



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 09:29 PM UTC

PDB ID : 7UDG / pdb_00007udg
Title : Integrin alpha IIB beta3 complex with lotrafiban
Authors : Lin, F.-Y.; Zhu, J.; Zhu, J.; Springer, T.A.
Deposited on : 2022-03-19
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

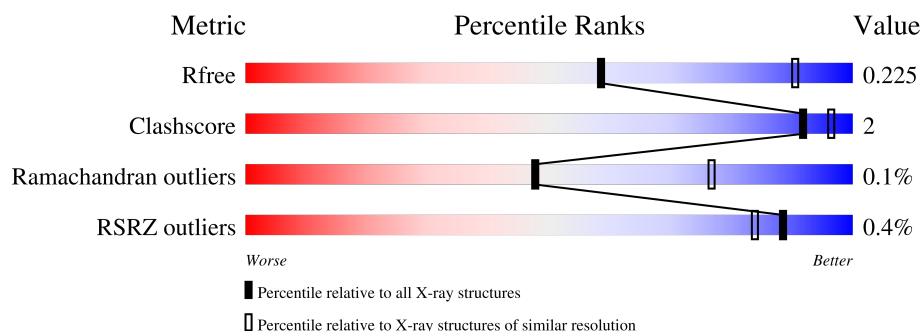
1 Overall quality at a glance

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)
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	457	96%
1	C	457	95%
2	B	472	93%
2	D	472	92%
3	E	221	95%
3	H	221	95%

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Mol	Chain	Length	Quality of chain
4	F	214	% 96% .
4	L	214	98% .

2 Entry composition

There are 14 unique types of molecules in this entry. The entry contains 21786 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 Integrin alpha-IIb heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	454	Total	C	N	O	S	0	4	0
			3499	2226	601	664	8			
1	C	453	Total	C	N	O	S	0	4	0
			3502	2224	604	666	8			

- Molecule 2 is a protein called Isoform Beta-3C of Integrin beta-3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	466	Total	C	N	O	S	4	3	0
			3610	2248	616	712	34			
2	D	471	Total	C	N	O	S	3	0	0
			3623	2255	619	715	34			

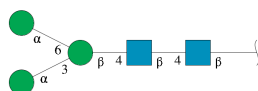
- Molecule 3 is a protein called 10E5 Fab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	214	Total	C	N	O	S	0	0	0
			1631	1035	264	326	6			
3	H	216	Total	C	N	O	S	0	0	0
			1642	1041	266	329	6			

- Molecule 4 is a protein called 10E5 Fab light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	F	214	Total	C	N	O	S	0	0	0
			1637	1019	268	341	9			
4	L	214	Total	C	N	O	S	0	0	0
			1637	1019	268	341	9			

- Molecule 5 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	G	5	Total	C	N	O	0	0	0
			61	34	2	25			

- Molecule 6 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



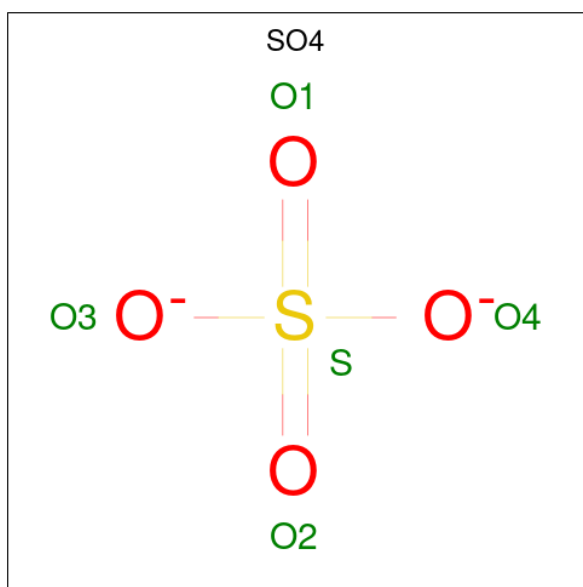
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
6	I	2	Total	C	N	O	0	0	0
			28	16	2	10			
6	K	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 7 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
7	J	4	Total	C	N	O	0	0	0
			50	28	2	20			

- Molecule 8 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	1	Total	O	S	0	0
			5	4	1		
8	A	1	Total	O	S	0	0
			5	4	1		
8	A	1	Total	O	S	0	0
			5	4	1		
8	A	1	Total	O	S	0	0
			5	4	1		
8	C	1	Total	O	S	0	0
			5	4	1		
8	C	1	Total	O	S	0	0
			5	4	1		
8	C	1	Total	O	S	0	0
			5	4	1		
8	C	1	Total	O	S	0	0
			5	4	1		
8	L	1	Total	O	S	0	0
			5	4	1		

- Molecule 9 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	4	Total	Ca	0	0
			4	4		
9	B	1	Total	Ca	0	0
			1	1		
9	C	4	Total	Ca	0	0
			4	4		

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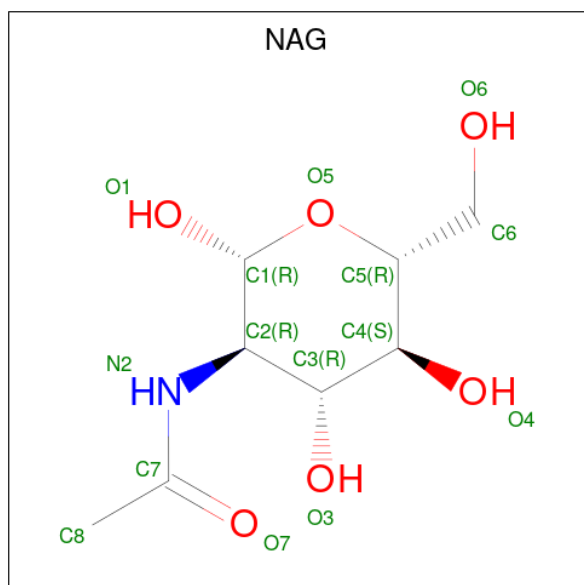
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	D	1	Total	Ca	0	0
			1	1		

- Molecule 10 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

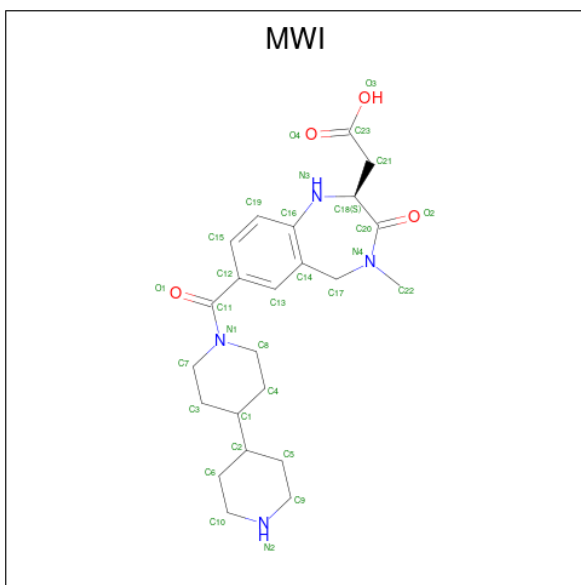
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	B	1	Total	Mg	0	0
			1	1		
10	D	1	Total	Mg	0	0
			1	1		

- Molecule 11 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



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

- Molecule 12 is Lotrafiban (CCD ID: MWI) (formula: C₂₃H₃₂N₄O₄) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
12	B	1	Total 31	C 23	N 4	O 4	0	0
12	D	1	Total 31	C 23	N 4	O 4	0	0

- Molecule 13 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
13	C	2	Total Cl 2 2	0	0
13	D	1	Total Cl 1 1	0	0

- Molecule 14 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
14	A	300	Total O 300 300	0	0
14	B	144	Total O 144 144	0	0
14	C	106	Total O 106 106	0	0
14	D	83	Total O 83 83	0	0
14	E	5	Total O 5 5	0	0

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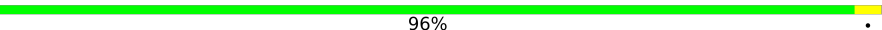
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
14	F	4	Total 4	O 4	0	0
14	H	22	Total 22	O 22	0	0
14	L	24	Total 24	O 24	0	0

3 Residue-property plots

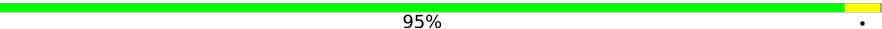
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Integrin alpha-IIb heavy chain

Chain A:  96%

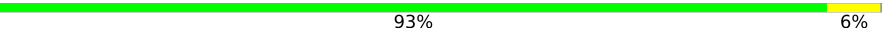


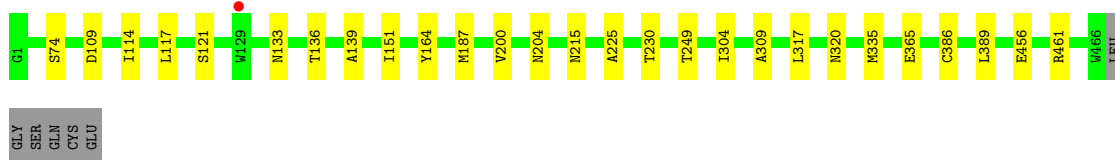
- Molecule 1: Integrin alpha-IIb heavy chain

Chain C:  95%

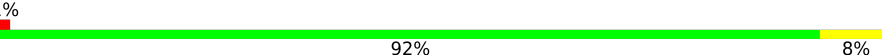


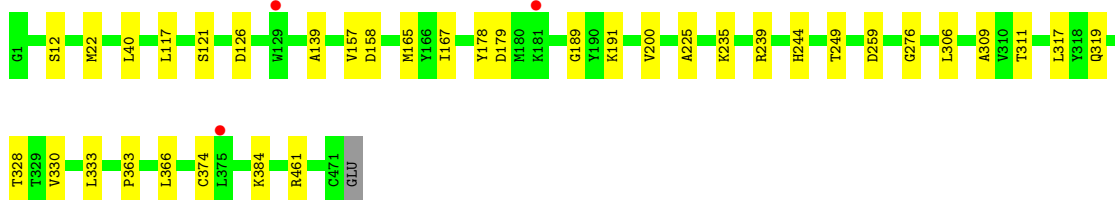
- Molecule 2: Isoform Beta-3C of Integrin beta-3

Chain B:  93% 6%

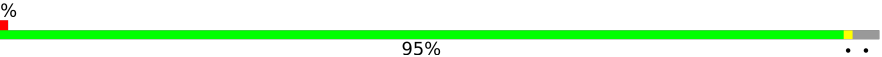


- Molecule 2: Isoform Beta-3C of Integrin beta-3

Chain D:  92% 8%



- Molecule 3: 10E5 Fab heavy chain

Chain E:  95%



- Molecule 3: 10E5 Fab heavy chain

Chain H: 95%



- Molecule 4: 10E5 Fab light chain

Chain F: 96%



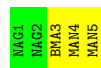
- Molecule 4: 10E5 Fab light chain

Chain L: 98%



- Molecule 5: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain G: 40%



- Molecule 6: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain I: 100%



- Molecule 6: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain K: 100%



- Molecule 7: alpha-D-mannopyranose-(1-3)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain J:  25% 75%

MAG1	MAG2	BMA3	MAN4
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4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	256.76Å 144.17Å 104.55Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.14 – 2.80 49.14 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.6 (49.14-2.80) 99.6 (49.14-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.05 (at 2.81Å)	Xtriage
Refinement program	PHENIX 1.11.1_2575	Depositor
R, R_{free}	0.213 , 0.235 (Not available) , 0.225	Depositor DCC
R_{free} test set	2000 reflections (1.08%)	wwPDB-VP
Wilson B-factor (Å ²)	74.2	Xtriage
Anisotropy	0.284	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 70.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	21786	wwPDB-VP
Average B, all atoms (Å ²)	103.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.10% 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: MWI, MG, BMA, NAG, CL, MAN, CA, SO4

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	0.18	0/3608	0.40	0/4918
1	C	0.15	0/3605	0.36	0/4912
2	B	0.16	0/3680	0.35	0/4989
2	D	0.14	0/3690	0.33	0/5003
3	E	0.14	0/1673	0.31	0/2290
3	H	0.13	0/1684	0.31	0/2305
4	F	0.13	0/1673	0.30	0/2269
4	L	0.14	0/1673	0.31	0/2269
All	All	0.15	0/21286	0.34	0/28955

There are no bond length outliers.

There are no bond angle outliers.

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	3499	0	3345	9	0
1	C	3502	0	3334	12	0
2	B	3610	0	3529	18	0
2	D	3623	0	3539	20	0
3	E	1631	0	1590	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	H	1642	0	1600	3	0
4	F	1637	0	1553	5	0
4	L	1637	0	1553	2	0
5	G	61	0	52	0	0
6	I	28	0	25	0	0
6	K	28	0	25	0	0
7	J	50	0	43	0	0
8	A	20	0	0	0	0
8	C	20	0	0	0	0
8	L	5	0	0	0	0
9	A	4	0	0	0	0
9	B	1	0	0	0	0
9	C	4	0	0	0	0
9	D	1	0	0	0	0
10	B	1	0	0	0	0
10	D	1	0	0	0	0
11	B	14	0	13	0	0
11	D	14	0	13	0	0
12	B	31	0	0	2	0
12	D	31	0	0	1	0
13	C	2	0	0	0	0
13	D	1	0	0	0	0
14	A	300	0	0	0	0
14	B	144	0	0	1	0
14	C	106	0	0	1	0
14	D	83	0	0	1	0
14	E	5	0	0	0	0
14	F	4	0	0	0	0
14	H	22	0	0	0	0
14	L	24	0	0	0	0
All	All	21786	0	20214	64	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 2.

All (64) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:3:LEU:O	1:C:405:GLN:NE2	2.31	0.63
2:B:456:GLU:OE2	2:B:461:ARG:NH1	2.35	0.59
2:B:139:ALA:HB2	2:B:200:VAL:HG11	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:230:THR:HG23	2:B:304:ILE:HG13	1.86	0.57
1:C:262:TRP:HB3	2:D:317:LEU:HD13	1.86	0.57
2:D:191:LYS:NZ	14:D:2102:HOH:O	2.38	0.55
2:D:12:SER:HB3	2:D:461:ARG:HD3	1.89	0.54
2:D:306:LEU:HB3	2:D:328:THR:HG22	1.91	0.53
2:D:121:SER:HB2	12:D:2005:MWI:C23	2.39	0.53
1:A:9:THR:HB	1:A:447:VAL:HB	1.91	0.52
1:C:1:LEU:N	14:C:605:HOH:O	2.41	0.52
2:D:126:ASP:OD1	2:D:126:ASP:N	2.36	0.52
1:C:319:ASP:OD1	2:D:384:LYS:NZ	2.43	0.51
2:B:121:SER:HB2	12:B:2004:MWI:C23	2.41	0.51
2:B:133:ASN:OD1	2:B:133:ASN:N	2.44	0.51
1:C:415:SER:HB2	1:C:439:ALA:HB1	1.93	0.50
2:D:117:LEU:HD11	2:D:225:ALA:HB1	1.93	0.49
3:E:149:LYS:NZ	4:F:180:THR:HG21	2.28	0.48
1:A:192:LEU:HD11	1:A:231:PHE:CD1	2.49	0.48
1:C:259:THR:HA	1:C:263:THR:HA	1.96	0.47
2:B:249:THR:HA	2:B:309:ALA:O	2.14	0.47
4:L:66:GLY:HA3	4:L:71:TYR:HA	1.97	0.47
1:A:314:MET:HE3	1:A:322:LEU:HD13	1.97	0.46
4:F:201:SER:OG	4:F:203:SER:O	2.27	0.46
1:A:394:GLY:HA2	1:A:399:LEU:HD23	1.97	0.46
2:D:235:LYS:HE3	2:D:276:GLY:O	2.15	0.46
1:C:83:THR:HB	1:C:116:LEU:HB2	1.99	0.45
2:D:22:MET:HG2	2:D:40:LEU:HD22	1.97	0.45
2:D:139:ALA:HB2	2:D:200:VAL:HG11	1.98	0.45
2:B:164:TYR:O	2:B:215:ASN:HB2	2.17	0.44
2:B:136:THR:HG23	2:B:204:ASN:HD21	1.83	0.44
1:A:285:MET:SD	2:B:320:ASN:HB3	2.57	0.44
1:C:107:CYS:HA	1:C:130:CYS:HA	1.99	0.44
4:F:29:ILE:O	4:F:32:ASN:ND2	2.40	0.44
4:L:4:MET:HE3	4:L:23:CYS:SG	2.58	0.44
2:B:365:GLU:H	2:B:365:GLU:CD	2.26	0.43
2:B:335:MET:HB2	14:B:2221:HOH:O	2.17	0.43
2:B:187[B]:MET:SD	2:B:215:ASN:HB3	2.59	0.43
1:C:278[A]:HIS:NE2	1:C:340:LEU:O	2.51	0.43
2:D:319:GLN:HA	2:D:330:VAL:HG21	2.00	0.43
2:B:74:SER:HB3	2:B:109:ASP:CG	2.43	0.43
1:C:288:TYR:OH	2:D:259:ASP:OD1	2.31	0.42
2:D:363:PRO:HG2	2:D:366:LEU:HB2	2.01	0.42
2:D:178:TYR:CG	2:D:179:ASP:N	2.87	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:262:TRP:HB3	2:B:317:LEU:HD13	2.00	0.42
2:D:239:ARG:O	2:D:244:HIS:NE2	2.40	0.42
3:E:172:PHE:CD2	4:F:164:THR:HG23	2.54	0.42
1:C:224:ASP:OD1	1:C:225:SER:N	2.46	0.42
3:H:69:THR:HB	3:H:82:GLN:HB3	2.02	0.42
3:H:91:THR:HG23	3:H:116:THR:HA	2.02	0.42
4:F:35:TRP:CZ3	4:F:88:CYS:HB3	2.55	0.41
1:A:167:CYS:HB3	1:A:185:ALA:HB1	2.03	0.41
2:D:158:ASP:O	2:D:189:GLY:HA2	2.20	0.41
3:H:51:ILE:O	3:H:53:PRO:HD3	2.20	0.41
2:B:386:CYS:HB3	2:B:389:LEU:HD11	2.03	0.41
2:D:165:MET:HE3	2:D:167:ILE:HG22	2.01	0.41
1:C:215:HIS:CE1	3:E:32:THR:HG22	2.56	0.41
1:A:107:CYS:HA	1:A:130:CYS:HA	2.02	0.41
2:B:114:ILE:O	2:B:151:ILE:HA	2.21	0.41
2:B:117:LEU:HD11	2:B:225:ALA:HB1	2.03	0.41
1:A:122:ALA:O	1:A:123:GLU:HB2	2.21	0.41
2:D:249:THR:HG22	2:D:309:ALA:HB3	2.03	0.40
2:D:311:THR:HA	2:D:333:LEU:O	2.21	0.40
2:B:121:SER:HB2	12:B:2004:MWI:O4	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	456/457 (100%)	438 (96%)	17 (4%)	1 (0%)	43	72
1	C	455/457 (100%)	434 (95%)	21 (5%)	0	100	100
2	B	467/472 (99%)	448 (96%)	19 (4%)	0	100	100
2	D	469/472 (99%)	454 (97%)	13 (3%)	2 (0%)	30	60

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	E	210/221 (95%)	196 (93%)	14 (7%)	0	100	100
3	H	212/221 (96%)	199 (94%)	13 (6%)	0	100	100
4	F	212/214 (99%)	199 (94%)	13 (6%)	0	100	100
4	L	212/214 (99%)	204 (96%)	7 (3%)	1 (0%)	24	55
All	All	2693/2728 (99%)	2572 (96%)	117 (4%)	4 (0%)	48	77

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	374	CYS
1	A	123	GLU
2	D	157	VAL
4	L	68	GLY

5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

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](#)

13 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	G	1	5,2	14,14,15	0.31	0	17,19,21	0.51	0
5	NAG	G	2	5	14,14,15	0.30	0	17,19,21	0.62	0
5	BMA	G	3	5	11,11,12	0.97	1 (9%)	15,15,17	0.93	0
5	MAN	G	4	5	11,11,12	1.27	1 (9%)	15,15,17	0.89	1 (6%)
5	MAN	G	5	5	11,11,12	1.42	1 (9%)	15,15,17	1.20	2 (13%)
6	NAG	I	1	6,2	14,14,15	0.46	0	17,19,21	0.42	0
6	NAG	I	2	6	14,14,15	0.46	0	17,19,21	0.43	0
7	NAG	J	1	2,7	14,14,15	0.34	0	17,19,21	0.75	1 (5%)
7	NAG	J	2	7	14,14,15	0.47	0	17,19,21	0.62	1 (5%)
7	BMA	J	3	7	11,11,12	1.03	0	15,15,17	0.84	0
7	MAN	J	4	7	11,11,12	1.24	2 (18%)	15,15,17	0.92	1 (6%)
6	NAG	K	1	6,2	14,14,15	0.60	0	17,19,21	0.55	0
6	NAG	K	2	6	14,14,15	0.52	0	17,19,21	0.55	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	G	1	5,2	-	0/6/23/26	0/1/1/1
5	NAG	G	2	5	-	0/6/23/26	0/1/1/1
5	BMA	G	3	5	-	2/2/19/22	0/1/1/1
5	MAN	G	4	5	-	0/2/19/22	0/1/1/1
5	MAN	G	5	5	-	2/2/19/22	0/1/1/1
6	NAG	I	1	6,2	-	0/6/23/26	0/1/1/1
6	NAG	I	2	6	-	3/6/23/26	0/1/1/1
7	NAG	J	1	2,7	-	0/6/23/26	0/1/1/1
7	NAG	J	2	7	-	0/6/23/26	0/1/1/1
7	BMA	J	3	7	-	1/2/19/22	0/1/1/1
7	MAN	J	4	7	-	0/2/19/22	0/1/1/1
6	NAG	K	1	6,2	-	2/6/23/26	0/1/1/1
6	NAG	K	2	6	-	4/6/23/26	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	G	5	MAN	C1-C2	3.50	1.60	1.52
5	G	4	MAN	O5-C5	2.92	1.49	1.43
5	G	3	BMA	C4-C5	2.50	1.58	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	J	4	MAN	O5-C5	2.42	1.48	1.43
7	J	4	MAN	C1-C2	2.34	1.57	1.52

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	J	1	NAG	C1-O5-C5	2.62	115.70	112.19
5	G	5	MAN	C1-C2-C3	2.57	113.39	109.64
7	J	4	MAN	O2-C2-C3	-2.31	105.36	110.15
5	G	4	MAN	O2-C2-C3	-2.28	105.43	110.15
7	J	2	NAG	C1-O5-C5	2.17	115.09	112.19
5	G	5	MAN	C1-O5-C5	2.11	115.02	112.19

There are no chirality outliers.

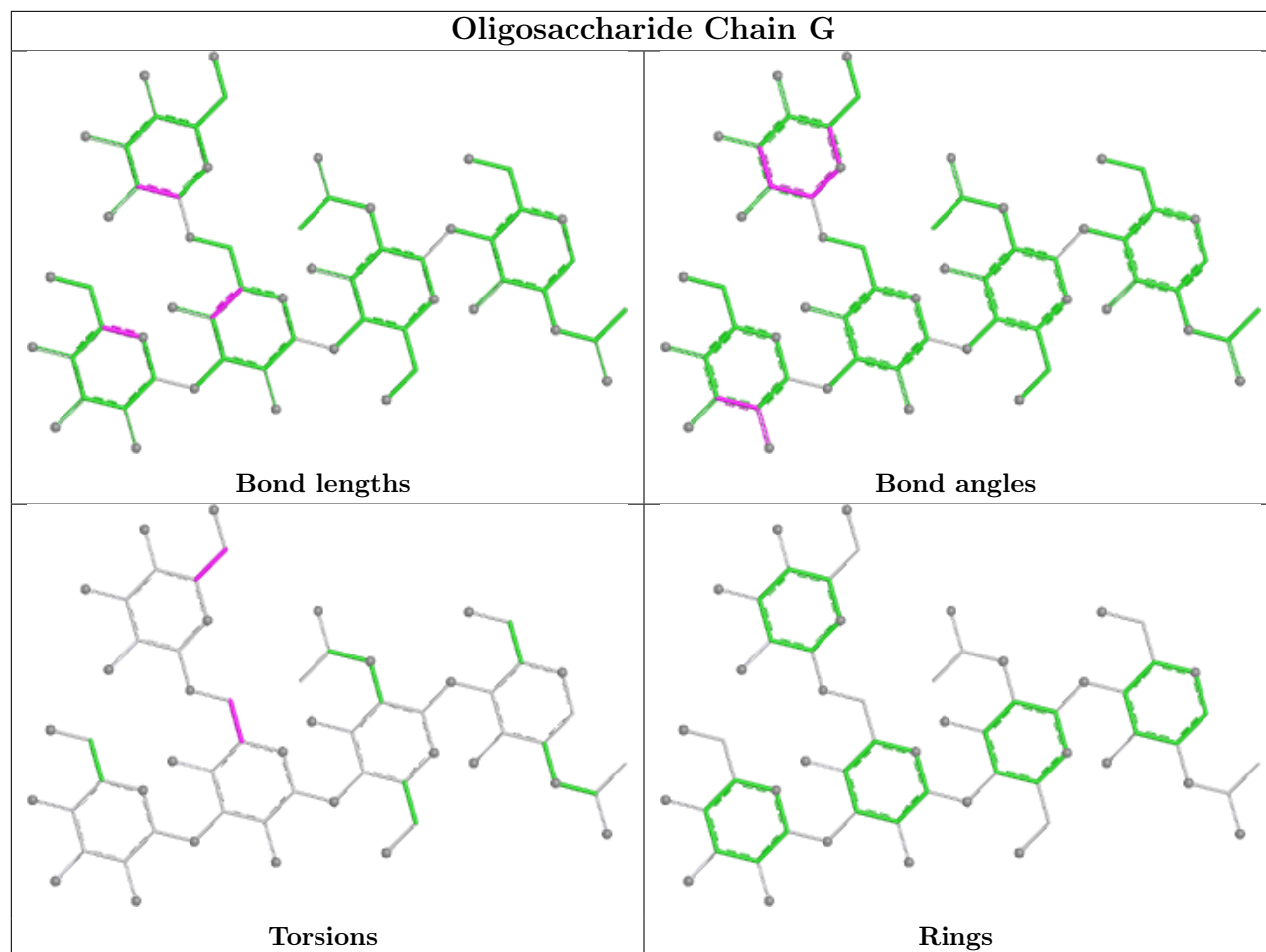
All (14) torsion outliers are listed below:

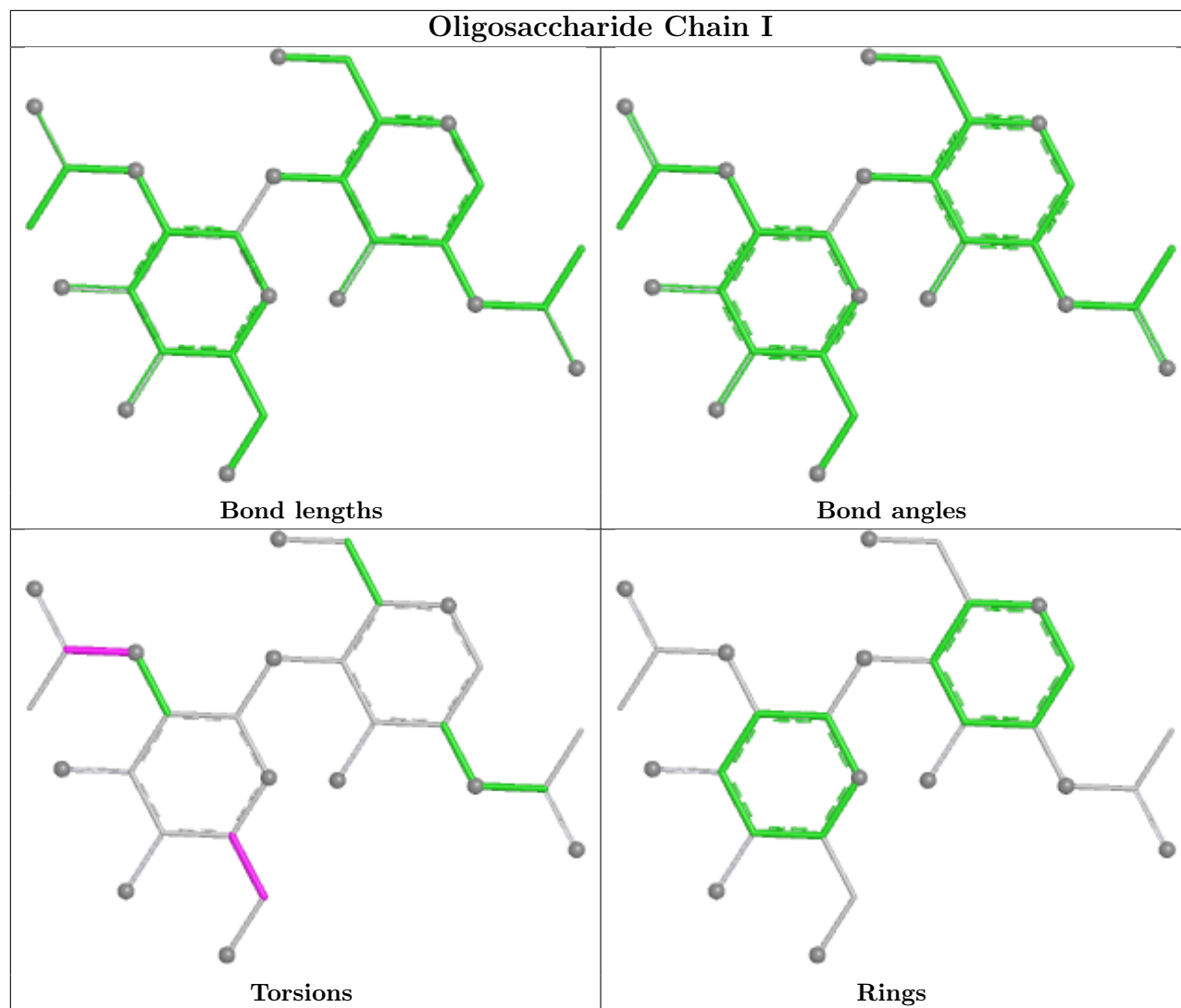
Mol	Chain	Res	Type	Atoms
5	G	5	MAN	O5-C5-C6-O6
5	G	5	MAN	C4-C5-C6-O6
6	I	2	NAG	C8-C7-N2-C2
6	I	2	NAG	O7-C7-N2-C2
6	K	2	NAG	C8-C7-N2-C2
6	K	2	NAG	O7-C7-N2-C2
6	K	2	NAG	O5-C5-C6-O6
6	K	2	NAG	C4-C5-C6-O6
7	J	3	BMA	O5-C5-C6-O6
6	K	1	NAG	O5-C5-C6-O6
5	G	3	BMA	O5-C5-C6-O6
6	I	2	NAG	O5-C5-C6-O6
6	K	1	NAG	C4-C5-C6-O6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

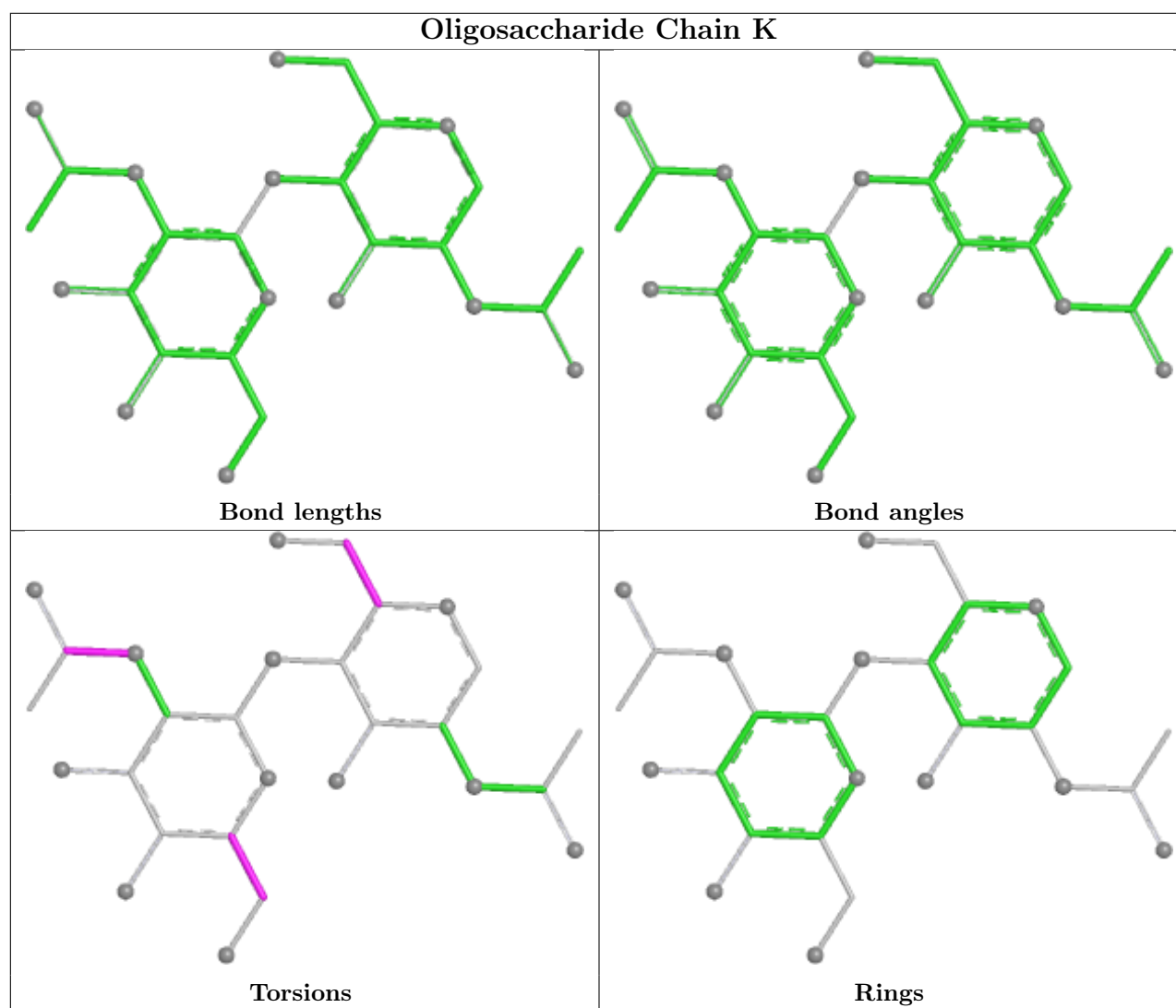
There are no ring outliers.

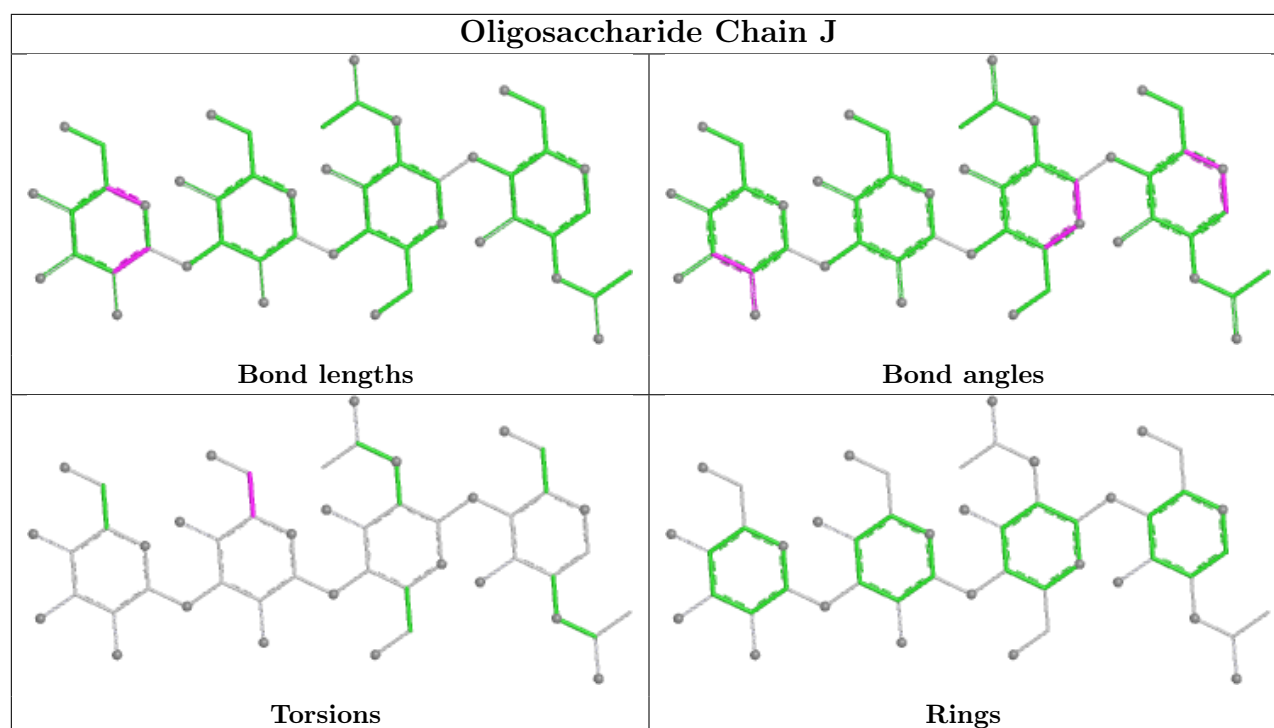
No monomer is involved in short contacts.

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](#)

Of 28 ligands modelled in this entry, 15 are monoatomic - leaving 13 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
8	SO4	C	510	-	4,4,4	0.61	0	6,6,6	0.22	0
12	MWI	D	2005	10	33,34,34	1.02	1 (3%)	42,48,48	1.28	4 (9%)
11	NAG	B	2003	2	14,14,15	0.67	0	17,19,21	0.84	1 (5%)
8	SO4	A	508	-	4,4,4	0.57	0	6,6,6	0.20	0
8	SO4	C	501	-	4,4,4	0.55	0	6,6,6	0.23	0
8	SO4	C	502	-	4,4,4	0.63	0	6,6,6	0.47	0
12	MWI	B	2004	10	33,34,34	0.95	0	42,48,48	1.31	4 (9%)
8	SO4	A	501	-	4,4,4	0.58	0	6,6,6	0.24	0
8	SO4	L	301	-	4,4,4	0.59	0	6,6,6	0.22	0
8	SO4	A	502	-	4,4,4	0.44	0	6,6,6	0.30	0
8	SO4	C	503	-	4,4,4	0.58	0	6,6,6	0.34	0
8	SO4	A	507	-	4,4,4	0.62	0	6,6,6	0.43	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
11	NAG	D	2003	2	14,14,15	0.42	0	17,19,21	0.71	1 (5%)

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
11	NAG	B	2003	2	-	2/6/23/26	0/1/1/1
12	MWI	D	2005	10	-	1/16/50/50	0/4/4/4
11	NAG	D	2003	2	-	0/6/23/26	0/1/1/1
12	MWI	B	2004	10	-	4/16/50/50	0/4/4/4

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
12	D	2005	MWI	C16-C14	2.07	1.42	1.40

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
12	B	2004	MWI	C22-N4-C20	5.58	123.53	118.58
12	D	2005	MWI	C22-N4-C20	4.68	122.73	118.58
11	B	2003	NAG	C1-O5-C5	2.88	116.04	112.19
12	D	2005	MWI	C12-C11-N1	2.74	122.08	118.66
12	B	2004	MWI	C12-C11-N1	2.60	121.91	118.66
11	D	2003	NAG	C1-O5-C5	2.51	115.54	112.19
12	B	2004	MWI	C19-C16-N3	-2.46	114.76	119.20
12	D	2005	MWI	C8-N1-C11	2.08	129.22	122.79
12	D	2005	MWI	C19-C16-N3	-2.08	115.45	119.20
12	B	2004	MWI	C14-C16-N3	2.05	125.39	121.69

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
12	B	2004	MWI	C3-C1-C2-C5
12	B	2004	MWI	C3-C1-C2-C6
12	B	2004	MWI	C4-C1-C2-C5
12	B	2004	MWI	C4-C1-C2-C6

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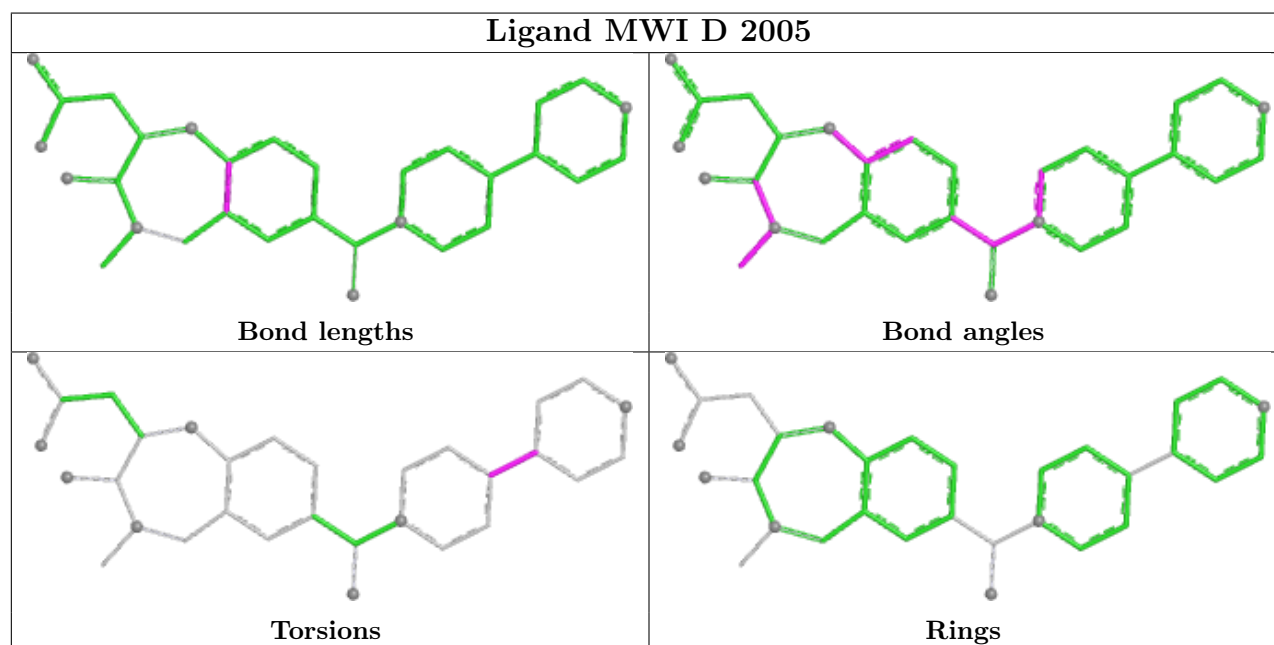
Mol	Chain	Res	Type	Atoms
11	B	2003	NAG	O5-C5-C6-O6
11	B	2003	NAG	C4-C5-C6-O6
12	D	2005	MWI	C3-C1-C2-C5

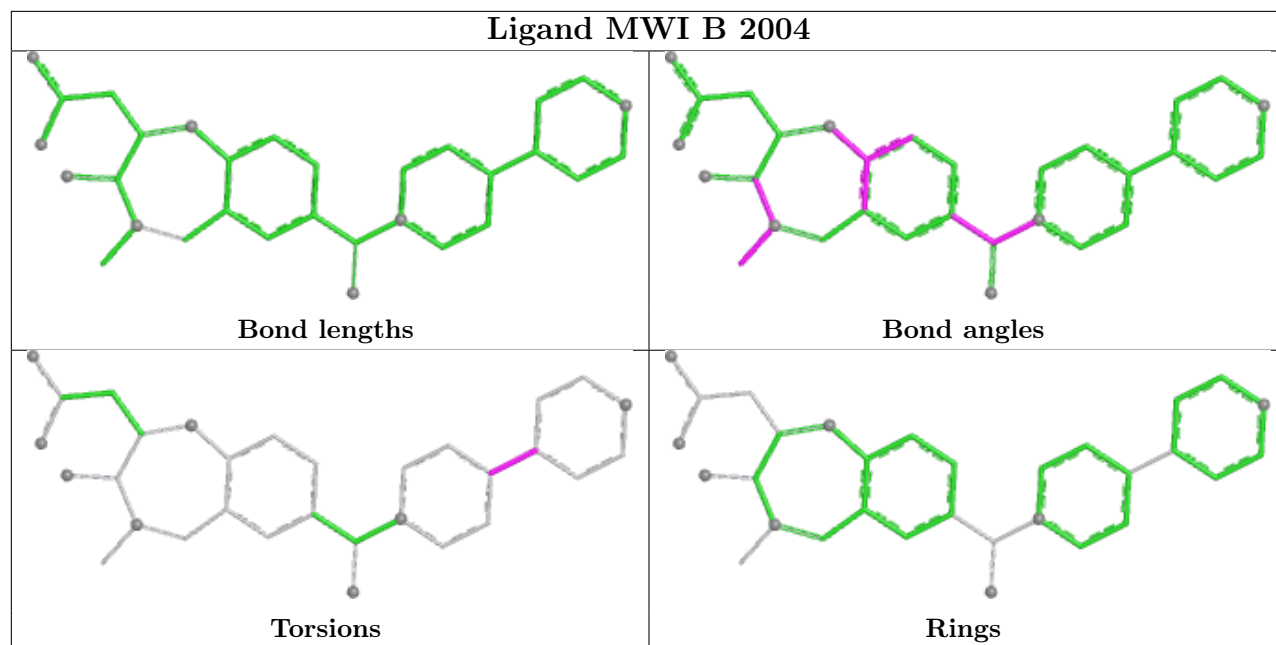
There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
12	D	2005	MWI	1	0
12	B	2004	MWI	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less 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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ > 2	OWAB(Å ²)	Q < 0.9
1	A	454/457 (99%)	-0.50	1 (0%) 91 88	34, 61, 90, 128	4 (0%)
1	C	453/457 (99%)	-0.31	0 100 100	34, 82, 114, 153	4 (0%)
2	B	466/472 (98%)	-0.20	1 (0%) 91 88	29, 93, 169, 191	4 (0%)
2	D	471/472 (99%)	-0.07	3 (0%) 85 80	60, 112, 158, 183	1 (0%)
3	E	214/221 (96%)	0.30	3 (1%) 73 64	100, 154, 225, 236	0
3	H	216/221 (97%)	-0.06	0 100 100	71, 115, 168, 184	0
4	F	214/214 (100%)	0.19	2 (0%) 81 74	103, 161, 218, 235	0
4	L	214/214 (100%)	-0.19	0 100 100	78, 107, 133, 174	0
All	All	2702/2728 (99%)	-0.16	10 (0%) 88 84	29, 99, 187, 236	13 (0%)

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	454	VAL	3.5
3	E	201	CYS	3.1
3	E	212	VAL	3.0
2	D	375	LEU	2.9
2	D	181	LYS	2.7
4	F	178	THR	2.3
2	B	129	TRP	2.3
3	E	203	VAL	2.3
4	F	115	VAL	2.3
2	D	129	TRP	2.1

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

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

6.3 Carbohydrates

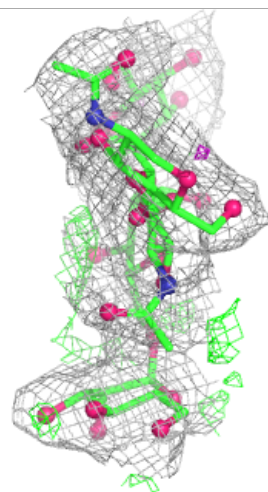
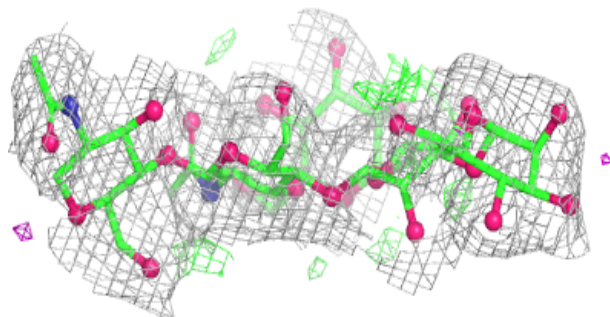
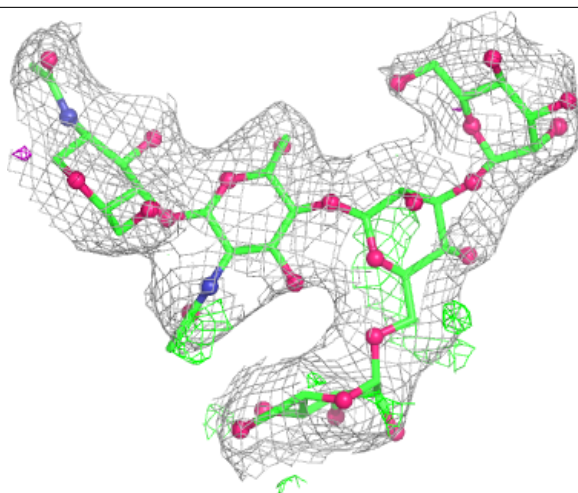
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	G	1	14/15	-	-	75,75,75,75	0
5	NAG	G	2	14/15	-	-	96,96,96,96	0
5	BMA	G	3	11/12	-	-	112,112,112,112	0
5	MAN	G	4	11/12	-	-	110,110,110,110	0
5	MAN	G	5	11/12	-	-	119,119,119,119	0
7	NAG	J	2	14/15	0.67	0.10	118,118,118,118	0
6	NAG	K	2	14/15	0.79	0.09	130,130,130,130	0
6	NAG	I	2	14/15	0.83	0.10	134,134,134,134	0
6	NAG	K	1	14/15	0.85	0.10	131,131,131,131	0
7	NAG	J	1	14/15	0.87	0.09	105,105,105,105	0
6	NAG	I	1	14/15	0.95	0.07	126,126,126,126	0
7	BMA	J	3	11/12	-	-	128,128,128,128	0
7	MAN	J	4	11/12	-	-	127,127,127,127	0

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

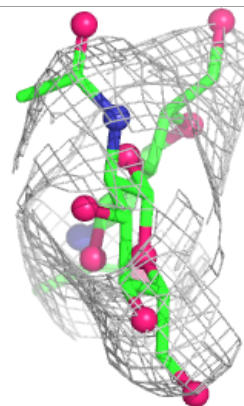
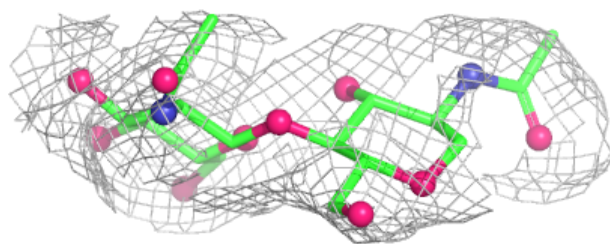
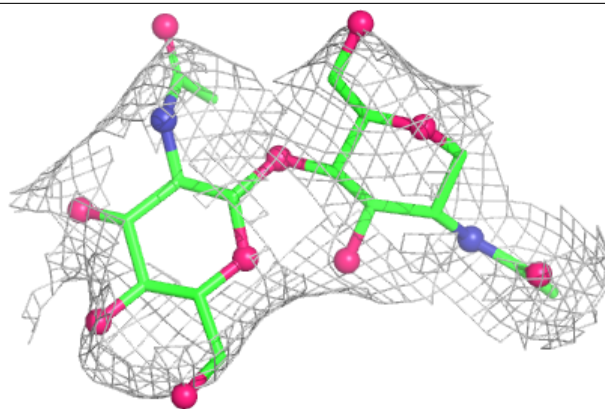
Electron density around Chain G:

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

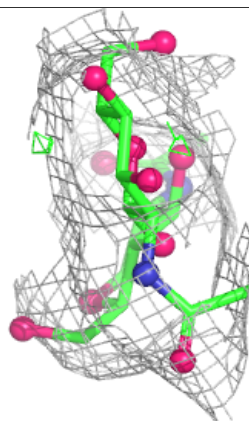
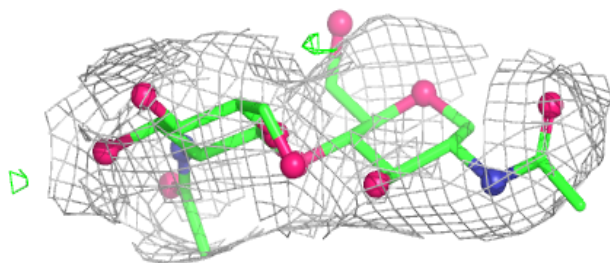
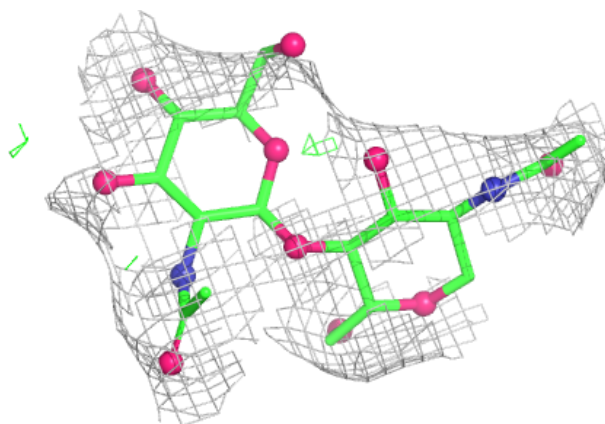


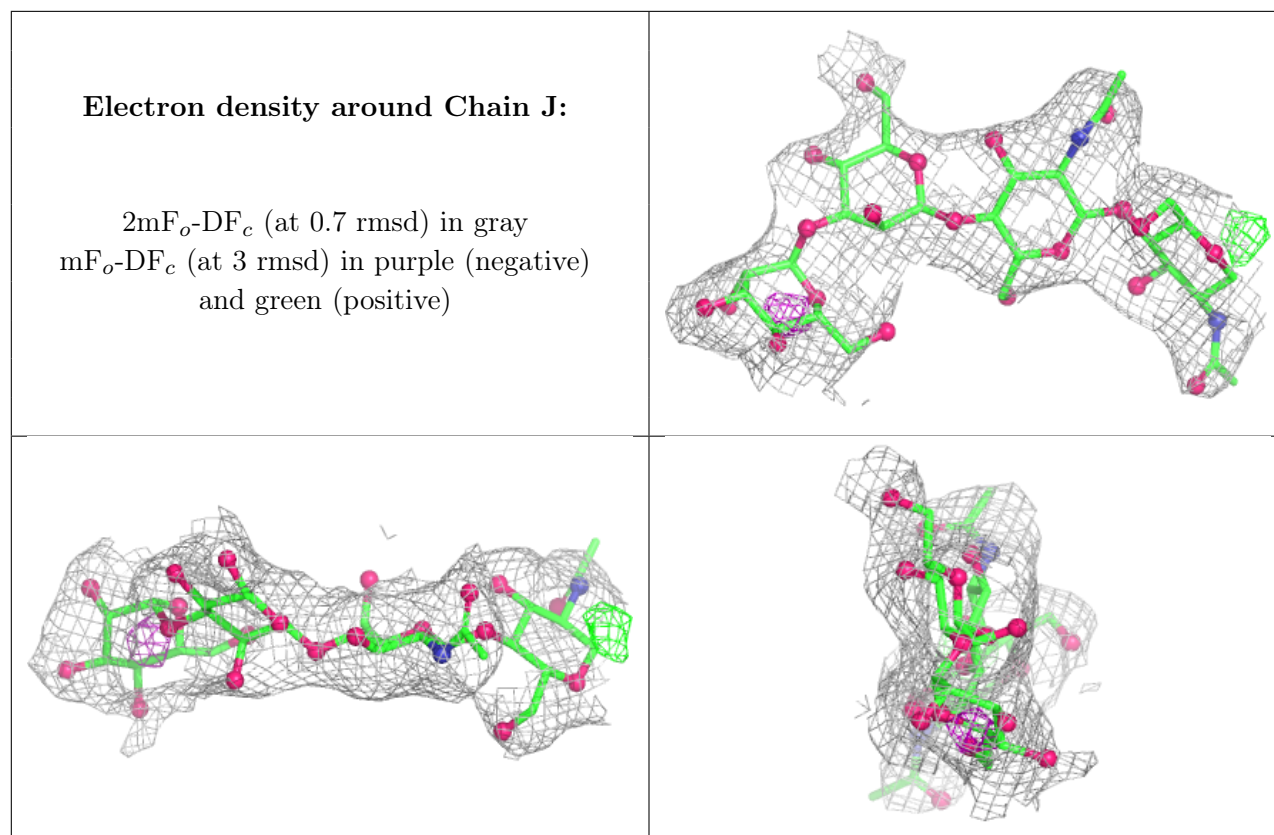
Electron density around Chain I:

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

**Electron density around Chain K:**

$2mF_o - DF_c$ (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
8	SO4	C	502	5/5	0.66	0.13	116,116,116,116	0
8	SO4	C	503	5/5	0.72	0.08	133,133,133,133	0
11	NAG	B	2003	14/15	0.73	0.10	135,135,135,135	0
11	NAG	D	2003	14/15	0.78	0.10	121,121,121,121	0
8	SO4	C	501	5/5	0.82	0.17	104,104,104,104	0
8	SO4	L	301	5/5	0.87	0.13	112,112,112,112	0
8	SO4	A	508	5/5	0.88	0.10	109,109,109,109	0
13	CL	D	2004	1/1	0.88	0.11	105,105,105,105	0
8	SO4	A	507	5/5	0.89	0.15	106,106,106,106	0
13	CL	C	505	1/1	0.89	0.08	110,110,110,110	0
8	SO4	C	510	5/5	0.89	0.07	120,120,120,120	0
8	SO4	A	501	5/5	0.90	0.26	87,87,87,87	0
8	SO4	A	502	5/5	0.94	0.06	80,80,80,80	0
12	MWI	D	2005	31/31	0.94	0.10	83,83,83,83	0

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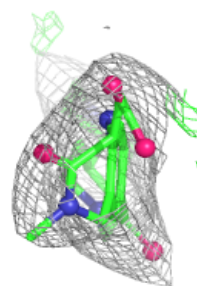
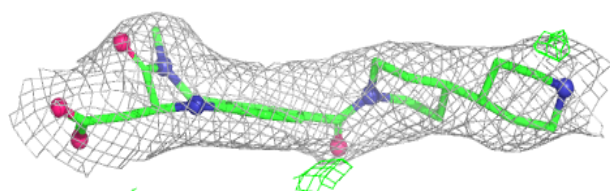
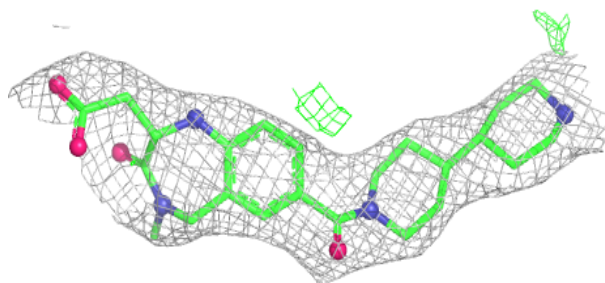
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
12	MWI	B	2004	31/31	0.95	0.09	61,61,61,61	0
13	CL	C	504	1/1	0.95	0.15	84,84,84,84	0
10	MG	D	2001	1/1	0.97	0.05	86,86,86,86	0
9	CA	C	506	1/1	0.98	0.04	111,111,111,111	0
9	CA	D	2002	1/1	0.98	0.03	72,72,72,72	0
9	CA	C	509	1/1	0.99	0.03	70,70,70,70	0
9	CA	A	505	1/1	0.99	0.04	50,50,50,50	0
9	CA	A	504	1/1	0.99	0.02	55,55,55,55	0
9	CA	C	507	1/1	0.99	0.03	91,91,91,91	0
9	CA	C	508	1/1	0.99	0.03	77,77,77,77	0
9	CA	B	2002	1/1	1.00	0.01	51,51,51,51	0
9	CA	A	503	1/1	1.00	0.03	72,72,72,72	0
9	CA	A	506	1/1	1.00	0.02	56,56,56,56	0
10	MG	B	2001	1/1	1.00	0.04	49,49,49,49	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.

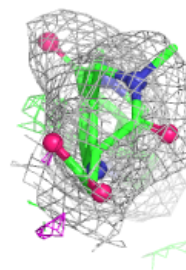
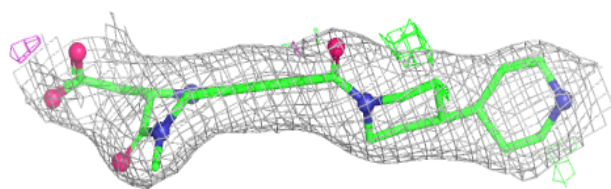
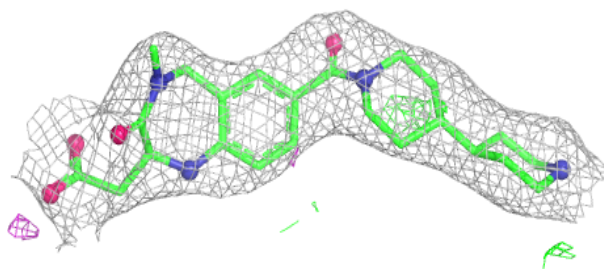
Electron density around MWI D 2005:

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 MWI B 2004:

$2mF_o - DF_c$ (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.