



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 05:31 PM UTC

PDB ID : 4Y4H / pdb_00004y4h
Title : Crystal structure of the mCD1d/GCK152/iNKTCR ternary complex
Authors : Zajonc, D.M.; Yu, E.D.
Deposited on : 2015-02-10
Resolution : 3.10 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

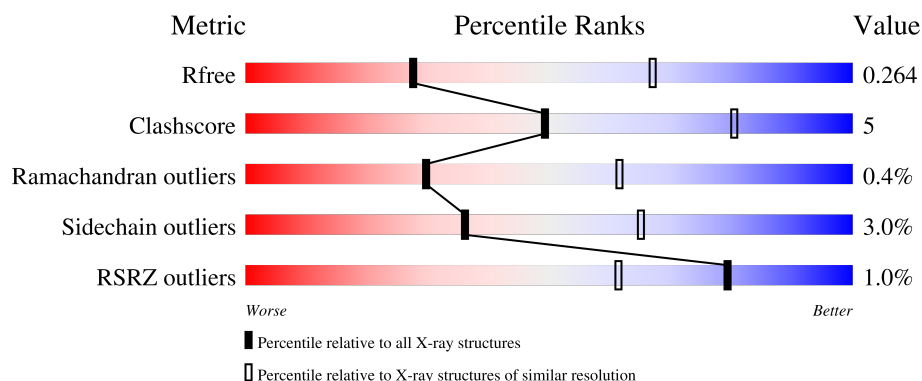
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

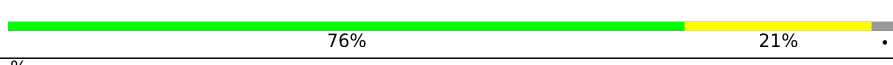
The reported resolution of this entry is 3.10 Å.

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.



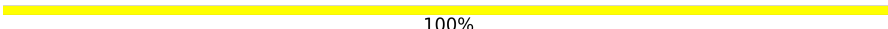
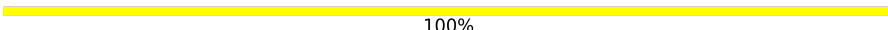
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1456 (3.10-3.10)
Clashscore	190562	1539 (3.10-3.10)
Ramachandran outliers	187476	1467 (3.10-3.10)
Sidechain outliers	187428	1467 (3.10-3.10)
RSRZ outliers	180081	1456 (3.10-3.10)

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	E	285	
2	B	99	
2	F	99	
3	C	209	

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Mol	Chain	Length	Quality of chain
3	G	209	 84% 12% .
4	D	241	 85% 13% .
4	H	241	 85% 13% .
5	I	2	 100%
5	J	2	 100%
5	K	2	 50% 50%

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 12468 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	267	Total	C	N	O	S	0	0	0
			2025	1287	339	386	13			
1	E	267	Total	C	N	O	S	0	0	0
			2032	1292	340	387	13			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	201	HIS	ASP	variant	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
E	201	HIS	ASP	variant	UNP P11609
E	280	HIS	-	expression tag	UNP P11609
E	281	HIS	-	expression tag	UNP P11609
E	282	HIS	-	expression tag	UNP P11609
E	283	HIS	-	expression tag	UNP P11609
E	284	HIS	-	expression tag	UNP P11609
E	285	HIS	-	expression tag	UNP P11609

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	96	Total	C	N	O	S	0	0	0
			723	459	122	136	6			
2	F	96	Total	C	N	O	S	0	0	0
			744	475	126	137	6			

- Molecule 3 is a protein called Chimeric TCR Valpha14/Jalpha18 chain (mouse variable do-

main/ human constant domain).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	202	Total	C	N	O	S	0	0	0
			1524	942	260	314	8			
3	G	202	Total	C	N	O	S	0	0	0
			1514	937	258	311	8			

- Molecule 4 is a protein called Chimeric TCR Vbeta8.2 chain (mouse variable domain/ human constant domain).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	239	Total	C	N	O	S	0	0	0
			1845	1157	328	354	6			
4	H	239	Total	C	N	O	S	0	0	0
			1839	1153	325	355	6			

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



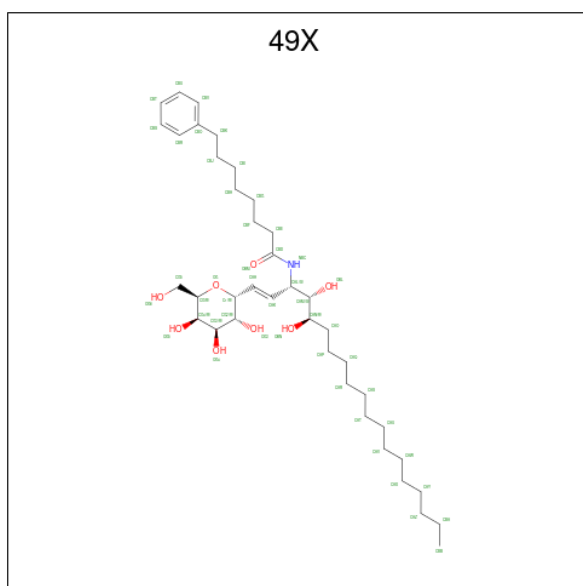
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	I	2	Total	C	N	O	0	0	0
			28	16	2	10			
5	J	2	Total	C	N	O	0	0	0
			28	16	2	10			
5	K	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 6 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



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

- Molecule 7 is (1R)-1,5-anhydro-1-[(1E,3S,4S,5R)-4,5-dihydroxy-3-[(8-phenyloctanoyl)amino]nonadec-1-en-1-yl]-D-galactitol (CCD ID: 49X) (formula: C₃₉H₆₇NO₈).

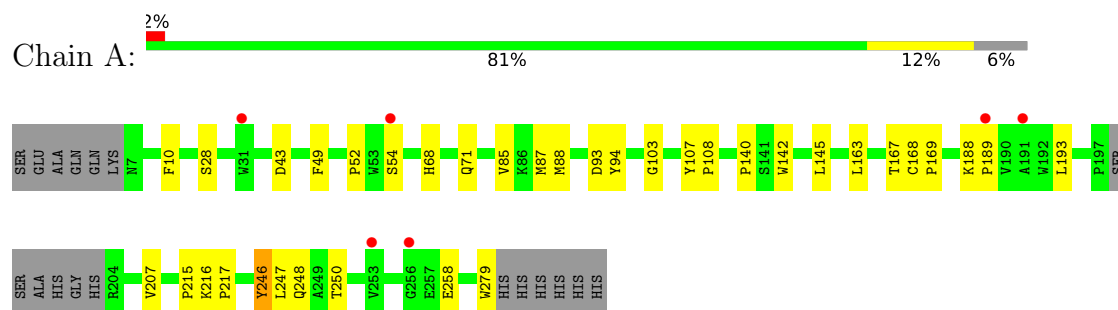


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	A	1	Total 48	C 39	N 1	O 8	0	0
7	E	1	Total 48	C 39	N 1	O 8	0	0

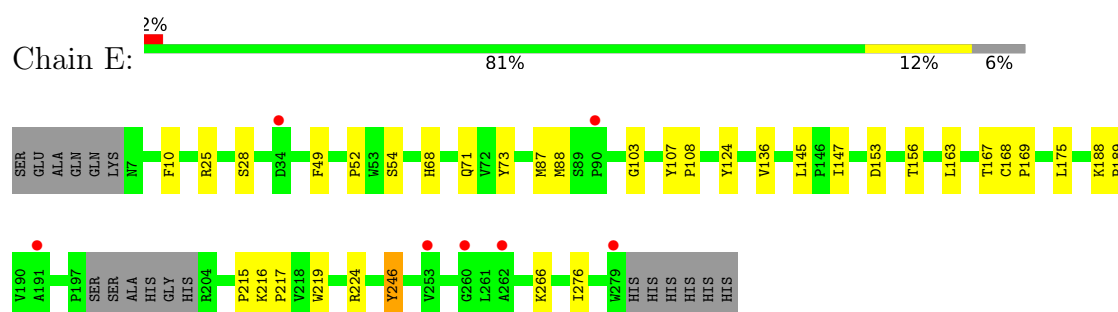
3 Residue-property plots [i](#)

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.

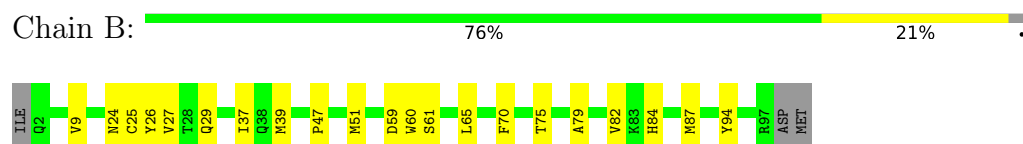
- Molecule 1: Antigen-presenting glycoprotein CD1d1



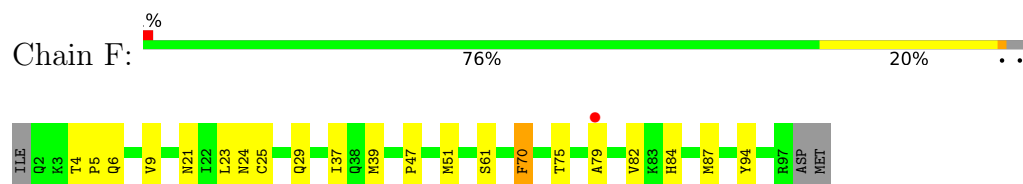
- Molecule 1: Antigen-presenting glycoprotein CD1d1



- Molecule 2: Beta-2-microglobulin

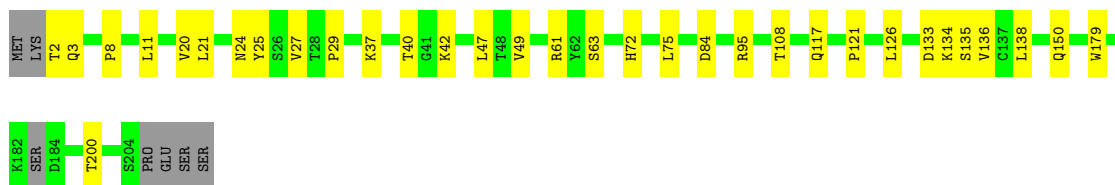


- Molecule 2: Beta-2-microglobulin



- Molecule 3: Chimeric TCR Valpha14/Jalpha18 chain (mouse variable domain/ human constant domain)

Chain C:  81% 16%



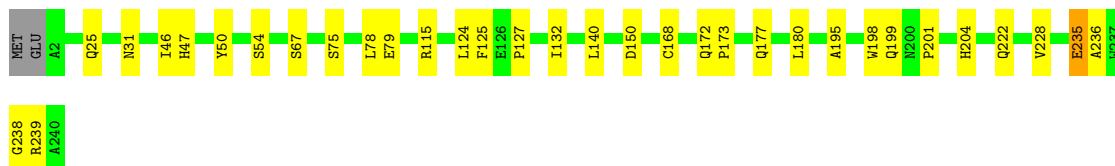
- Molecule 3: Chimeric TCR Valpha14/Jalpha18 chain (mouse variable domain/ human constant domain)

Chain G:  84% 12%



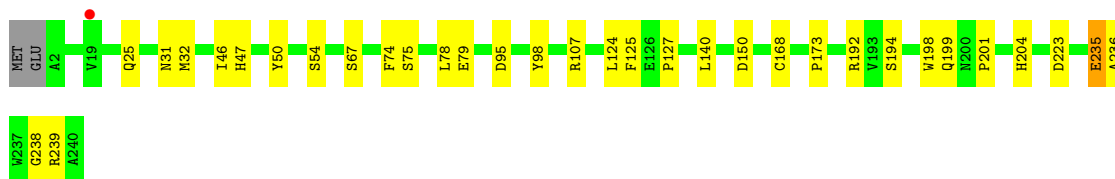
- Molecule 4: Chimeric TCR Vbeta8.2 chain (mouse variable domain/ human constant domain)

Chain D:  85% 13%

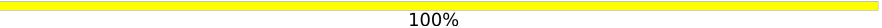


- Molecule 4: Chimeric TCR Vbeta8.2 chain (mouse variable domain/ human constant domain)

Chain H:  85% 13%



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

Chain I:  100%



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

Chain J:  100%

MAG1
MAG2

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

Chain K:  50%  50%

MAG1
MAG2

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	79.42Å 150.38Å 102.49Å 90.00° 96.38° 90.00°	Depositor
Resolution (Å)	38.17 – 3.10 38.17 – 3.10	Depositor EDS
% Data completeness (in resolution range)	98.6 (38.17-3.10) 98.6 (38.17-3.10)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.02 (at 3.12Å)	Xtriage
Refinement program	REFMAC 5.6.0104	Depositor
R, R_{free}	0.242 , 0.287 0.212 , 0.264	Depositor DCC
R_{free} test set	1334 reflections (3.08%)	wwPDB-VP
Wilson B-factor (Å ²)	83.9	Xtriage
Anisotropy	0.064	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 28.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	12468	wwPDB-VP
Average B, all atoms (Å ²)	74.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.24% 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: 49X, 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	0.80	0/2085	0.86	1/2854 (0.0%)
1	E	0.79	0/2092	0.87	0/2861
2	B	0.87	0/749	0.91	0/1031
2	F	0.83	0/770	0.85	0/1058
3	C	0.77	0/1551	0.84	0/2115
3	G	0.78	0/1541	0.82	1/2103 (0.0%)
4	D	0.79	0/1896	0.86	0/2593
4	H	0.81	0/1890	0.86	0/2585
All	All	0.80	0/12574	0.86	2/17200 (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	A	93	ASP	N-CA-C	5.51	117.46	108.76
3	G	135	SER	N-CA-C	5.17	115.87	108.74

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	2025	0	1794	23	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	2032	0	1812	24	0
2	B	723	0	612	9	0
2	F	744	0	663	11	0
3	C	1524	0	1414	16	0
3	G	1514	0	1397	18	0
4	D	1845	0	1713	15	0
4	H	1839	0	1700	18	0
5	I	28	0	25	0	0
5	J	28	0	25	0	0
5	K	28	0	25	0	0
6	A	14	0	13	0	0
6	E	28	0	26	1	0
7	A	48	0	67	3	0
7	E	48	0	67	4	0
All	All	12468	0	11353	129	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 5.

All (129) 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)
4:H:204:HIS:NE2	4:H:235:GLU:HG3	1.88	0.86
4:D:204:HIS:NE2	4:D:235:GLU:HG3	1.90	0.86
1:E:88:MET:HE1	1:E:145:LEU:HD23	1.64	0.79
1:A:168:CYS:HB3	1:A:169:PRO:HD3	1.67	0.75
1:A:88:MET:HE1	1:A:145:LEU:HD23	1.69	0.73
1:E:49:PHE:CD1	1:E:54:SER:HB2	2.24	0.73
1:A:49:PHE:HD1	1:A:54:SER:CB	2.02	0.72
3:G:8:PRO:O	3:G:108:THR:HG23	1.89	0.72
1:E:219:TRP:HB3	1:E:266:LYS:HB2	1.71	0.71
1:A:49:PHE:CD1	1:A:54:SER:HB2	2.26	0.70
3:C:61:ARG:NH2	3:C:84:ASP:OD2	2.23	0.69
1:E:49:PHE:HD1	1:E:54:SER:CB	2.09	0.66
1:E:168:CYS:HB3	1:E:169:PRO:HD3	1.76	0.66
1:A:258:GLU:HB3	1:A:279:TRP:CD1	2.32	0.65
3:G:61:ARG:NH2	3:G:84:ASP:OD2	2.28	0.64
3:C:8:PRO:O	3:C:108:THR:HG23	1.96	0.64
1:A:49:PHE:HD1	1:A:54:SER:HB2	1.61	0.63
3:C:2:THR:O	3:C:2:THR:HG23	1.99	0.62
1:E:49:PHE:HD1	1:E:54:SER:HB2	1.63	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:20:VAL:C	3:G:21:LEU:HD12	2.25	0.62
1:E:219:TRP:CE3	1:E:266:LYS:HG3	2.35	0.62
1:A:68:HIS:HA	1:A:71:GLN:OE1	1.99	0.61
3:C:20:VAL:C	3:C:21:LEU:HD12	2.26	0.60
4:H:201:PRO:HA	4:H:238:GLY:O	2.02	0.60
4:D:235:GLU:HG2	4:D:236:ALA:N	2.17	0.60
1:A:49:PHE:HD1	1:A:54:SER:HB3	1.68	0.59
1:A:107:TYR:HB3	1:A:108:PRO:HD2	1.85	0.58
4:D:201:PRO:HA	4:D:238:GLY:O	2.03	0.58
2:F:29:GLN:HA	2:F:61:SER:HB2	1.86	0.58
3:G:2:THR:O	3:G:2:THR:HG23	2.04	0.58
1:E:68:HIS:HA	1:E:71:GLN:OE1	2.04	0.56
1:A:215:PRO:HB2	1:A:217:PRO:HD2	1.88	0.56
2:B:84:HIS:HB3	2:B:87:MET:HE2	1.88	0.56
4:D:127:PRO:HD2	4:D:198:TRP:CZ2	2.41	0.55
3:C:126:LEU:O	3:C:135:SER:HB2	2.06	0.55
4:D:115:ARG:HB2	4:D:222:GLN:NE2	2.22	0.55
1:A:49:PHE:CD1	1:A:54:SER:CB	2.86	0.55
4:H:235:GLU:HG2	4:H:236:ALA:N	2.21	0.55
1:E:107:TYR:HB3	1:E:108:PRO:HD2	1.90	0.54
3:G:37:LYS:HB2	3:G:47:LEU:HD11	1.89	0.54
7:A:306:49X:H40	3:C:29:PRO:HB3	1.90	0.53
1:A:246:TYR:C	1:A:246:TYR:CD2	2.86	0.53
4:H:124:LEU:HD11	4:H:140:LEU:HD23	1.90	0.53
2:F:25:CYS:HB2	2:F:39:MET:HE3	1.89	0.53
4:H:199:GLN:HA	4:H:239:ARG:O	2.08	0.53
3:G:126:LEU:O	3:G:135:SER:HB2	2.09	0.52
2:B:25:CYS:HB2	2:B:39:MET:HE3	1.91	0.52
2:B:29:GLN:HA	2:B:61:SER:HB2	1.92	0.51
4:D:78:LEU:HD12	4:D:78:LEU:N	2.25	0.51
7:E:305:49X:NBC	7:E:305:49X:H30	2.26	0.51
3:C:117:GLN:HE21	3:C:117:GLN:HA	1.76	0.50
2:F:84:HIS:HB3	2:F:87:MET:HE2	1.93	0.50
4:H:204:HIS:NE2	4:H:235:GLU:CG	2.68	0.49
4:D:199:GLN:HA	4:D:239:ARG:O	2.12	0.49
2:F:37:ILE:HG12	2:F:82:VAL:HG22	1.95	0.49
2:F:37:ILE:HB	2:F:51:MET:HE1	1.95	0.49
2:F:9:VAL:HA	2:F:24:ASN:O	2.13	0.49
2:B:37:ILE:HG12	2:B:82:VAL:HG22	1.95	0.48
1:E:246:TYR:CD2	1:E:246:TYR:C	2.91	0.48
3:G:117:GLN:HE21	3:G:117:GLN:HA	1.78	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:31:ASN:OD1	4:H:50:TYR:HD1	1.97	0.47
4:H:78:LEU:HD12	4:H:78:LEU:N	2.29	0.47
2:B:26:TYR:HB2	2:B:65:LEU:CD1	2.44	0.47
3:G:32:HIS:CE1	4:H:98:TYR:CD1	3.03	0.47
4:D:177:GLN:HB3	4:D:180:LEU:HG	1.96	0.47
4:D:115:ARG:HB2	4:D:222:GLN:HE22	1.80	0.47
3:G:169:MET:HE1	4:H:194:SER:HB3	1.97	0.46
2:B:37:ILE:HB	2:B:51:MET:HE1	1.97	0.46
4:D:31:ASN:OD1	4:D:50:TYR:HD1	1.98	0.46
3:C:24:ASN:OD1	3:C:72:HIS:ND1	2.49	0.46
4:H:127:PRO:HD2	4:H:198:TRP:CZ2	2.50	0.46
2:B:9:VAL:HA	2:B:24:ASN:O	2.15	0.46
3:G:98:ALA:HB2	4:H:95:ASP:O	2.15	0.46
4:H:46:ILE:HG22	4:H:47:HIS:CD2	2.51	0.45
1:E:49:PHE:HD1	1:E:54:SER:HB3	1.78	0.45
1:A:85:VAL:HG11	1:A:94:TYR:CE1	2.52	0.45
7:E:305:49X:H40	3:G:29:PRO:CB	2.47	0.45
7:E:305:49X:H40	3:G:29:PRO:HB3	1.98	0.45
4:D:124:LEU:HD11	4:D:140:LEU:HD23	1.99	0.45
3:C:37:LYS:HB2	3:C:47:LEU:HD11	1.97	0.45
1:E:216:LYS:N	1:E:217:PRO:CD	2.80	0.45
2:F:21:ASN:HB3	2:F:70:PHE:CE2	2.52	0.45
1:A:87:MET:C	1:A:88:MET:HE2	2.43	0.44
3:C:136:VAL:HG12	3:C:179:TRP:HB3	2.00	0.44
1:E:215:PRO:HB2	1:E:217:PRO:HD2	1.98	0.44
1:A:207:VAL:HG22	1:A:250:THR:HG22	1.99	0.44
1:A:140:PRO:HB2	1:A:142:TRP:CD1	2.53	0.43
7:A:306:49X:H40	3:C:29:PRO:CB	2.47	0.43
1:E:49:PHE:CD1	1:E:54:SER:CB	2.90	0.43
4:D:46:ILE:HG22	4:D:47:HIS:CD2	2.54	0.43
1:A:163:LEU:HA	1:A:167:THR:HB	2.01	0.43
3:G:166:MET:HE3	4:H:192:ARG:O	2.18	0.43
1:A:10:PHE:O	1:A:103:GLY:HA3	2.19	0.43
1:E:163:LEU:HA	1:E:167:THR:HB	1.99	0.42
7:A:306:49X:H1	3:C:95:ARG:HD3	2.01	0.42
1:E:153:ASP:OD2	1:E:156:THR:OG1	2.36	0.42
3:G:136:VAL:HG12	3:G:179:TRP:HB3	2.01	0.42
2:B:79:ALA:HB2	2:B:94:TYR:CD2	2.54	0.42
1:E:73:TYR:CD1	7:E:305:49X:H56	2.54	0.42
1:E:188:LYS:HA	1:E:189:PRO:HD2	1.90	0.42
2:F:9:VAL:HG12	2:F:23:LEU:HD11	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:133:ASP:OD1	3:C:134:LYS:N	2.53	0.42
4:H:125:PHE:CD1	4:H:125:PHE:N	2.87	0.42
4:H:150:ASP:HB2	4:H:173:PRO:HG2	2.02	0.42
2:B:59:ASP:O	2:B:60:TRP:HB2	2.20	0.41
2:F:6:GLN:OE1	2:F:6:GLN:HA	2.21	0.41
2:F:79:ALA:HB2	2:F:94:TYR:CD2	2.55	0.41
4:D:125:PHE:CD1	4:D:125:PHE:N	2.89	0.41
1:E:87:MET:HB3	1:E:88:MET:HE2	2.02	0.41
1:E:124:TYR:CZ	1:E:136:VAL:HG11	2.56	0.41
1:A:188:LYS:HA	1:A:189:PRO:HD2	1.88	0.41
1:A:247:LEU:HD12	1:A:248:GLN:H	1.86	0.41
1:A:52:PRO:HB3	1:E:52:PRO:O	2.21	0.41
3:C:63:SER:O	3:C:75:LEU:HD12	2.21	0.41
1:E:25:ARG:HB3	6:E:302:NAG:H82	2.02	0.41
1:E:87:MET:C	1:E:88:MET:HE2	2.46	0.41
3:G:3:GLN:O	3:G:25:TYR:HA	2.21	0.41
3:G:21:LEU:HD12	3:G:21:LEU:N	2.36	0.41
3:G:41:GLY:CA	4:H:107:ARG:HH22	2.34	0.41
1:A:216:LYS:N	1:A:217:PRO:CD	2.84	0.41
4:H:32:MET:SD	4:H:74:PHE:HB2	2.61	0.41
1:A:168:CYS:HB3	1:A:169:PRO:CD	2.46	0.40
3:C:3:GLN:O	3:C:25:TYR:HA	2.20	0.40
1:E:10:PHE:O	1:E:103:GLY:HA3	2.21	0.40
3:C:121:PRO:HB2	3:C:200:THR:HG23	2.03	0.40
2:F:4:THR:HA	2:F:5:PRO:HD3	1.90	0.40
4:D:132:ILE:HG23	4:D:195:ALA:HB1	2.02	0.40
3:G:30:ASP:N	3:G:30:ASP:OD1	2.55	0.40
4:D:150:ASP:HB2	4:D:173:PRO:HG2	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	263/285 (92%)	248 (94%)	15 (6%)	0	100	100
1	E	263/285 (92%)	249 (95%)	14 (5%)	0	100	100
2	B	94/99 (95%)	91 (97%)	2 (2%)	1 (1%)	11	39
2	F	94/99 (95%)	91 (97%)	2 (2%)	1 (1%)	11	39
3	C	198/209 (95%)	183 (92%)	12 (6%)	3 (2%)	8	32
3	G	198/209 (95%)	181 (91%)	15 (8%)	2 (1%)	12	41
4	D	237/241 (98%)	230 (97%)	7 (3%)	0	100	100
4	H	237/241 (98%)	229 (97%)	8 (3%)	0	100	100
All	All	1584/1668 (95%)	1502 (95%)	75 (5%)	7 (0%)	30	61

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	G	150	GLN
3	C	42	LYS
3	C	150	GLN
2	B	47	PRO
2	F	47	PRO
3	G	132	SER
3	C	27	VAL

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	204/249 (82%)	200 (98%)	4 (2%)	48	72
1	E	206/249 (83%)	200 (97%)	6 (3%)	37	66
2	B	72/93 (77%)	69 (96%)	3 (4%)	26	58
2	F	78/93 (84%)	76 (97%)	2 (3%)	40	68
3	C	171/188 (91%)	167 (98%)	4 (2%)	44	70
3	G	168/188 (89%)	166 (99%)	2 (1%)	63	78
4	D	195/208 (94%)	186 (95%)	9 (5%)	24	55

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
4	H	194/208 (93%)	186 (96%)	8 (4%)	27 59
All	All	1288/1476 (87%)	1250 (97%)	38 (3%)	36 65

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	SER
1	A	43	ASP
1	A	193	LEU
1	A	246	TYR
2	B	27	VAL
2	B	70	PHE
2	B	75	THR
3	C	11	LEU
3	C	40	THR
3	C	49	VAL
3	C	138	LEU
4	D	25	GLN
4	D	54	SER
4	D	67	SER
4	D	75	SER
4	D	79	GLU
4	D	168	CYS
4	D	172	GLN
4	D	228	VAL
4	D	235	GLU
1	E	28	SER
1	E	147	ILE
1	E	175	LEU
1	E	224	ARG
1	E	246	TYR
1	E	276	ILE
2	F	70	PHE
2	F	75	THR
3	G	40	THR
3	G	49	VAL
4	H	25	GLN
4	H	54	SER
4	H	67	SER
4	H	75	SER
4	H	79	GLU
4	H	168	CYS

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Mol	Chain	Res	Type
4	H	223	ASP
4	H	235	GLU

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	186	GLN
1	A	229	GLN
2	B	13	HIS
3	C	31	ASN
3	C	117	GLN
4	D	41	HIS
4	D	208	GLN
4	D	210	GLN
1	E	7	ASN
1	E	186	GLN
1	E	229	GLN
2	F	13	HIS
2	F	38	GLN
3	G	31	ASN
3	G	38	GLN
3	G	117	GLN
4	H	37	GLN
4	H	71	GLN
4	H	164	HIS
4	H	230	GLN

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](#)

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	0.50	0	17,19,21	2.07	1 (5%)
5	NAG	I	2	5	14,14,15	0.68	0	17,19,21	1.02	1 (5%)
5	NAG	J	1	1,5	14,14,15	0.71	0	17,19,21	2.02	6 (35%)
5	NAG	J	2	5	14,14,15	0.57	0	17,19,21	1.59	3 (17%)
5	NAG	K	1	1,5	14,14,15	0.43	0	17,19,21	1.36	3 (17%)
5	NAG	K	2	5	14,14,15	0.60	0	17,19,21	1.17	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
5	NAG	I	1	1,5	-	2/6/23/26	0/1/1/1
5	NAG	I	2	5	-	2/6/23/26	0/1/1/1
5	NAG	J	1	1,5	-	2/6/23/26	0/1/1/1
5	NAG	J	2	5	-	2/6/23/26	0/1/1/1
5	NAG	K	1	1,5	-	0/6/23/26	0/1/1/1
5	NAG	K	2	5	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	I	1	NAG	C1-O5-C5	7.36	122.06	112.19
5	J	1	NAG	C6-C5-C4	4.38	123.77	113.02
5	J	1	NAG	C2-N2-C7	3.82	128.02	122.90
5	J	2	NAG	C4-C3-C2	3.48	116.11	111.02
5	J	1	NAG	C3-C4-C5	-3.45	103.97	110.23
5	K	1	NAG	C1-O5-C5	3.11	116.35	112.19
5	I	2	NAG	C4-C3-C2	3.04	115.47	111.02
5	J	2	NAG	O5-C1-C2	-3.00	106.65	111.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	J	2	NAG	O5-C5-C6	2.86	113.23	107.66
5	J	1	NAG	C1-C2-N2	2.58	114.50	110.43
5	K	1	NAG	C4-C3-C2	2.55	114.75	111.02
5	J	1	NAG	C4-C3-C2	2.42	114.56	111.02
5	K	1	NAG	O5-C1-C2	-2.36	107.64	111.29
5	J	1	NAG	O7-C7-C8	-2.24	118.06	122.05

There are no chirality outliers.

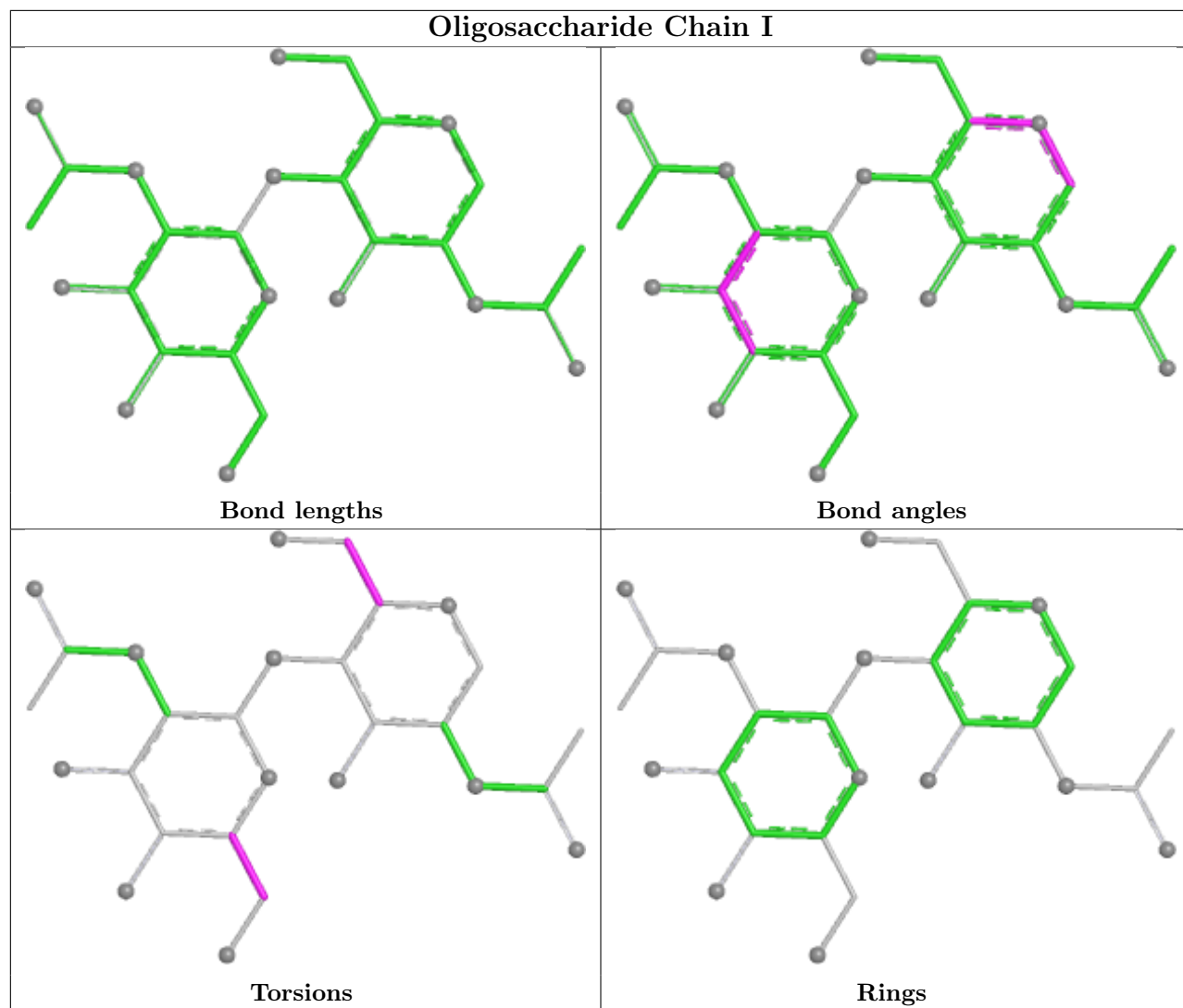
All (10) torsion outliers are listed below:

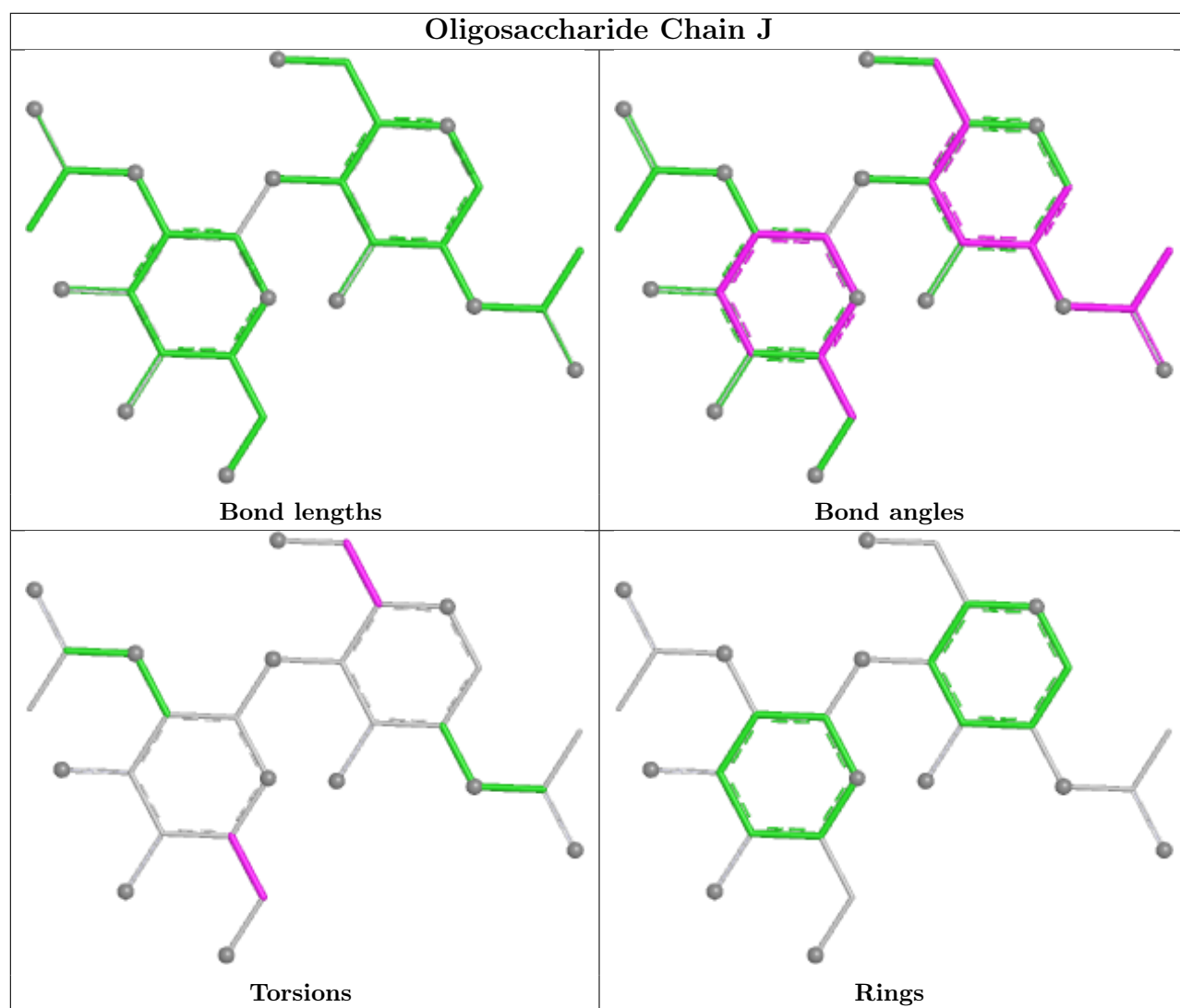
Mol	Chain	Res	Type	Atoms
5	J	1	NAG	O5-C5-C6-O6
5	J	1	NAG	C4-C5-C6-O6
5	J	2	NAG	O5-C5-C6-O6
5	I	1	NAG	C4-C5-C6-O6
5	J	2	NAG	C4-C5-C6-O6
5	K	2	NAG	O5-C5-C6-O6
5	K	2	NAG	C4-C5-C6-O6
5	I	1	NAG	O5-C5-C6-O6
5	I	2	NAG	O5-C5-C6-O6
5	I	2	NAG	C4-C5-C6-O6

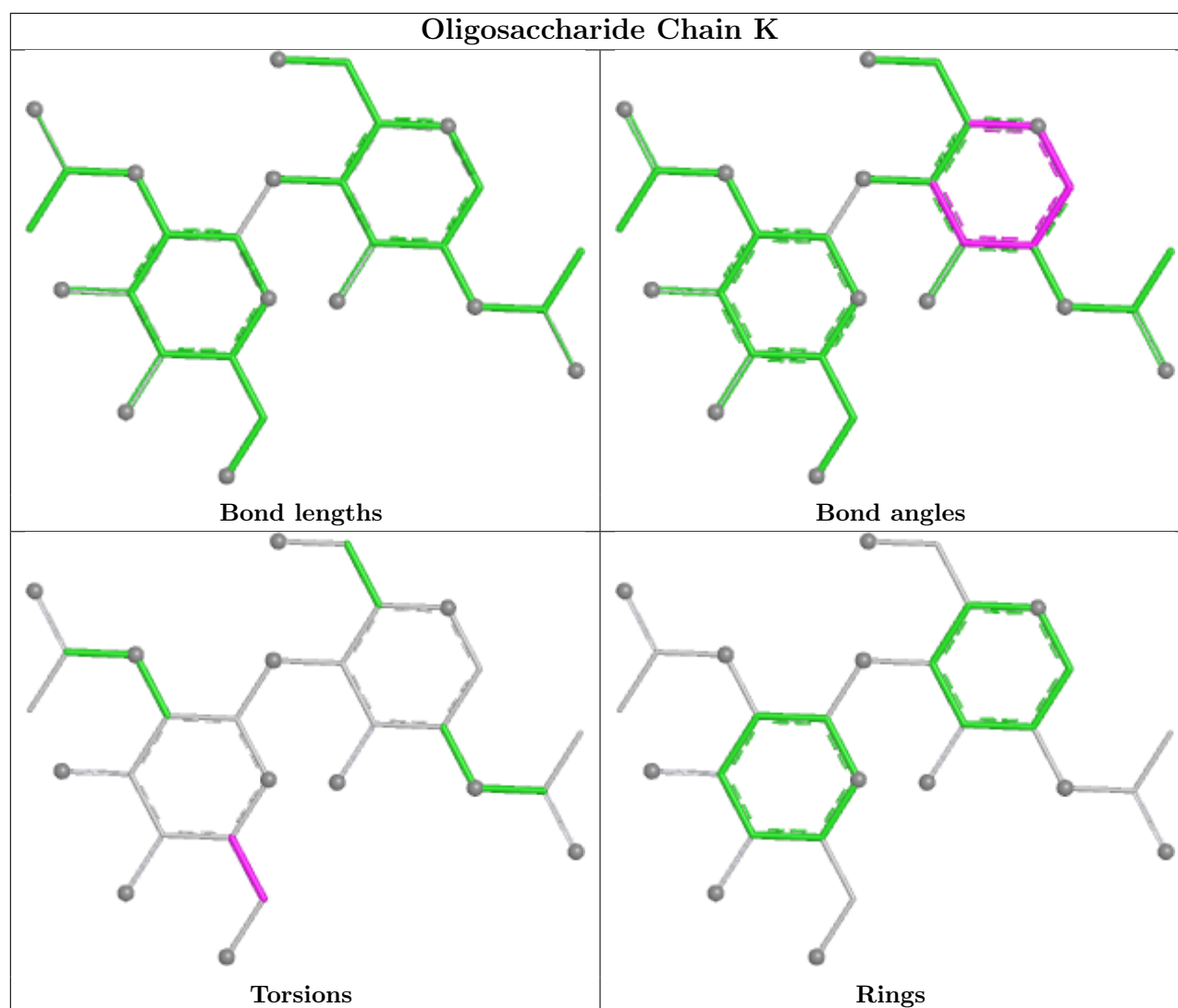
There are no ring outliers.

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](#)

5 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	$\# Z > 2$
7	49X	E	305	-	49,49,49	0.89	2 (4%)	53,60,60	1.10	5 (9%)
6	NAG	E	302	1	14,14,15	0.52	0	17,19,21	1.82	4 (23%)
6	NAG	A	301	1	14,14,15	0.51	0	17,19,21	1.24	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	NAG	E	301	1	14,14,15	0.73	0	17,19,21	1.14	1 (5%)
7	49X	A	306	-	49,49,49	0.94	3 (6%)	53,60,60	1.29	5 (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
7	49X	E	305	-	-	25/43/63/63	0/2/2/2
6	NAG	E	302	1	-	2/6/23/26	0/1/1/1
6	NAG	A	301	1	-	2/6/23/26	0/1/1/1
6	NAG	E	301	1	-	0/6/23/26	0/1/1/1
7	49X	A	306	-	-	21/43/63/63	0/2/2/2

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	E	305	49X	C1-CAH	-2.92	1.40	1.50
7	A	306	49X	CAH-CAK	2.88	1.40	1.32
7	A	306	49X	C1-CAH	-2.66	1.41	1.50
7	E	305	49X	CAH-CAK	2.63	1.40	1.32
7	A	306	49X	CAL-CAK	2.24	1.55	1.49

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	A	306	49X	CAP-CAO-CAN	-4.88	106.12	114.11
6	E	302	NAG	C1-O5-C5	4.59	118.34	112.19
7	A	306	49X	OBL-CAM-CAN	-3.78	100.33	108.93
6	A	301	NAG	O5-C5-C6	3.59	114.65	107.66
7	E	305	49X	OG-C1-CAH	3.35	118.93	108.68
7	A	306	49X	CG-OG-C1	-3.26	107.62	112.93
6	E	302	NAG	C3-C4-C5	-3.22	104.40	110.23
7	A	306	49X	OG-C1-CAH	2.96	117.72	108.68
6	E	301	NAG	C4-C3-C2	2.81	115.14	111.02
7	E	305	49X	CAO-CAN-CAM	-2.57	109.90	113.97
6	E	302	NAG	C4-C3-C2	-2.52	107.32	111.02
7	E	305	49X	OG6-CG5-CG	-2.37	103.28	111.33
7	E	305	49X	CG5-CG-CG4	2.28	118.61	113.02
7	A	306	49X	OG-CG-CG5	2.19	111.87	106.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	E	305	49X	CAP-CAO-CAN	-2.19	110.52	114.11
6	E	302	NAG	C1-C2-N2	2.08	113.72	110.43

There are no chirality outliers.

All (50) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	A	306	49X	OG-C1-CAH-CAK
7	A	306	49X	CG2-C1-CAH-CAK
7	A	306	49X	CAH-CAK-CAL-CAM
7	A	306	49X	CAH-CAK-CAL-NBC
7	E	305	49X	OG-C1-CAH-CAK
7	E	305	49X	CG2-C1-CAH-CAK
7	E	305	49X	OBN-CAN-CAO-CAP
7	E	305	49X	NBC-CAL-CAM-OBL
7	E	305	49X	CAH-CAK-CAL-CAM
7	E	305	49X	CAH-CAK-CAL-NBC
6	A	301	NAG	O5-C5-C6-O6
6	E	302	NAG	O5-C5-C6-O6
7	E	305	49X	OG-CG-CG5-OG6
6	A	301	NAG	C4-C5-C6-O6
7	A	306	49X	OG-CG-CG5-OG6
7	E	305	49X	CG4-CG-CG5-OG6
7	A	306	49X	CAR-CAS-CAT-CAU
7	E	305	49X	CBI-CBJ-CBK-CBO
6	E	302	NAG	C4-C5-C6-O6
7	A	306	49X	CBI-CBJ-CBK-CBO
7	E	305	49X	CBD-CBE-CBF-CBG
7	A	306	49X	CG4-CG-CG5-OG6
7	A	306	49X	CBD-CBE-CBF-CBG
7	A	306	49X	CAT-CAU-CAV-CAW
7	A	306	49X	CBF-CBG-CBH-CBI
7	E	305	49X	CAT-CAU-CAV-CAW
7	A	306	49X	CBE-CBF-CBG-CBH
7	E	305	49X	CAV-CAW-CAX-CAY
7	E	305	49X	CAX-CAY-CAZ-CBA
7	E	305	49X	CBE-CBF-CBG-CBH
7	E	305	49X	CAO-CAP-CAQ-CAR
7	A	306	49X	CAW-CAX-CAY-CAZ
7	E	305	49X	CBF-CBG-CBH-CBI
7	E	305	49X	CAP-CAQ-CAR-CAS
7	E	305	49X	CAU-CAV-CAW-CAX

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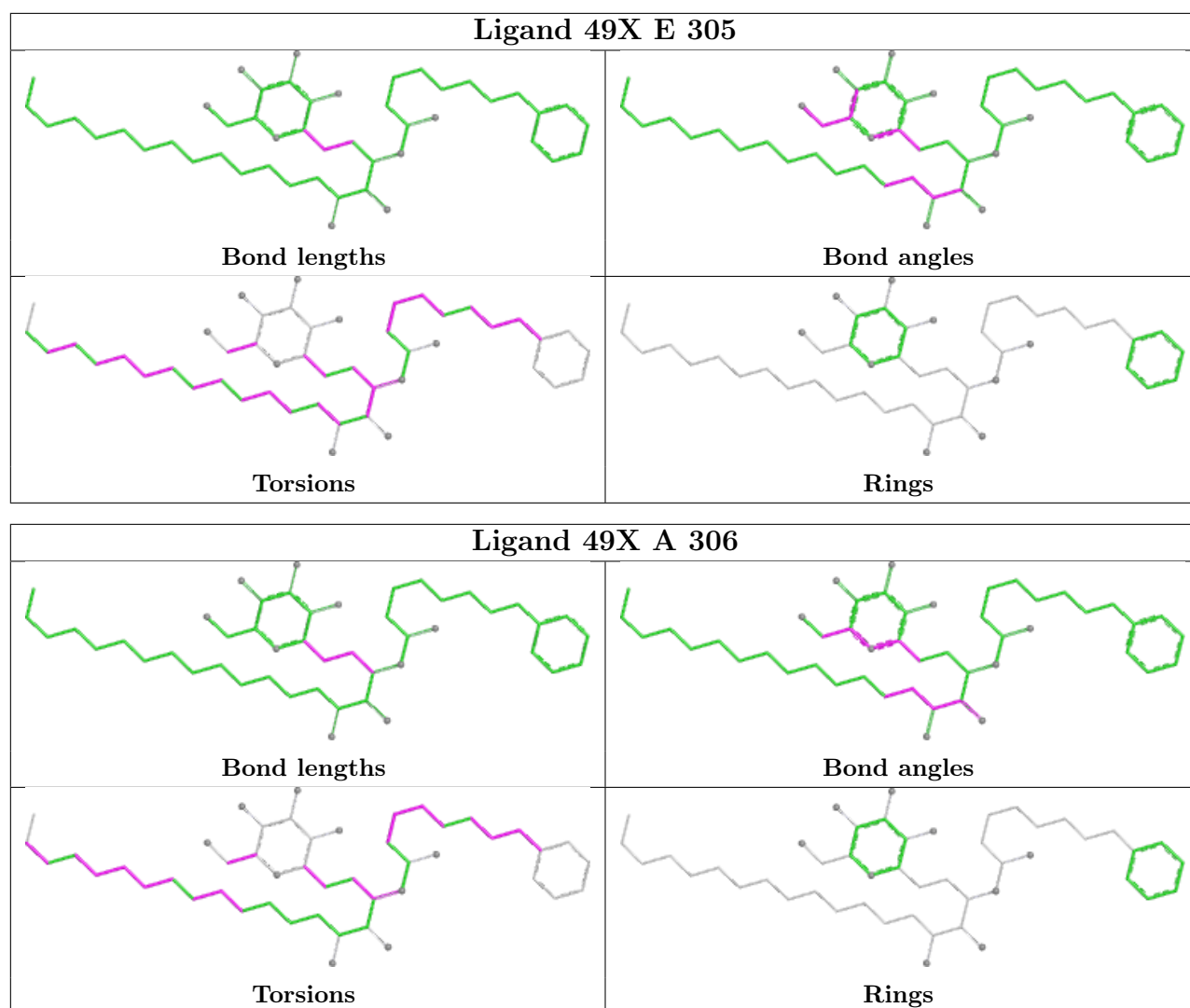
Mol	Chain	Res	Type	Atoms
7	E	305	49X	CBH-CBI-CBJ-CBK
7	E	305	49X	CAK-CAL-CAM-OBL
7	A	306	49X	CBH-CBI-CBJ-CBK
7	A	306	49X	CAQ-CAR-CAS-CAT
7	A	306	49X	CAV-CAW-CAX-CAY
7	A	306	49X	CAU-CAV-CAW-CAX
7	A	306	49X	CAY-CAZ-CBA-CBB
7	E	305	49X	CBJ-CBK-CBO-CBR
7	E	305	49X	CBJ-CBK-CBO-CBV
7	E	305	49X	CAR-CAS-CAT-CAU
7	E	305	49X	CAM-CAN-CAO-CAP
7	A	306	49X	CBJ-CBK-CBO-CBR
7	A	306	49X	CAK-CAL-NBC-CBD
7	E	305	49X	CAK-CAL-NBC-CBD
7	A	306	49X	CBJ-CBK-CBO-CBV

There are no ring outliers.

3 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	E	305	49X	4	0
6	E	302	NAG	1	0
7	A	306	49X	3	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	267/285 (93%)	0.10	6 (2%) 62 41	39, 79, 170, 204	0
1	E	267/285 (93%)	-0.09	7 (2%) 57 36	40, 76, 136, 169	0
2	B	96/99 (96%)	-0.13	0 100 100	47, 81, 111, 125	0
2	F	96/99 (96%)	-0.06	1 (1%) 79 61	59, 95, 118, 127	0
3	C	202/209 (96%)	-0.43	0 100 100	36, 57, 114, 148	0
3	G	202/209 (96%)	-0.47	1 (0%) 87 73	34, 57, 112, 153	0
4	D	239/241 (99%)	-0.44	0 100 100	40, 60, 108, 123	0
4	H	239/241 (99%)	-0.52	1 (0%) 88 76	36, 60, 97, 123	0
All	All	1608/1668 (96%)	-0.27	16 (0%) 79 61	34, 67, 128, 204	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	191	ALA	5.8
1	E	34	ASP	5.2
1	A	54	SER	3.4
1	E	279	TRP	3.1
1	A	253	VAL	3.0
1	E	191	ALA	2.9
1	E	262	ALA	2.9
1	A	189	PRO	2.5
2	F	79	ALA	2.4
1	A	31	TRP	2.4
1	E	260	GLY	2.4
3	G	2	THR	2.2
1	E	90	PRO	2.1
1	A	256	GLY	2.0
1	E	253	VAL	2.0
4	H	19	VAL	2.0

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

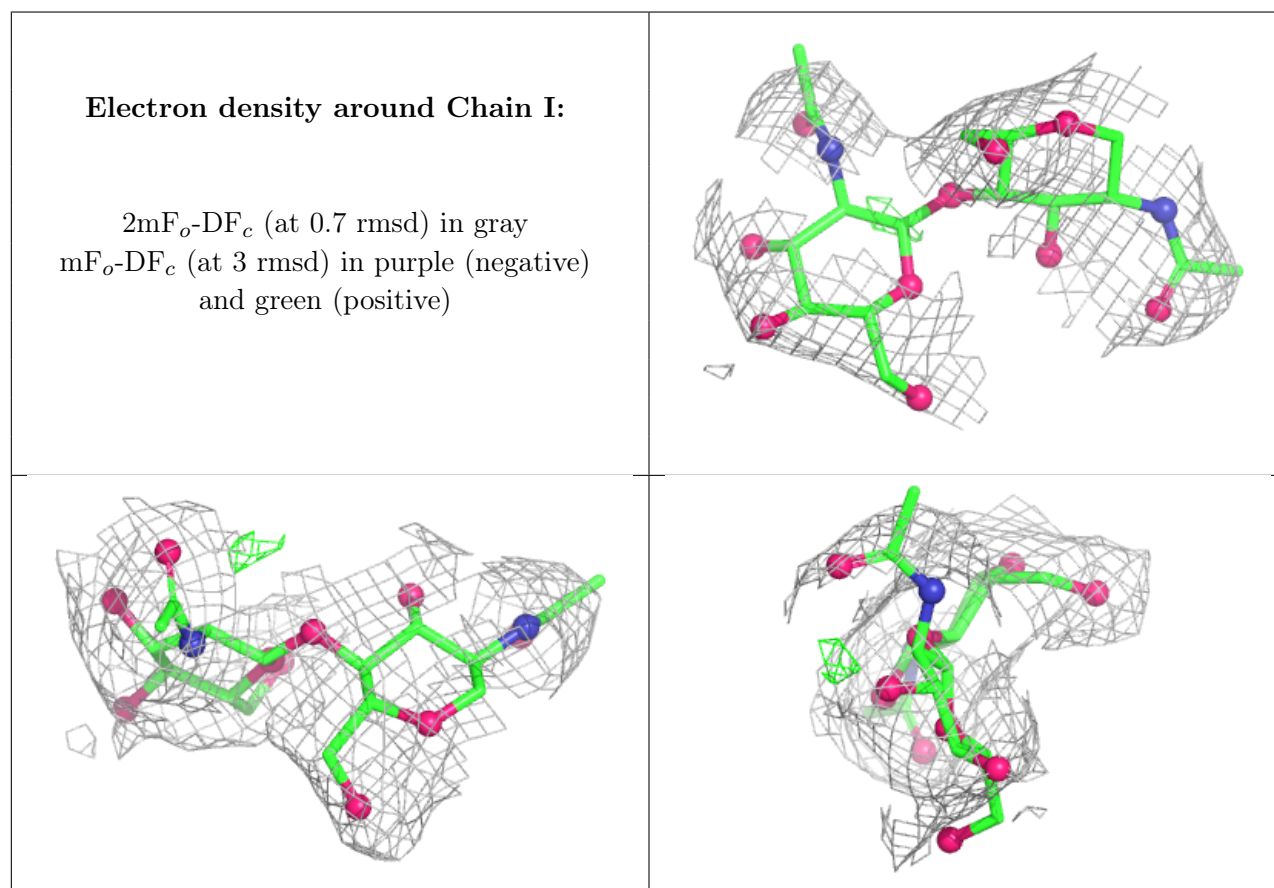
There are no non-standard protein/DNA/RNA residues in this entry.

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.

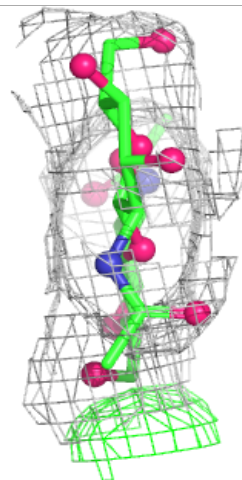
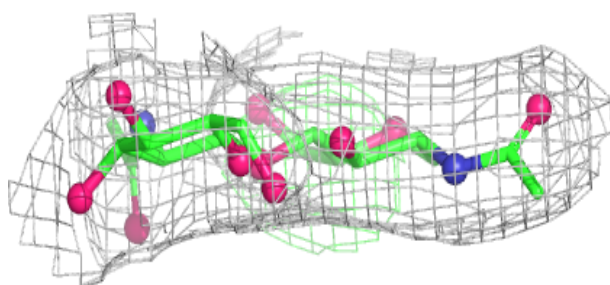
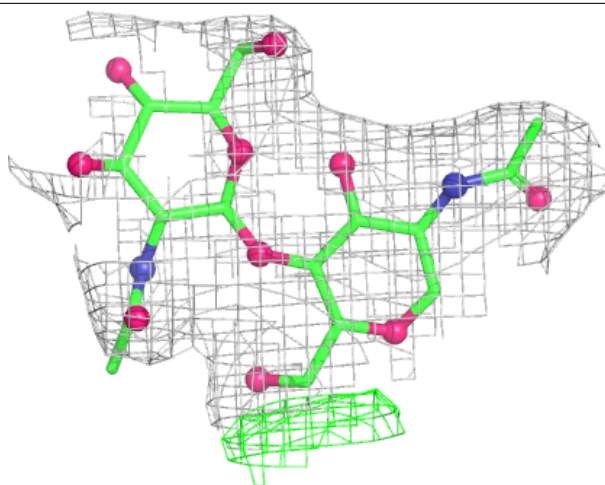
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	NAG	J	1	14/15	0.79	0.09	51,63,70,71	0
5	NAG	I	2	14/15	0.80	0.09	99,111,125,126	0
5	NAG	K	2	14/15	0.84	0.10	76,87,94,95	0
5	NAG	J	2	14/15	0.92	0.07	74,86,93,94	0
5	NAG	K	1	14/15	0.93	0.07	57,62,71,72	0
5	NAG	I	1	14/15	0.93	0.07	64,69,75,90	0

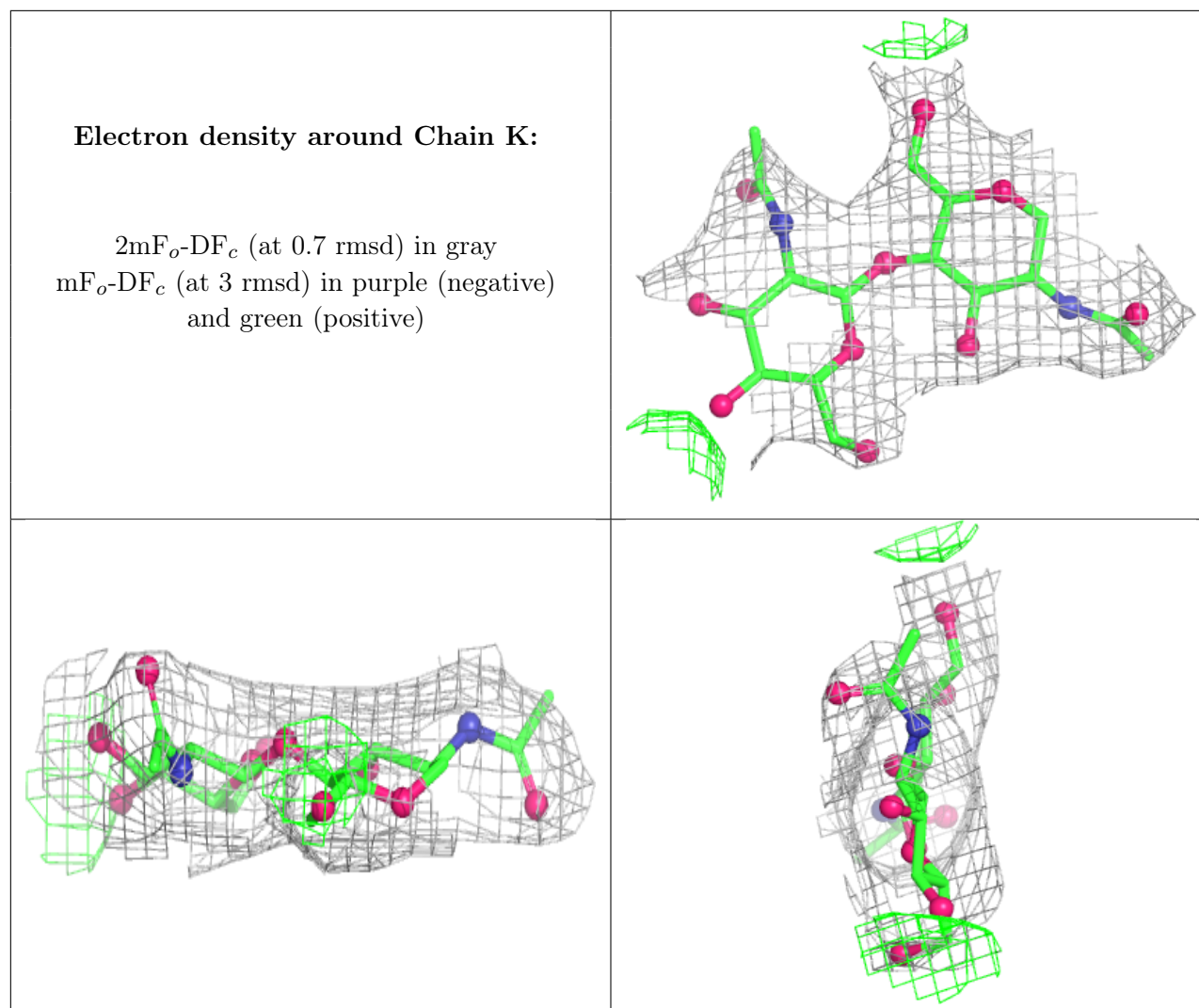
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.



Electron density around Chain 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 [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.

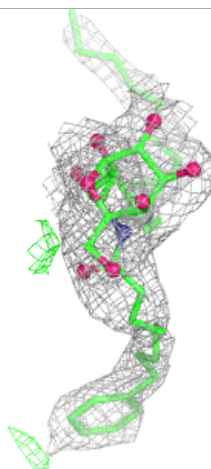
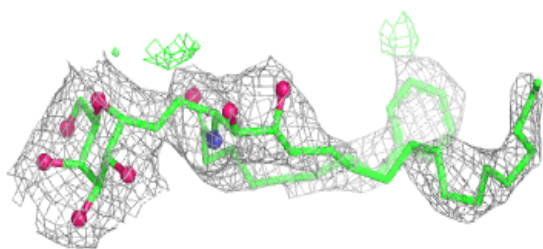
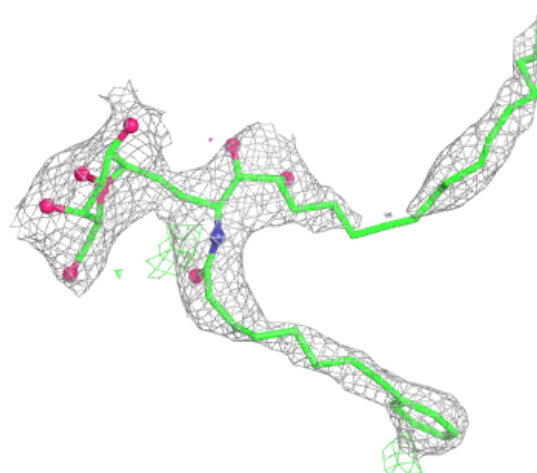
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
6	NAG	E	301	14/15	0.81	0.07	82,100,106,112	0
6	NAG	E	302	14/15	0.85	0.08	69,70,74,79	0
6	NAG	A	301	14/15	0.86	0.08	90,97,106,109	0
7	49X	A	306	48/48	0.95	0.10	46,57,75,77	0
7	49X	E	305	48/48	0.96	0.11	42,53,64,68	0

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.

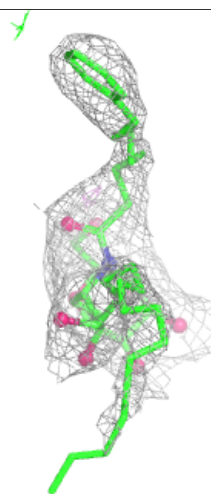
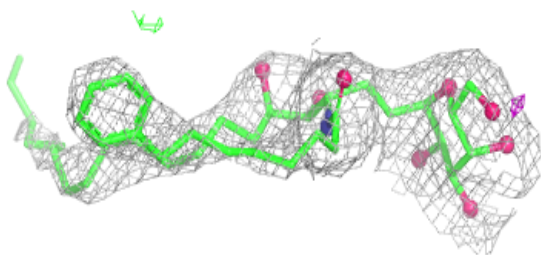
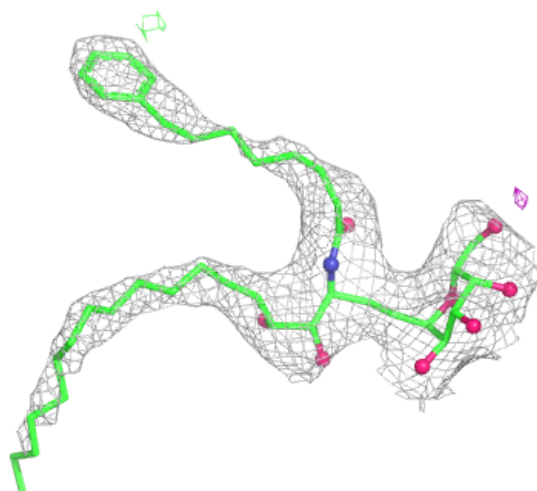
Electron density around 49X A 306:

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



Electron density around 49X E 305:

$2mF_o-DF_c$ (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.