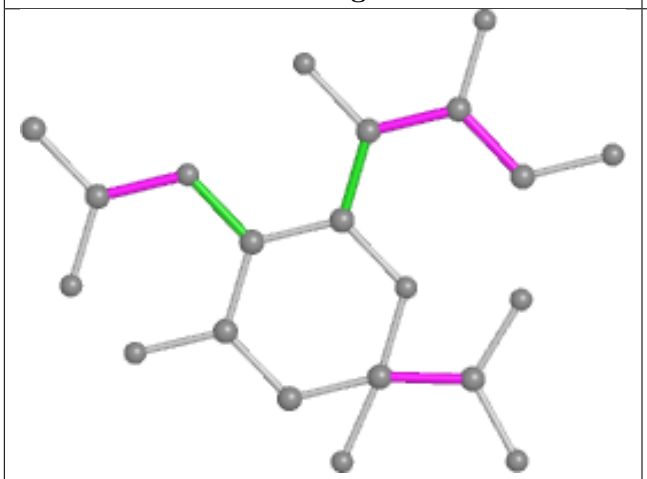
Ligand SIA C 405



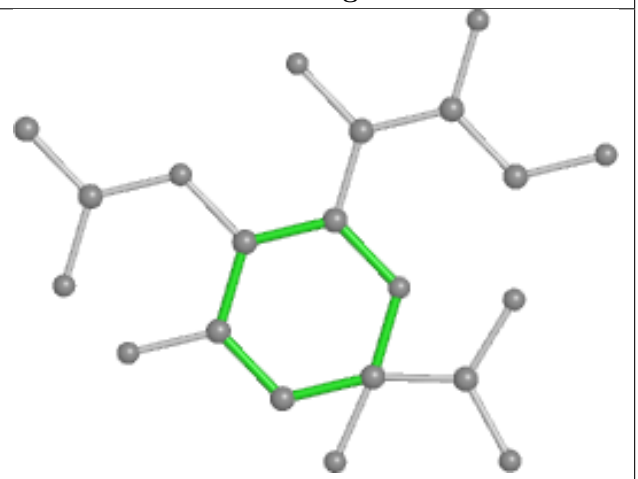
Bond lengths



Bond angles

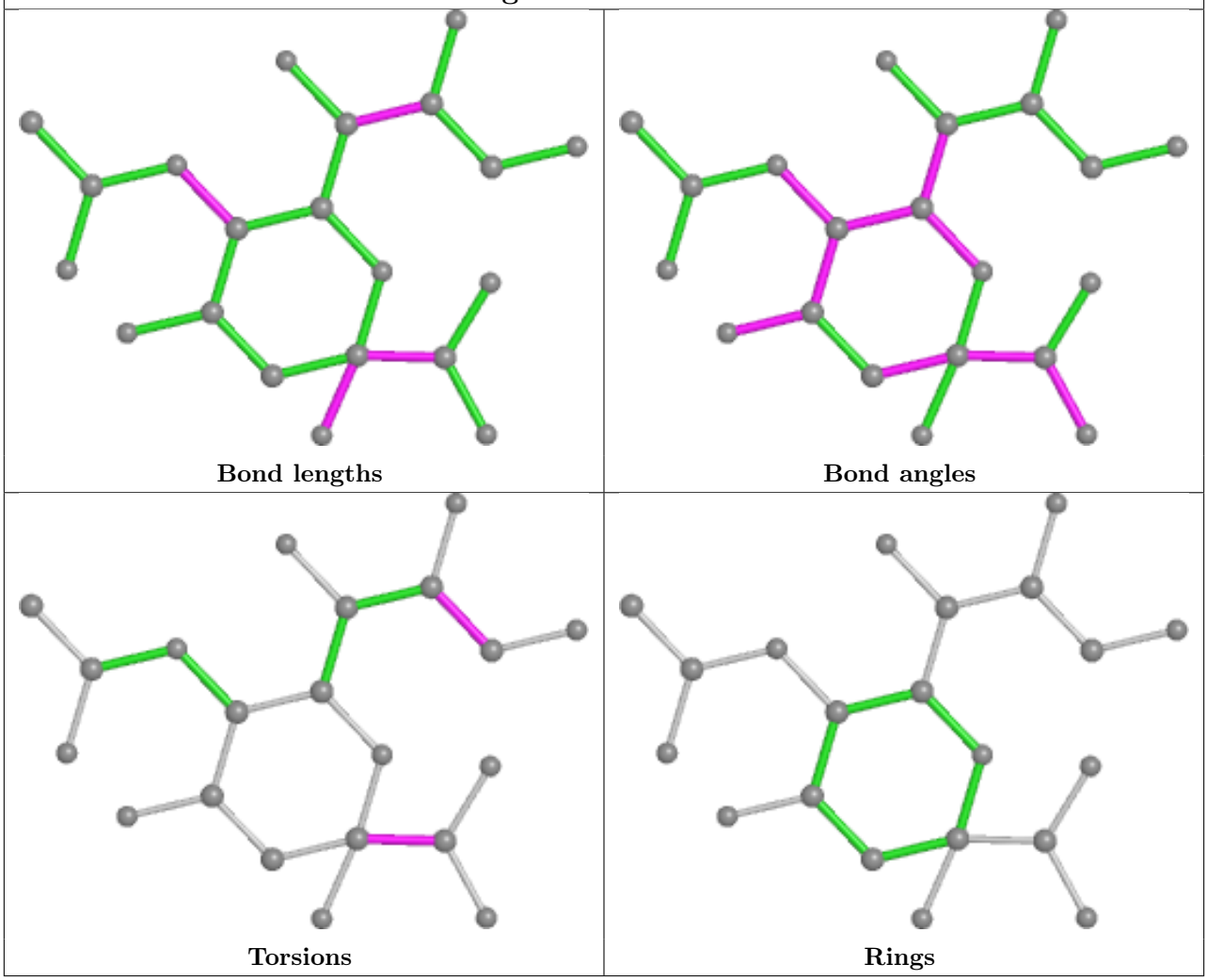


Torsions



Rings

Ligand SIA E 602



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 (i)

6.1 Protein, DNA and RNA chains (i)

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	318/323 (98%)	-0.23	0	100	100	26, 40, 62, 75	0
1	C	319/323 (98%)	-0.08	5 (1%)	72	52	27, 46, 66, 90	0
1	E	318/323 (98%)	-0.10	3 (0%)	84	69	29, 44, 59, 76	0
2	B	171/180 (95%)	0.19	1 (0%)	89	77	31, 54, 71, 80	0
2	D	168/180 (93%)	0.10	2 (1%)	79	61	32, 56, 73, 78	0
2	F	171/180 (95%)	0.00	1 (0%)	89	77	29, 49, 62, 84	0
All	All	1465/1509 (97%)	-0.05	12 (0%)	86	71	26, 46, 67, 90	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	220	ARG	3.0
1	E	222	GLN	2.8
2	F	59	THR	2.7
1	E	46	THR	2.6
1	C	222	GLN	2.4
2	D	2	LEU	2.3
1	C	219	ALA	2.2
1	C	326	LEU	2.2
1	C	325	GLU	2.1
2	D	168	LEU	2.1
1	E	297	LEU	2.1
2	B	143	ALA	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)

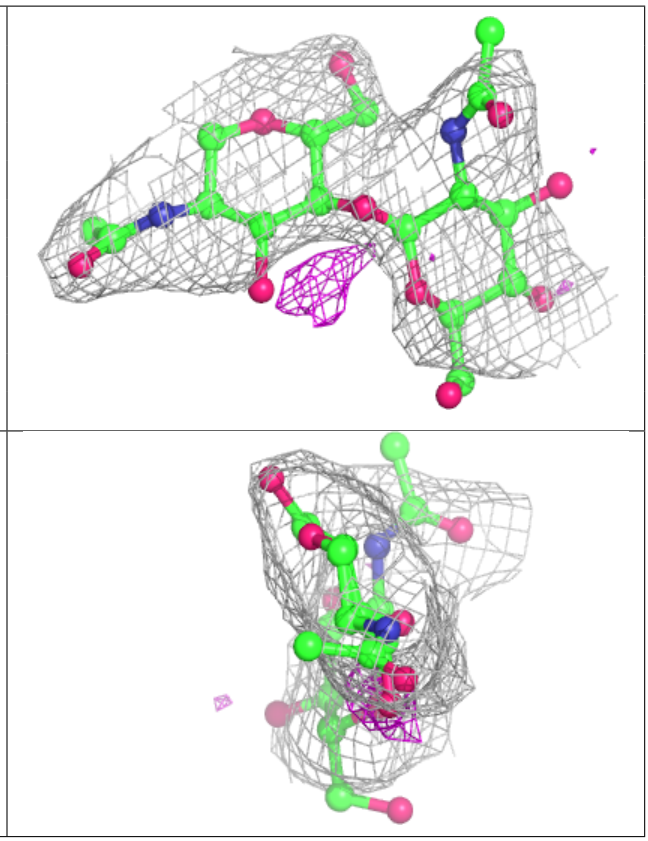
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	NAG	G	2	14/15	0.79	0.44	86,95,100,102	0
4	GAL	H	2	11/12	0.80	0.37	80,88,98,101	0
4	NAG	H	1	15/15	0.81	0.34	84,96,100,102	0
5	NAG	I	1	14/15	0.81	0.27	49,74,80,81	0
5	BMA	I	2	11/12	0.84	0.25	75,79,82,85	0
4	SIA	H	3	20/21	0.86	0.30	52,66,80,84	0
3	NAG	G	1	14/15	0.88	0.22	58,66,82,91	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.

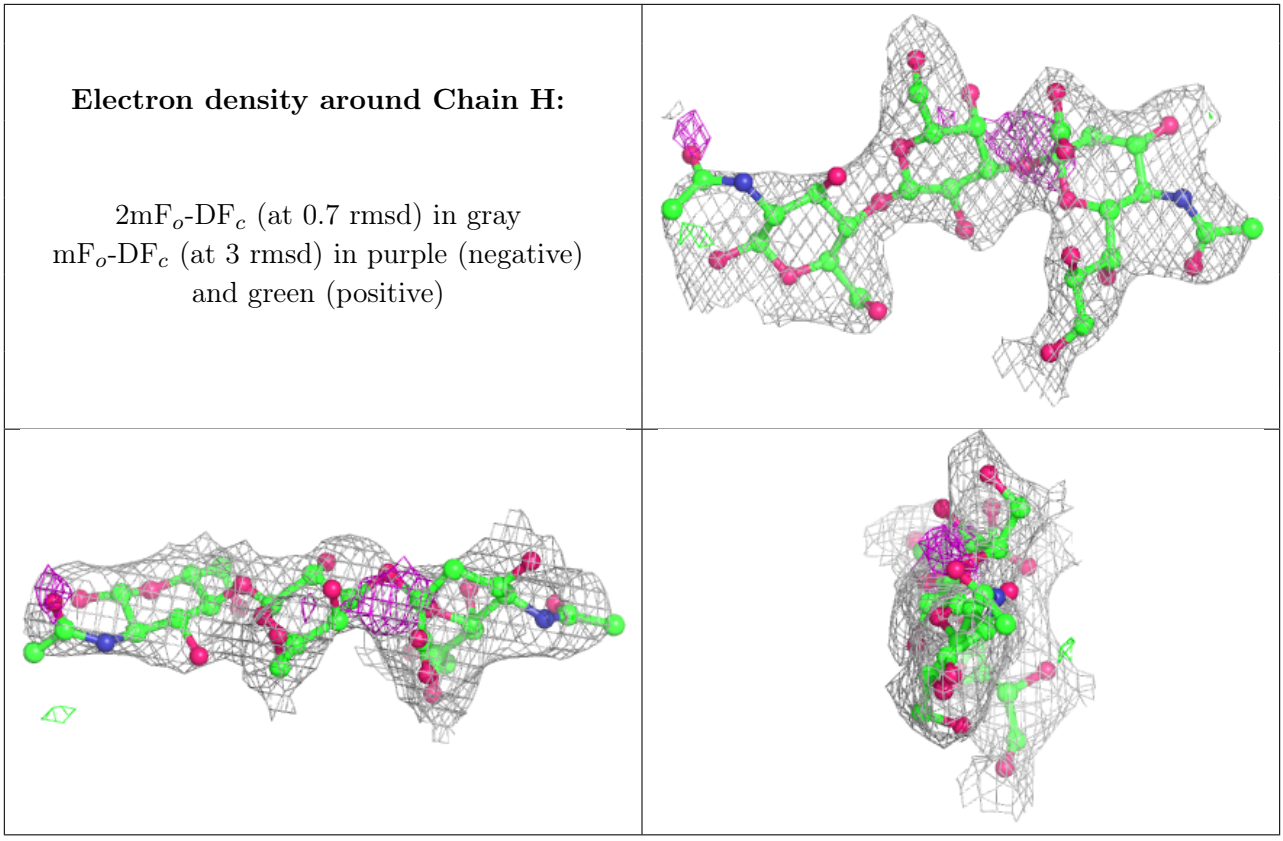
Electron density around Chain G:

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



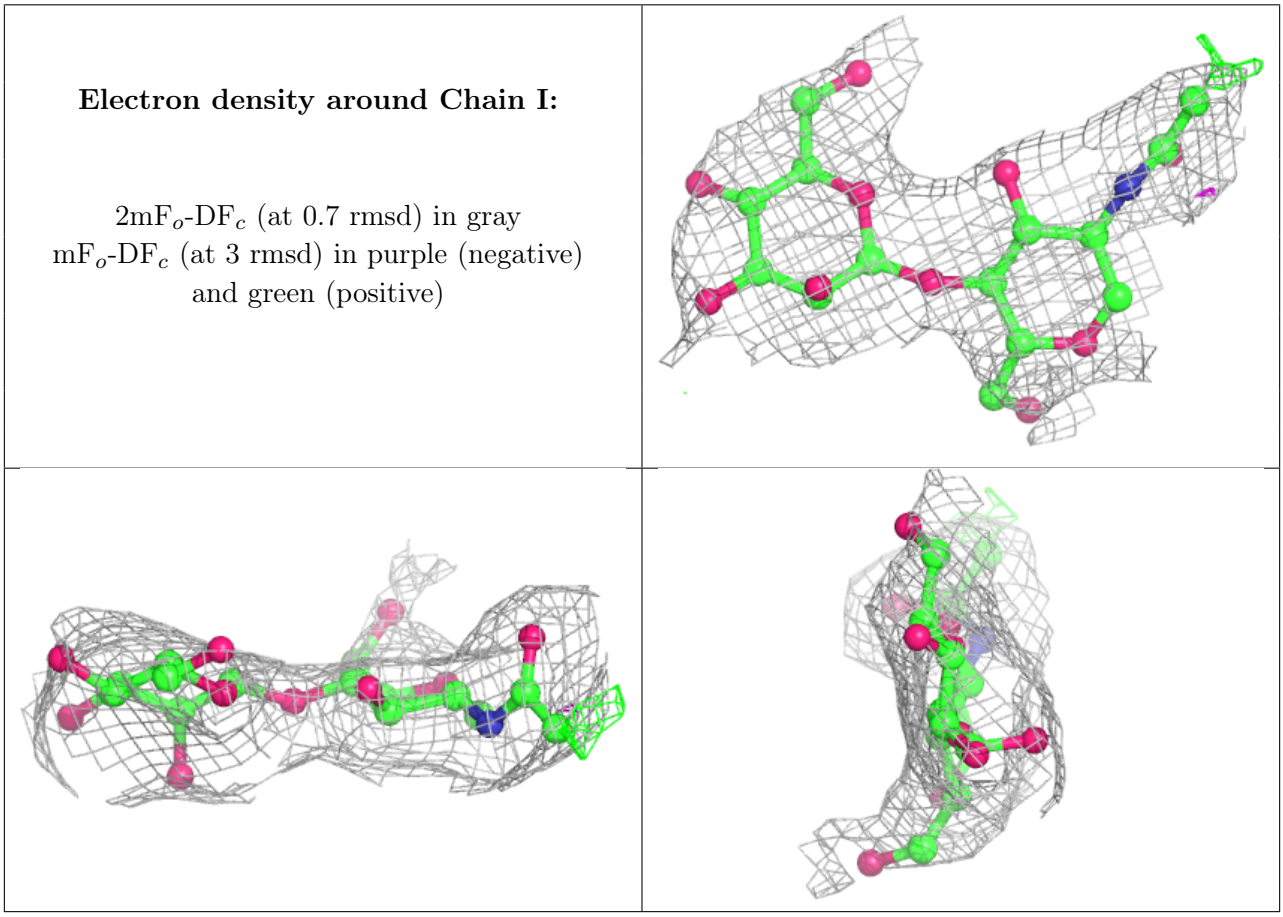
Electron density around Chain H:

$2mF_o$-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)

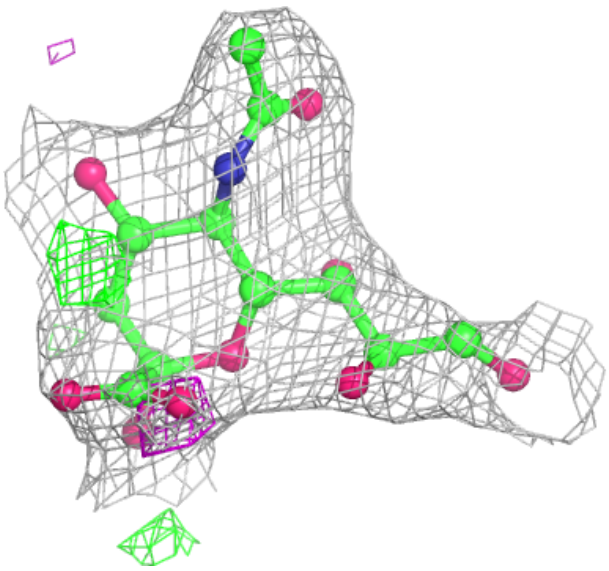
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

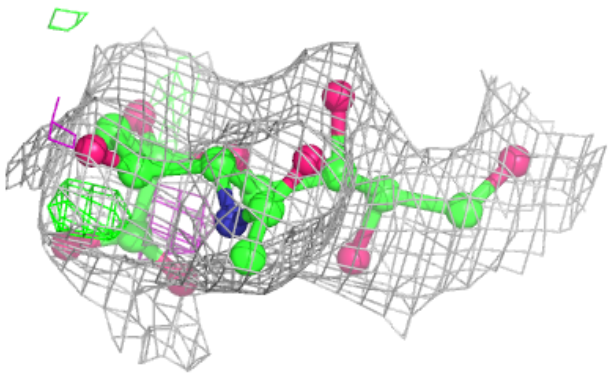
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
6	NAG	C	401	14/15	0.74	0.34	69,90,97,99	0
6	NAG	A	401	14/15	0.77	0.38	80,92,106,107	0
6	NAG	C	402	14/15	0.88	0.20	50,58,66,67	0
6	NAG	E	601	14/15	0.88	0.21	57,72,79,80	0
7	SIA	C	405	21/21	0.88	0.25	42,54,57,63	1
7	SIA	E	602	21/21	0.90	0.19	37,49,58,59	0

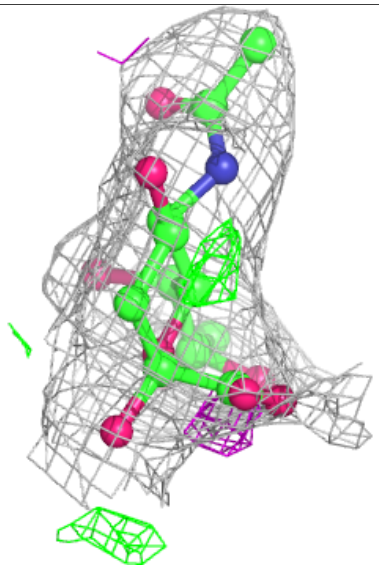
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around SIA C 405:

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

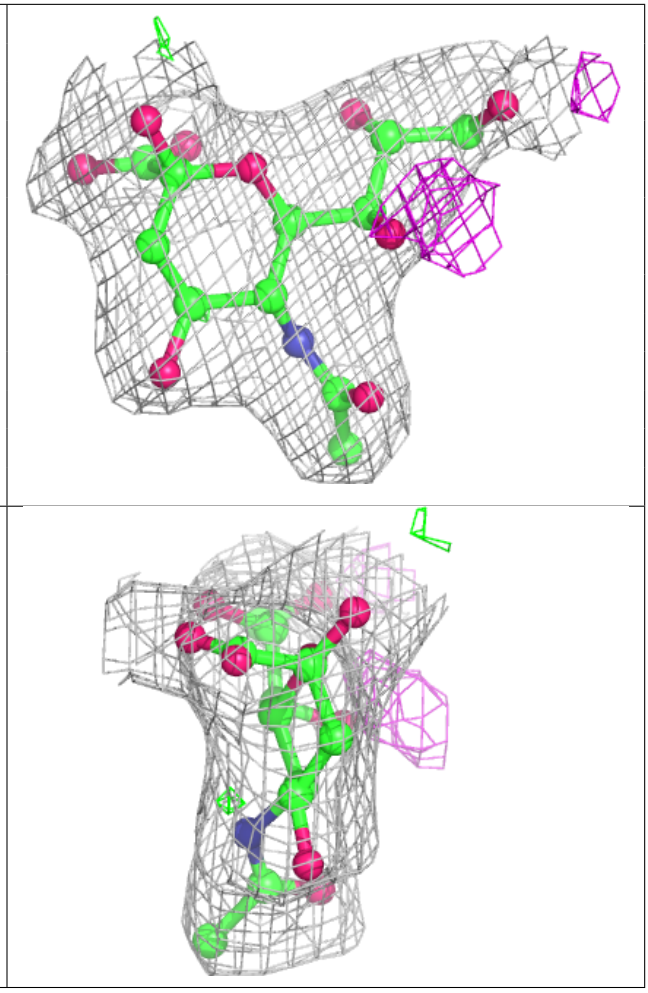






Electron density around SIA E 602:

$2mF_o$-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 (i)

There are no such residues in this entry.