



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 09:35 PM UTC

PDB ID : 6X4H / pdb_00006x4h
Title : Sortilin-Progranulin Interaction With Compound 24
Authors : Parthasarathy, G.; Soisson, S.M.; Klein, D.
Deposited on : 2020-05-22
Resolution : 2.90 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

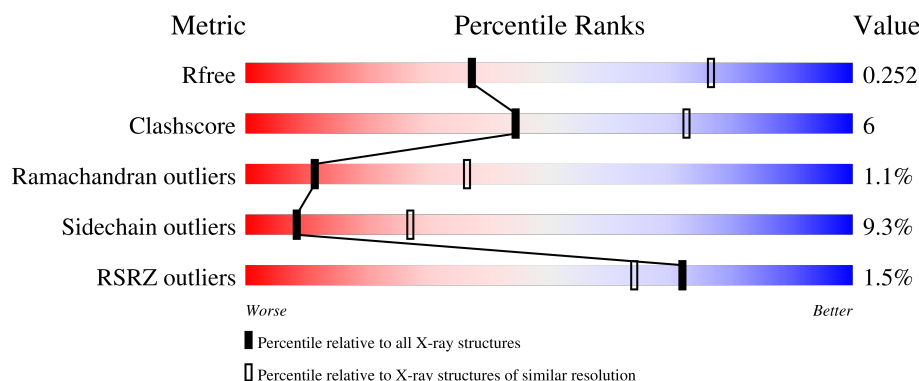
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

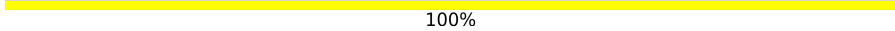
The reported resolution of this entry is 2.90 Å.

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	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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	663	 2% 76% 21% ..
2	B	2	 50% 50%
2	C	2	 100%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 5240 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 Sortilin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	654	Total	C	N	O	S	0	0	0
			5143	3241	867	1006	29			

There is a discrepancy between the modelled and reference sequences:

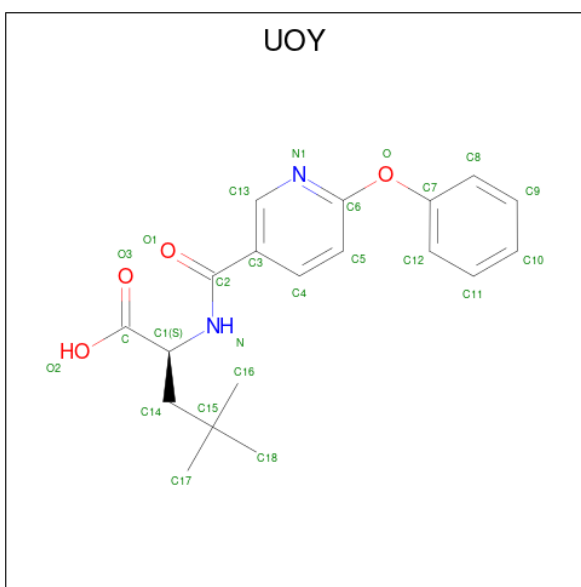
Chain	Residue	Modelled	Actual	Comment	Reference
A	617	MET	VAL	conflict	UNP Q99523

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



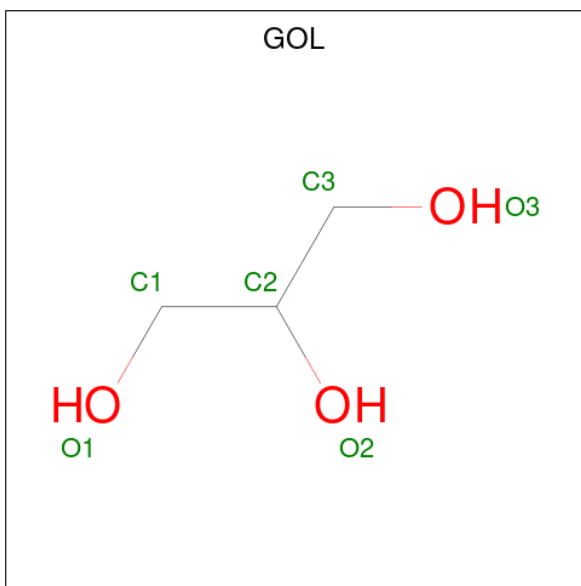
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	C	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 3 is 4-methyl-N-(6-phenoxy-pyridine-3-carbonyl)-L-leucine (CCD ID: UOY) (formula: C₁₉H₂₂N₂O₄) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			25	19	2	4		

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	10	Total	O	0	0
			10	10		

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	161.46Å 79.44Å 111.55Å 90.00° 127.00° 90.00°	Depositor
Resolution (Å)	32.29 – 2.90 32.29 – 2.90	Depositor EDS
% Data completeness (in resolution range)	97.6 (32.29-2.90) 97.8 (32.29-2.90)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.03 (at 2.90Å)	Xtriage
Refinement program	BUSTER 2.11.5	Depositor
R, R_{free}	0.201 , 0.253 0.208 , 0.252	Depositor DCC
R_{free} test set	1245 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	83.3	Xtriage
Anisotropy	0.446	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 61.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5240	wwPDB-VP
Average B, all atoms (Å ²)	92.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.70% 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: NAG, UOY, GOL

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	0.87	0/5262	1.33	34/7125 (0.5%)

There are no bond length outliers.

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	668	PHE	CA-C-N	7.09	130.74	120.38
1	A	668	PHE	C-N-CA	7.09	130.74	120.38
1	A	678	THR	CB-CA-C	6.72	120.82	109.80
1	A	363	THR	CA-C-N	6.69	129.25	120.28
1	A	363	THR	C-N-CA	6.69	129.25	120.28
1	A	470	PRO	N-CA-C	6.34	121.46	111.38
1	A	347	ASP	CA-CB-CG	6.17	118.77	112.60
1	A	614	LYS	CA-C-N	6.13	128.82	120.54
1	A	614	LYS	C-N-CA	6.13	128.82	120.54
1	A	320	ILE	N-CA-CB	5.86	117.75	112.06
1	A	643	ASP	CA-CB-CG	5.83	118.43	112.60
1	A	450	ASN	CA-CB-CG	5.69	118.29	112.60
1	A	97	PHE	CA-CB-CG	5.64	119.44	113.80
1	A	416	ASP	CA-CB-CG	5.64	118.24	112.60
1	A	651	ASN	CA-C-N	5.59	129.64	121.31
1	A	651	ASN	C-N-CA	5.59	129.64	121.31
1	A	191	PRO	CA-C-N	5.58	128.03	120.44
1	A	191	PRO	C-N-CA	5.58	128.03	120.44
1	A	705	LYS	N-CA-C	-5.53	106.58	113.55
1	A	455	VAL	N-CA-CB	5.42	119.43	111.52
1	A	155	VAL	CA-C-N	5.37	128.33	120.71
1	A	155	VAL	C-N-CA	5.37	128.33	120.71
1	A	519	ASP	CA-CB-CG	5.37	117.97	112.60
1	A	501	ILE	N-CA-C	-5.36	100.67	108.17

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	115	GLU	N-CA-C	-5.33	106.27	114.16
1	A	279	PHE	N-CA-C	5.16	118.47	110.17
1	A	136	GLU	N-CA-C	-5.12	105.89	111.82
1	A	679	THR	N-CA-C	5.05	117.06	109.23
1	A	61	ALA	CA-C-N	5.04	127.54	120.28
1	A	61	ALA	C-N-CA	5.04	127.54	120.28
1	A	459	GLY	CA-C-N	5.03	129.80	121.86
1	A	459	GLY	C-N-CA	5.03	129.80	121.86
1	A	231	ASP	CA-C-N	5.02	129.25	122.07
1	A	231	ASP	C-N-CA	5.02	129.25	122.07

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	5143	0	4933	57	0
2	B	28	0	25	0	0
2	C	28	0	25	0	0
3	A	25	0	0	0	0
4	A	6	0	8	0	0
5	A	10	0	0	0	0
All	All	5240	0	4991	57	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 6.

All (57) 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:A:416:ASP:OD1	1:A:418:THR:HG22	1.57	1.04
1:A:579:CYS:HG	1:A:618:CYS:HG	0.91	0.87
1:A:415:CYS:HG	1:A:425:CYS:HG	1.22	0.82
1:A:669:CYS:SG	1:A:707:CYS:HB3	2.28	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:513:VAL:HG11	1:A:526:THR:HG23	1.78	0.66
1:A:75:LEU:HD22	1:A:564:TRP:CD1	2.33	0.63
1:A:632:ILE:HG22	1:A:691:GLN:HG2	1.81	0.63
1:A:513:VAL:CG1	1:A:526:THR:HG23	2.29	0.62
1:A:95:THR:HG23	1:A:110:LYS:HB2	1.81	0.61
1:A:493:THR:HG21	1:A:550:ILE:HD13	1.82	0.60
1:A:134:ARG:HE	1:A:137:PHE:HE2	1.50	0.58
1:A:349:ARG:HB3	1:A:351:ILE:HD12	1.86	0.56
1:A:540:ALA:HB3	1:A:551:SER:HB2	1.87	0.56
1:A:72:PHE:HB3	1:A:75:LEU:HD12	1.87	0.55
1:A:579:CYS:HG	1:A:618:CYS:CB	2.20	0.54
1:A:286:ALA:HB3	1:A:289:ASP:HB3	1.89	0.54
1:A:125:THR:HG22	1:A:130:ASN:HA	1.89	0.54
1:A:206:LEU:HB2	1:A:223:VAL:HG21	1.90	0.53
1:A:536:PHE:HZ	1:A:539:LEU:HG	1.74	0.53
1:A:78:SER:HB3	1:A:96:THR:H	1.74	0.53
1:A:187:MET:HG2	1:A:199:ALA:HB2	1.92	0.51
1:A:281:PHE:CZ	1:A:328:VAL:HG11	2.45	0.51
1:A:75:LEU:HD21	1:A:79:VAL:HG13	1.93	0.50
1:A:323:ALA:CB	1:A:328:VAL:HG12	2.42	0.50
1:A:558:SER:HB3	1:A:562:SER:HB2	1.94	0.49
1:A:418:THR:HG23	1:A:463:ASP:H	1.77	0.49
1:A:281:PHE:HZ	1:A:328:VAL:HG11	1.77	0.49
1:A:632:ILE:CG2	1:A:691:GLN:HG2	2.42	0.48
1:A:667:GLU:O	1:A:671:TYR:HD1	1.97	0.47
1:A:377:LEU:HD11	1:A:589:ALA:HB1	1.95	0.47
1:A:72:PHE:HE1	1:A:566:SER:HB3	1.80	0.47
1:A:140:ALA:HB3	1:A:150:VAL:HB	1.97	0.47
1:A:351:ILE:HD13	1:A:676:HIS:CE1	2.50	0.47
1:A:699:GLU:HB3	1:A:701:LYS:HE2	1.96	0.46
1:A:293:ARG:HD2	1:A:306:MET:HE2	1.97	0.46
1:A:305:SER:HB3	1:A:676:HIS:NE2	2.30	0.46
1:A:69:GLN:HG2	1:A:70:HIS:N	2.32	0.45
1:A:321:LEU:HD11	1:A:331:HIS:HB2	1.98	0.45
1:A:380:VAL:HA	1:A:397:THR:O	2.17	0.45
1:A:634:LEU:HA	1:A:691:GLN:O	2.16	0.45
1:A:243:SER:H	1:A:246:ALA:HB3	1.82	0.44
1:A:442:PRO:HD2	1:A:458:HIS:CE1	2.52	0.44
1:A:493:THR:HG22	1:A:502:VAL:HB	1.98	0.44
1:A:606:LYS:HB2	1:A:631:SER:HB2	2.00	0.44
1:A:258:ASP:HB2	1:A:261:LYS:HB2	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:124:ILE:HB	1:A:127:LEU:HD12	1.99	0.43
1:A:395:MET:HE1	1:A:406:HIS:CE1	2.53	0.43
1:A:703:LEU:C	1:A:705:LYS:H	2.25	0.42
1:A:142:GLY:HA3	1:A:148:LYS:HB2	2.02	0.42
1:A:500:ILE:HG22	1:A:502:VAL:HG23	2.01	0.42
1:A:395:MET:HE1	1:A:406:HIS:CG	2.55	0.42
1:A:540:ALA:O	1:A:550:ILE:HD12	2.20	0.42
1:A:306:MET:HE1	1:A:671:TYR:HB2	2.02	0.41
1:A:185:THR:HG23	1:A:200:LEU:HB2	2.03	0.41
1:A:458:HIS:CD2	1:A:472:VAL:HG22	2.56	0.40
1:A:495:LEU:HB2	1:A:500:ILE:HB	2.04	0.40
1:A:394:THR:O	1:A:395:MET:HE2	2.22	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	648/663 (98%)	601 (93%)	40 (6%)	7 (1%)	11	36

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	73	ASP
1	A	337	ASP
1	A	242	GLY
1	A	130	ASN
1	A	241	ASN
1	A	204	ASN
1	A	79	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	570/579 (98%)	517 (91%)	53 (9%)	8 27

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	60	VAL
1	A	79	VAL
1	A	88	THR
1	A	95	THR
1	A	109	SER
1	A	135	THR
1	A	194	SER
1	A	200	LEU
1	A	217	GLU
1	A	233	THR
1	A	258	ASP
1	A	261	LYS
1	A	283	SER
1	A	296	VAL
1	A	305	SER
1	A	325	ASP
1	A	334	GLU
1	A	338	THR
1	A	352	VAL
1	A	361	LEU
1	A	365	THR
1	A	382	ILE
1	A	389	ASP
1	A	408	ARG
1	A	409	LYS
1	A	413	SER
1	A	426	SER
1	A	439	LEU
1	A	450	ASN
1	A	455	VAL

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Mol	Chain	Res	Type
1	A	460	SER
1	A	483	THR
1	A	493	THR
1	A	497	SER
1	A	502	VAL
1	A	504	ILE
1	A	537	THR
1	A	541	SER
1	A	566	SER
1	A	576	GLU
1	A	580	GLU
1	A	582	LYS
1	A	598	GLU
1	A	602	ILE
1	A	615	SER
1	A	617	MET
1	A	627	THR
1	A	661	LEU
1	A	669	CYS
1	A	676	HIS
1	A	678	THR
1	A	705	LYS
1	A	715	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	65	ASN
1	A	406	HIS
1	A	423	ASN
1	A	512	ASN
1	A	619	GLN
1	A	695	ASN

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 ⓘ

4 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$
2	NAG	B	1	1,2	14,14,15	0.43	0	17,19,21	0.57	0
2	NAG	B	2	2	14,14,15	0.33	0	17,19,21	0.99	2 (11%)
2	NAG	C	1	1,2	14,14,15	0.30	0	17,19,21	1.03	1 (5%)
2	NAG	C	2	2	14,14,15	0.32	0	17,19,21	1.03	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
2	NAG	B	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	B	2	2	-	1/6/23/26	0/1/1/1
2	NAG	C	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	C	2	2	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1	NAG	C1-O5-C5	3.42	116.77	112.19
2	C	2	NAG	C1-O5-C5	2.96	116.16	112.19
2	B	2	NAG	C1-O5-C5	2.54	115.59	112.19
2	B	2	NAG	C1-C2-N2	-2.16	107.02	110.43

There are no chirality outliers.

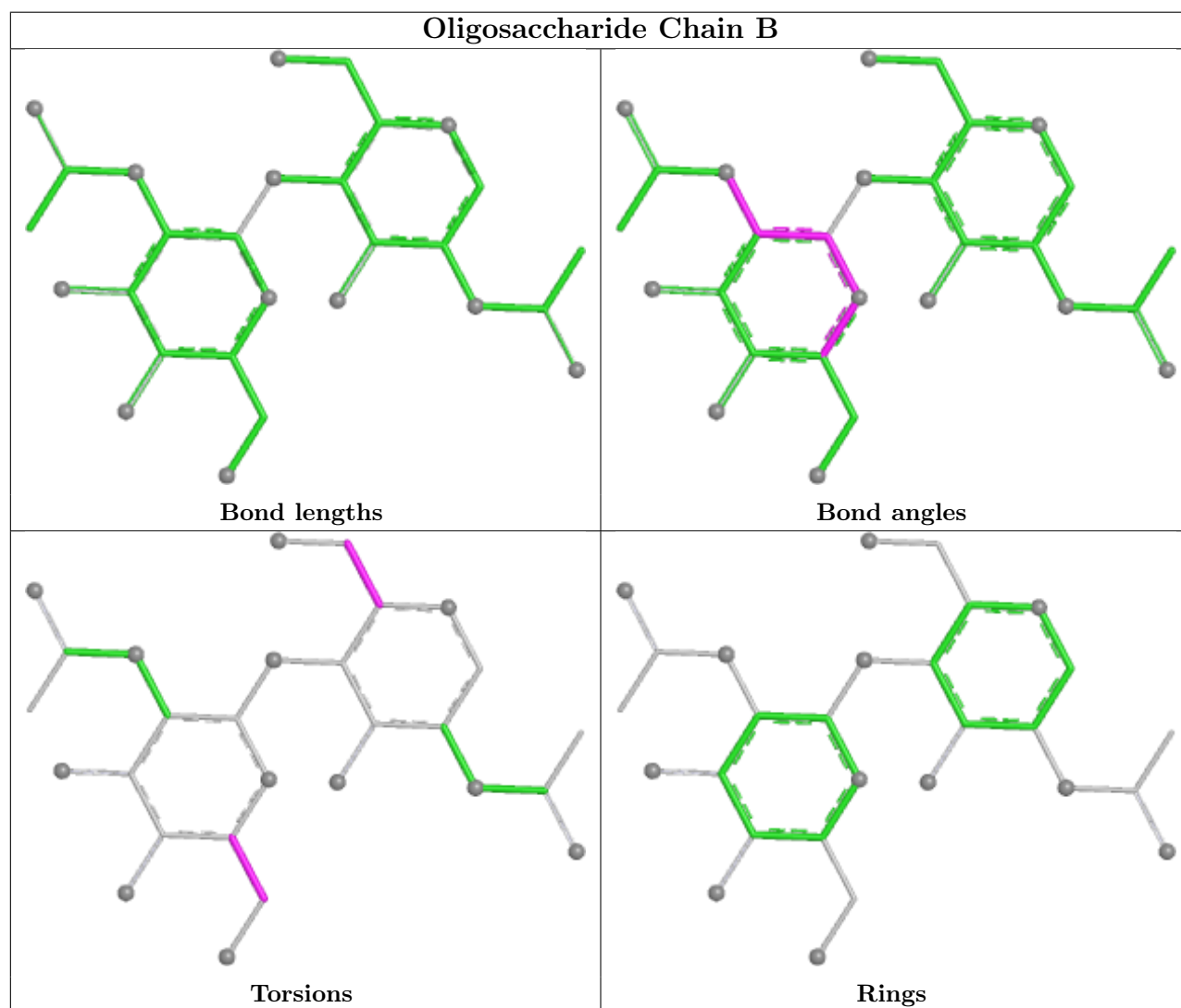
All (3) torsion outliers are listed below:

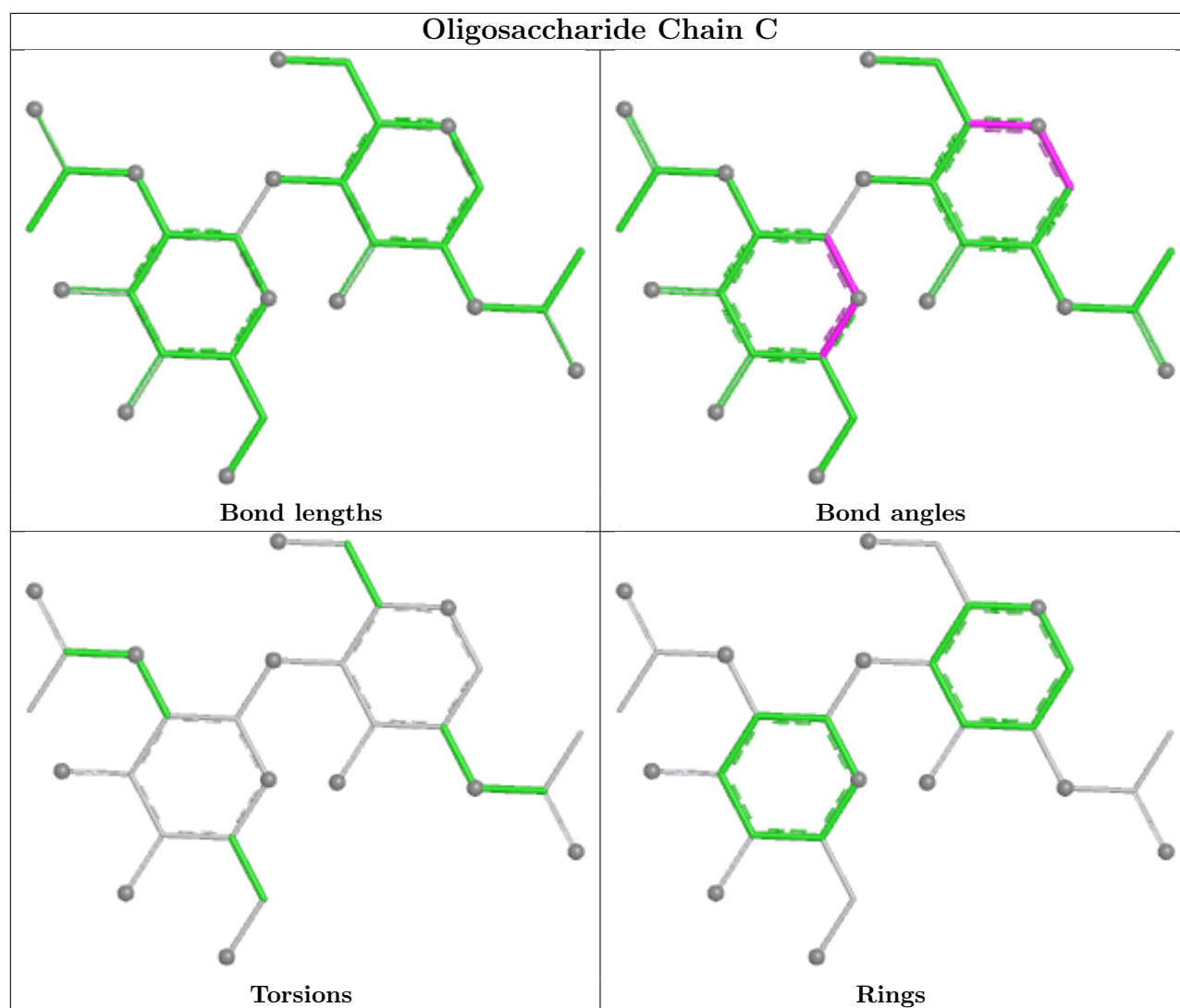
Mol	Chain	Res	Type	Atoms
2	B	2	NAG	O5-C5-C6-O6
2	B	1	NAG	C4-C5-C6-O6
2	B	1	NAG	O5-C5-C6-O6

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](#)

2 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	$\# Z > 2$
4	GOL	A	806	-	5,5,5	0.19	0	5,5,5	0.41	0
3	UOY	A	805	-	26,26,26	0.63	0	35,36,36	0.59	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
4	GOL	A	806	-	-	1/4/4/4	-
3	UOY	A	805	-	-	0/21/21/21	0/2/2/2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

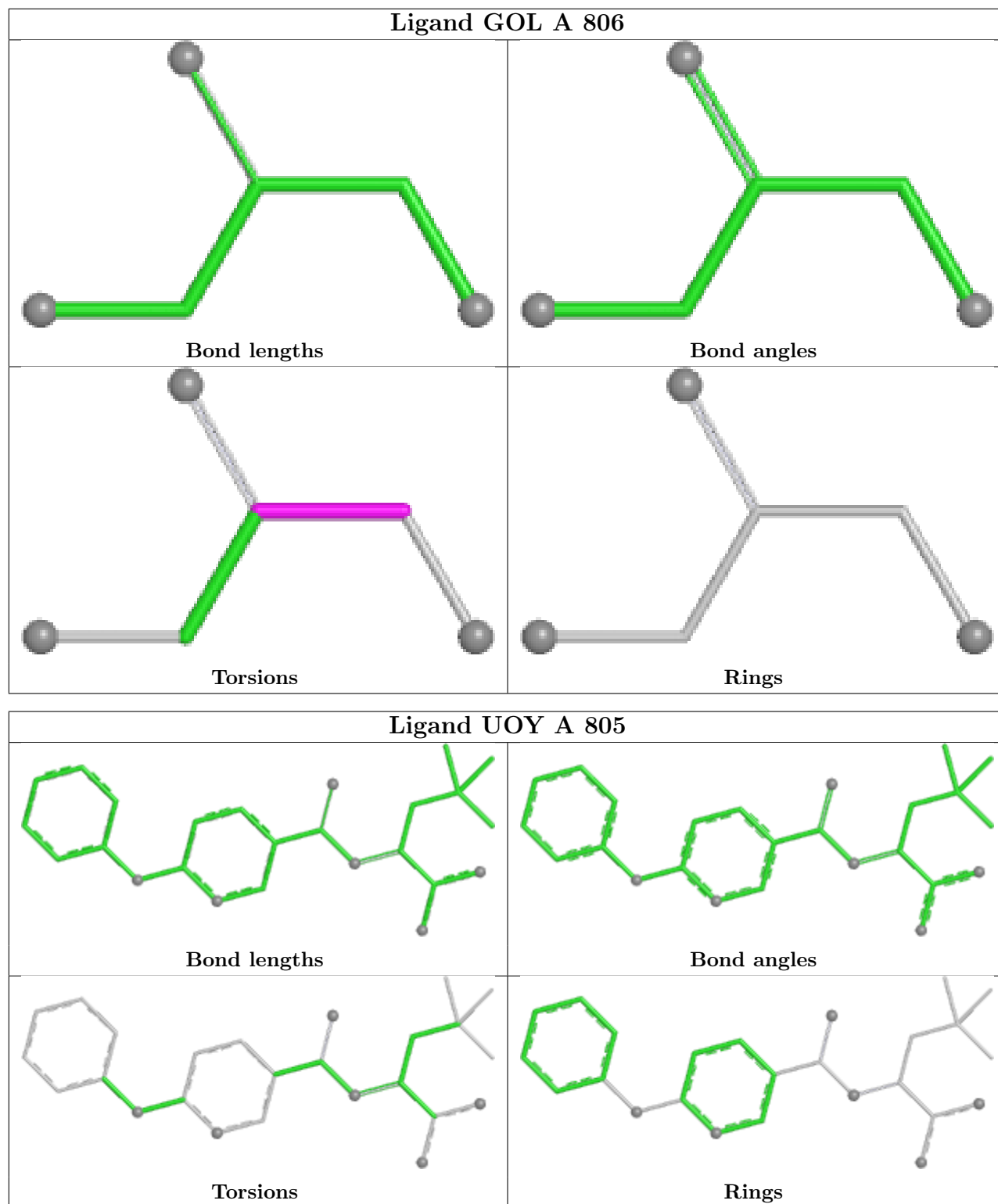
All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	806	GOL	O1-C1-C2-C3

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 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 ⓘ

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	654/663 (98%)	0.12	10 (1%) 72 64	52, 91, 132, 176	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	75	LEU	3.7
1	A	99	VAL	3.6
1	A	561	THR	3.1
1	A	677	LEU	2.6
1	A	700	VAL	2.5
1	A	676	HIS	2.5
1	A	697	VAL	2.4
1	A	696	PRO	2.4
1	A	638	GLU	2.3
1	A	286	ALA	2.2

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.

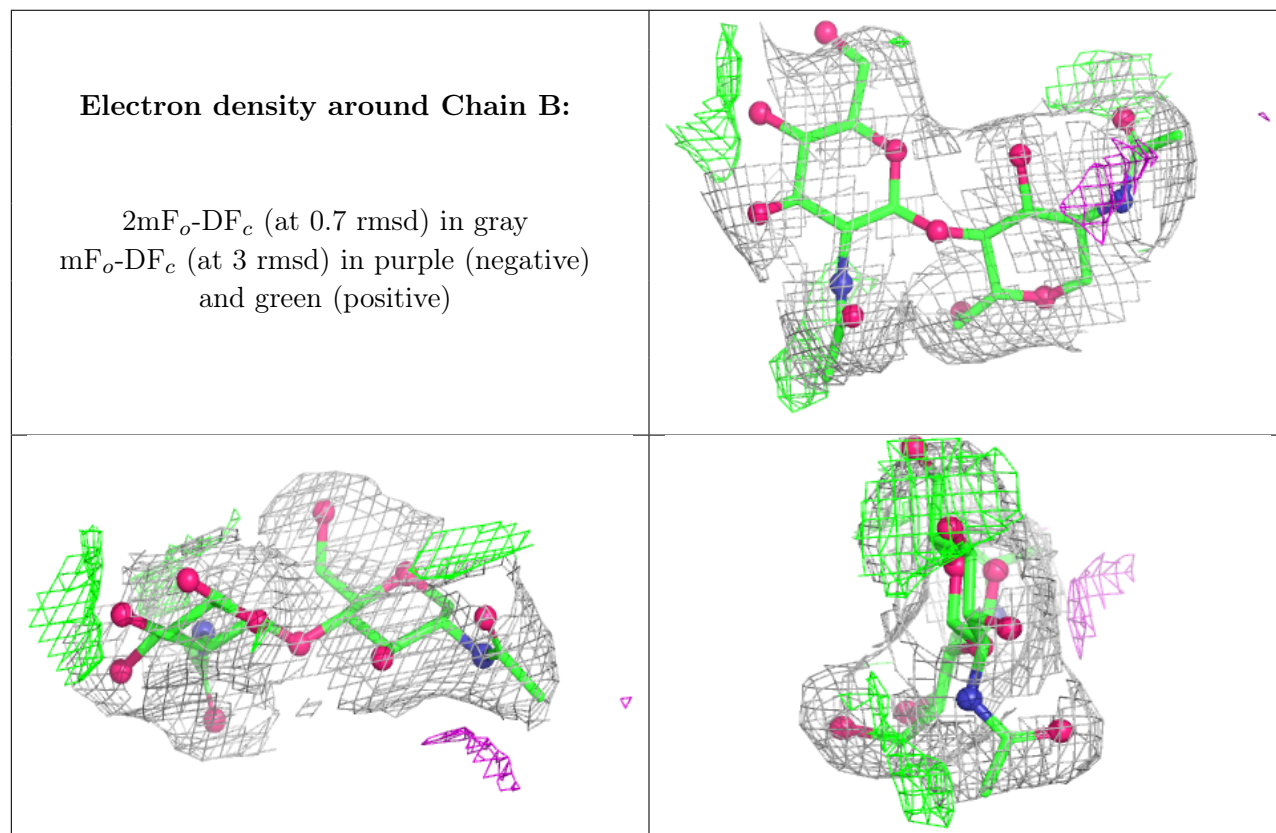
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	NAG	B	2	14/15	0.78	0.13	84,91,102,108	0
2	NAG	C	2	14/15	0.84	0.12	110,115,117,117	0

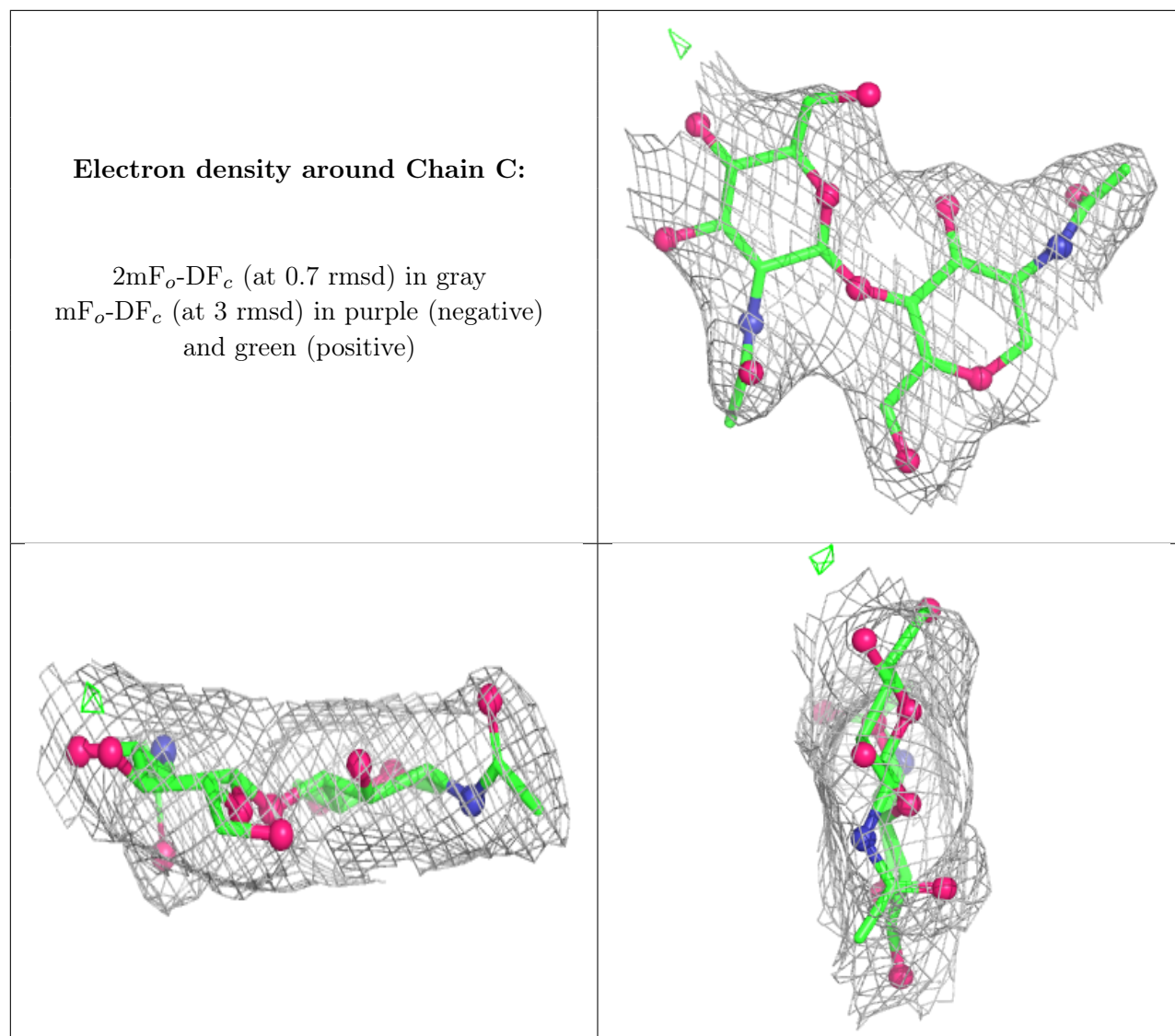
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	NAG	C	1	14/15	0.91	0.08	80,94,97,103	0
2	NAG	B	1	14/15	0.96	0.08	56,63,70,75	0

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.





6.4 Ligands [i](#)

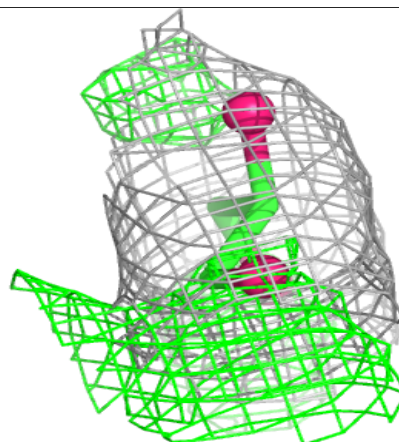
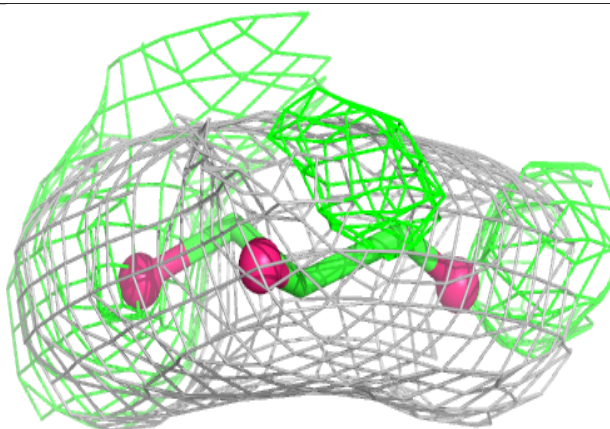
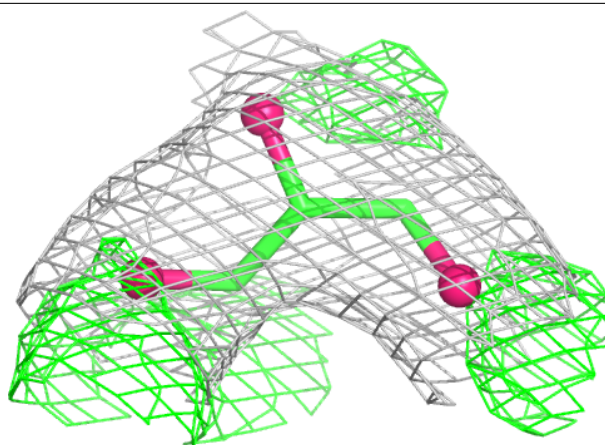
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	GOL	A	806	6/6	0.69	0.27	97,99,101,101	0
3	UOY	A	805	25/25	0.95	0.10	67,75,82,87	0

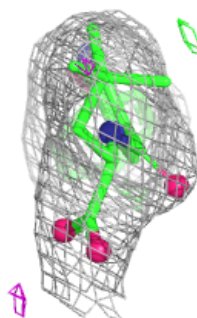
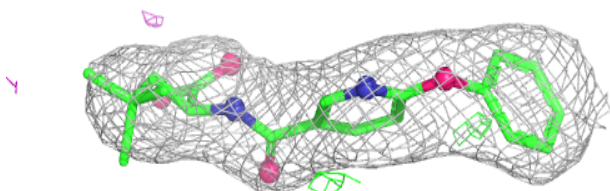
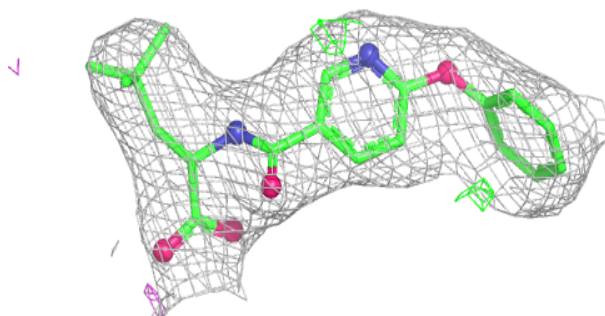
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around GOL A 806:

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

**Electron density around UOY A 805:**

$2mF_o-DF_c$ (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.