



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 12:33 PM UTC

PDB ID : 4WRL / pdb_00004wrl
Title : Structure of the human CSF-1:CSF-1R complex
Authors : Felix, J.; De Munck, S.; Elegheert, J.; Savvides, S.N.
Deposited on : 2014-10-24
Resolution : 2.80 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 ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

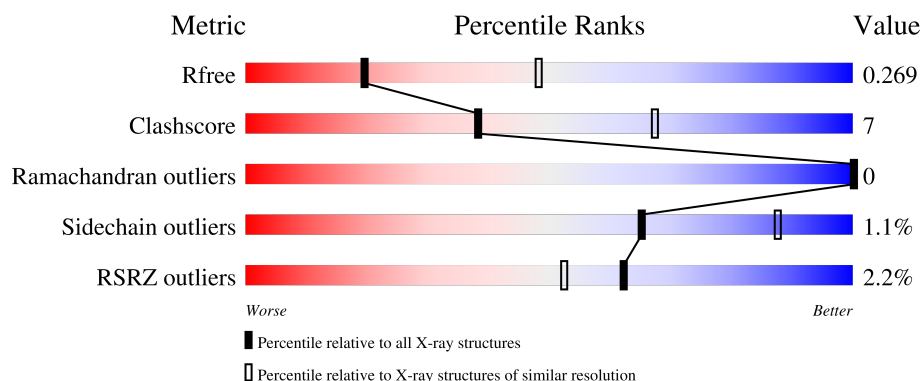
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

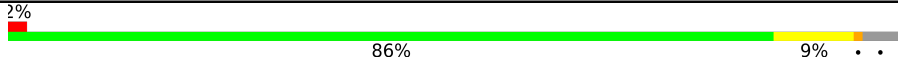
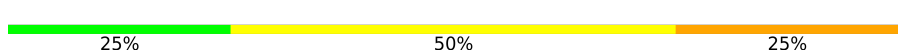
The reported resolution of this entry is 2.80 Å.

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}	180053	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	286	
1	C	286	
2	B	170	
2	D	170	
3	E	4	

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Mol	Chain	Length	Quality of chain
3	H	4	 25% 75%
4	F	2	 50% 50%
5	G	3	 100%
5	I	3	 33% 67%

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	SIA	H	3	-	-	X	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 6657 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 Macrophage colony-stimulating factor 1 receptor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	274	Total	C	N	O	S	0	0	0
			2074	1317	363	384	10			
1	C	277	Total	C	N	O	S	0	0	0
			2099	1333	366	390	10			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	240	GLN	ASN	engineered mutation	UNP P07333
A	297	GLY	-	expression tag	UNP P07333
A	298	THR	-	expression tag	UNP P07333
A	299	LYS	-	expression tag	UNP P07333
A	300	HIS	-	expression tag	UNP P07333
A	301	HIS	-	expression tag	UNP P07333
A	302	HIS	-	expression tag	UNP P07333
A	303	HIS	-	expression tag	UNP P07333
A	304	HIS	-	expression tag	UNP P07333
A	305	HIS	-	expression tag	UNP P07333
C	240	GLN	ASN	engineered mutation	UNP P07333
C	297	GLY	-	expression tag	UNP P07333
C	298	THR	-	expression tag	UNP P07333
C	299	LYS	-	expression tag	UNP P07333
C	300	HIS	-	expression tag	UNP P07333
C	301	HIS	-	expression tag	UNP P07333
C	302	HIS	-	expression tag	UNP P07333
C	303	HIS	-	expression tag	UNP P07333
C	304	HIS	-	expression tag	UNP P07333
C	305	HIS	-	expression tag	UNP P07333

- Molecule 2 is a protein called Macrophage colony-stimulating factor 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	141	Total	C	N	O	S	0	0	0
			1095	686	184	214	11			
2	D	141	Total	C	N	O	S	0	0	0
			1102	691	185	215	11			

There are 42 discrepancies between the modelled and reference sequences:

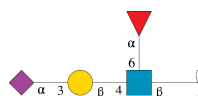
Chain	Residue	Modelled	Actual	Comment	Reference
B	-20	MET	-	initiating methionine	UNP P09603
B	-19	GLY	-	expression tag	UNP P09603
B	-18	SER	-	expression tag	UNP P09603
B	-17	SER	-	expression tag	UNP P09603
B	-16	HIS	-	expression tag	UNP P09603
B	-15	HIS	-	expression tag	UNP P09603
B	-14	HIS	-	expression tag	UNP P09603
B	-13	HIS	-	expression tag	UNP P09603
B	-12	HIS	-	expression tag	UNP P09603
B	-11	HIS	-	expression tag	UNP P09603
B	-10	SER	-	expression tag	UNP P09603
B	-9	SER	-	expression tag	UNP P09603
B	-8	GLY	-	expression tag	UNP P09603
B	-7	LEU	-	expression tag	UNP P09603
B	-6	VAL	-	expression tag	UNP P09603
B	-5	PRO	-	expression tag	UNP P09603
B	-4	ARG	-	expression tag	UNP P09603
B	-3	GLY	-	expression tag	UNP P09603
B	-2	SER	-	expression tag	UNP P09603
B	-1	HIS	-	expression tag	UNP P09603
B	0	MET	-	expression tag	UNP P09603
D	-20	MET	-	initiating methionine	UNP P09603
D	-19	GLY	-	expression tag	UNP P09603
D	-18	SER	-	expression tag	UNP P09603
D	-17	SER	-	expression tag	UNP P09603
D	-16	HIS	-	expression tag	UNP P09603
D	-15	HIS	-	expression tag	UNP P09603
D	-14	HIS	-	expression tag	UNP P09603
D	-13	HIS	-	expression tag	UNP P09603
D	-12	HIS	-	expression tag	UNP P09603
D	-11	HIS	-	expression tag	UNP P09603
D	-10	SER	-	expression tag	UNP P09603
D	-9	SER	-	expression tag	UNP P09603
D	-8	GLY	-	expression tag	UNP P09603
D	-7	LEU	-	expression tag	UNP P09603

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-6	VAL	-	expression tag	UNP P09603
D	-5	PRO	-	expression tag	UNP P09603
D	-4	ARG	-	expression tag	UNP P09603
D	-3	GLY	-	expression tag	UNP P09603
D	-2	SER	-	expression tag	UNP P09603
D	-1	HIS	-	expression tag	UNP P09603
D	0	MET	-	expression tag	UNP P09603

- Molecule 3 is an oligosaccharide called N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose.



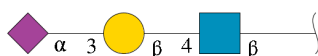
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	E	4	Total	C	N	O	0	0	0
			55	31	2	22			
3	H	4	Total	C	N	O	0	0	0
			55	31	2	22			

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



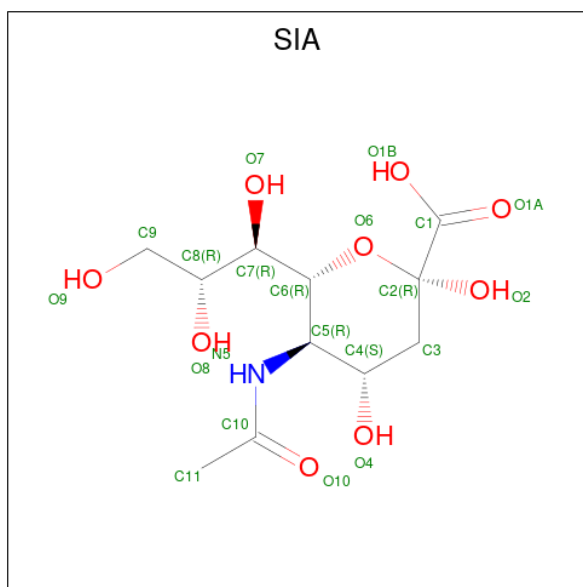
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	F	2	Total	C	N	O	0	0	0
			25	14	1	10			

- Molecule 5 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
5	G	3	Total	C	N	O	0	0	0
			45	25	2	18			
5	I	3	Total	C	N	O	0	0	0
			45	25	2	18			

- Molecule 6 is N-acetyl-alpha-neuraminic acid (CCD ID: SIA) (formula: $C_{11}H_{19}NO_9$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	N	O	0	0
			20	11	1	8		

- Molecule 7 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).

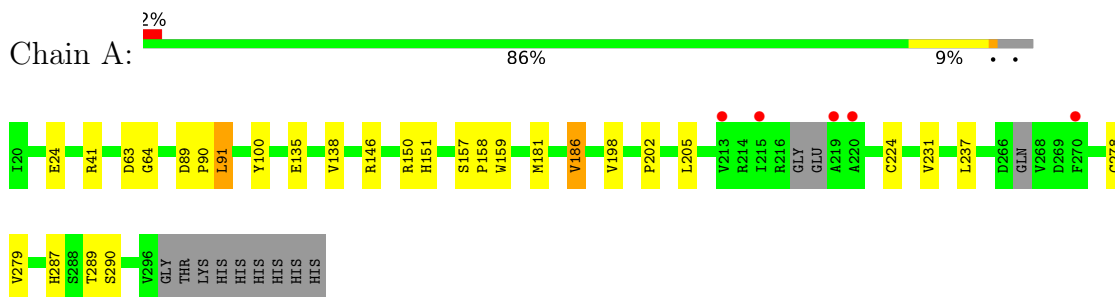


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	A	1	Total	C	N	O	0	0
			14	8	1	5		
7	C	1	Total	C	N	O	0	0
			14	8	1	5		
7	C	1	Total	C	N	O	0	0
			14	8	1	5		

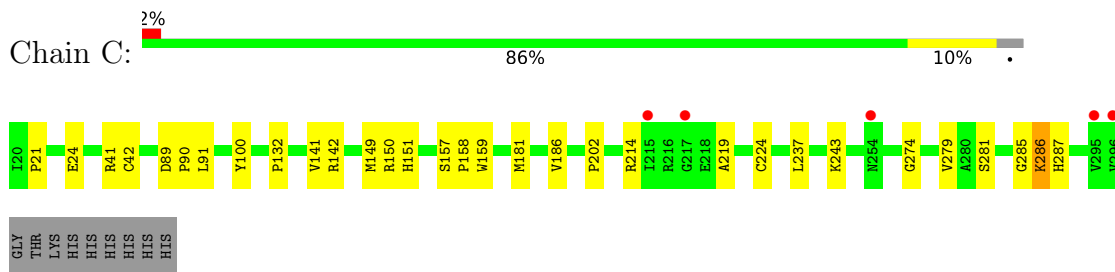
3 Residue-property plots [i](#)

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.

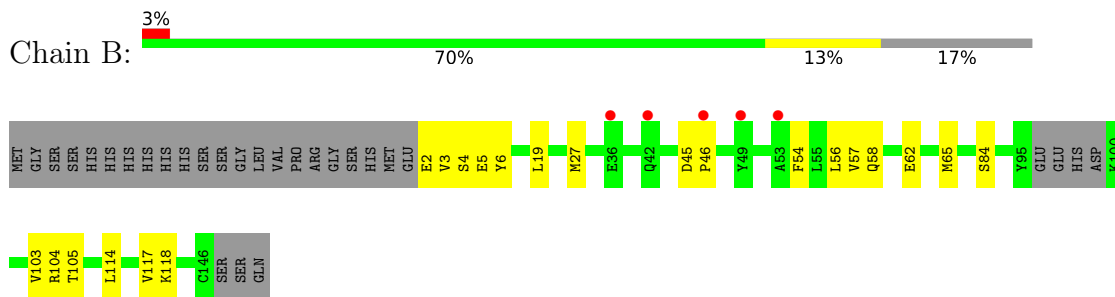
- Molecule 1: Macrophage colony-stimulating factor 1 receptor



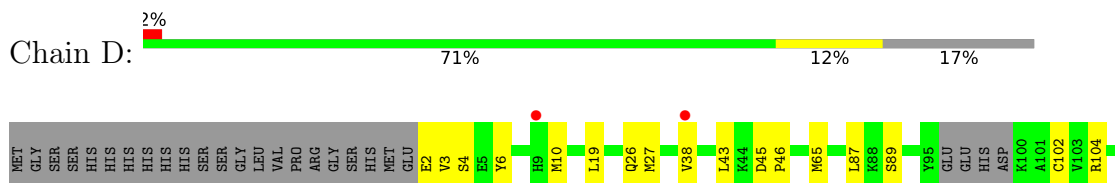
- Molecule 1: Macrophage colony-stimulating factor 1 receptor



- Molecule 2: Macrophage colony-stimulating factor 1



- Molecule 2: Macrophage colony-stimulating factor 1





- Molecule 3: N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 3: N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose



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



- Molecule 5: N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 5: N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



4 Data and refinement statistics

Property	Value	Source
Space group	I 41	Depositor
Cell constants a, b, c, α , β , γ	143.00Å 143.00Å 138.32Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.71 – 2.80 49.71 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.8 (49.71-2.80) 99.9 (49.71-2.80)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.66 (at 2.81Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, R_{free}	0.223 , 0.261 0.233 , 0.269	Depositor DCC
R_{free} test set	1706 reflections (4.73%)	wwPDB-VP
Wilson B-factor (Å ²)	82.2	Xtriage
Anisotropy	0.035	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 53.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.009 for l,-k,h 0.013 for -l,-k,-h 0.012 for -h,-l,-k 0.007 for -h,l,k 0.418 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	6657	wwPDB-VP
Average B, all atoms (Å ²)	100.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.24% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.44	0/2126	0.76	1/2905 (0.0%)
1	C	0.39	0/2152	0.76	1/2942 (0.0%)
2	B	0.29	0/1111	0.68	0/1504
2	D	0.33	0/1118	0.70	0/1513
All	All	0.38	0/6507	0.74	2/8864 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	274	GLY	N-CA-C	6.23	118.35	110.38
1	A	91	LEU	N-CA-C	-5.99	105.92	113.23

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2074	0	1999	18	0
1	C	2099	0	2038	23	0
2	B	1095	0	1018	18	1
2	D	1102	0	1033	15	1
3	E	55	0	47	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	H	55	0	47	11	0
4	F	25	0	21	2	0
5	G	45	0	38	3	0
5	I	45	0	38	3	0
6	A	20	0	17	2	0
7	A	14	0	13	0	0
7	C	28	0	26	1	0
All	All	6657	0	6335	90	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 7.

All (90) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:3:VAL:HG22	2:D:89:SER:O	1.73	0.88
6:A:407:SIA:O1A	4:F:2:GAL:O4	1.92	0.87
1:C:90:PRO:O	1:C:91:LEU:HB2	1.74	0.86
1:C:281:SER:CB	1:C:286:LYS:HG3	2.11	0.81
3:H:3:SIA:H113	3:H:3:SIA:H4	1.61	0.81
1:C:286:LYS:N	1:C:286:LYS:HD2	1.96	0.80
1:A:181:MET:HE2	1:A:186:VAL:HG11	1.66	0.78
1:A:90:PRO:O	1:A:91:LEU:HB2	1.83	0.77
2:B:104:ARG:HD2	2:B:105:THR:N	2.00	0.76
1:C:281:SER:HB3	1:C:286:LYS:HG3	1.68	0.75
3:H:3:SIA:H91	3:H:3:SIA:HN5	1.53	0.73
7:C:405:NAG:H82	7:C:405:NAG:O3	1.91	0.70
3:H:3:SIA:H6	3:H:3:SIA:C11	2.22	0.69
3:H:3:SIA:H91	3:H:3:SIA:N5	2.08	0.69
2:B:104:ARG:NE	2:B:105:THR:O	2.26	0.68
1:A:90:PRO:O	1:A:91:LEU:CB	2.43	0.65
1:C:89:ASP:OD1	1:C:90:PRO:HD2	1.96	0.64
1:C:281:SER:HB2	1:C:286:LYS:HG3	1.81	0.63
1:C:281:SER:HB2	1:C:286:LYS:HE2	1.79	0.63
1:A:89:ASP:OD1	1:A:90:PRO:HD2	1.98	0.63
2:B:104:ARG:HD2	2:B:105:THR:H	1.64	0.62
3:H:3:SIA:H6	3:H:3:SIA:H112	1.84	0.60
1:C:214:ARG:NH1	1:C:219:ALA:O	2.34	0.59
5:I:3:SIA:H6	5:I:3:SIA:O1B	2.02	0.59
3:H:1:NAG:H61	3:H:2:GAL:H2	1.85	0.59
1:A:89:ASP:OD1	1:A:90:PRO:CD	2.51	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:90:PRO:O	1:C:91:LEU:CB	2.46	0.58
1:A:290:SER:OG	5:G:1:NAG:O7	2.14	0.57
1:C:224:CYS:HB2	1:C:237:LEU:HD13	1.88	0.56
1:C:202:PRO:HB2	1:C:287:HIS:HB2	1.88	0.56
2:B:4:SER:HG	2:B:6:TYR:HD2	1.54	0.56
5:G:3:SIA:H4	5:G:3:SIA:O1B	2.06	0.55
3:H:3:SIA:H5	3:H:3:SIA:C9	2.35	0.55
2:B:2:GLU:O	2:B:3:VAL:C	2.48	0.55
1:C:89:ASP:OD1	1:C:90:PRO:CD	2.55	0.55
2:D:4:SER:HG	2:D:6:TYR:HD2	1.56	0.54
3:H:3:SIA:H91	3:H:3:SIA:C5	2.38	0.54
1:C:285:GLY:C	1:C:286:LYS:HD2	2.33	0.53
2:B:56:LEU:HD11	2:B:103:VAL:HG11	1.92	0.52
1:A:224:CYS:HB2	1:A:237:LEU:HD13	1.92	0.52
2:B:114:LEU:HD21	2:D:27:MET:HE1	1.92	0.51
2:B:104:ARG:CD	2:B:105:THR:N	2.73	0.51
1:A:202:PRO:HB2	1:A:287:HIS:HB2	1.93	0.51
2:D:4:SER:OG	2:D:6:TYR:CD2	2.64	0.50
6:A:407:SIA:C1	4:F:2:GAL:O4	2.60	0.50
2:B:104:ARG:CD	2:B:105:THR:O	2.60	0.49
1:C:150:ARG:HG2	1:C:151:HIS:CD2	2.47	0.49
3:H:3:SIA:C9	3:H:3:SIA:C5	2.90	0.49
1:C:100:TYR:HH	1:C:159:TRP:CD1	2.31	0.48
3:E:3:SIA:O4	3:E:3:SIA:O10	2.31	0.48
1:A:205:LEU:HD13	1:A:289:THR:HG22	1.96	0.47
1:A:100:TYR:HH	1:A:159:TRP:CD1	2.32	0.47
2:B:27:MET:HE1	2:D:114:LEU:HD21	1.97	0.47
5:G:2:GAL:H3	5:G:3:SIA:H32	1.35	0.47
1:A:150:ARG:HG2	1:A:151:HIS:CD2	2.50	0.46
2:B:4:SER:OG	2:B:5:GLU:N	2.47	0.46
1:A:89:ASP:HA	1:A:90:PRO:HD3	1.84	0.46
2:D:2:GLU:HB3	2:D:89:SER:HB2	1.97	0.46
1:A:63:ASP:OD1	1:A:64:GLY:N	2.49	0.45
2:D:19:LEU:HD13	2:D:118:LYS:HA	1.99	0.45
1:C:286:LYS:N	1:C:286:LYS:CD	2.73	0.45
2:B:19:LEU:HD13	2:B:118:LYS:HA	1.99	0.45
2:B:65:MET:HE1	2:B:117:VAL:HG21	2.00	0.44
1:C:24:GLU:OE2	1:C:41:ARG:NH1	2.41	0.43
2:D:65:MET:HE1	2:D:117:VAL:HG21	2.00	0.43
2:D:10:MET:HE3	2:D:87:LEU:HD23	2.00	0.43
2:D:26:GLN:HB2	2:D:111:LEU:HD21	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:58:GLN:HE22	2:B:84:SER:HB3	1.83	0.43
5:I:2:GAL:O2	5:I:3:SIA:H31	2.19	0.43
2:B:104:ARG:HD2	2:B:105:THR:O	2.19	0.42
1:C:181:MET:HE2	1:C:186:VAL:HG21	2.01	0.42
1:A:157:SER:HA	1:A:158:PRO:HD3	1.87	0.42
1:C:142:ARG:NH2	2:B:62:GLU:OE1	2.30	0.42
1:C:141:VAL:HG22	1:C:142:ARG:H	1.85	0.42
2:D:43:LEU:HD11	2:D:102:CYS:SG	2.59	0.42
1:C:100:TYR:CD1	1:C:132:PRO:HD3	2.55	0.42
2:B:45:ASP:HA	2:B:46:PRO:HD2	1.93	0.42
1:C:21:PRO:HB2	1:C:42:CYS:SG	2.60	0.41
2:D:38:VAL:HG23	2:D:43:LEU:HD13	2.02	0.41
1:A:24:GLU:OE1	1:A:41:ARG:NH1	2.44	0.41
2:B:54:PHE:O	2:B:57:VAL:HG22	2.20	0.41
1:C:157:SER:HA	1:C:158:PRO:HD3	1.89	0.41
3:H:2:GAL:H3	3:H:3:SIA:O1A	2.07	0.41
1:A:135:GLU:O	1:A:138:VAL:HG22	2.21	0.41
1:A:231:VAL:HG11	2:D:10:MET:SD	2.61	0.41
5:I:3:SIA:O1B	5:I:3:SIA:H4	2.20	0.41
2:D:45:ASP:HA	2:D:46:PRO:HD2	1.93	0.41
1:A:278:CYS:N	1:A:289:THR:O	2.44	0.41
3:H:3:SIA:H113	3:H:3:SIA:H6	1.98	0.41
2:D:2:GLU:O	2:D:2:GLU:HG2	2.20	0.40

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:104:ARG:CG	2:D:104:ARG:NH2[4_554]	2.03	0.17

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	268/286 (94%)	258 (96%)	10 (4%)	0	100	100
1	C	275/286 (96%)	259 (94%)	16 (6%)	0	100	100
2	B	137/170 (81%)	132 (96%)	5 (4%)	0	100	100
2	D	137/170 (81%)	132 (96%)	5 (4%)	0	100	100
All	All	817/912 (90%)	781 (96%)	36 (4%)	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	226/249 (91%)	222 (98%)	4 (2%)	51	82
1	C	230/249 (92%)	226 (98%)	4 (2%)	53	83
2	B	120/160 (75%)	120 (100%)	0	100	100
2	D	122/160 (76%)	122 (100%)	0	100	100
All	All	698/818 (85%)	690 (99%)	8 (1%)	65	87

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	146	ARG
1	A	186	VAL
1	A	198	VAL
1	A	279	VAL
1	C	149	MET
1	C	243	LYS
1	C	279	VAL
1	C	286	LYS

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	77	GLN
1	A	113	GLN
1	A	238	GLN
1	A	248	GLN
1	C	255	ASN
1	C	258	GLN
2	D	42	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

16 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	E	1	1,3	14,14,15	0.81	1 (7%)	17,19,21	1.20	3 (17%)
3	GAL	E	2	3	11,11,12	0.74	0	15,15,17	2.18	3 (20%)
3	SIA	E	3	3	20,20,21	0.55	0	21,28,31	0.83	2 (9%)
3	FUC	E	4	3	10,10,11	0.63	0	14,14,16	0.78	0
4	NAG	F	1	1,4	14,14,15	1.11	1 (7%)	17,19,21	0.92	0
4	GAL	F	2	4	11,11,12	2.36	3 (27%)	15,15,17	1.52	3 (20%)
5	NAG	G	1	1,5	14,14,15	0.45	0	17,19,21	1.54	3 (17%)
5	GAL	G	2	5	11,11,12	0.77	0	15,15,17	1.27	2 (13%)
5	SIA	G	3	5	20,20,21	0.54	0	21,28,31	0.84	2 (9%)
3	NAG	H	1	1,3	14,14,15	0.46	0	17,19,21	1.30	2 (11%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	GAL	H	2	3	11,11,12	1.32	2 (18%)	15,15,17	1.89	3 (20%)
3	SIA	H	3	3	20,20,21	0.55	0	21,28,31	0.83	2 (9%)
3	FUC	H	4	3	10,10,11	0.69	0	14,14,16	0.89	0
5	NAG	I	1	1,5	14,14,15	0.41	0	17,19,21	0.93	1 (5%)
5	GAL	I	2	5	11,11,12	1.26	1 (9%)	15,15,17	1.55	3 (20%)
5	SIA	I	3	5	20,20,21	0.55	0	21,28,31	0.83	2 (9%)

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	E	1	1,3	-	2/6/23/26	0/1/1/1
3	GAL	E	2	3	-	0/2/19/22	0/1/1/1
3	SIA	E	3	3	-	4/18/34/38	0/1/1/1
3	FUC	E	4	3	-	-	0/1/1/1
4	NAG	F	1	1,4	-	2/6/23/26	0/1/1/1
4	GAL	F	2	4	-	0/2/19/22	0/1/1/1
5	NAG	G	1	1,5	-	1/6/23/26	0/1/1/1
5	GAL	G	2	5	-	2/2/19/22	0/1/1/1
5	SIA	G	3	5	-	8/18/34/38	0/1/1/1
3	NAG	H	1	1,3	-	0/6/23/26	0/1/1/1
3	GAL	H	2	3	-	2/2/19/22	0/1/1/1
3	SIA	H	3	3	-	10/18/34/38	0/1/1/1
3	FUC	H	4	3	-	-	0/1/1/1
5	NAG	I	1	1,5	-	2/6/23/26	0/1/1/1
5	GAL	I	2	5	-	0/2/19/22	0/1/1/1
5	SIA	I	3	5	-	8/18/34/38	0/1/1/1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	F	2	GAL	C4-C3	5.40	1.66	1.52
4	F	2	GAL	O3-C3	4.20	1.53	1.43
4	F	1	NAG	O5-C1	-3.77	1.37	1.43
4	F	2	GAL	C2-C3	3.41	1.57	1.52
3	E	1	NAG	C1-C2	2.90	1.56	1.52
5	I	2	GAL	C2-C3	2.89	1.56	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	H	2	GAL	C2-C3	-2.63	1.48	1.52
3	H	2	GAL	O3-C3	2.04	1.48	1.43

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	H	2	GAL	C1-O5-C5	4.91	118.77	112.19
3	E	2	GAL	C1-O5-C5	4.78	118.59	112.19
3	E	2	GAL	C1-C2-C3	4.65	116.41	109.64
3	H	1	NAG	O4-C4-C5	3.97	119.11	109.32
5	I	2	GAL	C1-C2-C3	3.95	115.39	109.64
3	E	2	GAL	O3-C3-C2	-3.85	102.20	110.05
5	G	1	NAG	C1-O5-C5	3.76	117.22	112.19
4	F	2	GAL	O4-C4-C3	3.08	117.63	110.38
4	F	2	GAL	O3-C3-C4	3.00	117.45	110.38
5	G	2	GAL	C1-C2-C3	2.96	113.95	109.64
5	I	2	GAL	O3-C3-C4	-2.92	103.49	110.38
5	G	2	GAL	C1-O5-C5	2.76	115.89	112.19
5	G	1	NAG	O4-C4-C3	2.75	116.87	110.38
5	G	1	NAG	C2-N2-C7	2.55	126.32	122.90
5	I	2	GAL	C1-O5-C5	2.49	115.52	112.19
3	H	2	GAL	O3-C3-C2	2.45	115.05	110.05
3	H	1	NAG	C1-O5-C5	2.42	115.43	112.19
4	F	2	GAL	C2-C3-C4	2.38	115.05	110.86
5	I	1	NAG	C2-N2-C7	2.38	126.09	122.90
3	E	1	NAG	C3-C4-C5	-2.28	106.10	110.23
3	H	2	GAL	C3-C4-C5	-2.24	106.18	110.23
3	H	3	SIA	O1B-C1-C2	2.23	118.52	112.71
5	I	3	SIA	O1B-C1-C2	2.23	118.52	112.71
5	G	3	SIA	O1B-C1-C2	2.22	118.50	112.71
3	E	3	SIA	O1B-C1-C2	2.22	118.47	112.71
3	E	1	NAG	O4-C4-C5	2.07	114.42	109.32
3	E	1	NAG	C1-O5-C5	2.06	114.94	112.19
3	H	3	SIA	O1A-C1-C2	-2.03	118.47	122.85
5	I	3	SIA	O1A-C1-C2	-2.03	118.47	122.85
3	E	3	SIA	O1A-C1-C2	-2.01	118.51	122.85
5	G	3	SIA	O1A-C1-C2	-2.01	118.51	122.85

There are no chirality outliers.

All (41) torsion outliers are listed below:

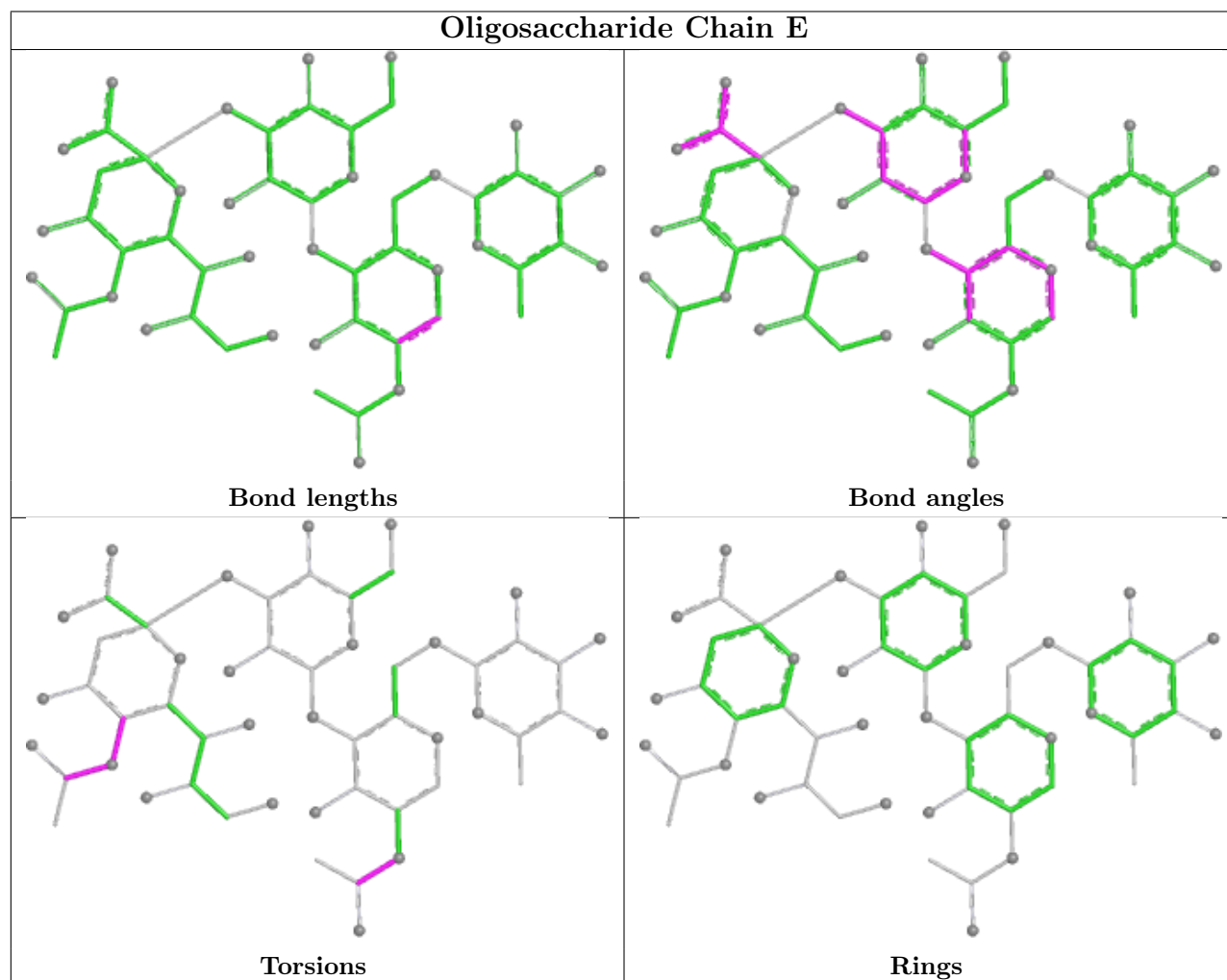
Mol	Chain	Res	Type	Atoms
3	E	3	SIA	C4-C5-N5-C10
3	H	3	SIA	C4-C5-N5-C10
3	H	3	SIA	O8-C8-C9-O9
3	H	3	SIA	C11-C10-N5-C5
3	H	3	SIA	O10-C10-N5-C5
5	G	1	NAG	C3-C2-N2-C7
5	G	3	SIA	O1A-C1-C2-O6
5	G	3	SIA	C5-C6-C7-O7
5	G	3	SIA	O6-C6-C7-O7
5	I	3	SIA	C11-C10-N5-C5
5	I	3	SIA	O10-C10-N5-C5
3	H	3	SIA	C7-C8-C9-O9
3	H	2	GAL	C4-C5-C6-O6
4	F	1	NAG	O5-C5-C6-O6
3	H	2	GAL	O5-C5-C6-O6
3	E	1	NAG	C8-C7-N2-C2
3	E	1	NAG	O7-C7-N2-C2
3	E	3	SIA	C11-C10-N5-C5
4	F	1	NAG	C4-C5-C6-O6
5	G	2	GAL	O5-C5-C6-O6
3	E	3	SIA	O10-C10-N5-C5
5	I	3	SIA	C6-C5-N5-C10
3	E	3	SIA	C6-C5-N5-C10
5	G	3	SIA	C7-C8-C9-O9
5	G	3	SIA	O8-C8-C9-O9
5	I	1	NAG	C3-C2-N2-C7
3	H	3	SIA	C6-C7-C8-O8
5	I	1	NAG	O5-C5-C6-O6
3	H	3	SIA	C6-C7-C8-C9
5	I	3	SIA	C4-C5-N5-C10
5	I	3	SIA	C6-C7-C8-C9
5	G	3	SIA	C5-C6-C7-C8
5	I	3	SIA	O1A-C1-C2-O6
3	H	3	SIA	O7-C7-C8-O8
5	I	3	SIA	O7-C7-C8-C9
5	I	3	SIA	O7-C7-C8-O8
5	G	3	SIA	O1B-C1-C2-O6
3	H	3	SIA	O7-C7-C8-C9
3	H	3	SIA	C6-C5-N5-C10
5	G	2	GAL	C4-C5-C6-O6
5	G	3	SIA	O6-C6-C7-C8

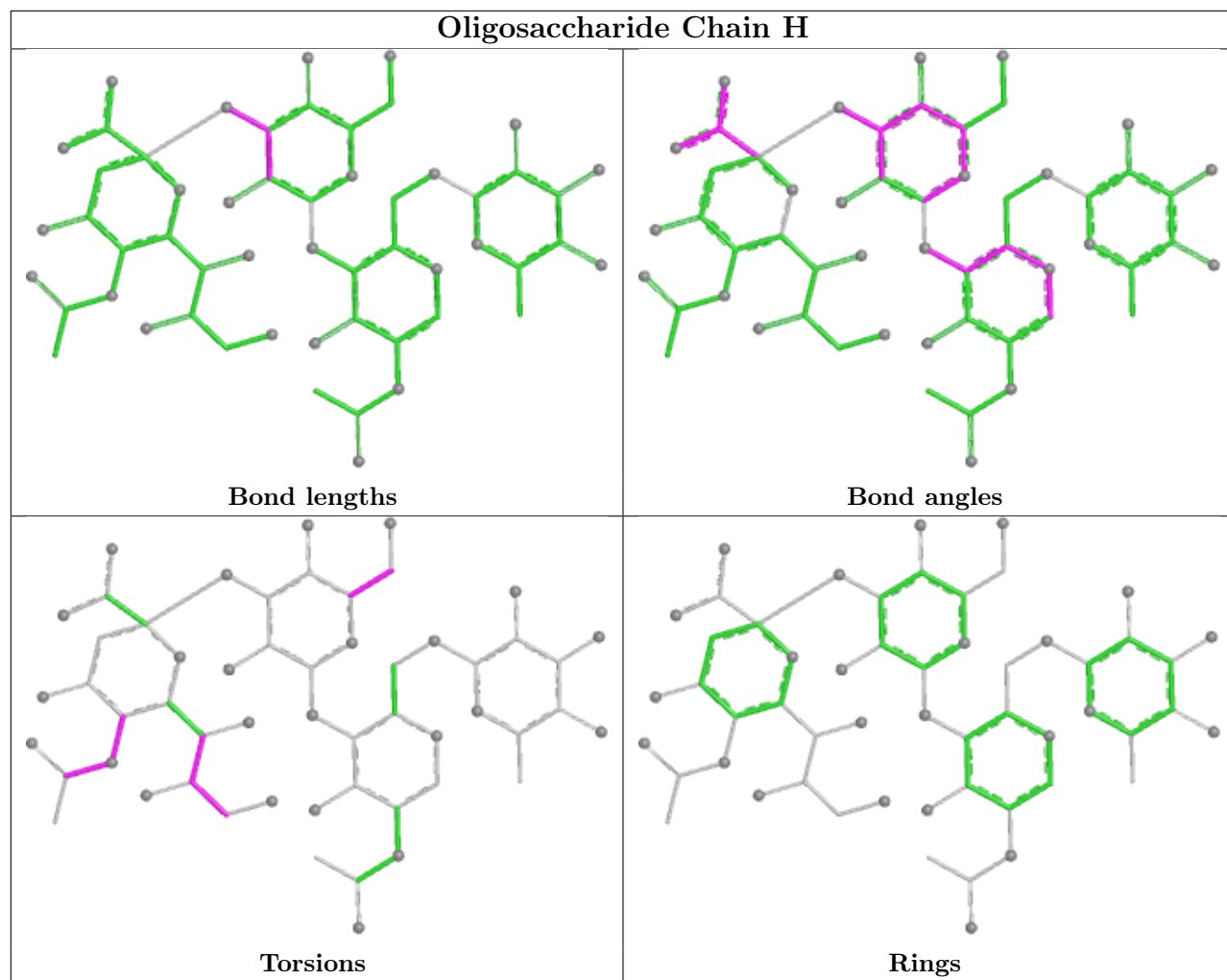
There are no ring outliers.

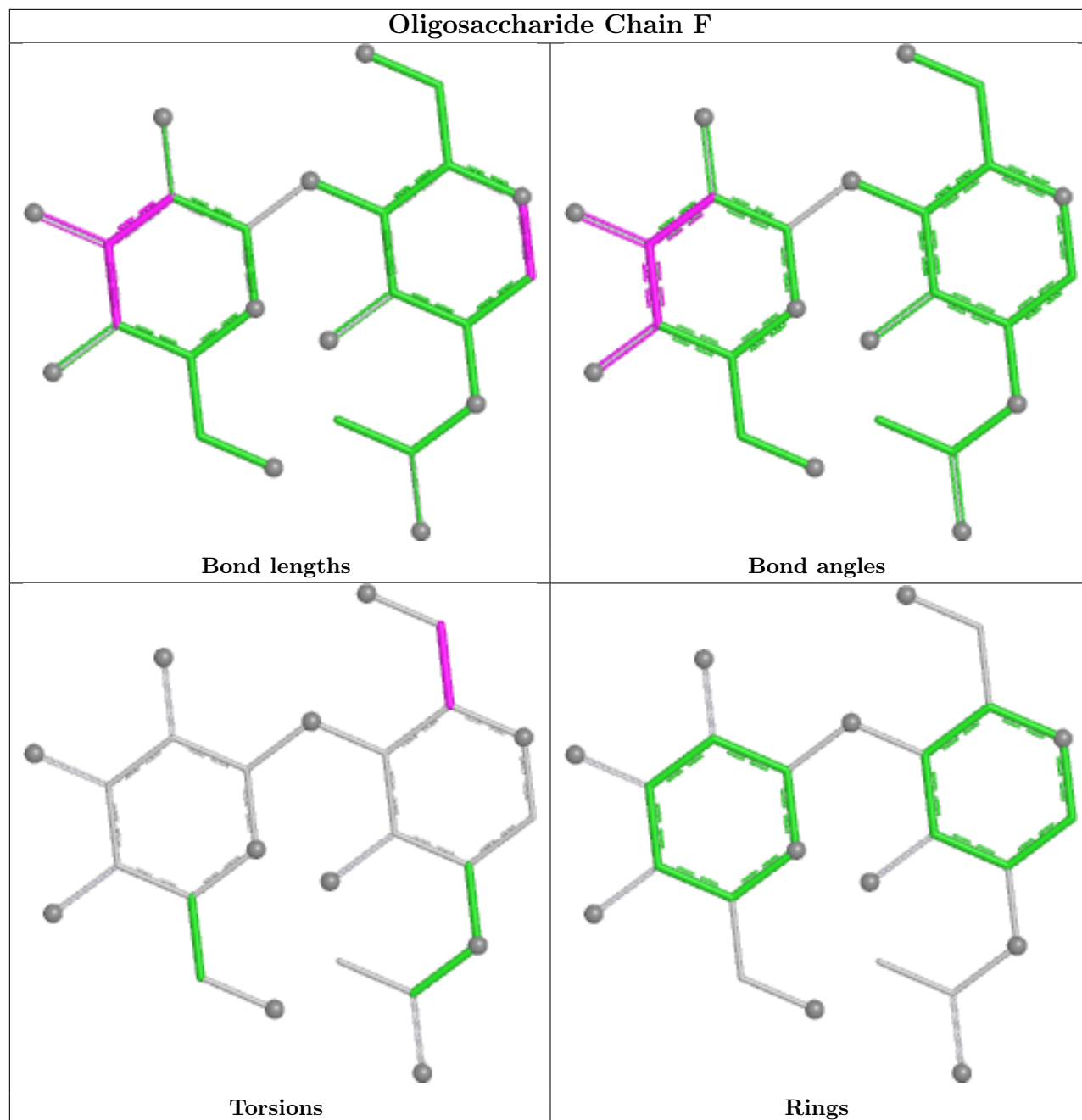
10 monomers are involved in 20 short contacts:

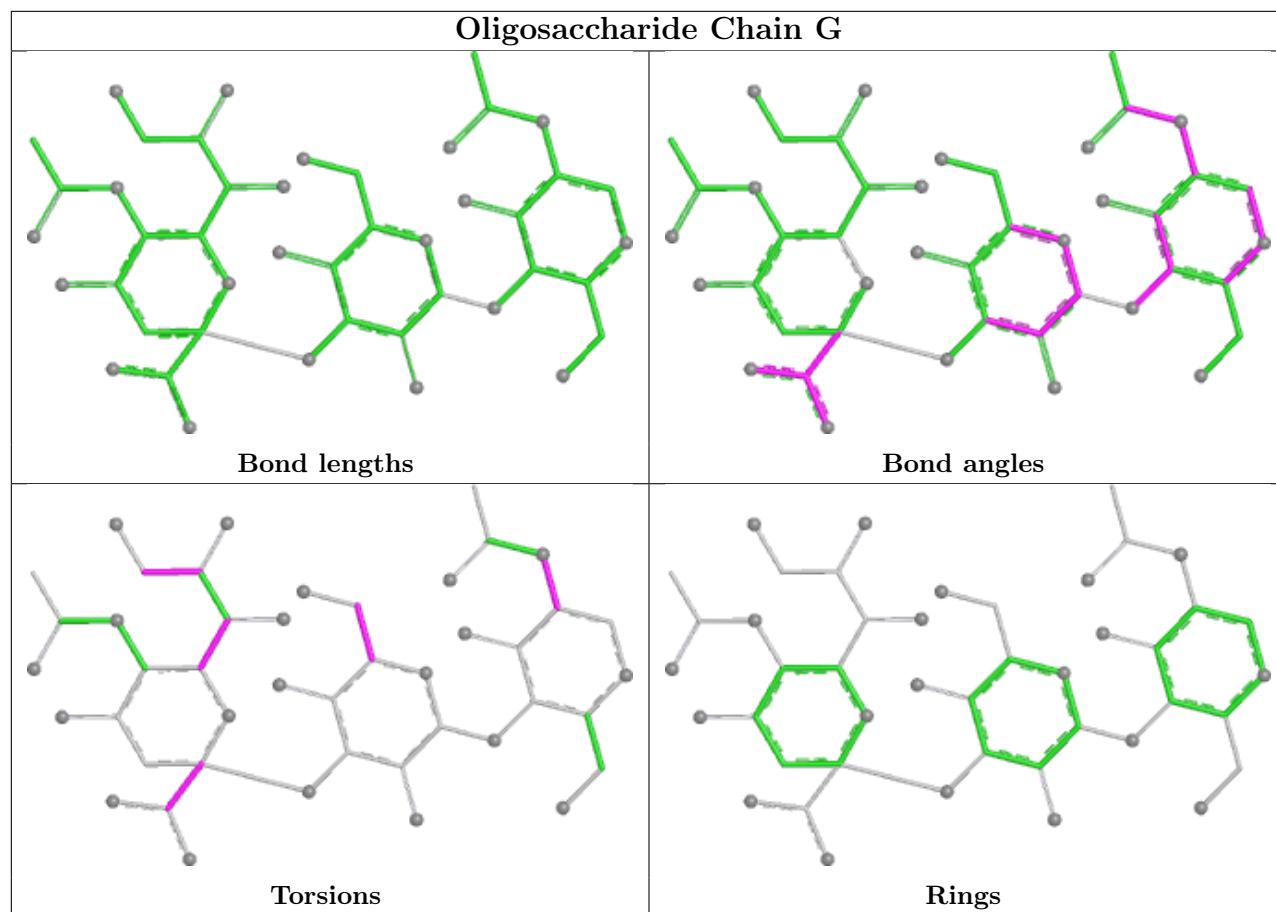
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	H	3	SIA	10	0
5	I	2	GAL	1	0
5	G	1	NAG	1	0
5	G	3	SIA	2	0
4	F	2	GAL	2	0
5	I	3	SIA	3	0
3	E	3	SIA	1	0
3	H	2	GAL	2	0
5	G	2	GAL	1	0
3	H	1	NAG	1	0

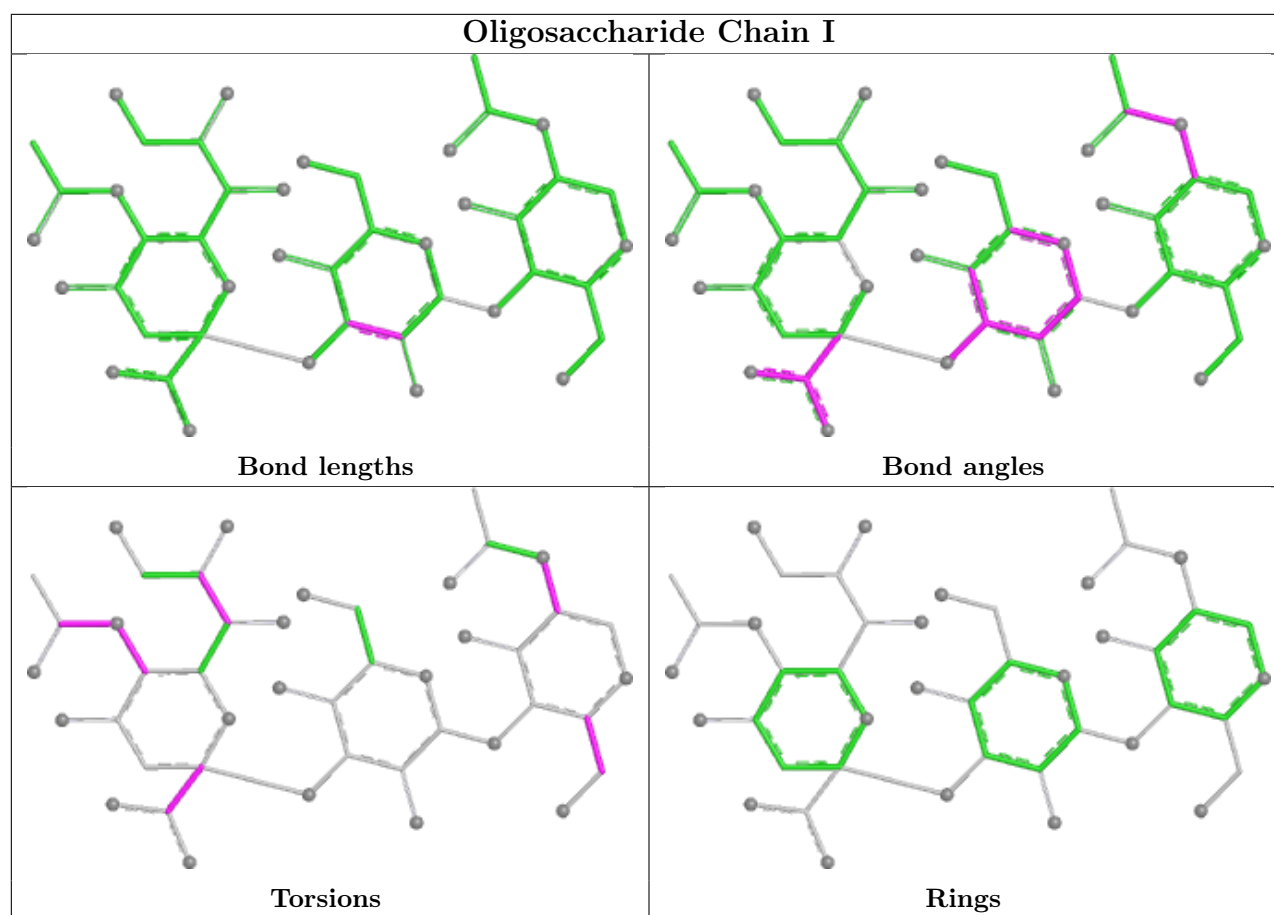
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.











5.6 Ligand geometry [i](#)

4 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	SIA	A	407	-	20,20,21	0.54	0	21,28,31	0.84	2 (9%)
7	NAG	A	408	1	14,14,15	0.34	0	17,19,21	0.47	0
7	NAG	C	405	1	14,14,15	0.35	0	17,19,21	1.50	3 (17%)
7	NAG	C	406	1	14,14,15	0.28	0	17,19,21	0.36	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
6	SIA	A	407	-	-	6/18/34/38	0/1/1/1
7	NAG	A	408	1	-	0/6/23/26	0/1/1/1
7	NAG	C	405	1	-	5/6/23/26	0/1/1/1
7	NAG	C	406	1	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	C	405	NAG	C2-N2-C7	4.15	128.46	122.90
7	C	405	NAG	C4-C3-C2	-2.72	107.04	111.02
6	A	407	SIA	O1B-C1-C2	2.23	118.52	112.71
6	A	407	SIA	O1A-C1-C2	-2.02	118.49	122.85
7	C	405	NAG	O5-C1-C2	-2.01	108.17	111.29

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	407	SIA	C5-C6-C7-C8
6	A	407	SIA	C5-C6-C7-O7
6	A	407	SIA	O6-C6-C7-C8
6	A	407	SIA	O6-C6-C7-O7
7	C	405	NAG	C3-C2-N2-C7
7	C	405	NAG	C8-C7-N2-C2
7	C	405	NAG	O7-C7-N2-C2
6	A	407	SIA	O8-C8-C9-O9
7	C	406	NAG	O5-C5-C6-O6
7	C	406	NAG	C4-C5-C6-O6
7	C	405	NAG	O5-C5-C6-O6
6	A	407	SIA	C7-C8-C9-O9
7	C	405	NAG	C4-C5-C6-O6

There are no ring outliers.

2 monomers are involved in 3 short contacts:

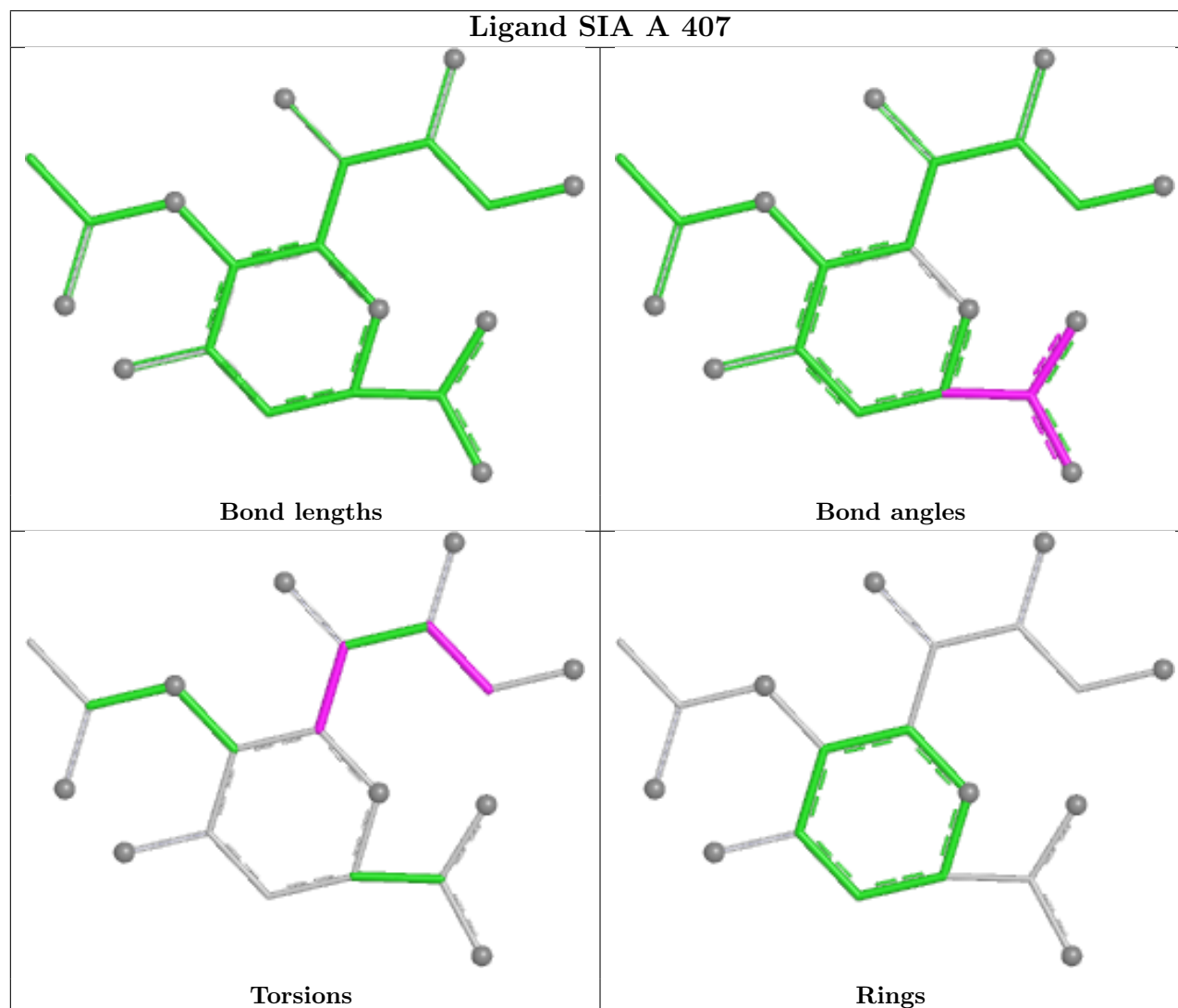
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	407	SIA	2	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	C	405	NAG	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 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 than 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.



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	274/286 (95%)	-0.13	5 (1%) 67 58	48, 83, 146, 188	0
1	C	277/286 (96%)	-0.15	5 (1%) 67 58	44, 84, 137, 168	0
2	B	141/170 (82%)	0.06	5 (3%) 47 38	50, 104, 168, 217	0
2	D	141/170 (82%)	-0.02	3 (2%) 63 54	53, 102, 169, 189	0
All	All	833/912 (91%)	-0.08	18 (2%) 62 52	44, 92, 157, 217	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	270	PHE	4.4
2	D	144	ALA	4.3
1	C	217	GLY	3.8
2	D	9	HIS	3.1
2	B	53	ALA	2.9
1	C	296	VAL	2.8
1	A	215	ILE	2.7
1	A	213	VAL	2.5
2	B	36	GLU	2.4
1	A	219	ALA	2.3
2	B	49	TYR	2.3
1	C	215	ILE	2.2
2	D	38	VAL	2.2
1	C	295	VAL	2.2
2	B	46	PRO	2.2
1	A	220	ALA	2.1
1	C	254	ASN	2.1
2	B	42	GLN	2.0

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

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

6.3 Carbohydrates ⓘ

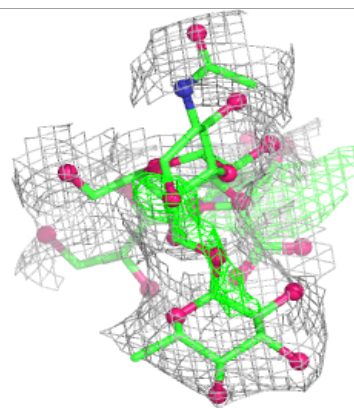
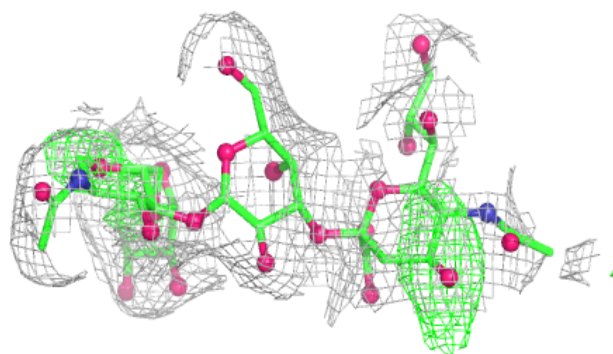
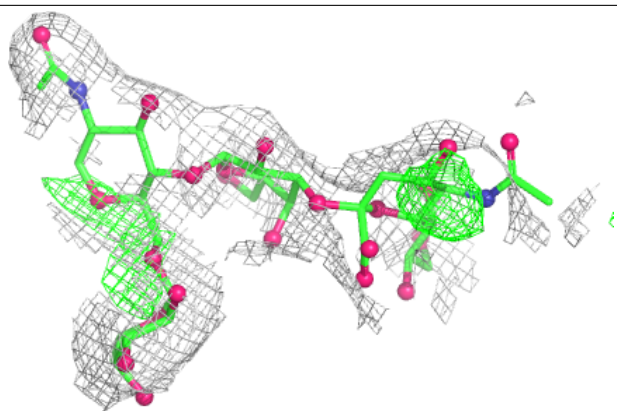
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	SIA	E	3	20/21	0.88	0.12	152,169,185,190	0
5	SIA	I	3	20/21	0.89	0.12	179,210,225,232	0
5	GAL	G	2	11/12	0.90	0.09	184,203,208,208	0
5	NAG	I	1	14/15	0.90	0.12	138,151,169,170	0
3	SIA	H	3	20/21	0.90	0.10	151,180,189,191	0
3	FUC	E	4	10/11	0.91	0.14	124,133,141,145	0
5	GAL	I	2	11/12	0.91	0.13	134,142,186,189	0
5	SIA	G	3	20/21	0.91	0.11	163,182,200,201	0
3	GAL	E	2	11/12	0.93	0.08	130,150,161,173	0
3	GAL	H	2	11/12	0.94	0.11	124,144,161,172	0
3	NAG	E	1	14/15	0.94	0.09	113,120,145,155	0
5	NAG	G	1	14/15	0.94	0.10	146,151,162,179	0
3	NAG	H	1	14/15	0.94	0.11	98,107,134,136	0
3	FUC	H	4	10/11	0.95	0.10	137,147,151,160	0
4	NAG	F	1	14/15	0.95	0.10	130,153,158,166	0
4	GAL	F	2	11/12	0.95	0.08	181,198,224,227	0

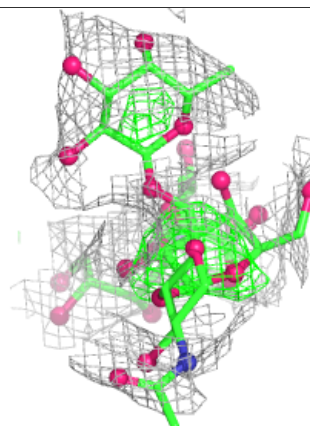
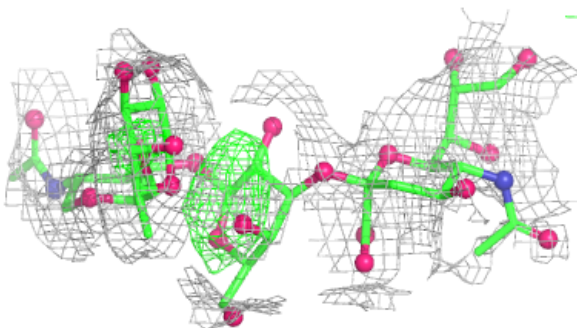
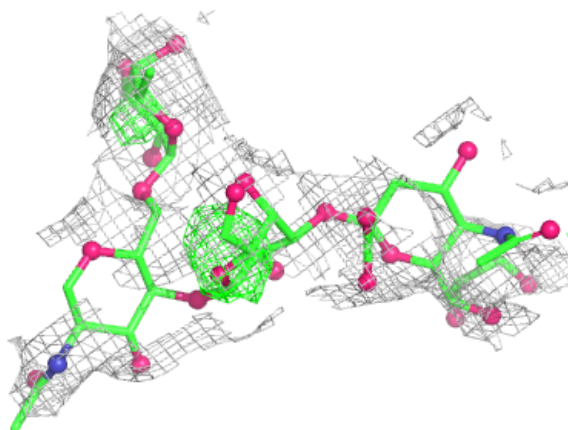
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain E:

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

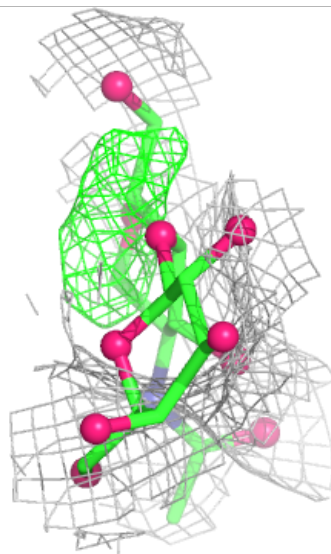
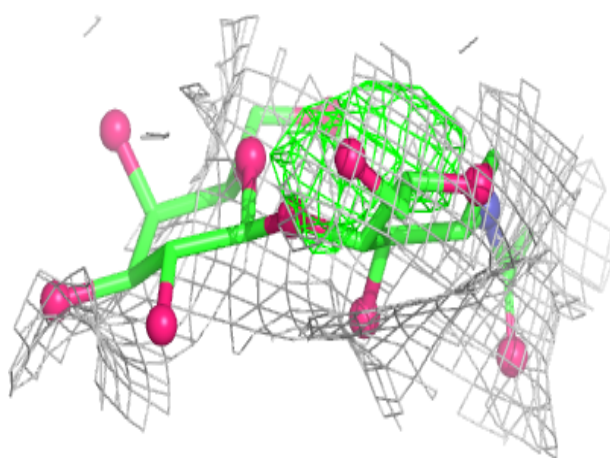
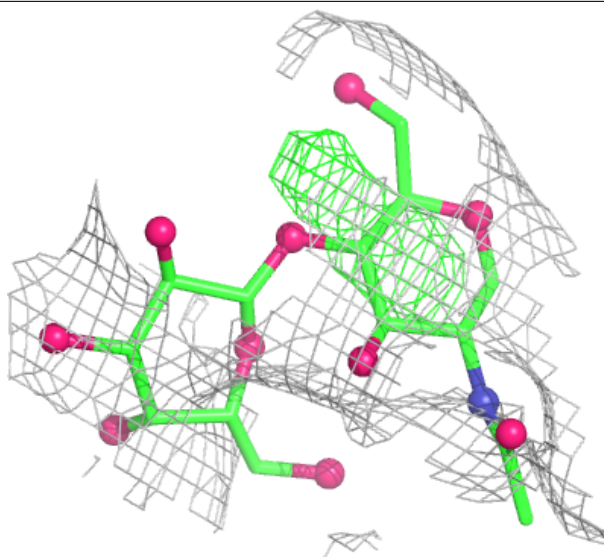
**Electron density around Chain H:**

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



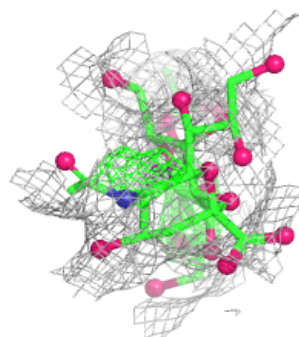
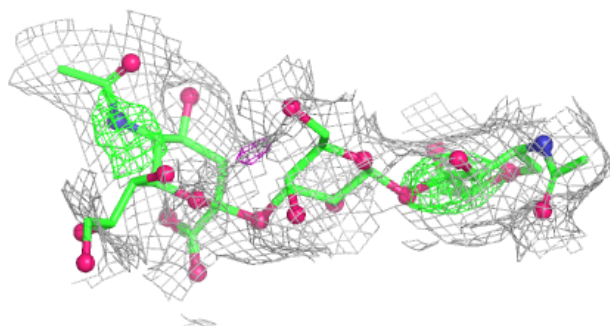
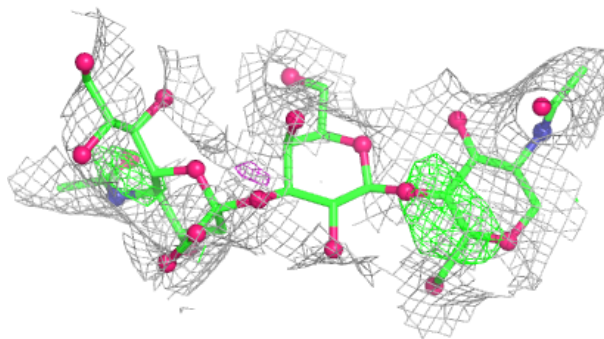
Electron density around Chain F:

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

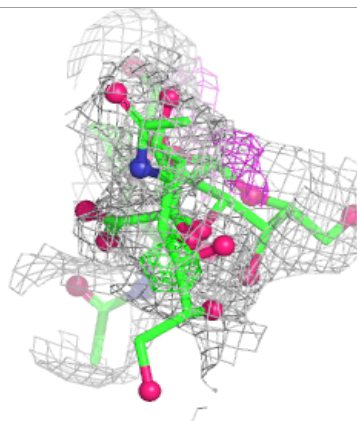
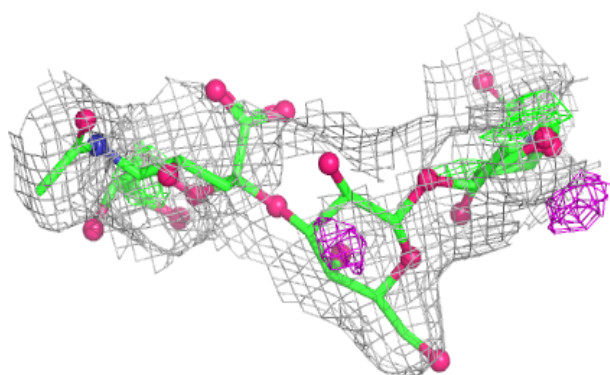
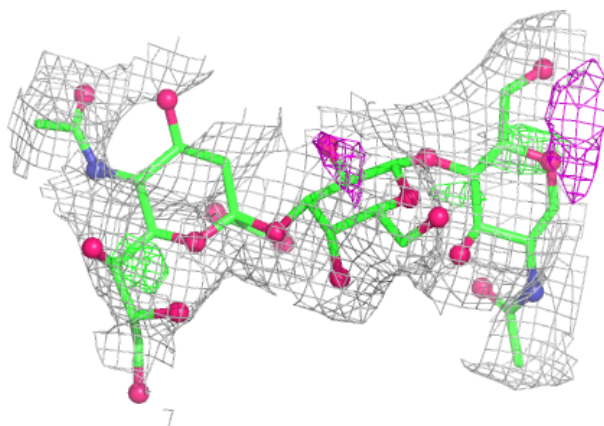


Electron density around Chain G:

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

**Electron density around Chain 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

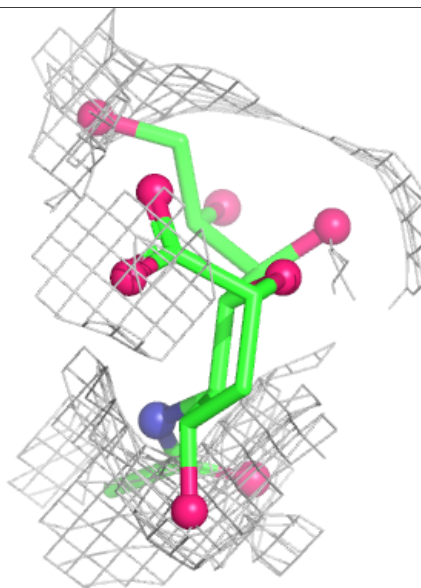
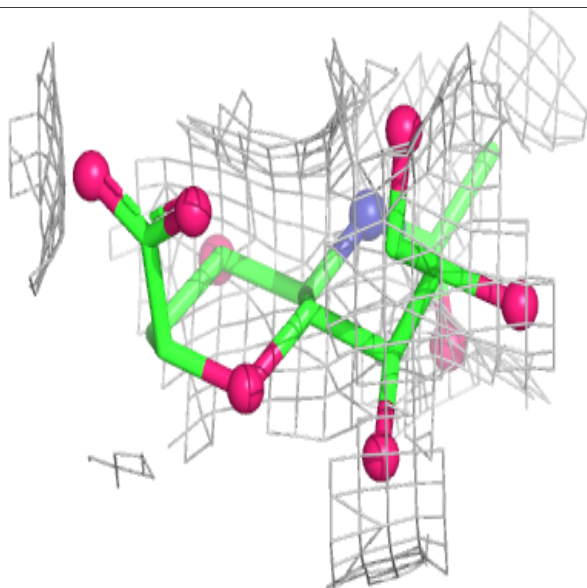
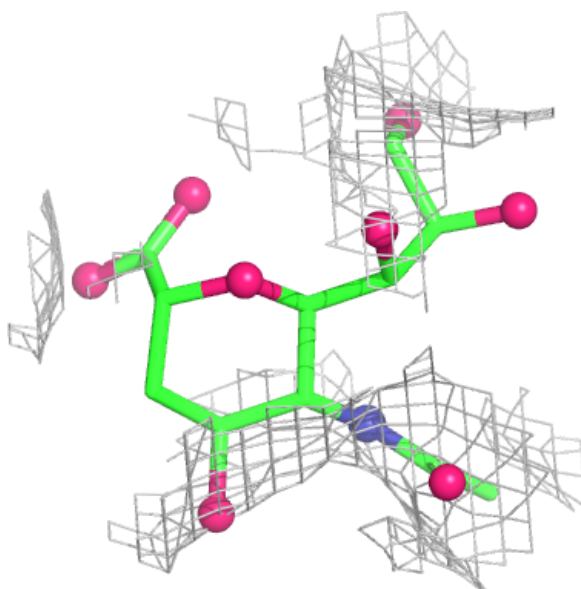
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists 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	SIA	A	407	20/21	0.87	0.09	182,203,218,221	0
7	NAG	A	408	14/15	0.94	0.13	150,162,168,170	0
7	NAG	C	405	14/15	0.94	0.07	145,154,167,175	0
7	NAG	C	406	14/15	0.95	0.11	123,140,142,145	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 A 407:

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.