



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 28, 2024 – 08:31 AM EDT

PDB ID : 4ZGA
Title : Structural basis for inhibition of human autotaxin by four novel compounds
Authors : Stein, A.J.; Bain, G.; Hutchinson, J.H.; Evans, J.F.
Deposited on : 2015-04-22
Resolution : 2.60 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 : 4.02b-467
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 1.20.1
EDS : 3.0
buster-report : 1.1.7 (2018)
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4 : 9.0.003 (Gargrove)
Density-Fitness : 1.0.11
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.39

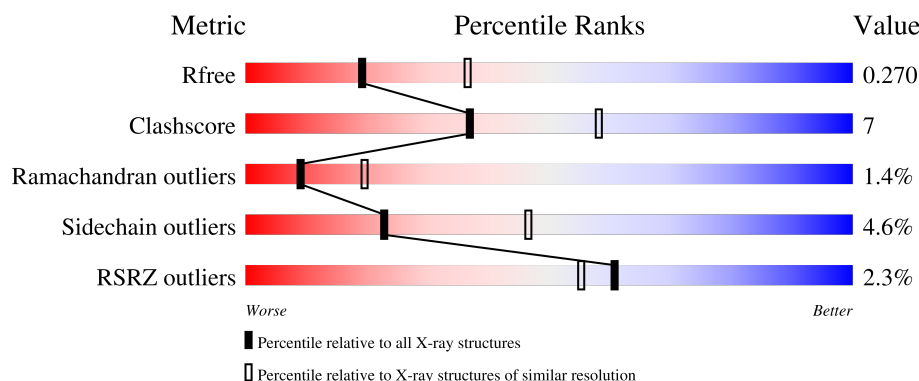
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION


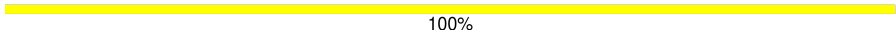
The reported resolution of this entry is 2.60 Å.

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}	164625	3775 (2.60-2.60)
Clashscore	180529	4181 (2.60-2.60)
Ramachandran outliers	177936	4129 (2.60-2.60)
Sidechain outliers	177891	4129 (2.60-2.60)
RSRZ outliers	164620	3775 (2.60-2.60)

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	863	
2	B	2	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
7	ACD	A	909	-	-	X	-

2 Entry composition [i](#)

There are 9 unique types of molecules in this entry. The entry contains 5870 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 Ectonucleotide pyrophosphatase/phosphodiesterase family member 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	740	Total	C	N	O	S	0	3	0
			5763	3668	979	1071	45			

- 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			

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Zn	0	0
			1	1		

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

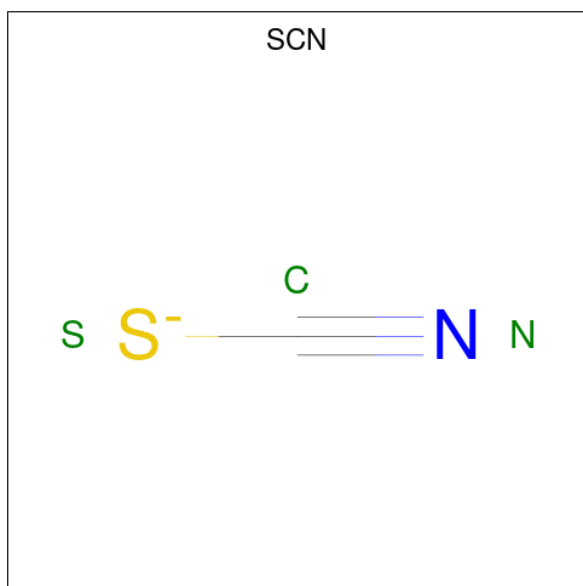
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	Ca	0	0
			2	2		

- Molecule 5 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



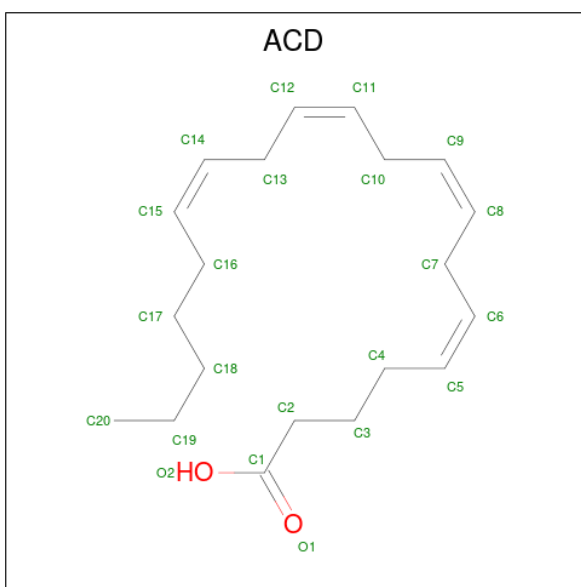
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			4	2	2		

- Molecule 6 is THIOCYANATE ION (three-letter code: SCN) (formula: CNS).



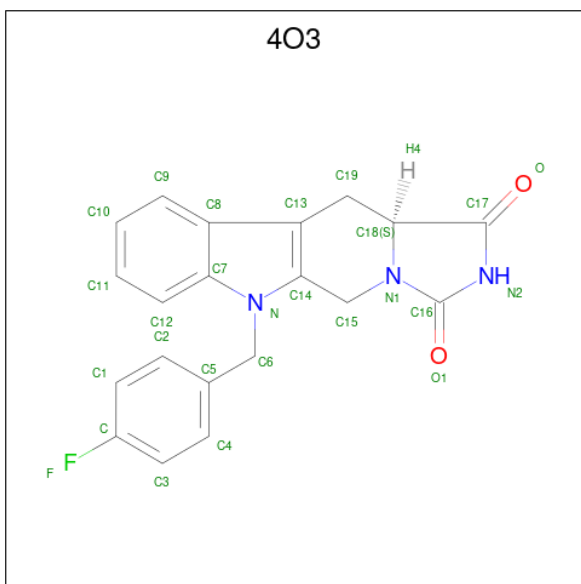
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	N	S	0	0
			3	1	1	1		
6	A	1	Total	C	N	S	0	0
			3	1	1	1		

- Molecule 7 is ARACHIDONIC ACID (three-letter code: ACD) (formula: C₂₀H₃₂O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	A	1	Total	C	O		0	0
			22	20	2			

- Molecule 8 is (11aS)-6-(4-fluorobenzyl)-5,6,11,11a-tetrahydro-1H-imidazo[1',5':1,6]pyrido[3,4-b]indole-1,3(2H)-dione (three-letter code: 4O3) (formula: C₂₀H₁₆FN₃O₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
8	A	1	Total	C	F	N	O	0	0
			26	20	1	3	2		

- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	18	Total	O	0	0
			18	18		

4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	53.67Å 64.00Å 70.50Å 99.44° 108.39° 98.61°	Depositor
Resolution (Å)	49.74 – 2.60 49.74 – 2.60	Depositor EDS
% Data completeness (in resolution range)	92.3 (49.74-2.60) 92.7 (49.74-2.60)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.81 (at 2.58Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.208 , 0.274 0.212 , 0.270	Depositor DCC
R_{free} test set	1218 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	65.0	Xtriage
Anisotropy	0.065	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 52.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5870	wwPDB-VP
Average B, all atoms (Å ²)	69.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.94% 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: CA, EDO, SCN, 4O3, ACD, NAG, ZN

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	0.58	0/5937	0.75	1/8089 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	199	SER	C-N-CD	5.45	139.85	128.40

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	5763	0	5294	63	0
2	B	28	0	25	0	0
3	A	1	0	0	0	0
4	A	2	0	0	0	0
5	A	4	0	6	2	0
6	A	6	0	0	0	0
7	A	22	0	31	13	0
8	A	26	0	16	0	0
9	A	18	0	0	2	0
All	All	5870	0	5372	78	0

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

All (78) 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)
7:A:909:ACD:H132	7:A:909:ACD:H9	1.44	1.00
7:A:909:ACD:C12	7:A:909:ACD:H161	2.11	0.81
7:A:909:ACD:H15	7:A:909:ACD:H192	1.64	0.78
7:A:909:ACD:H15	7:A:909:ACD:C19	2.16	0.74
7:A:909:ACD:H6	7:A:909:ACD:C2	2.21	0.71
1:A:191:LYS:O	1:A:195:CYS:HB2	1.92	0.69
1:A:450:ARG:HG3	1:A:451:ARG:HG3	1.76	0.68
7:A:909:ACD:H132	7:A:909:ACD:C9	2.22	0.67
1:A:848:GLU:O	1:A:851:THR:HB	1.95	0.66
1:A:166:LEU:HD13	1:A:516[B]:LEU:HD23	1.79	0.65
1:A:147:ASP:OD2	1:A:344:LYS:NZ	2.30	0.64
1:A:547:GLU:OE2	1:A:601:ARG:NH1	2.30	0.64
1:A:223:PRO:HA	1:A:226:HIS:CE1	2.35	0.61
1:A:549:THR:HG21	1:A:608:THR:H	1.65	0.61
7:A:909:ACD:H9	7:A:909:ACD:C13	2.27	0.61
1:A:770:SER:HB2	1:A:794:PHE:CZ	2.39	0.58
1:A:169:PHE:CZ	1:A:306:PHE:CE2	2.92	0.57
1:A:81:LYS:O	1:A:83:TYR:N	2.38	0.56
1:A:273:THR:HG22	1:A:275:PHE:H	1.70	0.56
1:A:145:VAL:HG22	1:A:188:ASN:OD1	2.06	0.56
1:A:103:CYS:O	1:A:137:VAL:HG21	2.06	0.56
1:A:191:LYS:O	1:A:195:CYS:CB	2.54	0.55
1:A:767:HIS:CD2	1:A:797:PRO:HA	2.42	0.55
7:A:909:ACD:C19	7:A:909:ACD:C15	2.85	0.55
7:A:909:ACD:H161	7:A:909:ACD:H12	1.87	0.54
1:A:695:ASP:HA	1:A:698:LEU:HD23	1.88	0.54
1:A:344:LYS:HA	1:A:349:HIS:ND1	2.22	0.53
1:A:622:TYR:HA	1:A:629:PRO:HA	1.90	0.53
7:A:909:ACD:C9	7:A:909:ACD:C13	2.85	0.53
1:A:501:VAL:HG21	1:A:534:LEU:HD11	1.90	0.52
1:A:695:ASP:HA	1:A:698:LEU:CD2	2.39	0.52
5:A:905:EDO:H21	9:A:1009:HOH:O	2.10	0.52
7:A:909:ACD:H6	7:A:909:ACD:H22	1.90	0.51
1:A:448:VAL:HG22	1:A:454:VAL:CG2	2.41	0.51
1:A:79:LEU:HD21	1:A:275:PHE:HB2	1.93	0.51
1:A:425:LEU:O	1:A:428:HIS:N	2.33	0.50
1:A:477:PHE:O	1:A:478:ASP:C	2.51	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:A:909:ACD:C2	7:A:909:ACD:C6	2.86	0.49
1:A:448:VAL:HG22	1:A:454:VAL:HG23	1.95	0.48
1:A:122:GLU:C	1:A:124:CYS:H	2.16	0.48
1:A:81:LYS:C	1:A:83:TYR:N	2.67	0.48
1:A:124:CYS:C	1:A:130:CYS:SG	2.92	0.48
1:A:186:MET:O	1:A:190:GLU:HB2	2.13	0.47
1:A:549:THR:CG2	1:A:607:ARG:HA	2.44	0.47
1:A:170:SER:HA	1:A:357:VAL:O	2.15	0.47
1:A:223:PRO:HA	1:A:226:HIS:NE2	2.30	0.47
1:A:511:ASN:HB3	1:A:528:HIS:CD2	2.50	0.47
1:A:233:MET:HG3	1:A:242:PHE:HB3	1.97	0.47
1:A:532:ASN:HB3	1:A:538:ASN:ND2	2.30	0.47
1:A:549:THR:HG23	1:A:606:TYR:O	2.14	0.47
1:A:82:SER:OG	1:A:277:SER:OG	2.27	0.47
1:A:632:THR:O	1:A:734:ILE:HA	2.15	0.47
1:A:418:ASP:O	1:A:420:HIS:CE1	2.68	0.46
1:A:81:LYS:C	1:A:83:TYR:H	2.19	0.46
1:A:432:ARG:HD3	1:A:824:THR:O	2.16	0.45
1:A:709:PHE:O	1:A:712:VAL:N	2.49	0.45
1:A:143:HIS:HB3	1:A:145:VAL:HG23	1.99	0.45
1:A:618:PHE:CD2	1:A:682:PHE:CD2	3.05	0.45
1:A:275:PHE:N	1:A:275:PHE:CD1	2.85	0.45
1:A:656:ASP:OD1	1:A:658:ARG:HD3	2.17	0.45
1:A:780:GLN:HA	1:A:781:PRO:HD2	1.86	0.45
1:A:623:SER:HB2	1:A:630:LEU:HD21	1.99	0.44
1:A:210:THR:HG23	1:A:214:LEU:HD22	1.99	0.44
1:A:699:VAL:HG23	1:A:702:MET:CE	2.48	0.44
1:A:226:HIS:CE1	1:A:228:ILE:HB	2.53	0.44
1:A:192:LEU:HD23	1:A:491:TYR:CG	2.53	0.44
1:A:741:TYR:N	1:A:748:ASP:OD2	2.47	0.44
1:A:424:TYR:CG	1:A:429:LEU:HD23	2.54	0.43
5:A:905:EDO:C1	9:A:1009:HOH:O	2.65	0.43
1:A:794:PHE:CD1	1:A:796:LEU:HG	2.54	0.43
1:A:113:ASN:HD22	1:A:114:GLU:N	2.16	0.42
1:A:122:GLU:C	1:A:124:CYS:N	2.73	0.42
7:A:909:ACD:H6	7:A:909:ACD:H21	1.99	0.42
1:A:846:TYR:CZ	1:A:850:LEU:HD21	2.55	0.42
1:A:617:ASP:HB3	1:A:682:PHE:CZ	2.55	0.42
1:A:87:CYS:O	1:A:88:HIS:C	2.58	0.42
1:A:123:ASP:O	1:A:127:ARG:CB	2.69	0.40
1:A:740:ASP:OD2	1:A:743:TYR:HA	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	727/863 (84%)	658 (90%)	59 (8%)	10 (1%)	9	19

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	114	GLU
1	A	85	SER
1	A	106	ASP
1	A	478	ASP
1	A	782	ALA
1	A	82	SER
1	A	108	CYS
1	A	134	TYR
1	A	729	ASN
1	A	123	ASP

5.3.2 Protein sidechains [i](#)

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	610/780 (78%)	582 (95%)	28 (5%)	23	46

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	66	LEU
1	A	97	THR
1	A	107	ARG
1	A	113	ASN
1	A	122	GLU
1	A	141	GLU
1	A	145	VAL
1	A	170	SER
1	A	171	VAL
1	A	214	LEU
1	A	233	MET
1	A	277	SER
1	A	394	ARG
1	A	416	LYS
1	A	453	HIS
1	A	487	VAL
1	A	493	SER
1	A	494	THR
1	A	520	LYS
1	A	557	MET
1	A	597	LEU
1	A	627	LEU
1	A	654	ARG
1	A	658	ARG
1	A	668	LEU
1	A	698	LEU
1	A	835	SER
1	A	840	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	113	ASN
1	A	116	ASN
1	A	538	ASN
1	A	717	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

2 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	2,1	14,14,15	1.00	0	17,19,21	1.69	4 (23%)
2	NAG	B	2	2	14,14,15	0.93	1 (7%)	17,19,21	1.87	5 (29%)

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	2,1	-	1/6/23/26	0/1/1/1
2	NAG	B	2	2	-	2/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	2	NAG	O5-C1	-2.32	1.39	1.43

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1	NAG	O5-C1-C2	-4.06	105.01	111.29
2	B	2	NAG	O5-C1-C2	-4.05	105.03	111.29
2	B	2	NAG	C4-C3-C2	3.89	116.71	111.02
2	B	2	NAG	O3-C3-C4	-2.96	103.40	110.38
2	B	1	NAG	C4-C3-C2	2.63	114.87	111.02

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	2	NAG	C1-O5-C5	-2.60	108.70	112.19
2	B	1	NAG	O6-C6-C5	-2.56	102.61	111.33
2	B	1	NAG	C6-C5-C4	-2.44	107.04	113.02
2	B	2	NAG	O5-C5-C4	-2.34	105.12	110.83

There are no chirality outliers.

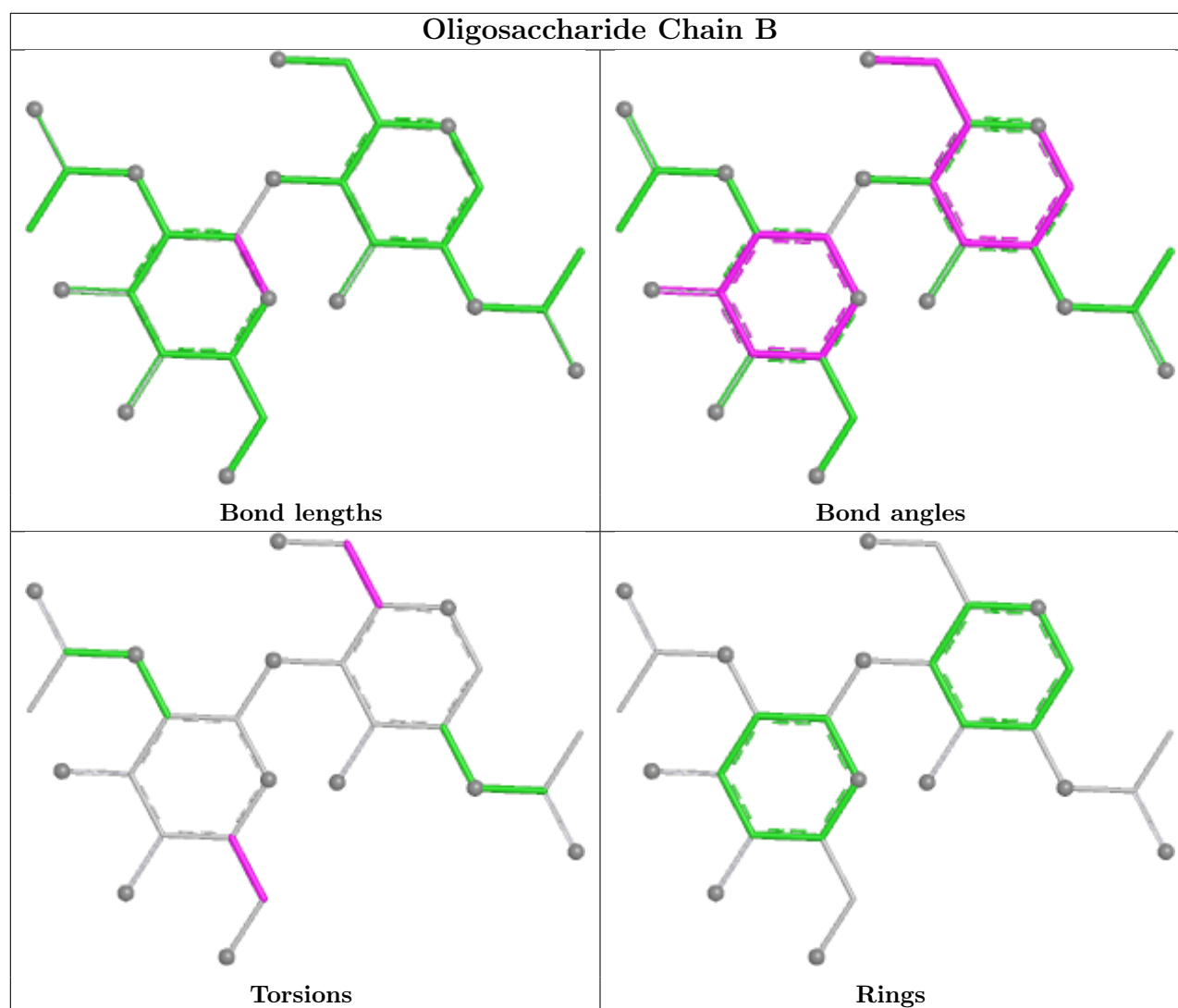
All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	1	NAG	O5-C5-C6-O6
2	B	2	NAG	O5-C5-C6-O6
2	B	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](#)

Of 8 ligands modelled in this entry, 3 are monoatomic - leaving 5 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	4O3	A	910	-	28,30,30	1.69	8 (28%)	28,45,45	4.10	12 (42%)
6	SCN	A	907	-	1,2,2	0.32	0	0,1,1	-	-
5	EDO	A	905	3	3,3,3	0.53	0	2,2,2	0.23	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
7	ACD	A	909	-	21,21,21	1.00	2 (9%)	21,21,21	1.20	1 (4%)
6	SCN	A	906	-	1,2,2	0.53	0	0,1,1	-	-

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	EDO	A	905	3	-	0/1/1/1	-
8	4O3	A	910	-	-	0/4/28/28	0/5/5/5
7	ACD	A	909	-	-	9/19/19/19	-

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	A	909	ACD	O2-C1	-3.55	1.19	1.30
8	A	910	4O3	C14-C13	3.28	1.46	1.38
8	A	910	4O3	C18-C17	-3.12	1.46	1.51
8	A	910	4O3	C13-C8	3.01	1.46	1.41
8	A	910	4O3	O1-C16	2.67	1.27	1.23
8	A	910	4O3	O-C17	2.66	1.28	1.23
8	A	910	4O3	C16-N1	-2.65	1.32	1.36
8	A	910	4O3	C17-N2	-2.65	1.34	1.37
7	A	909	ACD	O1-C1	2.57	1.30	1.22
8	A	910	4O3	C19-C13	-2.32	1.48	1.51

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	A	910	4O3	C18-C17-N2	15.08	120.45	106.65
8	A	910	4O3	C17-C18-N1	-7.43	92.67	102.18
8	A	910	4O3	O-C17-C18	-7.15	118.26	126.53
8	A	910	4O3	N2-C16-N1	6.14	113.20	107.30
8	A	910	4O3	C13-C19-C18	-6.02	105.64	111.77
7	A	909	ACD	C3-C2-C1	-4.32	103.23	114.51
8	A	910	4O3	O-C17-N2	-3.20	121.16	124.93
8	A	910	4O3	C14-N-C7	3.12	112.84	106.63
8	A	910	4O3	C12-C7-C8	-3.01	116.85	120.94
8	A	910	4O3	C14-C15-N1	-2.38	107.34	112.21
8	A	910	4O3	C4-C5-C2	2.21	121.52	118.23
8	A	910	4O3	C10-C9-C8	-2.17	117.98	120.91

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	A	910	4O3	O1-C16-N1	-2.06	122.59	125.45

There are no chirality outliers.

All (9) torsion outliers are listed below:

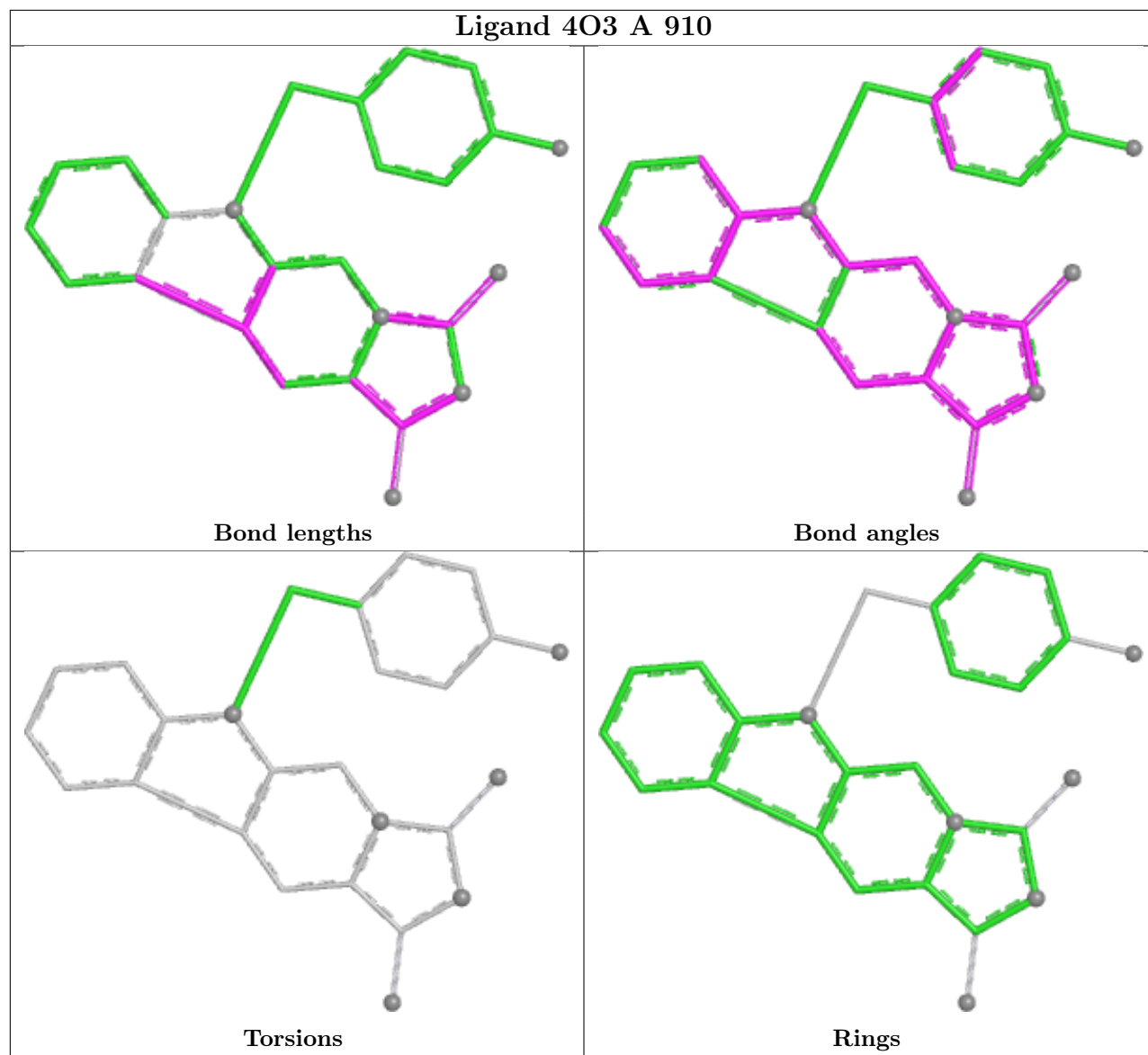
Mol	Chain	Res	Type	Atoms
7	A	909	ACD	C1-C2-C3-C4
7	A	909	ACD	C6-C7-C8-C9
7	A	909	ACD	C9-C10-C11-C12
7	A	909	ACD	C11-C12-C13-C14
7	A	909	ACD	C3-C4-C5-C6
7	A	909	ACD	O1-C1-C2-C3
7	A	909	ACD	C12-C13-C14-C15
7	A	909	ACD	O2-C1-C2-C3
7	A	909	ACD	C16-C17-C18-C19

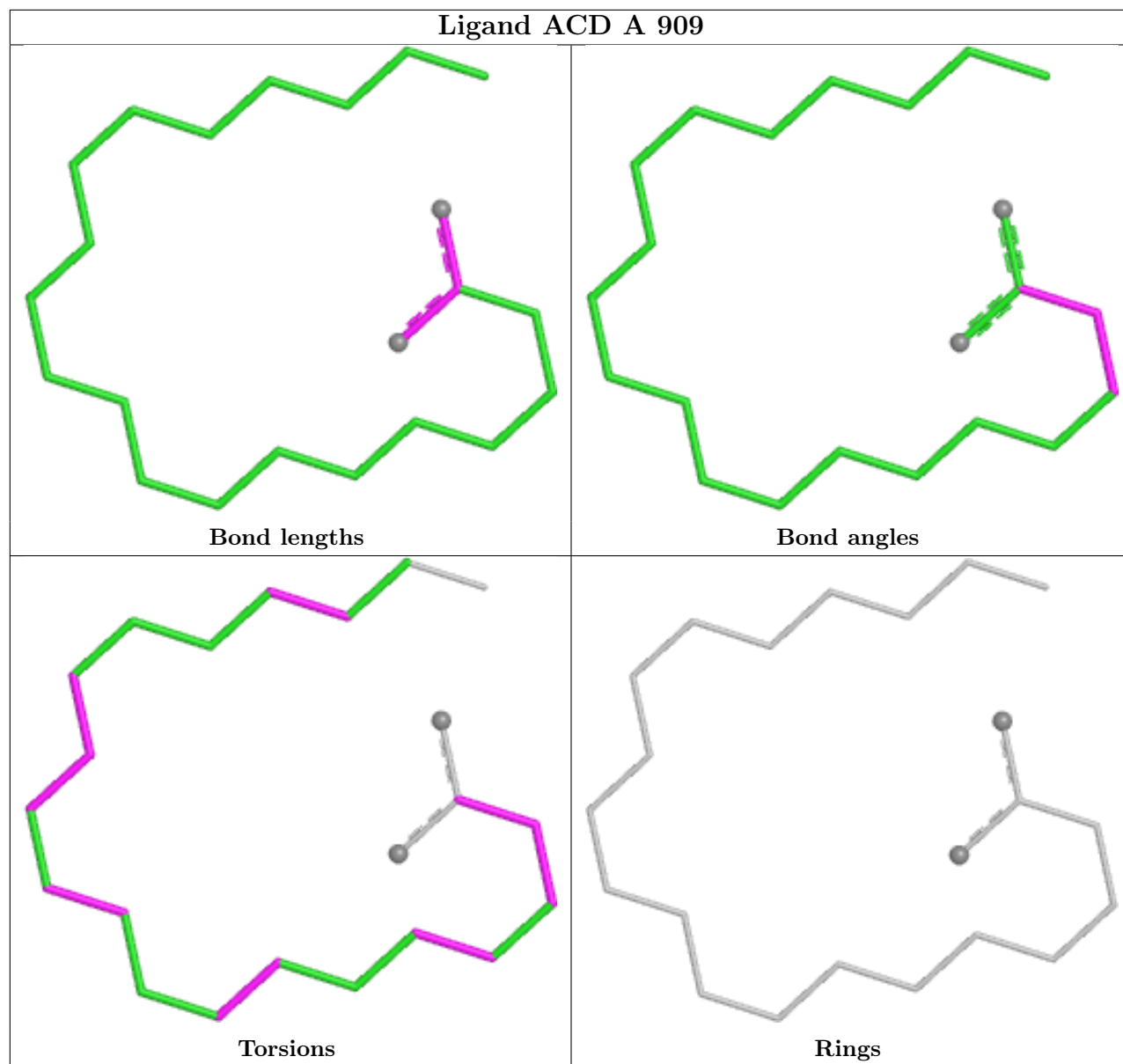
There are no ring outliers.

2 monomers are involved in 15 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	905	EDO	2	0
7	A	909	ACD	13	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	740/863 (85%)	-0.00	17 (2%) 61 55	35, 68, 99, 112	3 (0%)

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	453	HIS	4.0
1	A	66	LEU	3.6
1	A	218	ALA	3.5
1	A	397	PHE	3.2
1	A	342	GLY	2.9
1	A	135	GLN	2.8
1	A	557	MET	2.6
1	A	217	LEU	2.6
1	A	91	ASP	2.6
1	A	553	TYR	2.6
1	A	106	ASP	2.6
1	A	471	PHE	2.5
1	A	357	VAL	2.4
1	A	321	PHE	2.3
1	A	592	THR	2.2
1	A	126	ALA	2.2
1	A	739	PHE	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.

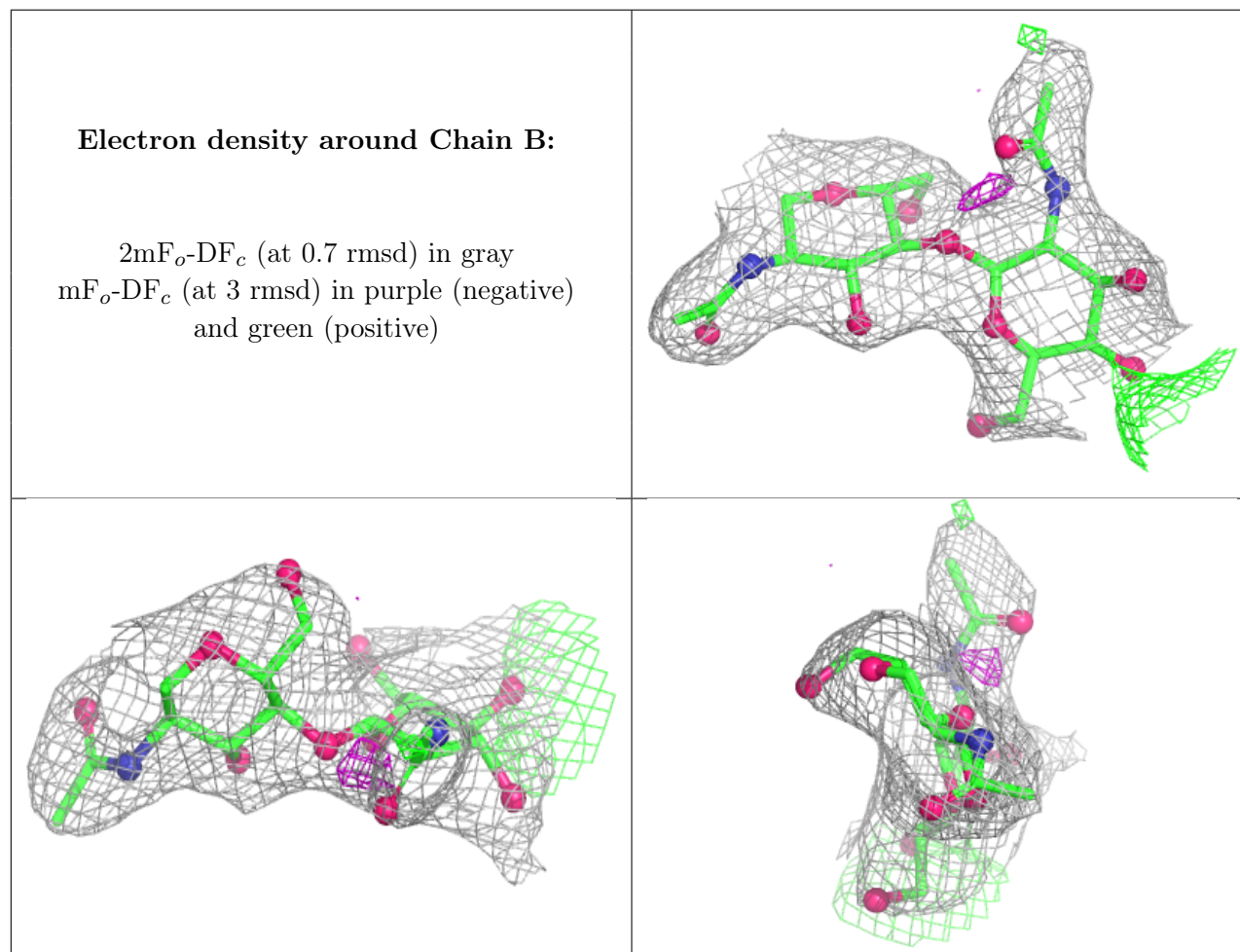
6.3 Carbohydrates [i](#)

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	0.67	0.15	69,79,91,97	0
2	NAG	B	1	14/15	0.94	0.09	46,55,65,67	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.

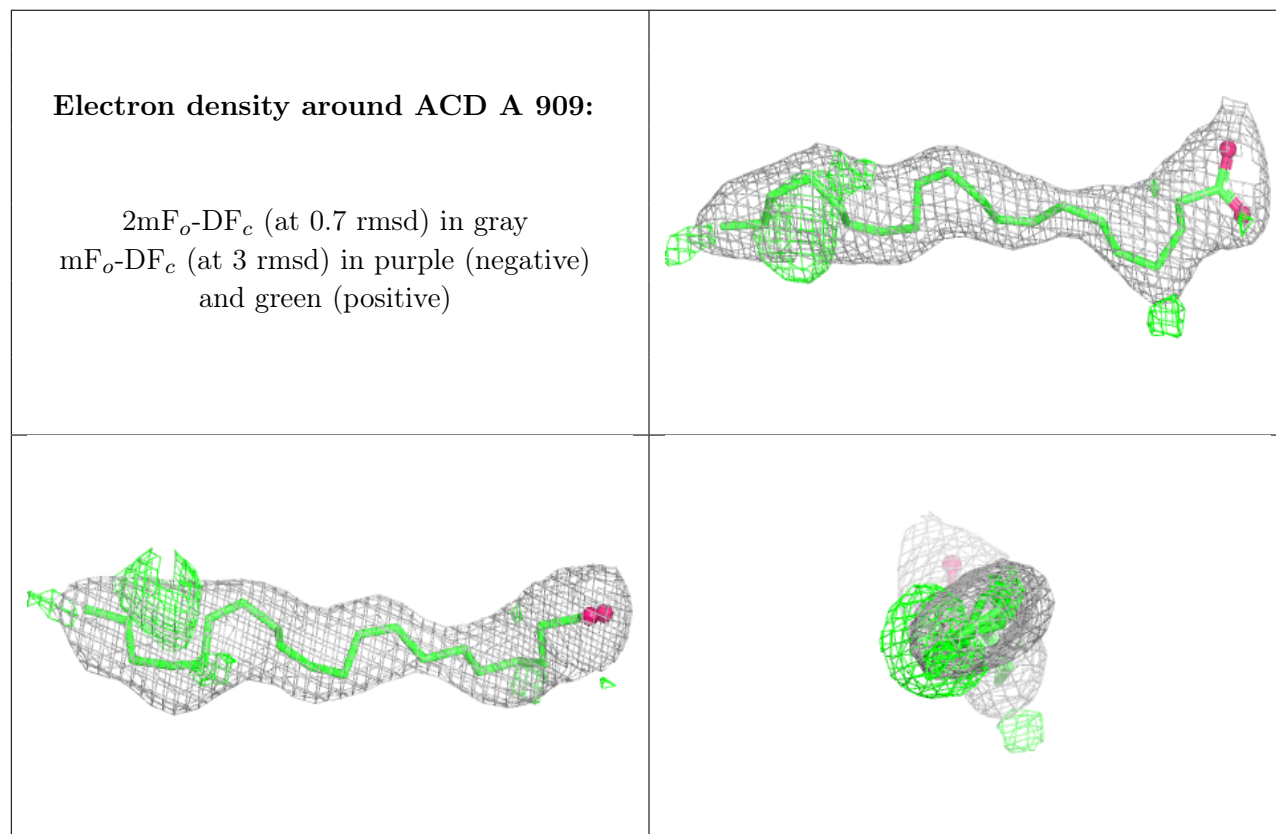
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	EDO	A	905	4/4	0.85	0.18	65,65,66,72	0

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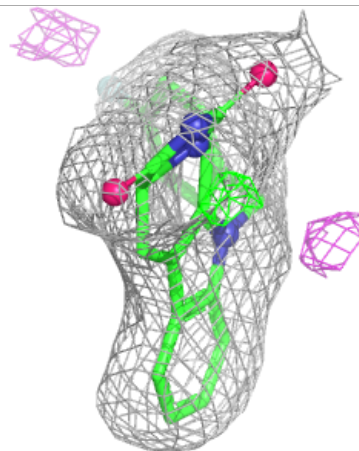
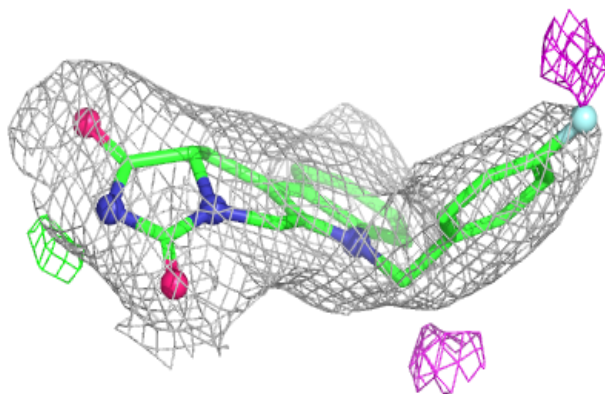
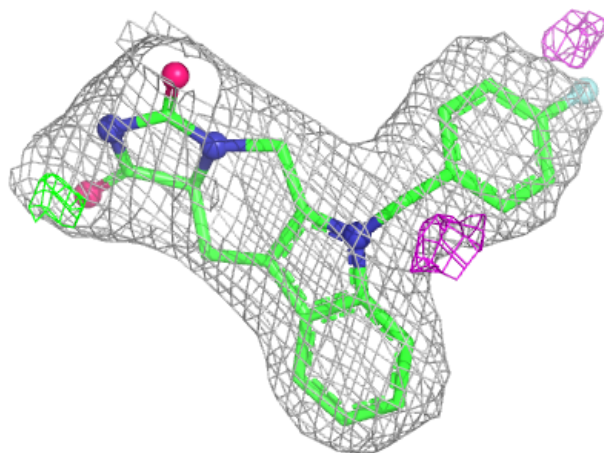
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
7	ACD	A	909	22/22	0.86	0.15	43,54,60,62	0
6	SCN	A	906	3/3	0.89	0.14	61,61,75,92	0
6	SCN	A	907	3/3	0.91	0.10	71,71,78,85	0
8	4O3	A	910	26/26	0.93	0.08	50,58,65,67	0
4	CA	A	908	1/1	0.98	0.05	65,65,65,65	0
3	ZN	A	903	1/1	0.99	0.06	65,65,65,65	0
4	CA	A	904	1/1	0.99	0.05	61,61,61,61	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 4O3 A 910:

$2mF_o-DF_c$ (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.