



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 3, 2021 – 04:32 AM EDT

PDB ID : 3KPF
Title : X-ray structure of the mutant Lys300Met of polyamine oxidase from Zea mays
Authors : Fiorillo, A.; Ilari, A.; Tavladoraki, P.
Deposited on : 2009-11-16
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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.23.2
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.23.2

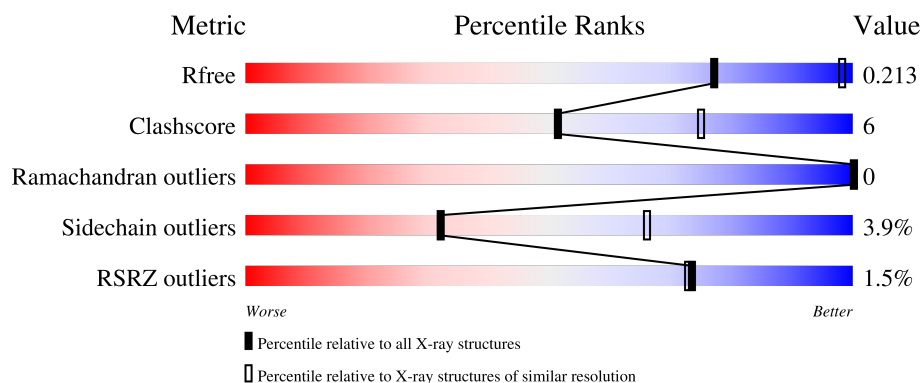
1 Overall quality at a glance

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}	130704	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (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	478	<div> <div>2%</div> <div>81%</div> <div>16%</div> <div>..</div> </div>
1	B	478	<div> <div>%</div> <div>85%</div> <div>12%</div> <div>.</div> </div>
2	C	2	<div> <div>100%</div> </div>

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 7861 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 Polyamine oxidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	467	Total	C	N	O	S	22	7	0
			3782	2417	633	714	18			
1	B	464	Total	C	N	O	S	24	3	0
			3746	2394	634	702	16			

There are 14 discrepancies between the modelled and reference sequences:

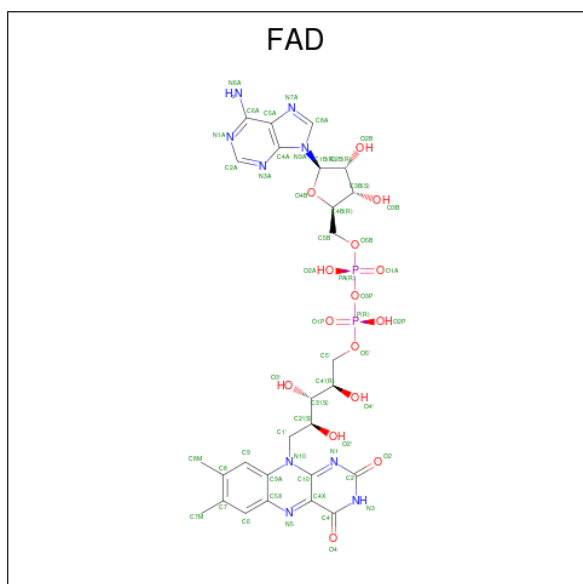
Chain	Residue	Modelled	Actual	Comment	Reference
A	300	MET	LYS	engineered mutation	UNP O64411
A	473	HIS	-	expression tag	UNP O64411
A	474	HIS	-	expression tag	UNP O64411
A	475	HIS	-	expression tag	UNP O64411
A	476	HIS	-	expression tag	UNP O64411
A	477	HIS	-	expression tag	UNP O64411
A	478	HIS	-	expression tag	UNP O64411
B	300	MET	LYS	engineered mutation	UNP O64411
B	473	HIS	-	expression tag	UNP O64411
B	474	HIS	-	expression tag	UNP O64411
B	475	HIS	-	expression tag	UNP O64411
B	476	HIS	-	expression tag	UNP O64411
B	477	HIS	-	expression tag	UNP O64411
B	478	HIS	-	expression tag	UNP O64411

- 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	C	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 3 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $\text{C}_{27}\text{H}_{33}\text{N}_9\text{O}_{15}\text{P}_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 53	C 27	N 9	O 15	P 2	0	0
3	B	1	Total 53	C 27	N 9	O 15	P 2	0	0

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $\text{C}_8\text{H}_{15}\text{NO}_6$).

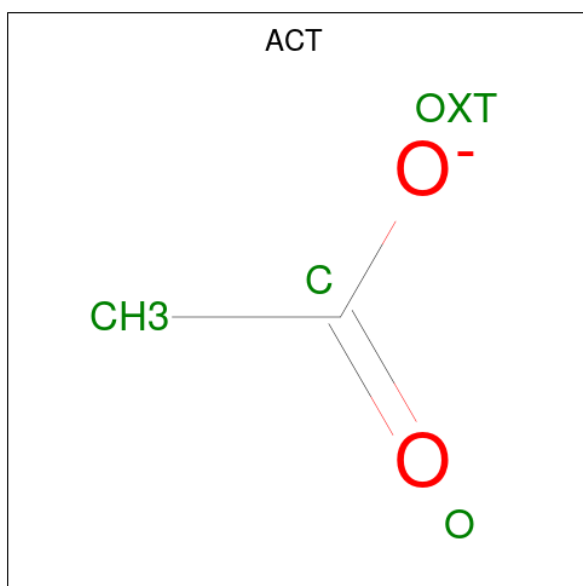


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 5 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

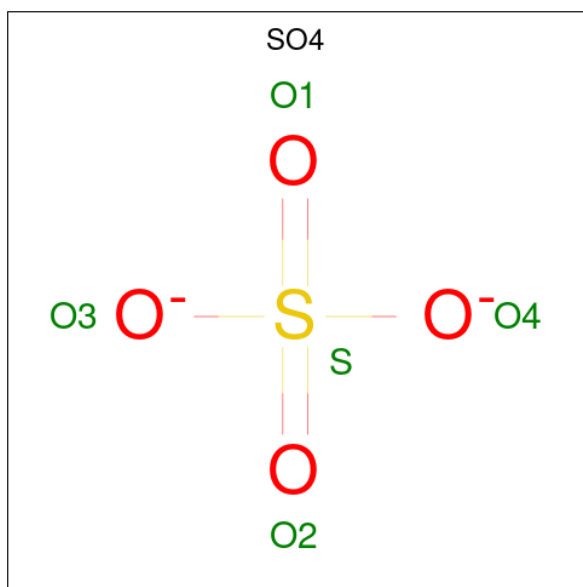
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Cl	0	0
			1	1		

- Molecule 6 is ACETATE ION (three-letter code: ACT) (formula: C₂H₃O₂).



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

- Molecule 7 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	B	1	Total O S 5 4 1	0	0

- Molecule 8 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	93	Total O 93 93	0	0
8	B	78	Total O 78 78	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	138.61Å 138.61Å 189.47Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 2.90 26.19 – 2.90	Depositor EDS
% Data completeness (in resolution range)	96.9 (50.00-2.90) 97.0 (26.19-2.90)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.54 (at 2.89Å)	Xtriage
Refinement program	REFMAC 5.5.0043	Depositor
R, R_{free}	0.186 , 0.223 0.178 , 0.213	Depositor DCC
R_{free} test set	2316 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	29.8	Xtriage
Anisotropy	0.517	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 38.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.024 for -h,-k,l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7861	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.55% 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: CL, SO4, NAG, FAD, ACT

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	0.51	2/3896 (0.1%)	0.58	0/5280
1	B	0.48	0/3848	0.57	0/5213
All	All	0.49	2/7744 (0.0%)	0.57	0/10493

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	464	LYS	CG-CD	-5.18	1.34	1.52
1	A	99	GLU	C-N	5.05	1.45	1.34

There are no bond angle outliers.

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	3782	0	3688	53	0
1	B	3746	0	3658	41	0
2	C	28	0	25	0	0
3	A	53	0	31	0	0
3	B	53	0	31	1	0
4	A	14	0	13	0	0
5	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	A	4	0	3	0	0
6	B	4	0	3	0	0
7	B	5	0	0	0	0
8	A	93	0	0	3	0
8	B	78	0	0	4	0
All	All	7861	0	7452	84	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 (84) 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:454:LEU:HA	1:A:457[B]:CYS:SG	1.84	1.17
1:B:298:TYR:HA	1:B:350:THR:HG22	1.39	1.00
1:A:412:ARG:HG2	1:A:412:ARG:HH11	1.30	0.97
1:B:412:ARG:HG3	1:B:412:ARG:HH11	1.30	0.96
1:A:454:LEU:CA	1:A:457[B]:CYS:SG	2.58	0.91
1:A:453:ILE:O	1:A:457[B]:CYS:SG	2.32	0.87
1:A:352:GLU:HG3	8:B:514:HOH:O	1.84	0.77
1:A:366:LYS:HD2	1:A:385:ALA:HB3	1.69	0.74
1:A:351:ASP:OD1	1:B:176:ARG:NH2	2.17	0.74
1:A:412:ARG:HH11	1:A:412:ARG:CG	2.03	0.71
1:A:454:LEU:O	1:A:457[B]:CYS:SG	2.52	0.68
1:B:255:SER:HB3	1:B:257:TYR:CE1	2.30	0.67
1:B:349:VAL:O	1:B:350:THR:HG23	1.97	0.65
1:A:177:VAL:HG11	1:B:176:ARG:HD2	1.78	0.65
1:B:349:VAL:O	1:B:350:THR:CG2	2.45	0.64
1:A:250:LYS:HG3	1:A:256:VAL:HG22	1.79	0.64
1:B:412:ARG:HG3	1:B:412:ARG:NH1	2.06	0.63
1:B:6[A]:ARG:NH1	1:B:260:ASP:OD2	2.32	0.62
1:A:454:LEU:C	1:A:457[B]:CYS:SG	2.79	0.62
1:A:353:GLU:OE1	8:A:546:HOH:O	2.16	0.61
1:A:98:LYS:HG3	1:A:102:GLY:O	2.02	0.60
1:A:156:THR:HG21	8:A:572:HOH:O	2.01	0.60
1:A:374:ARG:HG3	1:A:382:VAL:CG2	2.32	0.59
1:A:176:ARG:HD2	8:B:518:HOH:O	2.04	0.58
1:B:6[B]:ARG:HD3	1:B:28:ILE:HG23	1.86	0.58
1:B:410:VAL:HG13	1:B:414:GLU:HG3	1.85	0.58
1:A:394:TRP:HE1	1:B:152:ASN:ND2	2.02	0.57
1:A:412:ARG:HG2	1:A:412:ARG:NH1	2.11	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:349:VAL:C	1:B:350:THR:HG23	2.29	0.54
1:A:400:LYS:HB3	1:B:151:PRO:HG3	1.90	0.53
1:B:411:ASN:OD1	1:B:414:GLU:HG2	2.07	0.53
1:B:95:ASN:HB3	1:B:319:PHE:HB3	1.91	0.52
1:A:30:ASP:HB2	8:A:491:HOH:O	2.09	0.52
1:A:302:PHE:HB2	1:A:389:LEU:HB3	1.91	0.52
1:B:78:SER:O	1:B:81:LYS:HG3	2.11	0.51
1:B:218:LYS:HG3	1:B:227:VAL:HG23	1.93	0.51
1:B:335:PHE:HB3	1:B:338:GLN:HB2	1.92	0.51
1:A:152:ASN:ND2	1:B:394:TRP:HE1	2.09	0.50
1:A:458:ALA:O	1:A:461:LYS:HE3	2.13	0.49
1:A:136:ASP:OD1	1:A:136:ASP:N	2.46	0.48
1:B:431:HIS:CD2	8:B:511:HOH:O	2.67	0.48
1:A:352:GLU:OE2	1:B:326:ARG:HB3	2.14	0.48
1:A:138:MET:HB3	1:B:292:GLN:NE2	2.28	0.47
1:B:412:ARG:NH1	1:B:412:ARG:CG	2.74	0.47
1:B:370:MET:HE1	1:B:385:ALA:HB2	1.96	0.46
1:B:154:PRO:HB3	1:B:159:ASP:HB3	1.98	0.46
1:A:220:ASP:HB3	1:A:223:SER:HB3	1.98	0.46
1:B:131:HIS:CE1	1:B:135[A]:ARG:HH21	2.34	0.45
1:A:158:VAL:O	1:A:162:VAL:HG23	2.15	0.45
1:B:351:ASP:CG	1:B:352:GLU:N	2.70	0.45
1:B:351:ASP:CG	1:B:352:GLU:H	2.18	0.45
1:A:394:TRP:HE1	1:B:152:ASN:HD22	1.64	0.45
1:A:50:ALA:HB1	1:A:304:LYS:HE3	1.98	0.45
1:A:105:ASP:HB3	1:A:107[A]:ASP:OD2	2.16	0.45
1:B:241:LYS:NZ	8:B:505:HOH:O	2.49	0.44
1:A:156:THR:HG22	1:A:159:ASP:H	1.82	0.44
1:A:99:GLU:HG3	1:A:323:SER:HA	2.00	0.44
1:A:359:GLN:NE2	1:B:156:THR:HG22	2.32	0.44
1:A:20:ALA:HB3	1:A:216:TYR:OH	2.17	0.44
1:A:462:MET:HA	1:A:462:MET:HE3	2.00	0.44
1:A:171:PHE:CZ	1:A:183:THR:HB	2.53	0.43
1:A:353:GLU:O	1:A:357:ILE:HD12	2.17	0.43
1:A:110:GLN:HE21	1:A:114:GLU:HG3	1.83	0.43
1:B:131:HIS:CG	1:B:135[A]:ARG:HE	2.37	0.43
1:A:241:LYS:HA	1:A:279:LYS:O	2.19	0.42
1:B:300:MET:HB3	1:B:346:LEU:HD11	2.01	0.42
1:A:400:LYS:HA	1:A:400:LYS:HD2	1.92	0.42
1:A:307:ARG:HG2	1:A:383:PRO:HG3	2.03	0.41
1:A:358:GLU:HG2	1:A:392:ARG:HG2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:241:LYS:HB2	1:B:248:THR:HB	2.03	0.41
1:B:5:PRO:HD2	1:B:258:SER:O	2.20	0.41
1:A:98:LYS:O	1:A:99:GLU:C	2.59	0.41
1:A:219:THR:HG22	1:A:226:ILE:HA	2.03	0.41
1:A:332:TRP:HH2	1:A:369:ILE:HG23	1.84	0.41
1:A:406:TRP:HA	1:A:407:PRO:HD3	1.94	0.41
1:A:417:GLN:HG2	1:A:467:VAL:HG11	2.03	0.41
1:B:420:ALA:HA	1:B:421:PRO:HD3	1.94	0.41
1:B:58:ALA:HA	3:B:500:FAD:N5	2.36	0.41
1:A:118:SER:O	1:A:122[A]:MET:HG3	2.21	0.40
1:B:49:PHE:O	1:B:50:ALA:C	2.59	0.40
1:A:151:PRO:HG3	1:B:400:LYS:HB3	2.03	0.40
1:A:451:ALA:O	1:A:455:ILE:HG13	2.22	0.40
1:B:431:HIS:CD2	1:B:431:HIS:H	2.39	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	472/478 (99%)	451 (96%)	21 (4%)	0	100	100
1	B	465/478 (97%)	441 (95%)	24 (5%)	0	100	100
All	All	937/956 (98%)	892 (95%)	45 (5%)	0	100	100

There are no Ramachandran outliers to report.

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	408/410 (100%)	390 (96%)	18 (4%)	28	61
1	B	401/410 (98%)	386 (96%)	15 (4%)	34	68
All	All	809/820 (99%)	776 (96%)	33 (4%)	32	64

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	THR
1	A	6	ARG
1	A	59	ASN
1	A	107[A]	ASP
1	A	107[B]	ASP
1	A	115	LEU
1	A	156	THR
1	A	205	TYR
1	A	241	LYS
1	A	243	SER
1	A	248	THR
1	A	293	PHE
1	A	308	LYS
1	A	360	GLN
1	A	361	SER
1	A	384	ASP
1	A	412	ARG
1	A	418	LEU
1	B	6[A]	ARG
1	B	6[B]	ARG
1	B	59	ASN
1	B	103	VAL
1	B	105	ASP
1	B	115	LEU
1	B	243	SER
1	B	293	PHE
1	B	308	LYS
1	B	364	GLN
1	B	412	ARG
1	B	418	LEU
1	B	419	ARG
1	B	461	LYS
1	B	462	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	45	HIS
1	A	110	GLN
1	A	131	HIS
1	A	152	ASN
1	A	359	GLN
1	A	360	GLN
1	A	411	ASN
1	A	431	HIS
1	B	110	GLN
1	B	152	ASN
1	B	359	GLN
1	B	360	GLN
1	B	431	HIS

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	C	1	1,2	14,14,15	0.57	0	17,19,21	1.69	2 (11%)
2	NAG	C	2	2	14,14,15	0.69	0	17,19,21	1.39	2 (11%)

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	C	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	C	2	2	-	2/6/23/26	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	5.24	119.30	112.19
2	C	2	NAG	C4-C3-C2	4.51	117.63	111.02
2	C	1	NAG	C4-C3-C2	3.28	115.83	111.02
2	C	2	NAG	O5-C5-C6	2.00	110.34	107.20

There are no chirality outliers.

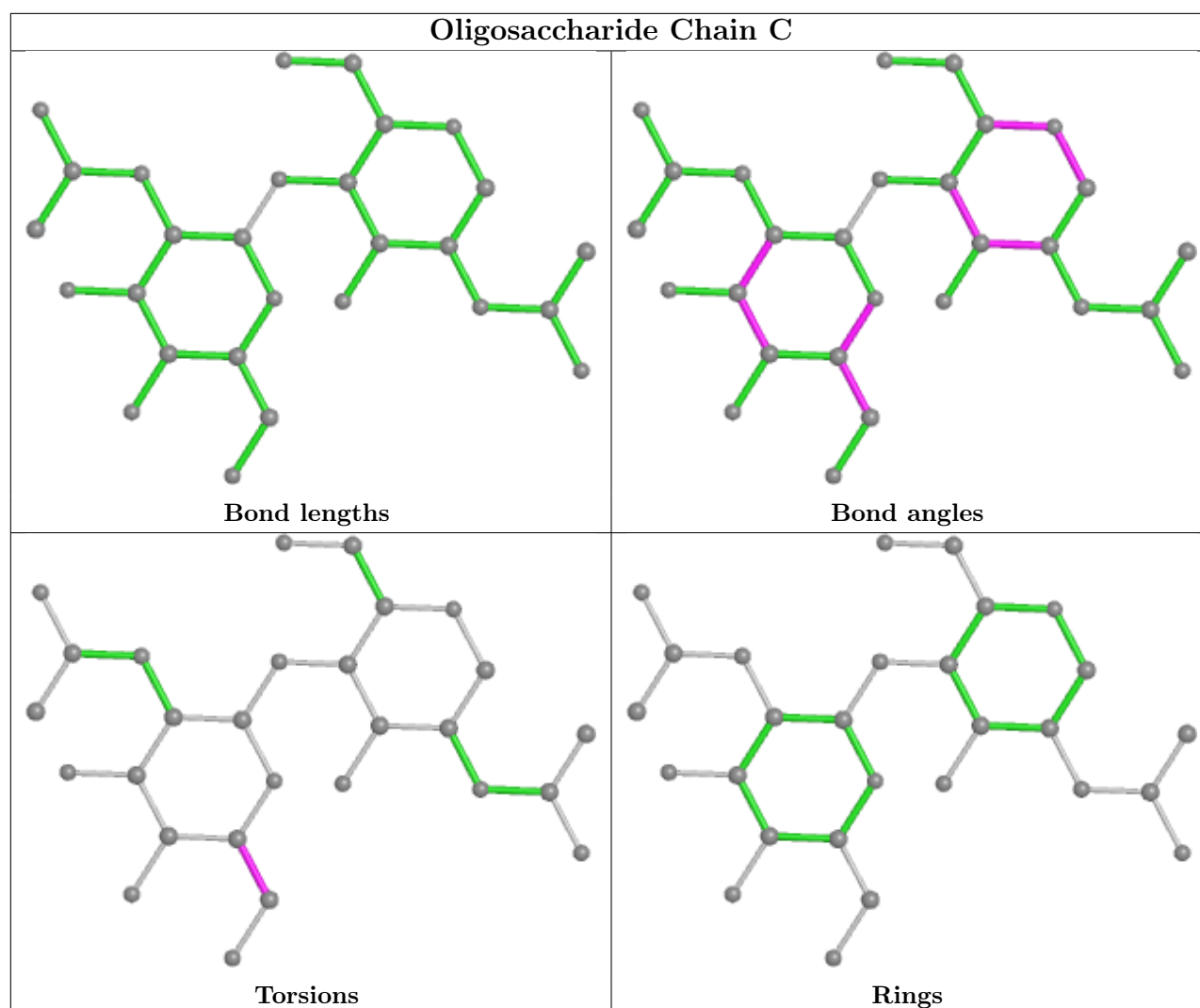
All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	2	NAG	O5-C5-C6-O6
2	C	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 7 ligands modelled in this entry, 1 is monoatomic - leaving 6 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	$\# Z > 2$
4	NAG	A	600	1	14,14,15	0.52	0	17,19,21	1.39	2 (11%)
3	FAD	A	500	-	51,58,58	1.45	7 (13%)	60,89,89	1.73	12 (20%)
7	SO4	B	479	-	4,4,4	0.16	0	6,6,6	0.17	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	ACT	B	480	-	1,3,3	1.89	0	0,3,3	-	-
3	FAD	B	500	-	51,58,58	1.42	6 (11%)	60,89,89	1.78	8 (13%)
6	ACT	A	480	-	1,3,3	1.75	0	0,3,3	-	-

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	NAG	A	600	1	-	0/6/23/26	0/1/1/1
3	FAD	B	500	-	-	8/30/50/50	0/6/6/6
3	FAD	A	500	-	-	5/30/50/50	0/6/6/6

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	500	FAD	C10-N1	4.49	1.39	1.33
3	A	500	FAD	C10-N1	4.27	1.38	1.33
3	A	500	FAD	C4X-N5	4.23	1.39	1.33
3	B	500	FAD	C4X-N5	4.19	1.39	1.33
3	A	500	FAD	C2A-N3A	4.19	1.38	1.32
3	B	500	FAD	C2A-N3A	3.77	1.38	1.32
3	A	500	FAD	C4-N3	3.61	1.39	1.33
3	B	500	FAD	C4-N3	3.61	1.39	1.33
3	A	500	FAD	C2A-N1A	2.86	1.39	1.33
3	B	500	FAD	C5X-N5	2.61	1.39	1.35
3	B	500	FAD	C2A-N1A	2.52	1.38	1.33
3	A	500	FAD	C5X-N5	2.40	1.39	1.35
3	A	500	FAD	C1'-N10	2.14	1.50	1.48

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	500	FAD	C4-N3-C2	7.03	121.08	115.14
3	B	500	FAD	N3A-C2A-N1A	-6.22	118.96	128.68
3	A	500	FAD	C4-N3-C2	6.07	120.27	115.14
3	A	500	FAD	N3A-C2A-N1A	-5.18	120.58	128.68
3	B	500	FAD	C10-C4X-N5	-4.60	118.08	121.26
3	A	500	FAD	C4X-N5-C5X	4.41	121.17	116.77
3	A	500	FAD	C10-C4X-N5	-4.10	118.42	121.26

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	500	FAD	C4X-N5-C5X	4.04	120.81	116.77
4	A	600	NAG	C3-C4-C5	-3.32	104.32	110.24
3	B	500	FAD	C4-C4X-C10	-3.23	117.81	119.95
3	A	500	FAD	C1'-N10-C9A	3.22	120.83	118.29
4	A	600	NAG	C1-O5-C5	2.97	116.22	112.19
3	B	500	FAD	C4X-C4-N3	-2.79	119.61	123.43
3	A	500	FAD	C4X-C4-N3	-2.77	119.64	123.43
3	B	500	FAD	P-O3P-PA	-2.77	123.32	132.83
3	A	500	FAD	C1B-N9A-C4A	-2.60	122.08	126.64
3	A	500	FAD	C4A-C5A-N7A	-2.51	106.79	109.40
3	A	500	FAD	C5'-C4'-C3'	-2.50	107.37	112.20
3	A	500	FAD	C4-C4X-N5	2.27	121.19	118.60
3	A	500	FAD	P-O3P-PA	-2.17	125.38	132.83
3	A	500	FAD	C5X-C9A-N10	2.12	119.25	117.72
3	B	500	FAD	C4-C4X-N5	2.02	120.91	118.60

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	500	FAD	C5B-O5B-PA-O1A
3	B	500	FAD	O3'-C3'-C4'-C5'
3	B	500	FAD	C2'-C3'-C4'-C5'
3	B	500	FAD	C2'-C3'-C4'-O4'
3	A	500	FAD	O3'-C3'-C4'-C5'
3	B	500	FAD	O3'-C3'-C4'-O4'
3	B	500	FAD	C5B-O5B-PA-O3P
3	B	500	FAD	C5B-O5B-PA-O2A
3	A	500	FAD	C2'-C3'-C4'-C5'
3	A	500	FAD	C2'-C3'-C4'-O4'
3	A	500	FAD	O4B-C4B-C5B-O5B
3	B	500	FAD	O4B-C4B-C5B-O5B
3	A	500	FAD	O3'-C3'-C4'-O4'

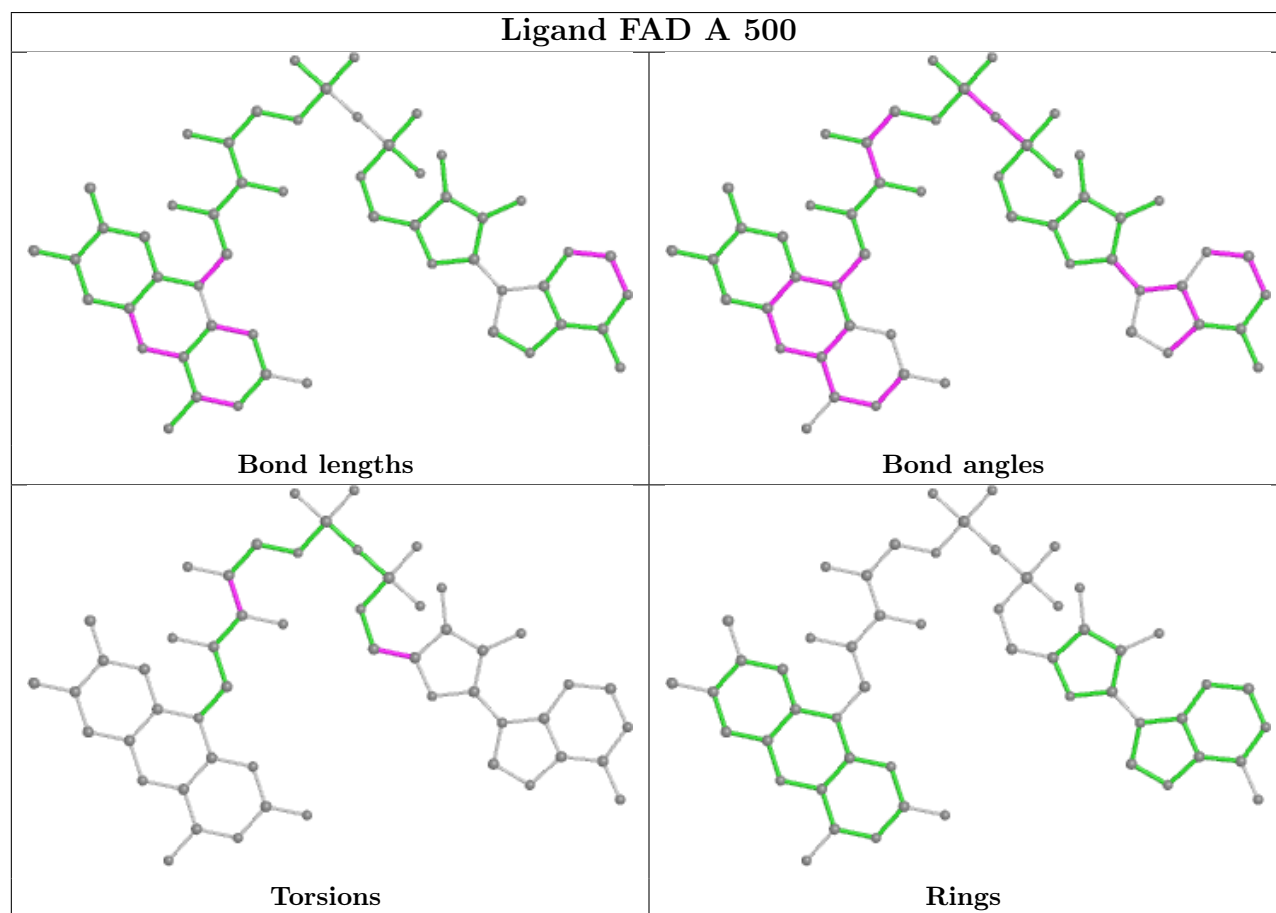
There are no ring outliers.

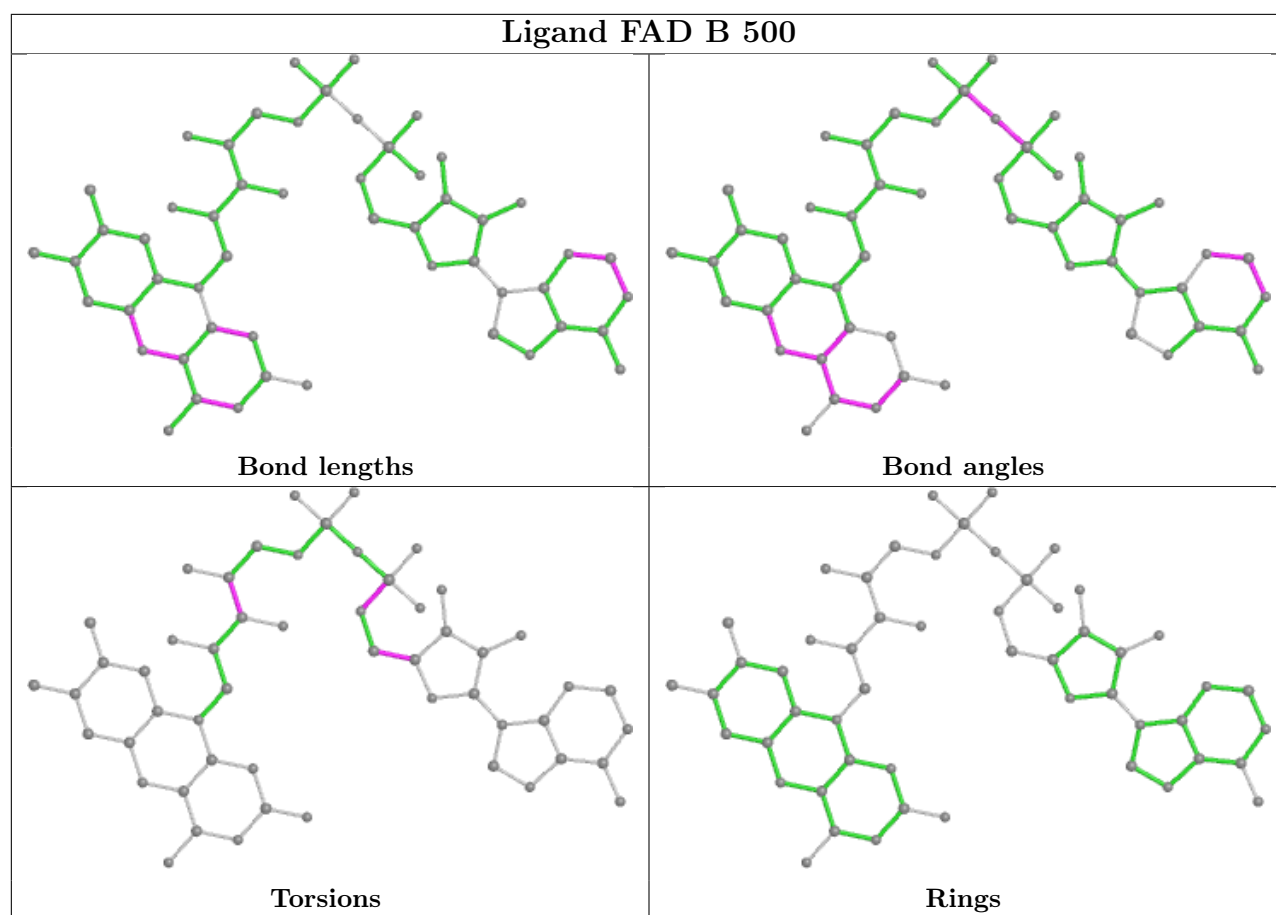
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	500	FAD	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths,

bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [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	467/478 (97%)	-0.45	9 (1%) 66 65	14, 23, 35, 55	19 (4%)
1	B	464/478 (97%)	-0.49	5 (1%) 80 80	13, 23, 37, 50	16 (3%)
All	All	931/956 (97%)	-0.47	14 (1%) 73 73	13, 23, 36, 55	35 (3%)

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	221	ASP	4.2
1	B	466	HIS	3.2
1	A	468	GLN	3.1
1	A	466	HIS	2.6
1	A	220	ASP	2.6
1	B	467	VAL	2.5
1	B	223	SER	2.3
1	A	381	ASP	2.2
1	A	107[A]	ASP	2.2
1	B	222	LYS	2.1
1	A	78	SER	2.1
1	A	223	SER	2.1
1	B	105	ASP	2.1
1	A	467	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.

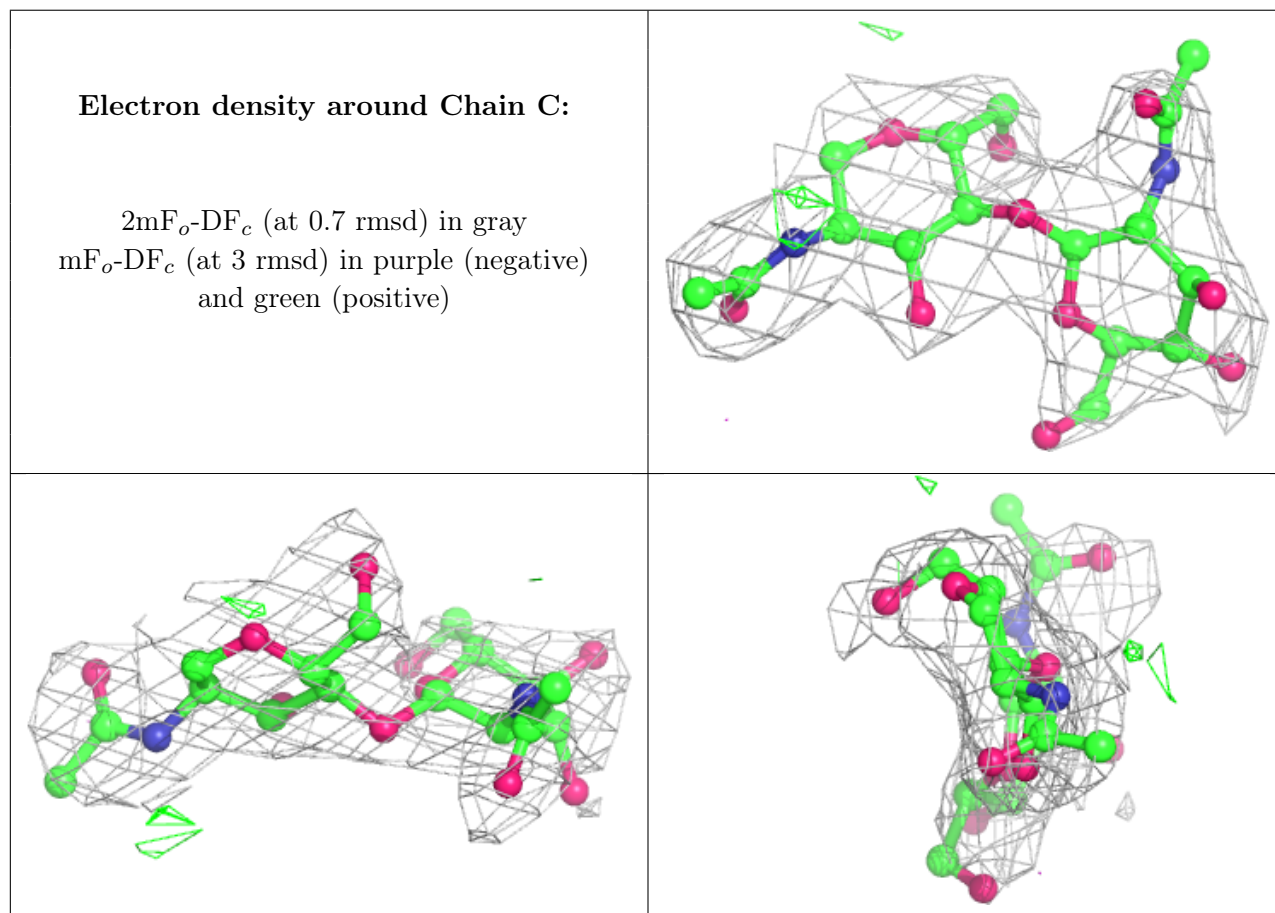
6.3 Carbohydrates [i](#)

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(Å ²)	Q<0.9
2	NAG	C	2	14/15	0.86	0.54	40,67,81,82	0
2	NAG	C	1	14/15	0.94	0.26	30,43,53,54	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 ⓘ

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	CL	A	479	1/1	0.89	0.16	52,52,52,52	0
6	ACT	B	480	4/4	0.91	0.26	30,48,53,58	0
6	ACT	A	480	4/4	0.95	0.23	24,42,43,46	0

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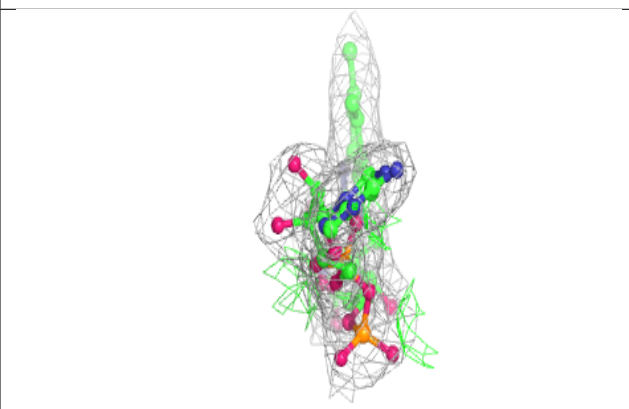
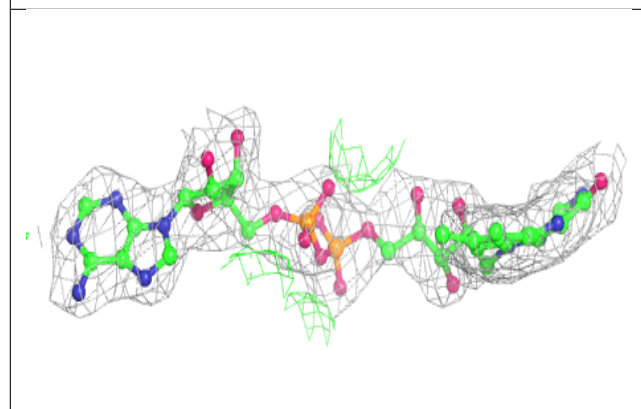
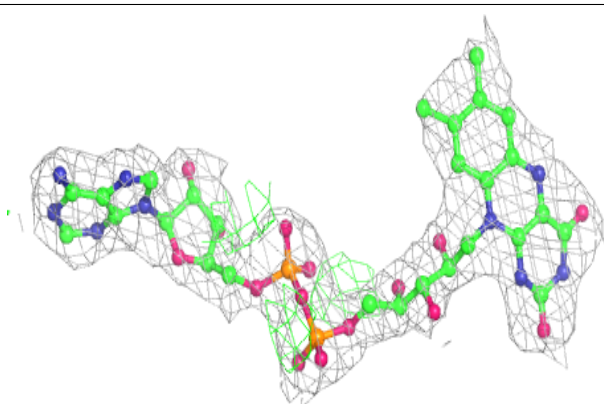
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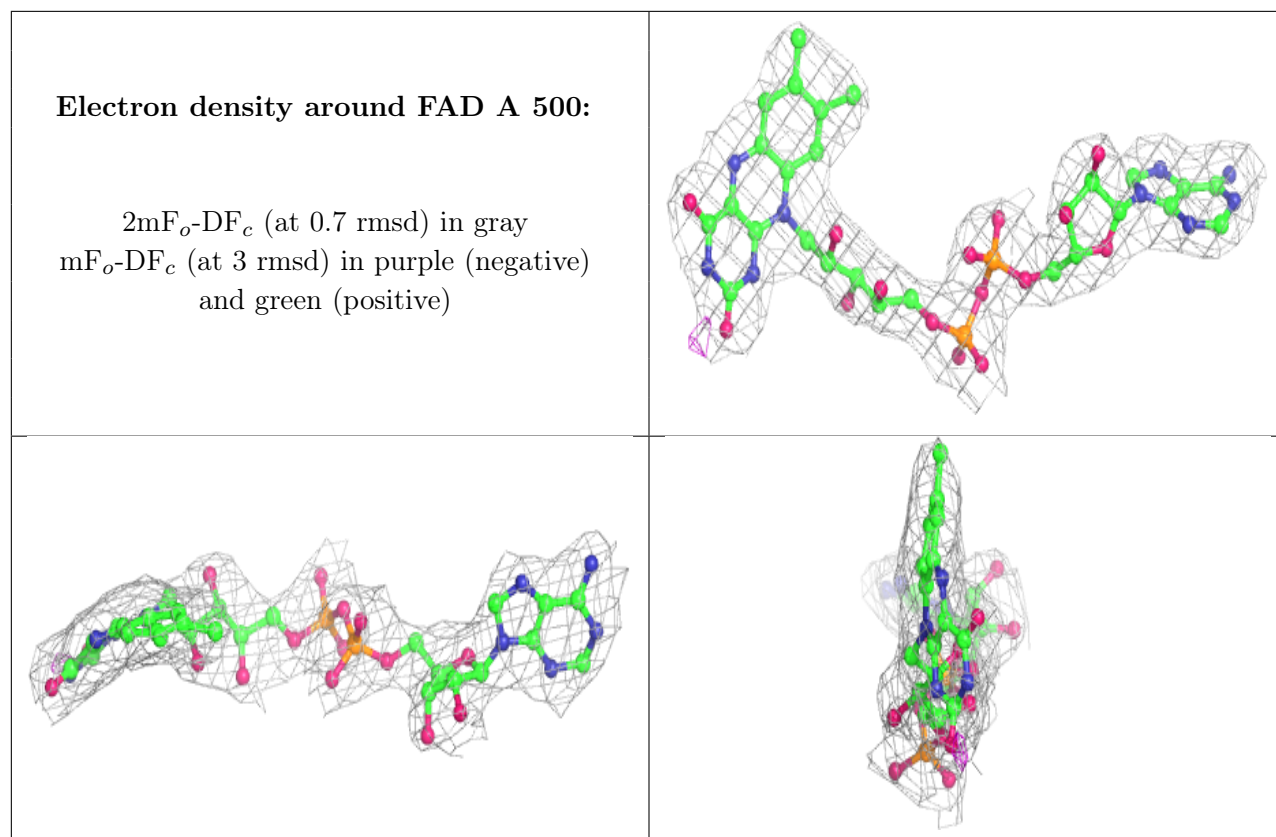
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	NAG	A	600	14/15	0.95	0.31	30,38,45,48	0
7	SO4	B	479	5/5	0.97	0.10	32,36,46,48	5
3	FAD	B	500	53/53	0.99	0.11	9,17,24,27	0
3	FAD	A	500	53/53	0.99	0.11	8,15,23,25	0

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 FAD B 500:

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.