



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 15, 2026 – 02:00 PM UTC

PDB ID : 4ELM / pdb_00004elm
Title : Crystal structure of the mouse CD1d-lysosulfatide-Hy19.3 TCR complex
Authors : Girardi, E.; Maricic, I.; Wang, J.; Mac, T.T.; Iyer, P.; Kumar, V.; Zajonc, D.M.
Deposited on : 2012-04-11
Resolution : 3.48 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

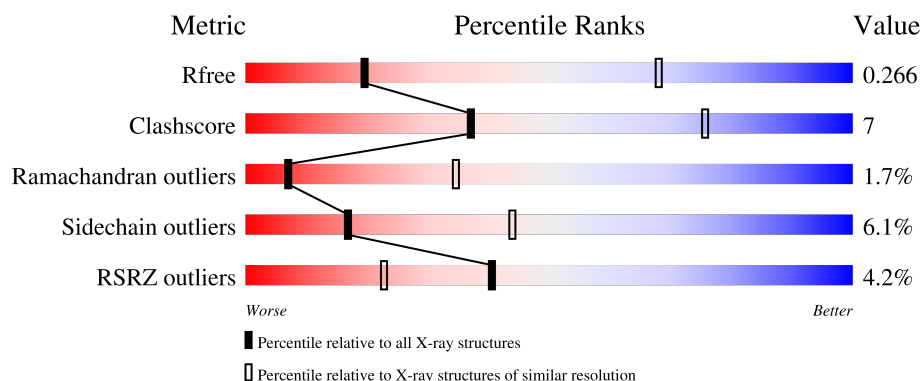
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.48 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	1083 (3.52-3.44)
Clashscore	190562	1139 (3.52-3.44)
Ramachandran outliers	187476	1111 (3.52-3.44)
Sidechain outliers	187428	1112 (3.52-3.44)
RSRZ outliers	180081	1082 (3.52-3.44)

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	285	
1	C	285	
2	B	99	
2	D	99	
3	E	208	

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Mol	Chain	Length	Quality of chain
3	G	208	5% 65% 16% • 18%
4	F	244	8% 77% 13% •• 8%
4	H	244	6% 71% 13% •• 14%
5	I	2	50% 50%
6	J	4	25% 75%

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 11360 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 Antigen-presenting glycoprotein CD1d1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	266	Total	C	N	O	S	0	0	0
			2047	1310	343	382	12			
1	C	269	Total	C	N	O	S	0	0	0
			2093	1340	358	383	12			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	201	HIS	ASP	SEE REMARK 999	UNP P11609
A	280	HIS	-	expression tag	UNP P11609
A	281	HIS	-	expression tag	UNP P11609
A	282	HIS	-	expression tag	UNP P11609
A	283	HIS	-	expression tag	UNP P11609
A	284	HIS	-	expression tag	UNP P11609
A	285	HIS	-	expression tag	UNP P11609
C	201	HIS	ASP	SEE REMARK 999	UNP P11609
C	280	HIS	-	expression tag	UNP P11609
C	281	HIS	-	expression tag	UNP P11609
C	282	HIS	-	expression tag	UNP P11609
C	283	HIS	-	expression tag	UNP P11609
C	284	HIS	-	expression tag	UNP P11609
C	285	HIS	-	expression tag	UNP P11609

- Molecule 2 is a protein called Beta-2 microglobulin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	98	Total	C	N	O	S	0	0	0
			747	471	126	143	7			
2	D	99	Total	C	N	O	S	0	0	0
			777	495	129	146	7			

- Molecule 3 is a protein called Hy19.3 TCR alpha chain (mouse variable domain, human

constant domain).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	179	Total	C	N	O	S	0	0	0
			1234	772	217	239	6			
3	G	171	Total	C	N	O	S	0	0	0
			1185	735	207	236	7			

- Molecule 4 is a protein called Hy19.3 TCR beta chain (mouse variable domain, human constant domain).

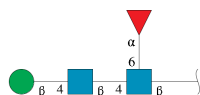
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	F	224	Total	C	N	O	S	0	0	0
			1578	1007	266	296	9			
4	H	211	Total	C	N	O	S	0	0	0
			1459	929	247	274	9			

- Molecule 5 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
5	I	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 6 is an oligosaccharide called beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose.



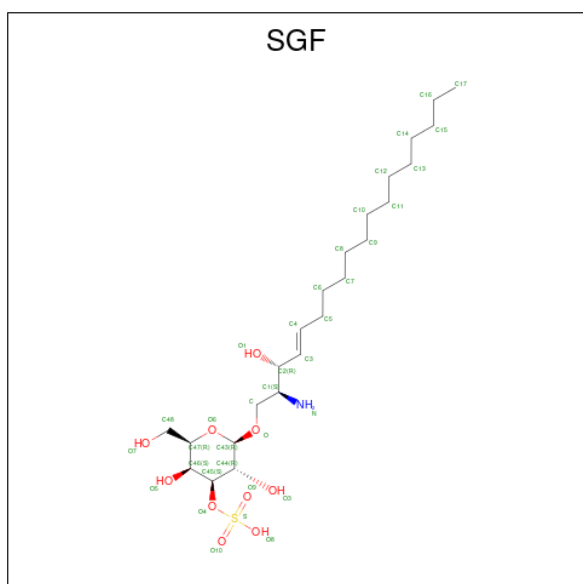
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
6	J	4	Total	C	N	O	0	0	0
			49	28	2	19			

- Molecule 7 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



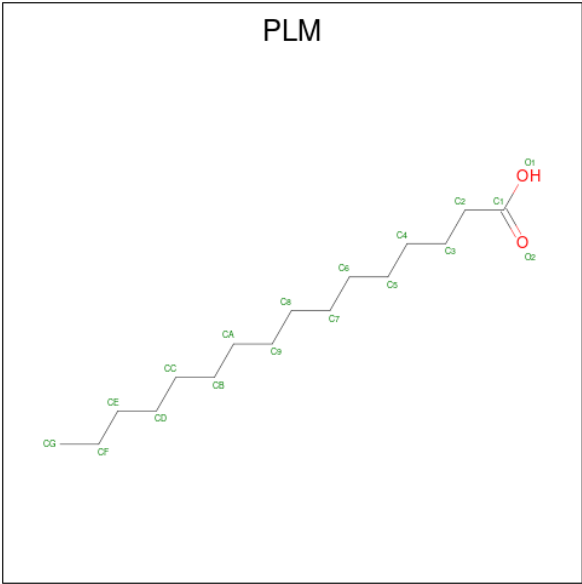
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	A	1	Total	C	N	O	0	0
			13	8	1	4		
7	A	1	Total	C	N	O	0	0
			14	8	1	5		
7	C	1	Total	C	N	O	0	0
			14	8	1	5		
7	C	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 8 is (2S,3R,4E)-2-amino-3-hydroxyoctadec-4-en-1-yl 3-O-sulfo-beta-D-galactopyranoside (CCD ID: SGF) (formula: C₂₄H₄₇NO₁₀S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
8	A	1	Total	C	N	O	S	0	0
			36	24	1	10	1		
8	C	1	Total	C	N	O	S	0	0
			36	24	1	10	1		

- Molecule 9 is PALMITIC ACID (CCD ID: PLM) (formula: C₁₆H₃₂O₂).

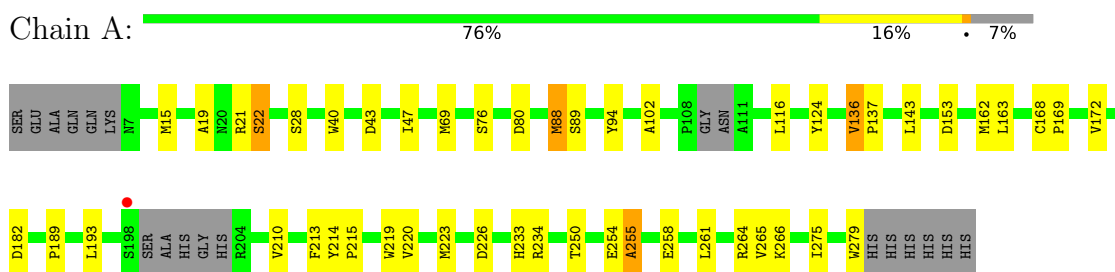


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	A	1	Total	C	O	0	0
			18	16	2		
9	C	1	Total	C	O	0	0
			18	16	2		

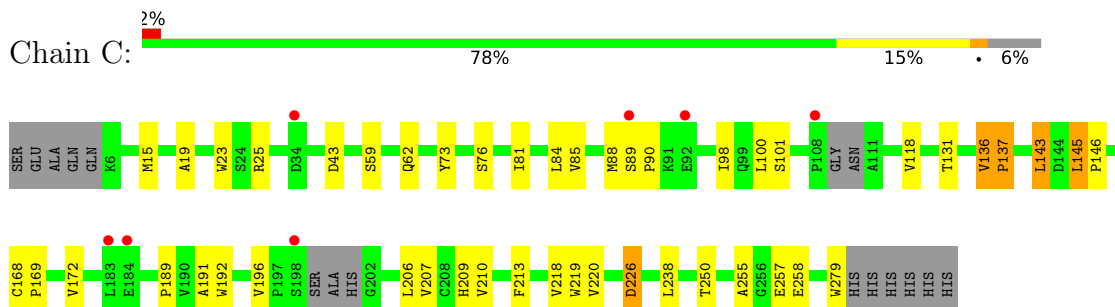
3 Residue-property plots

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.

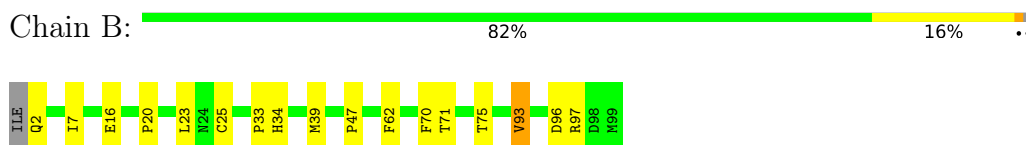
- Molecule 1: Antigen-presenting glycoprotein CD1d1



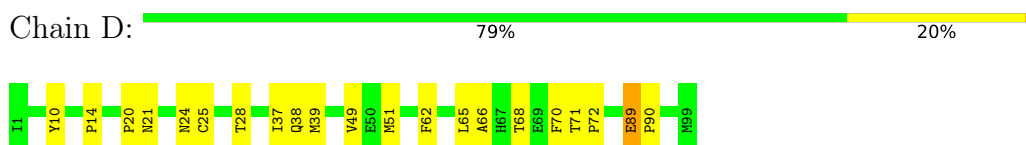
- Molecule 1: Antigen-presenting glycoprotein CD1d1



- Molecule 2: Beta-2 microglobulin

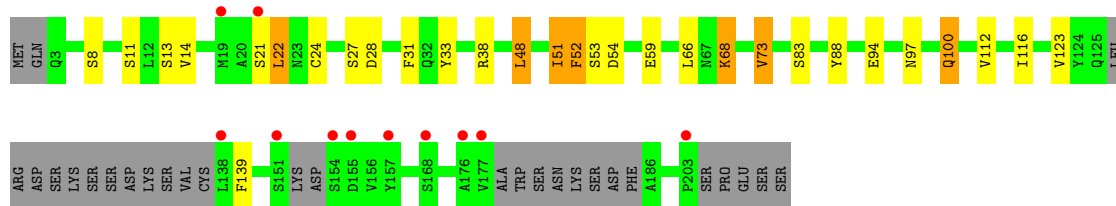


- Molecule 2: Beta-2 microglobulin

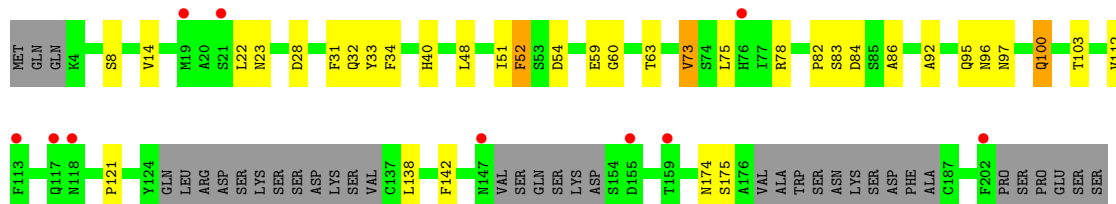


- Molecule 3: Hy19.3 TCR alpha chain (mouse variable domain, human constant domain)

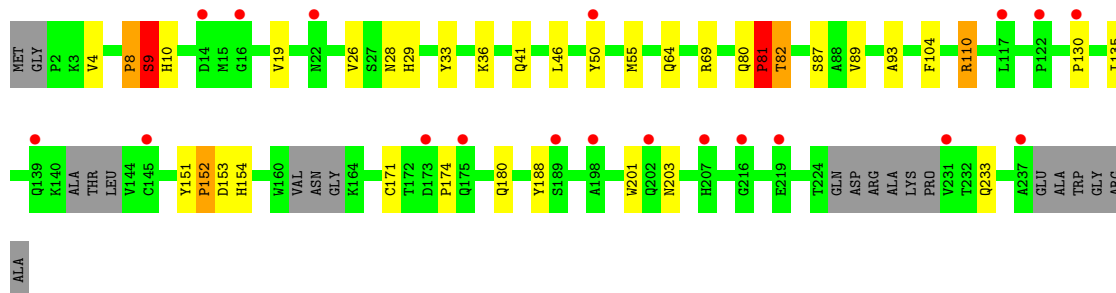




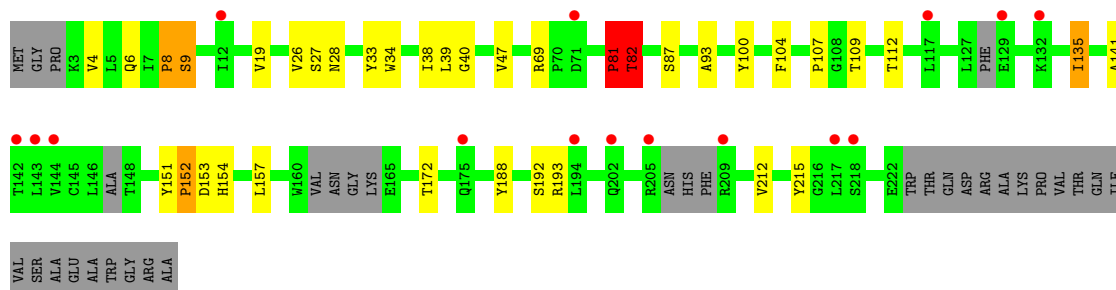
- Molecule 3: Hy19.3 TCR alpha chain (mouse variable domain, human constant domain)



- Molecule 4: Hy19.3 TCR beta chain (mouse variable domain, human constant domain)



- Molecule 4: Hy19.3 TCR beta chain (mouse variable domain, human constant domain)



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





- Molecule 6: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain J:  25% 75%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	98.51Å 127.00Å 104.35Å 90.00° 110.53° 90.00°	Depositor
Resolution (Å)	92.26 – 3.48 92.26 – 3.48	Depositor EDS
% Data completeness (in resolution range)	99.0 (92.26-3.48) 99.0 (92.26-3.48)	Depositor EDS
R_{merge}	0.24	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.63 (at 3.49Å)	Xtriage
Refinement program	REFMAC 5.6.0104	Depositor
R, R_{free}	0.209 , 0.265 0.208 , 0.266	Depositor DCC
R_{free} test set	1561 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	73.3	Xtriage
Anisotropy	0.076	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 70.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	11360	wwPDB-VP
Average B, all atoms (Å ²)	82.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 16.32% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

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

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	0.69	0/2108	0.78	0/2882
1	C	0.70	0/2155	0.80	0/2943
2	B	0.72	0/773	0.77	0/1061
2	D	0.74	0/803	0.82	0/1099
3	E	0.63	0/1263	0.83	1/1727 (0.1%)
3	G	0.63	0/1211	0.81	0/1654
4	F	0.67	0/1623	0.86	6/2230 (0.3%)
4	H	0.66	0/1501	0.86	4/2063 (0.2%)
All	All	0.68	0/11437	0.82	11/15659 (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
4	F	0	1

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	F	81	PRO	CA-C-N	6.83	133.99	121.70
4	F	81	PRO	C-N-CA	6.83	133.99	121.70
3	E	54	ASP	N-CA-C	5.96	120.28	113.19
4	H	8	PRO	CA-C-N	5.96	132.42	121.70
4	H	8	PRO	C-N-CA	5.96	132.42	121.70
4	H	81	PRO	CA-C-N	5.71	131.98	121.70
4	H	81	PRO	C-N-CA	5.71	131.98	121.70
4	F	233	GLN	N-CA-C	5.66	116.03	108.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	F	9	SER	N-CA-C	-5.59	95.36	111.00
4	F	180	GLN	CA-C-N	5.16	124.83	119.56
4	F	180	GLN	C-N-CA	5.16	124.83	119.56

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
4	F	81	PRO	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	2047	0	1853	24	0
1	C	2093	0	1929	29	0
2	B	747	0	631	8	0
2	D	777	0	703	13	0
3	E	1234	0	986	17	0
3	G	1185	0	953	21	0
4	F	1578	0	1247	22	0
4	H	1459	0	1123	20	0
5	I	28	0	25	0	0
6	J	49	0	43	0	0
7	A	27	0	24	0	0
7	C	28	0	26	1	0
8	A	36	0	47	0	0
8	C	36	0	47	5	0
9	A	18	0	31	0	0
9	C	18	0	31	0	0
All	All	11360	0	9699	149	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 7.

All (149) 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:33:TYR:HD1	3:G:52:PHE:HB3	1.11	1.10
3:G:33:TYR:HD1	3:G:52:PHE:CB	1.68	1.07
3:G:33:TYR:CD1	3:G:52:PHE:HB3	1.90	1.06
4:F:110:ARG:HH11	4:F:110:ARG:HG2	1.23	1.01
4:H:8:PRO:HA	4:H:9:SER:HB3	1.50	0.94
3:E:24:CYS:HB3	3:E:73:VAL:HG23	1.56	0.87
4:H:135:ILE:HG12	4:H:141:ALA:HB2	1.59	0.85
4:F:81:PRO:HA	4:F:82:THR:HG23	1.59	0.84
2:D:20:PRO:HA	2:D:71:THR:HG22	1.61	0.82
4:H:8:PRO:CA	4:H:9:SER:HB3	2.09	0.82
2:B:20:PRO:HA	2:B:71:THR:HG22	1.62	0.80
1:C:168:CYS:HB3	1:C:169:PRO:HD3	1.64	0.79
3:G:33:TYR:CD1	3:G:52:PHE:CB	2.59	0.77
1:A:168:CYS:HB3	1:A:169:PRO:HD3	1.72	0.72
4:F:110:ARG:HH11	4:F:110:ARG:CG	2.02	0.70
3:E:97:ASN:HB3	3:E:100:GLN:NE2	2.07	0.70
4:F:50:TYR:HB3	4:F:55:MET:HE3	1.74	0.69
3:E:33:TYR:HB3	3:E:52:PHE:CD1	2.29	0.66
3:G:97:ASN:HB3	3:G:100:GLN:NE2	2.11	0.66
4:F:33:TYR:HB2	4:F:93:ALA:HB3	1.79	0.63
1:C:118:VAL:HG11	8:C:507:SGF:H17	1.79	0.63
4:H:81:PRO:HA	4:H:82:THR:HG23	1.81	0.63
4:F:64:GLN:HB2	4:F:80:GLN:O	1.99	0.62
1:A:189:PRO:HB3	1:A:213:PHE:HB3	1.80	0.62
1:A:219:TRP:HB3	1:A:266:LYS:HB2	1.81	0.62
4:F:110:ARG:HG2	4:F:110:ARG:NH1	2.03	0.60
3:E:33:TYR:HD2	3:E:52:PHE:HA	1.66	0.60
1:A:254:GLU:HG3	1:A:255:ALA:N	2.17	0.59
1:C:84:LEU:O	1:C:88:MET:HG2	2.02	0.59
1:C:89:SER:N	1:C:90:PRO:HA	2.18	0.58
3:E:97:ASN:HB3	3:E:100:GLN:HE21	1.67	0.58
3:G:33:TYR:HD1	3:G:52:PHE:HB2	1.64	0.57
1:C:15:MET:HG2	2:D:62:PHE:HE2	1.68	0.57
3:E:14:VAL:O	3:E:112:VAL:HA	2.05	0.57
1:C:189:PRO:HB3	1:C:213:PHE:HB3	1.86	0.57
4:H:6:GLN:HB2	4:H:107:PRO:HD2	1.87	0.56
1:C:25:ARG:HB3	7:C:502:NAG:H82	1.87	0.56
2:D:21:ASN:HB3	2:D:70:PHE:CE1	2.41	0.55
1:A:254:GLU:HG3	1:A:255:ALA:H	1.71	0.54
1:C:145:LEU:HB3	1:C:146:PRO:HD3	1.89	0.54
4:H:33:TYR:HB2	4:H:93:ALA:HB3	1.90	0.54
4:H:93:ALA:HA	4:H:104:PHE:O	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:F:80:GLN:O	4:F:80:GLN:HG3	2.08	0.53
2:D:37:ILE:HB	2:D:51:MET:HE1	1.91	0.53
1:A:168:CYS:O	1:A:172:VAL:HG23	2.09	0.53
1:C:73:TYR:HE1	8:C:507:SGF:H26	1.74	0.52
4:H:112:THR:HG21	4:H:152:PRO:HG2	1.91	0.52
1:C:168:CYS:O	1:C:172:VAL:HG23	2.10	0.52
4:F:153:ASP:HB3	4:F:188:TYR:CD1	2.44	0.52
1:C:136:VAL:HG22	1:C:137:PRO:HD2	1.91	0.51
2:B:20:PRO:CA	2:B:71:THR:HG22	2.36	0.51
3:E:123:VAL:HG22	3:E:139:PHE:HD1	1.76	0.51
3:E:66:LEU:HD21	3:E:68:LYS:HD2	1.91	0.51
4:F:110:ARG:CG	4:F:110:ARG:NH1	2.67	0.51
2:D:25:CYS:HB2	2:D:39:MET:HE3	1.92	0.50
1:C:207:VAL:HG22	1:C:250:THR:HG22	1.92	0.50
1:A:69:MET:HE1	1:A:163:LEU:HD11	1.93	0.50
1:C:218:VAL:HG22	1:C:219:TRP:H	1.77	0.50
1:A:88:MET:HA	1:A:88:MET:HE3	1.94	0.50
3:G:138:LEU:HD11	3:G:175:SER:HB2	1.94	0.50
1:A:233:HIS:HB2	1:A:250:THR:OG1	2.12	0.50
1:A:162:MET:HE3	4:H:100:TYR:CD1	2.47	0.49
1:C:19:ALA:HB3	1:C:23:TRP:HB3	1.95	0.49
3:G:14:VAL:O	3:G:112:VAL:HA	2.13	0.49
4:H:27:SER:O	4:H:28:ASN:HB2	2.13	0.49
1:A:265:VAL:HB	1:A:275:ILE:HB	1.93	0.49
1:A:193:LEU:HD23	1:A:279:TRP:HE3	1.77	0.49
4:F:130:PRO:HG2	4:F:201:TRP:CE2	2.48	0.49
1:A:40:TRP:CG	1:A:47:ILE:HG13	2.48	0.48
4:H:9:SER:H	4:H:109:THR:HG23	1.78	0.48
1:A:21:ARG:O	1:A:22:SER:CB	2.62	0.48
3:G:40:HIS:CD2	3:G:86:ALA:HB2	2.48	0.48
2:D:25:CYS:HB3	2:D:66:ALA:HB3	1.96	0.48
4:H:8:PRO:CB	4:H:9:SER:HB3	2.44	0.47
4:F:36:LYS:HB2	4:F:46:LEU:HD11	1.96	0.47
3:E:123:VAL:HG22	3:E:139:PHE:CD1	2.49	0.47
1:C:143:LEU:HD21	8:C:507:SGF:H11	1.97	0.47
3:G:8:SER:HB2	3:G:23:ASN:HB2	1.95	0.47
1:C:59:SER:OG	1:C:62:GLN:HG3	2.14	0.47
1:C:258:GLU:HB3	1:C:279:TRP:CD1	2.50	0.47
2:B:96:ASP:O	2:B:97:ARG:HB3	2.14	0.47
1:A:15:MET:HG2	2:B:62:PHE:HE2	1.79	0.47
4:H:157:LEU:HA	4:H:212:VAL:HA	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:33:PRO:HD3	2:B:62:PHE:CE1	2.49	0.46
4:F:81:PRO:HA	4:F:82:THR:CG2	2.36	0.46
4:H:38:ILE:C	4:H:40:GLY:H	2.23	0.46
1:C:136:VAL:CG2	1:C:137:PRO:HD2	2.46	0.46
4:F:151:TYR:CD1	4:F:152:PRO:HB3	2.51	0.46
3:G:28:ASP:HB3	3:G:31:PHE:HD2	1.79	0.46
2:D:89:GLU:HG3	2:D:90:PRO:HD2	1.98	0.46
1:C:98:ILE:HD11	8:C:507:SGF:H9	1.98	0.45
3:E:48:LEU:HD23	3:E:48:LEU:HA	1.76	0.45
4:F:201:TRP:C	4:F:203:ASN:H	2.25	0.45
3:G:60:GLY:O	3:G:78:ARG:NH1	2.50	0.45
1:C:218:VAL:HG22	1:C:219:TRP:N	2.32	0.45
1:A:223:MET:HE2	1:A:264:ARG:HE	1.82	0.45
1:C:191:ALA:HA	1:C:209:HIS:O	2.16	0.45
1:C:192:TRP:CE3	2:D:14:PRO:HG3	2.51	0.45
1:A:258:GLU:HA	1:A:261:LEU:HD12	1.99	0.45
2:D:20:PRO:CA	2:D:71:THR:HG22	2.42	0.45
4:H:153:ASP:HB3	4:H:188:TYR:CD1	2.52	0.45
1:A:19:ALA:O	1:A:94:TYR:HB3	2.16	0.45
1:C:145:LEU:HB3	1:C:146:PRO:CD	2.47	0.44
3:G:33:TYR:CD1	3:G:52:PHE:HB2	2.45	0.44
4:F:93:ALA:HA	4:F:104:PHE:O	2.18	0.44
1:C:81:ILE:O	1:C:85:VAL:HG23	2.17	0.44
1:C:226:ASP:OD1	1:C:226:ASP:N	2.45	0.44
4:H:151:TYR:CD1	4:H:152:PRO:HB3	2.53	0.44
3:G:31:PHE:HB3	3:G:92:ALA:HB1	1.99	0.44
1:C:76:SER:HB3	8:C:507:SGF:H32	1.99	0.43
1:A:124:TYR:CZ	1:A:136:VAL:HG21	2.53	0.43
3:G:142:PHE:CE2	3:G:174:ASN:HB3	2.54	0.43
3:E:94:GLU:HB3	3:E:97:ASN:HD22	1.83	0.43
3:G:48:LEU:HD13	3:G:75:LEU:HD21	1.99	0.43
1:A:214:TYR:CG	1:A:215:PRO:HA	2.54	0.43
1:C:168:CYS:HB3	1:C:169:PRO:CD	2.43	0.43
2:B:7:ILE:HB	2:B:93:VAL:HG11	2.01	0.43
3:G:33:TYR:HB3	3:G:52:PHE:CD1	2.54	0.43
3:G:121:PRO:HB3	3:G:142:PHE:HB3	2.01	0.43
3:E:38:ARG:HD3	3:E:88:TYR:CE2	2.53	0.43
4:H:135:ILE:H	4:H:135:ILE:HG13	1.66	0.42
3:E:51:ILE:HB	3:E:66:LEU:HD22	2.00	0.42
4:H:172:THR:HG23	4:H:192:SER:HB2	2.02	0.42
4:F:89:VAL:HG22	4:F:110:ARG:HG3	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:210:VAL:HG21	1:A:220:VAL:HG21	2.02	0.42
1:A:124:TYR:OH	1:A:136:VAL:HG21	2.20	0.42
4:F:8:PRO:CA	4:F:9:SER:HB3	2.50	0.42
4:H:154:HIS:ND1	4:H:215:TYR:HB2	2.34	0.42
2:B:25:CYS:HB2	2:B:39:MET:HE3	2.02	0.41
3:G:95:GLN:O	3:G:96:ASN:HB2	2.20	0.41
4:F:10:HIS:HB3	4:F:154:HIS:ND1	2.35	0.41
4:F:151:TYR:HA	4:F:152:PRO:HA	1.95	0.41
3:E:28:ASP:HB3	3:E:31:PHE:HD2	1.84	0.41
3:G:82:PRO:C	3:G:84:ASP:H	2.28	0.41
1:A:102:ALA:HB2	1:A:116:LEU:HG	2.02	0.41
1:C:210:VAL:HG21	1:C:220:VAL:HG21	2.03	0.41
3:G:34:PHE:CG	3:G:73:VAL:HG21	2.54	0.41
3:E:66:LEU:HD13	3:E:73:VAL:HG13	2.02	0.41
2:D:49:VAL:HG22	2:D:68:THR:HB	2.03	0.41
2:D:71:THR:HA	2:D:72:PRO:HD2	1.92	0.41
3:E:21:SER:C	3:E:22:LEU:HD23	2.45	0.41
3:E:51:ILE:C	3:E:53:SER:H	2.29	0.41
4:F:26:VAL:O	4:F:29:HIS:HB2	2.21	0.40
1:C:238:LEU:HB3	2:D:10:TYR:CZ	2.55	0.40
2:D:24:ASN:HB3	2:D:65:LEU:HD11	2.04	0.40
1:A:153:ASP:OD1	1:A:153:ASP:C	2.64	0.40
2:B:23:LEU:HB2	2:B:70:PHE:CD1	2.57	0.40
4:F:8:PRO:HA	4:F:9:SER:HB3	2.03	0.40
4:H:34:TRP:HB2	4:H:47:VAL:HG12	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	260/285 (91%)	252 (97%)	3 (1%)	5 (2%)	6	32

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	263/285 (92%)	252 (96%)	9 (3%)	2 (1%)	16	49
2	B	96/99 (97%)	86 (90%)	9 (9%)	1 (1%)	12	44
2	D	97/99 (98%)	91 (94%)	6 (6%)	0	100	100
3	E	171/208 (82%)	146 (85%)	23 (14%)	2 (1%)	10	41
3	G	163/208 (78%)	142 (87%)	20 (12%)	1 (1%)	21	53
4	F	216/244 (88%)	181 (84%)	27 (12%)	8 (4%)	2	21
4	H	201/244 (82%)	173 (86%)	22 (11%)	6 (3%)	3	25
All	All	1467/1672 (88%)	1323 (90%)	119 (8%)	25 (2%)	7	34

All (25) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	22	SER
4	F	82	THR
4	F	152	PRO
4	H	81	PRO
4	H	152	PRO
4	F	9	SER
4	H	9	SER
1	A	88	MET
1	A	255	ALA
1	C	255	ALA
4	F	41	GLN
4	F	87	SER
4	H	39	LEU
4	H	87	SER
3	E	52	PHE
4	F	8	PRO
4	F	28	ASN
4	H	82	THR
1	A	137	PRO
1	C	137	PRO
1	A	89	SER
4	F	174	PRO
3	G	83	SER
3	E	116	ILE
2	B	47	PRO

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	208/249 (84%)	199 (96%)	9 (4%)	26	52
1	C	215/249 (86%)	204 (95%)	11 (5%)	21	48
2	B	75/93 (81%)	70 (93%)	5 (7%)	15	41
2	D	83/93 (89%)	80 (96%)	3 (4%)	31	57
3	E	104/185 (56%)	92 (88%)	12 (12%)	5	24
3	G	104/185 (56%)	94 (90%)	10 (10%)	8	30
4	F	131/217 (60%)	125 (95%)	6 (5%)	24	50
4	H	118/217 (54%)	111 (94%)	7 (6%)	18	45
All	All	1038/1488 (70%)	975 (94%)	63 (6%)	17	44

All (63) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	SER
1	A	43	ASP
1	A	76	SER
1	A	80	ASP
1	A	136	VAL
1	A	143	LEU
1	A	182	ASP
1	A	226	ASP
1	A	234	ARG
1	C	43	ASP
1	C	100	LEU
1	C	101	SER
1	C	131	THR
1	C	136	VAL
1	C	143	LEU
1	C	145	LEU
1	C	196	VAL
1	C	206	LEU
1	C	226	ASP

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Mol	Chain	Res	Type
1	C	257	GLU
2	B	2	GLN
2	B	16	GLU
2	B	34	HIS
2	B	75	THR
2	B	93	VAL
2	D	28	THR
2	D	38	GLN
2	D	89	GLU
3	E	8	SER
3	E	11	SER
3	E	13	SER
3	E	22	LEU
3	E	27	SER
3	E	48	LEU
3	E	51	ILE
3	E	59	GLU
3	E	68	LYS
3	E	73	VAL
3	E	83	SER
3	E	100	GLN
4	F	4	VAL
4	F	19	VAL
4	F	69	ARG
4	F	110	ARG
4	F	135	ILE
4	F	171	CYS
3	G	22	LEU
3	G	32	GLN
3	G	51	ILE
3	G	52	PHE
3	G	54	ASP
3	G	59	GLU
3	G	63	THR
3	G	73	VAL
3	G	100	GLN
3	G	103	THR
4	H	4	VAL
4	H	19	VAL
4	H	26	VAL
4	H	69	ARG
4	H	82	THR

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Mol	Chain	Res	Type
4	H	135	ILE
4	H	193	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	7	ASN
1	A	62	GLN
1	A	117	HIS
1	A	186	GLN
1	A	248	GLN
1	C	7	ASN
1	C	60	ASN
1	C	154	GLN
2	B	8	GLN
2	D	13	HIS
3	E	39	GLN
3	E	40	HIS
4	F	6	GLN
4	F	10	HIS
4	F	22	ASN
4	F	37	GLN
3	G	30	ASN
3	G	40	HIS
3	G	76	HIS
4	H	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 ⓘ

6 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	$\# Z > 2$
5	NAG	I	1	1,5	14,14,15	0.56	0	17,19,21	0.70	0
5	NAG	I	2	5	14,14,15	0.53	0	17,19,21	1.94	3 (17%)
6	NAG	J	1	1,6	14,14,15	0.60	0	17,19,21	0.79	0
6	NAG	J	2	6	14,14,15	0.62	0	17,19,21	2.29	5 (29%)
6	BMA	J	3	6	11,11,12	0.53	0	15,15,17	1.15	3 (20%)
6	FUC	J	4	6	10,10,11	0.60	0	14,14,16	0.89	1 (7%)

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
5	NAG	I	1	1,5	-	1/6/23/26	0/1/1/1
5	NAG	I	2	5	-	4/6/23/26	0/1/1/1
6	NAG	J	1	1,6	-	1/6/23/26	0/1/1/1
6	NAG	J	2	6	-	5/6/23/26	0/1/1/1
6	BMA	J	3	6	-	2/2/19/22	0/1/1/1
6	FUC	J	4	6	-	-	0/1/1/1

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	J	2	NAG	C2-N2-C7	5.29	129.99	122.90
5	I	2	NAG	C1-O5-C5	5.16	119.10	112.19
6	J	2	NAG	C4-C3-C2	3.64	116.35	111.02
6	J	2	NAG	C8-C7-N2	3.59	122.06	116.12
6	J	2	NAG	C3-C4-C5	3.45	116.48	110.23
5	I	2	NAG	C2-N2-C7	3.17	127.15	122.90
5	I	2	NAG	C3-C4-C5	3.06	115.78	110.23
6	J	3	BMA	C3-C4-C5	2.43	114.64	110.23
6	J	3	BMA	C2-C3-C4	2.30	114.90	110.86

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	J	4	FUC	O5-C5-C6	2.24	112.26	107.40
6	J	2	NAG	O5-C1-C2	-2.16	107.94	111.29
6	J	3	BMA	C1-O5-C5	-2.11	109.35	112.19

There are no chirality outliers.

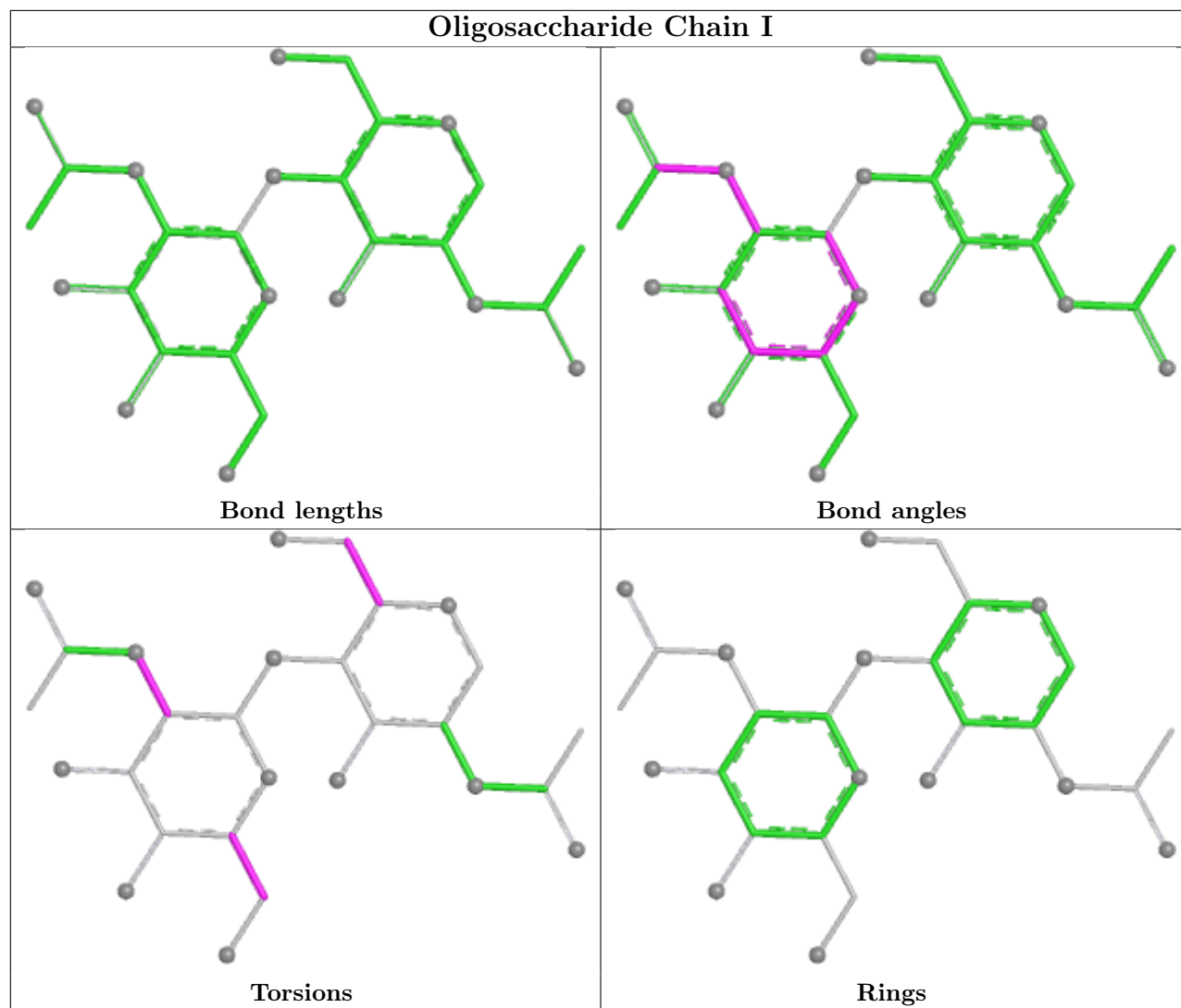
All (13) torsion outliers are listed below:

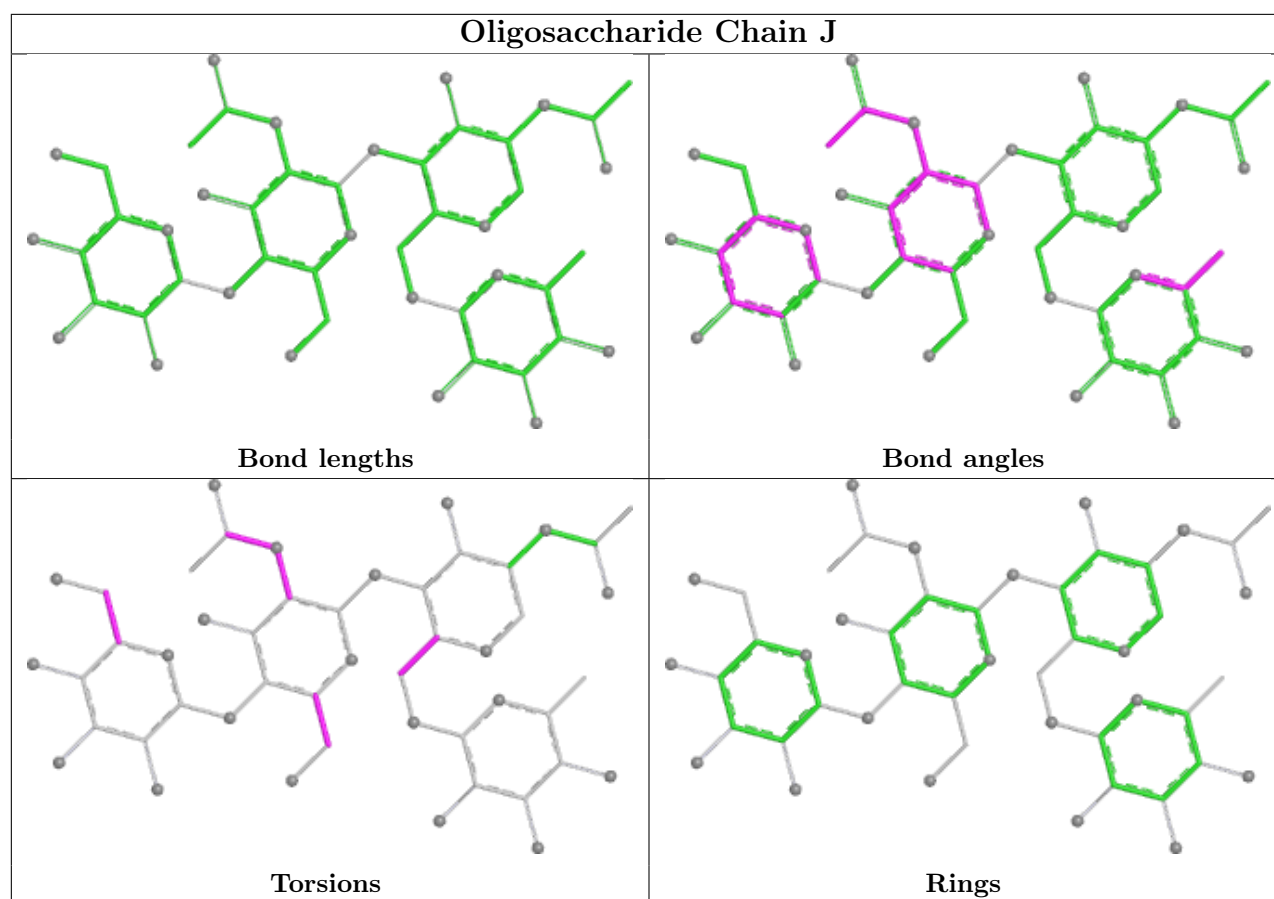
Mol	Chain	Res	Type	Atoms
5	I	2	NAG	O5-C5-C6-O6
6	J	2	NAG	C4-C5-C6-O6
6	J	3	BMA	O5-C5-C6-O6
6	J	3	BMA	C4-C5-C6-O6
6	J	2	NAG	C8-C7-N2-C2
6	J	2	NAG	O7-C7-N2-C2
6	J	2	NAG	O5-C5-C6-O6
5	I	2	NAG	C4-C5-C6-O6
6	J	2	NAG	C1-C2-N2-C7
5	I	1	NAG	C4-C5-C6-O6
6	J	1	NAG	O5-C5-C6-O6
5	I	2	NAG	C1-C2-N2-C7
5	I	2	NAG	C3-C2-N2-C7

There are no ring outliers.

No monomer is involved in short contacts.

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](#)

8 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	$\# Z > 2$
8	SGF	C	507	-	35,36,36	1.10	2 (5%)	40,46,46	1.20	6 (15%)
9	PLM	C	508	-	17,17,17	0.61	0	17,17,17	0.63	0
8	SGF	A	305	-	35,36,36	1.12	2 (5%)	40,46,46	1.17	5 (12%)
7	NAG	A	304	1	14,14,15	0.50	0	17,19,21	1.12	1 (5%)
7	NAG	C	502	1	14,14,15	0.57	0	17,19,21	1.34	2 (11%)
7	NAG	A	303	1	13,13,15	0.59	0	12,17,21	0.86	0
7	NAG	C	501	1	14,14,15	0.55	0	17,19,21	0.62	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
9	PLM	A	306	-	17,17,17	0.57	0	17,17,17	0.66	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
8	SGF	C	507	-	-	9/31/51/51	0/1/1/1
9	PLM	C	508	-	-	11/15/15/15	-
8	SGF	A	305	-	-	14/31/51/51	0/1/1/1
7	NAG	A	304	1	-	0/6/23/26	0/1/1/1
7	NAG	C	502	1	-	2/6/23/26	0/1/1/1
7	NAG	A	303	1	-	0/6/19/26	0/1/1/1
7	NAG	C	501	1	-	2/6/23/26	0/1/1/1
9	PLM	A	306	-	-	11/15/15/15	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	A	305	SGF	O4-S	-4.36	1.44	1.57
8	C	507	SGF	O4-S	-4.17	1.44	1.57
8	C	507	SGF	C2-C3	2.33	1.53	1.50
8	A	305	SGF	C2-C3	2.02	1.53	1.50

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	A	304	NAG	C1-O5-C5	3.48	116.85	112.19
8	C	507	SGF	C-O-C43	-3.28	106.76	113.80
7	C	502	NAG	C1-O5-C5	3.20	116.48	112.19
8	A	305	SGF	C-O-C43	-2.92	107.53	113.80
8	C	507	SGF	O-C-C1	2.87	115.00	108.74
8	A	305	SGF	C45-O4-S	-2.82	112.27	119.04
7	C	502	NAG	C1-C2-N2	-2.80	106.02	110.43
8	A	305	SGF	C2-C3-C4	-2.76	118.98	124.69
8	A	305	SGF	O-C43-C44	2.52	112.09	108.27
8	C	507	SGF	C2-C3-C4	-2.24	120.05	124.69
8	A	305	SGF	O-C-C1	2.21	113.57	108.74
8	C	507	SGF	C45-O4-S	-2.15	113.88	119.04
8	C	507	SGF	O8-S-O4	2.13	111.27	106.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	C	507	SGF	O-C43-C44	2.09	111.45	108.27

There are no chirality outliers.

All (49) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
8	A	305	SGF	C-C1-C2-C3
7	C	502	NAG	O5-C5-C6-O6
8	C	507	SGF	C44-C43-O-C
7	C	502	NAG	C4-C5-C6-O6
9	A	306	PLM	C1-C2-C3-C4
8	C	507	SGF	O6-C43-O-C
8	A	305	SGF	C7-C8-C9-C10
8	C	507	SGF	C11-C10-C9-C8
9	C	508	PLM	C3-C4-C5-C6
9	A	306	PLM	C5-C6-C7-C8
9	A	306	PLM	C6-C7-C8-C9
8	A	305	SGF	C5-C6-C7-C8
8	C	507	SGF	C11-C12-C13-C14
9	A	306	PLM	C8-C9-CA-CB
8	A	305	SGF	C9-C10-C11-C12
9	A	306	PLM	C3-C4-C5-C6
8	C	507	SGF	C4-C5-C6-C7
9	A	306	PLM	CB-CC-CD-CE
9	C	508	PLM	CA-CB-CC-CD
9	C	508	PLM	CB-CC-CD-CE
8	A	305	SGF	C11-C12-C13-C14
8	A	305	SGF	C12-C13-C14-C15
9	C	508	PLM	C4-C5-C6-C7
8	C	507	SGF	C10-C11-C12-C13
8	C	507	SGF	C12-C13-C14-C15
8	A	305	SGF	C4-C5-C6-C7
8	C	507	SGF	C6-C7-C8-C9
9	A	306	PLM	CA-CB-CC-CD
9	C	508	PLM	C9-CA-CB-CC
9	C	508	PLM	C8-C9-CA-CB
9	C	508	PLM	CD-CE-CF-CG
8	C	507	SGF	C7-C8-C9-C10
9	C	508	PLM	C5-C6-C7-C8
7	C	501	NAG	C4-C5-C6-O6
8	A	305	SGF	C-C1-C2-O1
8	A	305	SGF	C11-C10-C9-C8

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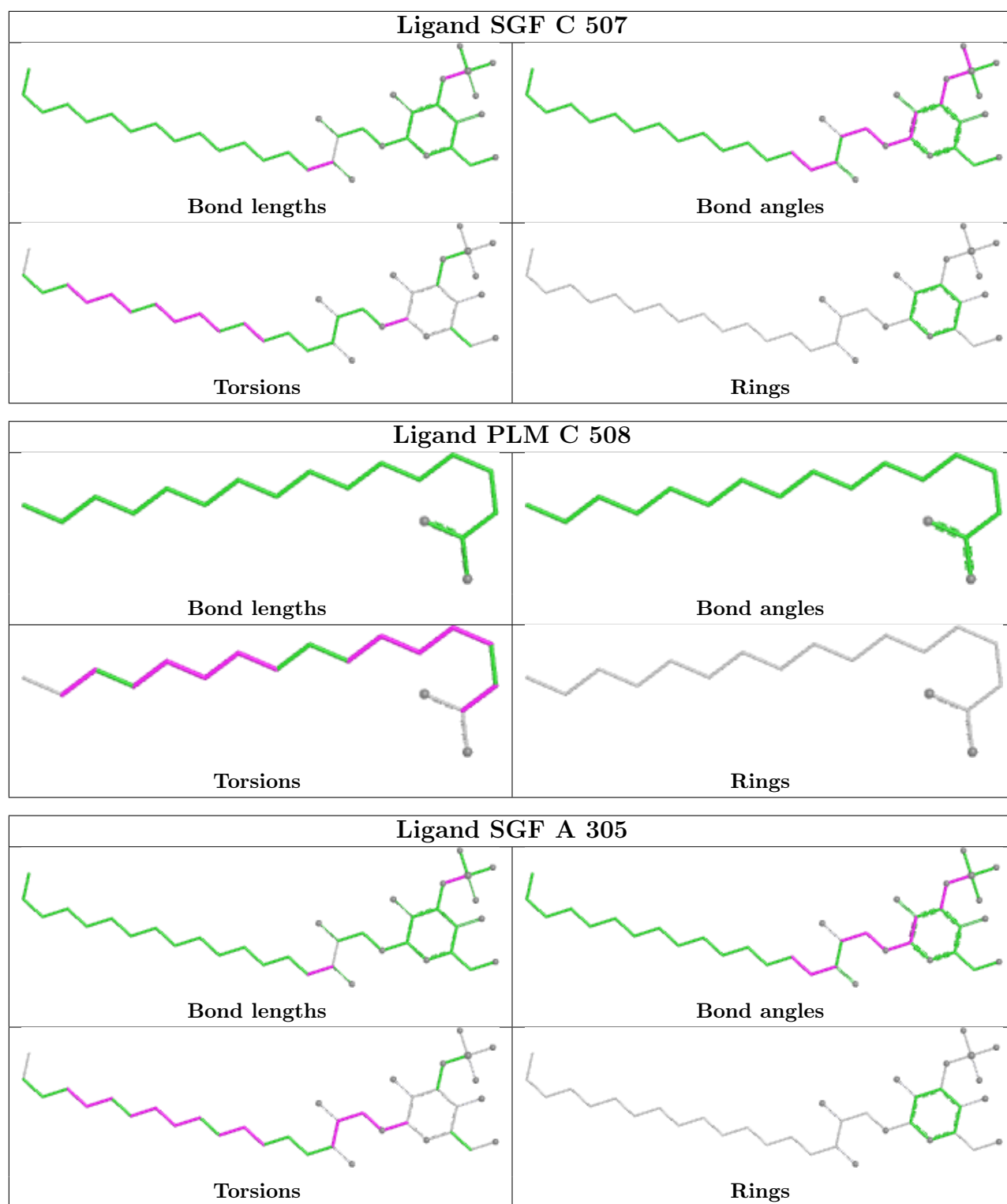
Mol	Chain	Res	Type	Atoms
9	A	306	PLM	C9-CA-CB-CC
9	A	306	PLM	C2-C3-C4-C5
9	A	306	PLM	O2-C1-C2-C3
9	A	306	PLM	O1-C1-C2-C3
9	C	508	PLM	C2-C3-C4-C5
8	A	305	SGF	C1-C-O-C43
8	A	305	SGF	C44-C43-O-C
8	A	305	SGF	N-C1-C2-O1
7	C	501	NAG	O5-C5-C6-O6
8	A	305	SGF	O6-C43-O-C
9	C	508	PLM	O1-C1-C2-C3
8	A	305	SGF	O-C-C1-N
9	C	508	PLM	O2-C1-C2-C3

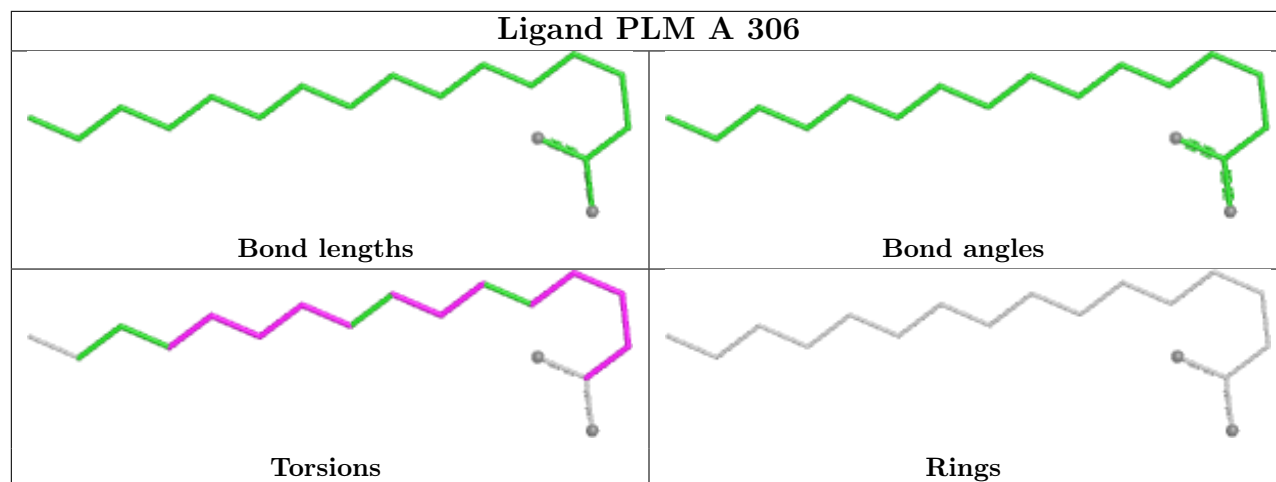
There are no ring outliers.

2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	C	507	SGF	5	0
7	C	502	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	266/285 (93%)	-0.21	1 (0%) 88 73	46, 61, 99, 152	0
1	C	269/285 (94%)	-0.25	7 (2%) 57 34	38, 54, 95, 133	0
2	B	98/99 (98%)	-0.14	0 100 100	55, 85, 108, 117	0
2	D	99/99 (100%)	-0.52	0 100 100	36, 49, 72, 78	0
3	E	179/208 (86%)	0.29	11 (6%) 27 17	40, 62, 171, 193	0
3	G	171/208 (82%)	0.47	10 (5%) 29 17	52, 81, 174, 203	0
4	F	224/244 (91%)	0.46	19 (8%) 16 11	40, 101, 169, 221	0
4	H	211/244 (86%)	0.56	15 (7%) 22 15	53, 127, 189, 236	0
All	All	1517/1672 (90%)	0.11	63 (4%) 40 23	36, 71, 162, 236	0

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	F	237	ALA	4.3
3	E	176	ALA	4.3
4	H	142	THR	3.9
3	E	155	ASP	3.5
3	G	113	PHE	3.4
3	E	154	SER	3.4
3	E	19	MET	3.3
3	G	202	PHE	3.2
4	F	198	ALA	3.1
4	F	16	GLY	2.9
3	E	138	LEU	2.9
1	C	108	PRO	2.8
4	F	130	PRO	2.8
4	F	231	VAL	2.8
4	H	129	GLU	2.8
3	G	118	ASN	2.8

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Mol	Chain	Res	Type	RSRZ
3	G	147	ASN	2.8
3	G	19	MET	2.7
4	F	122	PRO	2.7
3	G	21	SER	2.7
4	H	205	ARG	2.7
3	E	203	PRO	2.7
3	G	159	THR	2.7
4	H	12	ILE	2.7
4	H	209	ARG	2.7
4	H	117	LEU	2.6
3	E	21	SER	2.6
3	E	157	TYR	2.6
3	G	117	GLN	2.6
4	F	202	GLN	2.6
3	G	155	ASP	2.6
4	F	22	ASN	2.6
4	F	207	HIS	2.5
1	C	184	GLU	2.4
3	E	151	SER	2.4
4	F	145	CYS	2.4
1	C	183	LEU	2.4
4	H	218	SER	2.3
4	H	202	GLN	2.3
1	C	198	SER	2.3
3	G	76	HIS	2.3
4	F	216	GLY	2.3
4	F	189	SER	2.3
4	F	14	ASP	2.3
4	H	71	ASP	2.3
4	H	194	LEU	2.2
4	H	144	VAL	2.2
3	E	168	SER	2.2
4	H	143	LEU	2.2
4	H	217	LEU	2.2
4	F	219	GLU	2.2
3	E	177	VAL	2.2
4	H	175	GLN	2.1
4	F	173	ASP	2.1
4	F	50	TYR	2.1
4	F	139	GLN	2.1
1	C	92	GLU	2.1
4	F	175	GLN	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	34	ASP	2.1
1	A	198	SER	2.0
4	H	132	LYS	2.0
1	C	89	SER	2.0
4	F	117	LEU	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](#)

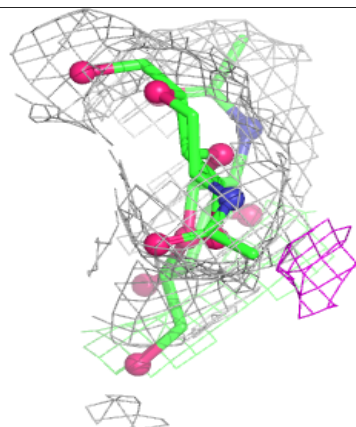
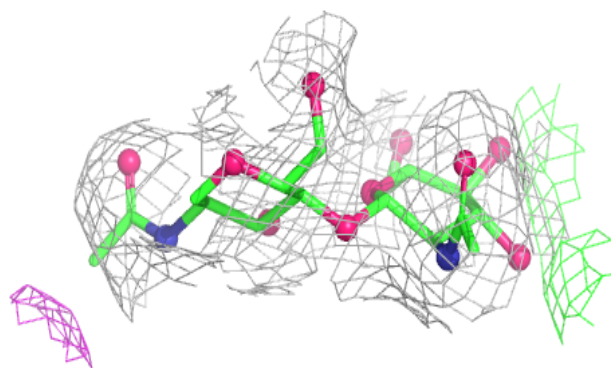
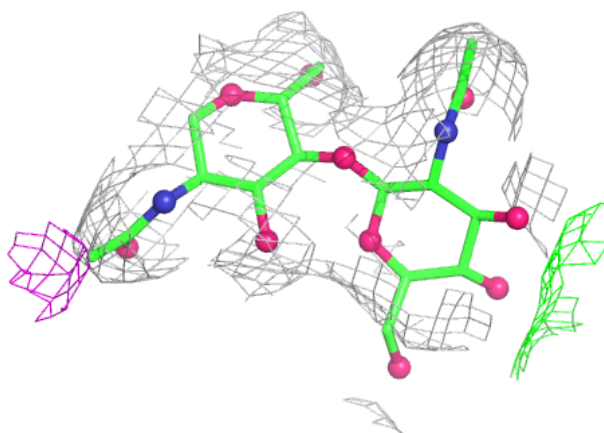
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	BMA	J	3	11/12	0.70	0.12	105,111,116,117	0
5	NAG	I	2	14/15	0.82	0.12	72,76,85,85	0
6	NAG	J	2	14/15	0.86	0.12	82,86,95,99	0
5	NAG	I	1	14/15	0.94	0.08	60,62,67,69	0
6	NAG	J	1	14/15	0.95	0.09	63,67,77,79	0
6	FUC	J	4	10/11	0.95	0.14	72,76,80,84	0

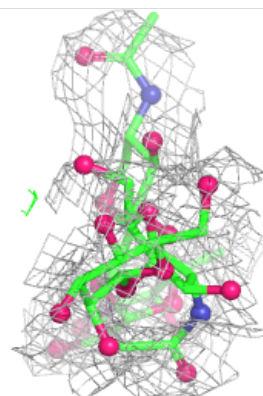
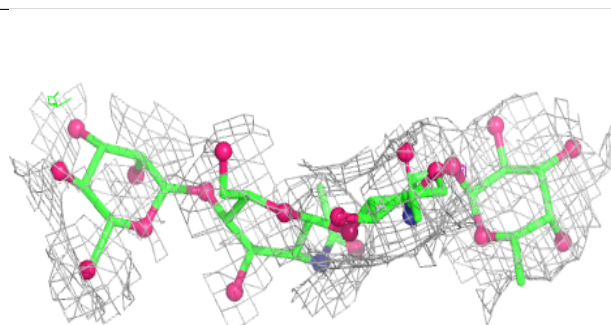
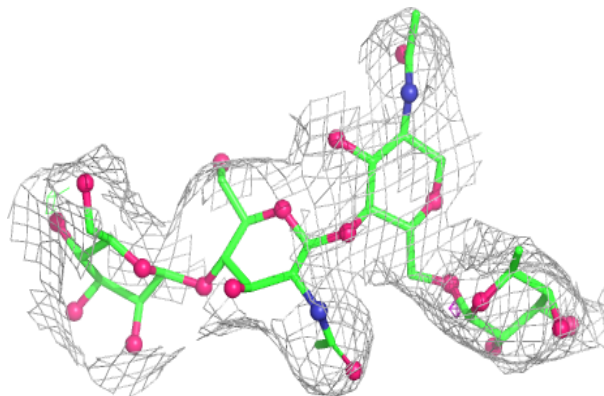
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain I:

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

**Electron density around Chain J:**

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

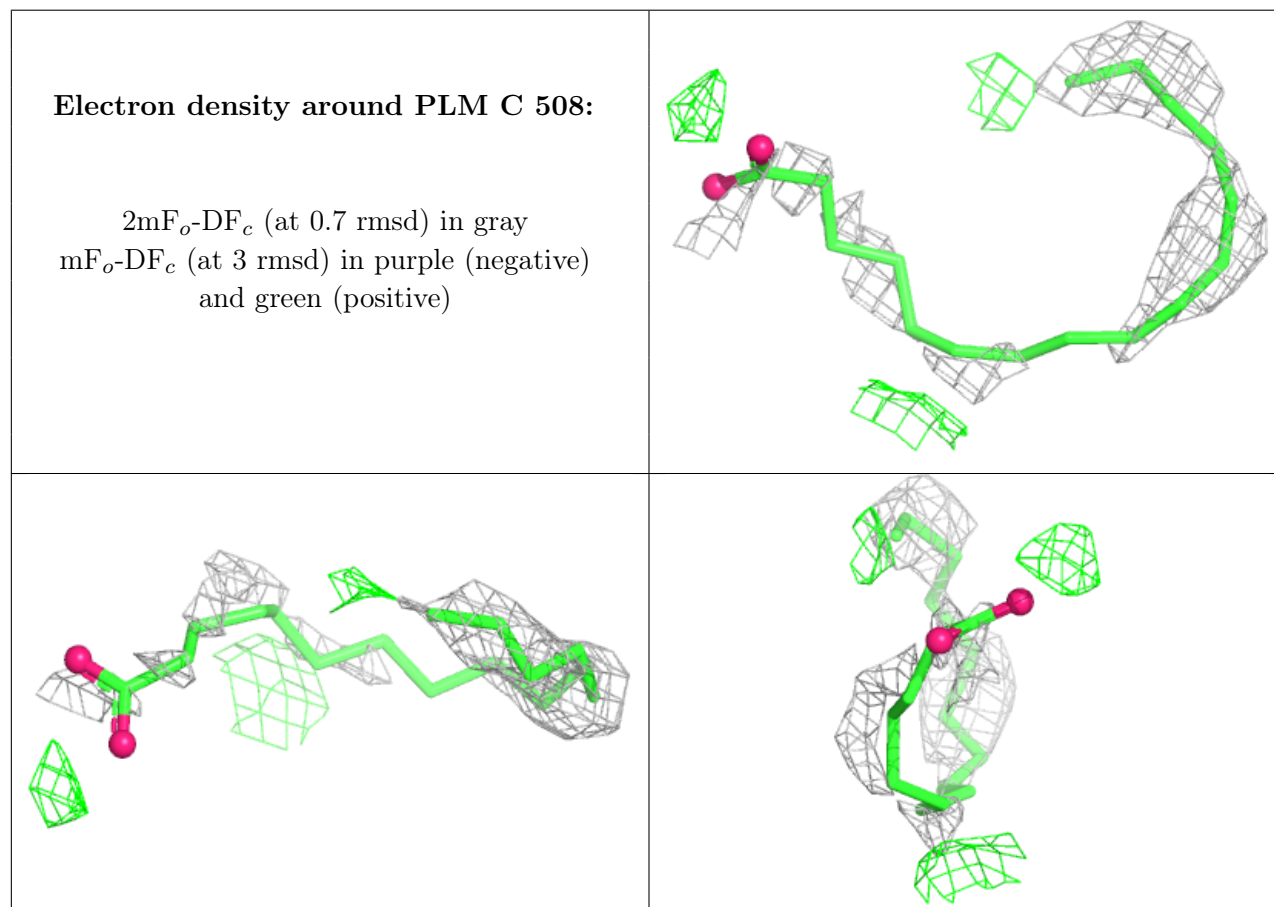


6.4 Ligands ⓘ

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.

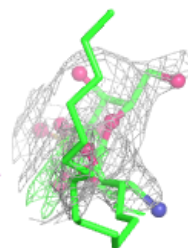
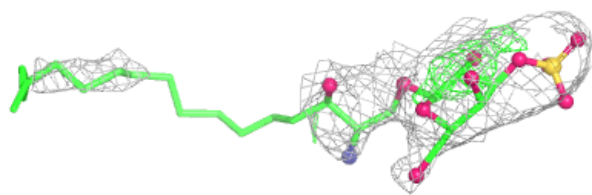
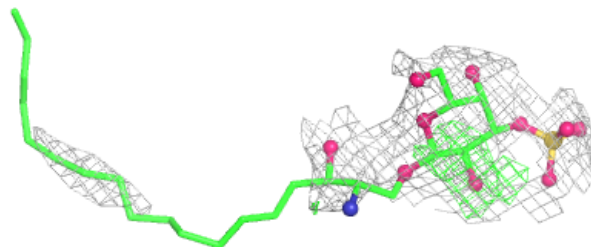
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
7	NAG	C	501	14/15	0.73	0.13	83,88,98,99	0
7	NAG	A	303	13/15	0.75	0.13	102,109,118,118	0
7	NAG	A	304	14/15	0.81	0.12	77,83,86,90	0
7	NAG	C	502	14/15	0.89	0.10	53,58,61,62	0
9	PLM	C	508	18/18	0.90	0.26	52,69,98,100	0
8	SGF	A	305	36/36	0.92	0.25	169,222,273,280	0
9	PLM	A	306	18/18	0.93	0.19	64,69,87,87	0
8	SGF	C	507	36/36	0.93	0.25	126,173,216,221	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.



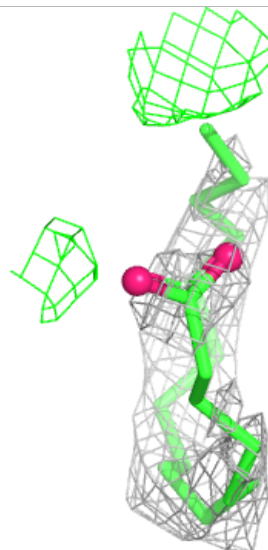
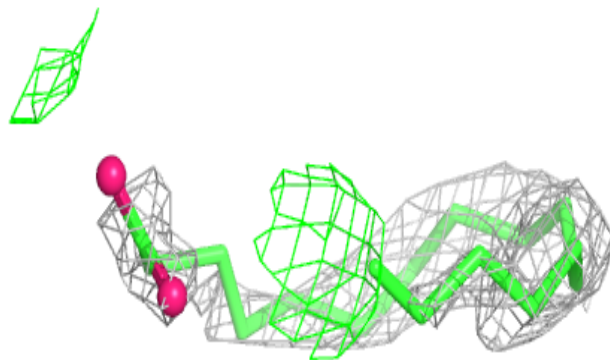
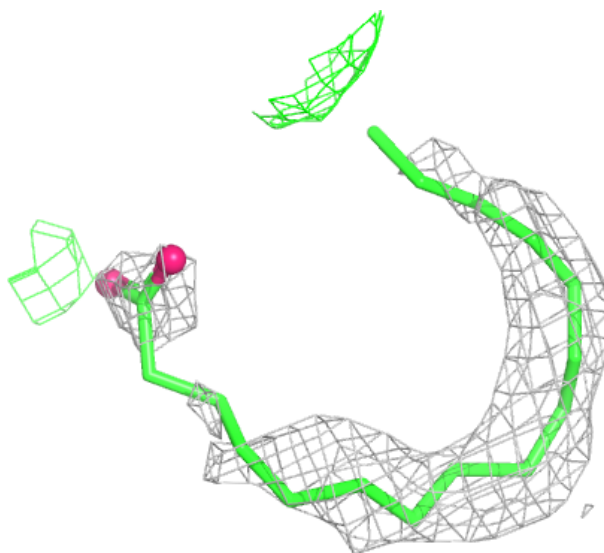
Electron density around SGF A 305:

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



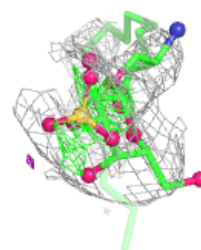
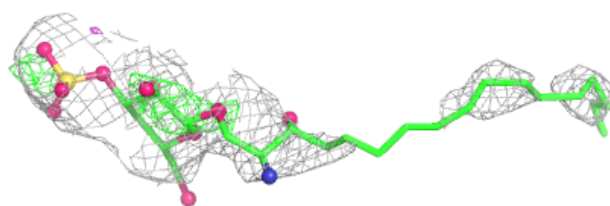
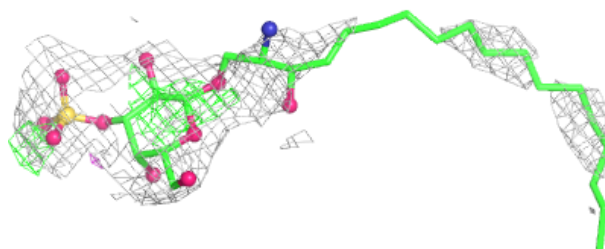
Electron density around PLM A 306:

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



Electron density around SGF C 507:

$2mF_o-DF_c$ (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.