WORLDWIDE wwPDB PROTEIN DATA BANK

Full wwPDB X-ray Structure Validation Report (i)

Dec 15, 2022 – 04:04 pm GMT

PDB ID : 7OWM

Title : HsNMT1 in complex with both MyrCoA and HCPA substrate peptide GKQN-SKLR

Authors : Dian, C.; Giglione, C.; Meinnel, T.

Deposited on : 2021-06-18

Resolution : 1.50 Å(reported)

The following versions of software and data (see references (i)) were used in the production of this report:

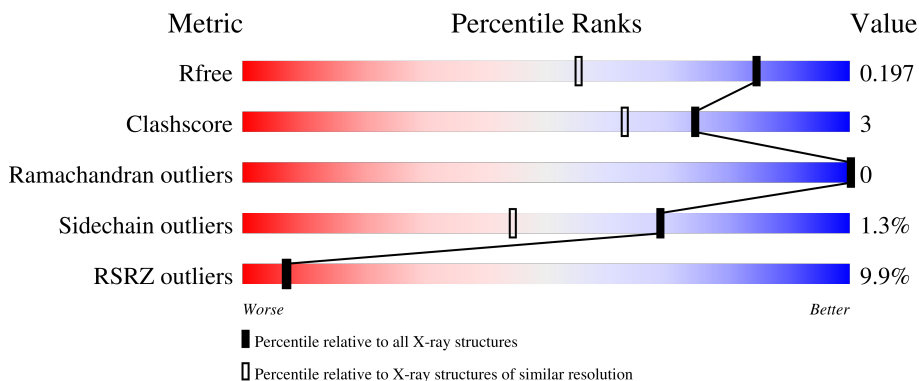|		
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.31.3
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.31.3

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.50 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	Percentile Ranks	Value
Rfree		0.197
Clashscore		3
Ramachandran outliers		0
Sidechain outliers		1.3%
RSRZ outliers		9.9%

Worse   *Better*

▐ Percentile relative to all X-ray structures

▯ Percentile relative to X-ray structures of similar resolution

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	2936 (1.50-1.50)
Clashscore	141614	3144 (1.50-1.50)
Ramachandran outliers	138981	3066 (1.50-1.50)
Sidechain outliers	138945	3064 (1.50-1.50)
RSRZ outliers	127900	2884 (1.50-1.50)

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 <=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	402	8% ... 92% ... 6% ·
1	B	402	10% ... 88% ... 7% ·
2	C	8	75% ... 62% ... 38%

2 Entry composition ⓘ

There are 5 unique types of molecules in this entry. The entry contains 7658 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 Glycylpeptide N-tetradecanoyltransferase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	392	Total	C	N	O	S	0	3	0
			3176	2056	531	572	17			
1	B	387	Total	C	N	O	S	0	11	0
			3185	2060	535	573	17			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	95	GLY	-	expression tag	UNP P30419
A	96	GLY	-	expression tag	UNP P30419
A	97	SER	-	expression tag	UNP P30419
A	98	GLU	-	expression tag	UNP P30419
B	95	GLY	-	expression tag	UNP P30419
B	96	GLY	-	expression tag	UNP P30419
B	97	SER	-	expression tag	UNP P30419
B	98	GLU	-	expression tag	UNP P30419

- Molecule 2 is a protein called Neuron-specific calcium-binding protein hippocalcin.
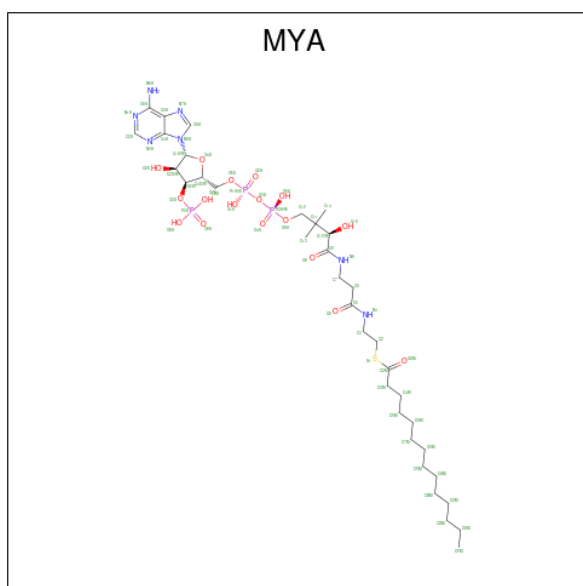
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	8	Total	C	N	O	0	2	0
			72	43	15	14			

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).

GOL

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total 6	C 3	O 3	0	0
3	B	1	Total 6	C 3	O 3	0	0
3	B	1	Total 6	C 3	O 3	0	0
3	B	1	Total 12	C 6	O 6	0	1

- Molecule 4 is TETRADECANOYL-COA (three-letter code: MYA) (formula: $C_{35}H_{62}N_7O_{17}P_3S$) (labeled as "Ligand of Interest" by depositor).

MYA

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
4	B	1	Total 63	C 35	N 7	O 17	P 3	S 1	0	0
4	C	1	Total 126	C 70	N 14	O 34	P 6	S 2	0	1

- Molecule 5 is water.

| Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	495	Total 495	O 495	0	0
5	B	495	Total 495	O 495	0	1
5	C	16	Total 16	O 16	0	1

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.
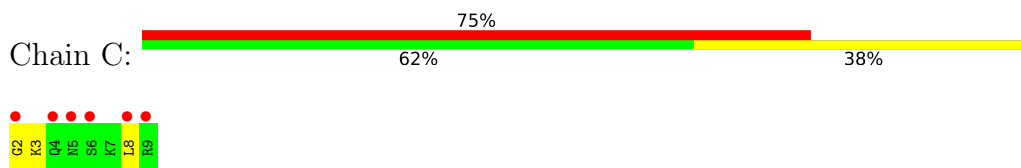
• Molecule 1: Glycylpeptide N-tetradecanoyltransferase 1

Chain A:

• Molecule 1: Glycylpeptide N-tetradecanoyltransferase 1

Chain B:
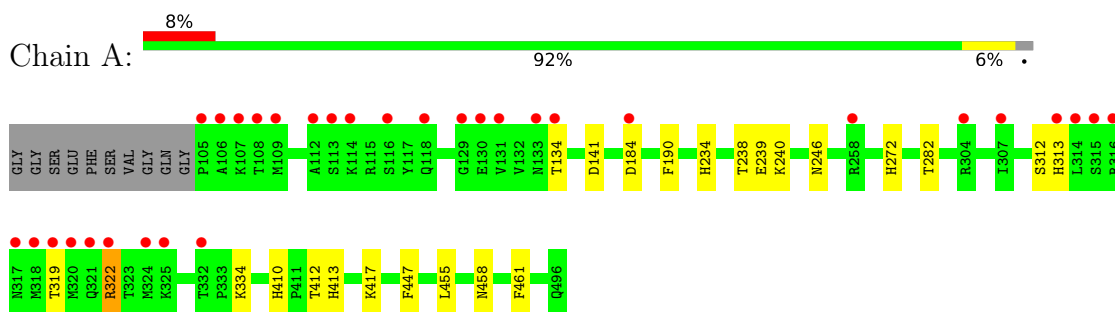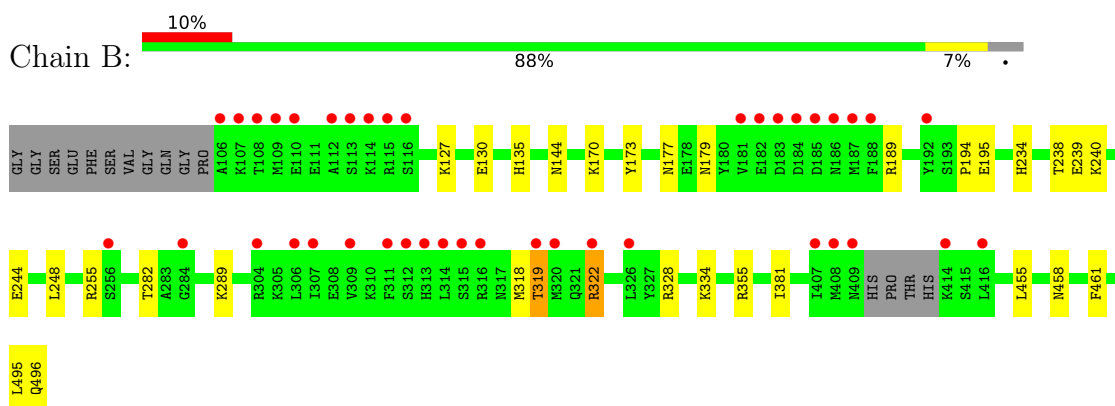
• Molecule 2: Neuron-specific calcium-binding protein hippocalcin

Chain C:

4 Data and refinement statistics ⓘ

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α, β, γ	79.88Å 178.38Å 58.26Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.77 – 1.50 48.77 – 1.50	Depositor EDS
% Data completeness (in resolution range)	96.0 (48.77-1.50) 96.0 (48.77-1.50)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.56 (at 1.50Å)	Xtriage
Refinement program	PHENIX 1.17_3644	Depositor
R, R_{free}	0.154 , 0.196 0.155 , 0.197	Depositor DCC
R_{free} test set	6514 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å^2)	16.7	Xtriage
Anisotropy	0.018	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 46.3	EDS
L-test for twinning[2]	$< |L| > = 0.50, < L^2 > = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.97	EDS
Total number of atoms	7658	wwPDB-VP
Average B, all atoms (Å^2)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 16.41% of the height of the origin peak. No significant pseudotranslation is detected.*

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[1]Intensities estimated from amplitudes.
[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: GOL, MYA

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	0.34	0/3265	0.51	0/4442
1	B	0.34	0/3270	0.52	0/4448
2	C	0.32	0/71	0.65	0/89
All	All	0.34	0/6606	0.52	0/8979

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 ⓘ

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	3176	0	3118	16	0
1	B	3185	0	3102	23	0
2	C	72	0	74	7	0
3	A	6	0	8	0	0
3	B	24	0	32	2	0
4	B	63	0	58	1	0
4	C	126	0	116	2	0
5	A	495	0	0	3	0
5	B	495	0	0	5	1
5	C	16	0	0	1	0
All	All	7658	0	6508	42	1

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 3.

All (42) 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:282:THR:HG1	2:C:2[A]:GLY:N	1.86	0.74
1:A:417:LYS:HG2	1:A:447:PHE:HA	1.72	0.69
1:B:173[B]:TYR:CE2	1:B:194:PRO:HG3	2.29	0.67
2:C:2[A]:GLY:HA3	4:C:501[A]:MYA:C2M	2.26	0.66
1:B:170:LYS:HA	1:B:173[B]:TYR:CE2	2.33	0.63
1:A:234:HIS:HD2	1:A:239:GLU:HG3	1.64	0.62
2:C:2[A]:GLY:N	5:C:601:HOH:O	2.32	0.62
1:A:312:SER:HB3	2:C:8:LEU:HD12	1.84	0.59
1:A:410:HIS:HD2	1:A:412:THR:H	1.49	0.58
2:C:2[A]:GLY:HA3	4:C:501[A]:MYA:S1	2.45	0.57
1:B:282:THR:HG23	1:B:495:LEU:HD12	1.87	0.56
1:B:319:THR:HB	1:B:322:ARG:H	1.72	0.54
1:B:328[B]:ARG:NH1	5:B:613:HOH:O	2.36	0.54
1:B:238:THR:HG22	1:B:240:LYS:HG2	1.89	0.54
1:A:410:HIS:CD2	1:A:413:HIS:H	2.27	0.53
1:B:244:GLU:OE2	3:B:504[A]:GOL:O3	2.25	0.51
1:B:458:ASN:HA	1:B:461:PHE:CE2	2.46	0.50
1:A:234:HIS:CD2	1:A:239:GLU:HG3	2.46	0.50
1:B:144[B]:ASN:OD1	5:B:601:HOH:O	2.19	0.49
1:B:495:LEU:HD13	5:B:638:HOH:O	2.13	0.49
1:A:458:ASN:HA	1:A:461:PHE:CE2	2.48	0.48
1:A:246:ASN:OD1	2:C:2[B]:GLY:HA2	2.13	0.48
1:B:130:GLU:O	1:B:289:LYS:HE3	2.13	0.48
1:B:195:GLU:HB3	1:B:381:ILE:HD11	1.95	0.48
1:B:170:LYS:HA	1:B:173[B]:TYR:CD2	2.48	0.47
1:A:238:THR:HG22	1:A:240:LYS:HG2	1.97	0.47
1:B:135:HIS:HD2	5:B:968:HOH:O	1.98	0.47
1:A:334:LYS:NZ	5:A:717:HOH:O	2.45	0.47
1:A:272:HIS:HE1	5:A:889:HOH:O	1.96	0.46
1:B:177[B]:ASN:O	1:B:189:ARG:HD2	2.16	0.45
1:B:127:LYS:O	1:B:130:GLU:HB3	2.16	0.45
1:A:410:HIS:CD2	1:A:412:THR:H	2.30	0.45
1:A:190:PHE:CE2	2:C:3[A]:LYS:HB3	2.52	0.44
1:B:240:LYS:HE3	1:B:240:LYS:HB3	1.64	0.44
1:B:355[A]:ARG:HG3	5:B:893:HOH:O	2.17	0.44
1:A:141:ASP:OD2	5:A:701:HOH:O	2.21	0.44
1:B:318:MET:HG2	1:B:322:ARG:HH21	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:179:ASN:O	1:B:255:ARG:NH2	2.52	0.43
1:B:234:HIS:HD2	1:B:239[B]:GLU:HG3	1.84	0.43
1:B:248:LEU:HG	4:B:501:MYA:H14	2.01	0.42
1:A:319:THR:HG23	1:A:322:ARG:H	1.85	0.41
1:B:496:GLN:HE22	3:B:504[A]:GOL:H12	1.85	0.41

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:B:780:HOH:O	5:B:928:HOH:O[1_556]	2.13	0.07

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	393/402 (98%)	382 (97%)	11 (3%)	0	100	100
1	B	394/402 (98%)	387 (98%)	7 (2%)	0	100	100
2	C	7/8 (88%)	7 (100%)	0	0	100	100
All	All	794/812 (98%)	776 (98%)	18 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	346/362 (96%)	341 (99%)	5 (1%)	67	42
1	B	342/362 (94%)	338 (99%)	4 (1%)	71	48
2	C	7/7 (100%)	7 (100%)	0	100	100
All	All	695/731 (95%)	686 (99%)	9 (1%)	69	44

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	134	THR
1	A	184	ASP
1	A	313	HIS
1	A	322	ARG
1	A	455	LEU
1	B	319	THR
1	B	322	ARG
1	B	334	LYS
1	B	455	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	133	ASN
1	A	272	HIS
1	A	410	HIS
1	B	135	HIS
1	B	391	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 ⓘ

There are no monosaccharides in this entry.

5.6 Ligand geometry ⓘ

8 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
3	GOL	B	504[A]	-	5,5,5	0.95	0	5,5,5	1.06	0
4	MYA	B	501	-	50,65,65	1.22	5 (10%)	65,91,91	1.39	7 (10%)
3	GOL	B	504[B]	-	5,5,5	0.80	0	5,5,5	1.01	0
4	MYA	C	501[A]	-	50,65,65	1.26	6 (12%)	65,91,91	1.42	10 (15%)
3	GOL	B	503	-	5,5,5	0.88	0	5,5,5	1.01	0
3	GOL	A	601	-	5,5,5	0.80	0	5,5,5	0.85	0
4	MYA	C	501[B]	2	50,65,65	1.27	5 (10%)	65,91,91	1.48	12 (18%)
3	GOL	B	502	-	5,5,5	0.84	0	5,5,5	0.90	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
3	GOL	B	504[A]	-	-	3/4/4/4	-
4	MYA	B	501	-	-	0/44/80/80	0/3/3/3
3	GOL	B	504[B]	-	-	2/4/4/4	-
4	MYA	C	501[A]	-	-	0/44/80/80	0/3/3/3
3	GOL	B	503	-	-	0/4/4/4	-
3	GOL	A	601	-	-	0/4/4/4	-
4	MYA	C	501[B]	2	-	1/44/80/80	0/3/3/3
3	GOL	B	502	-	-	0/4/4/4	-

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	501[B]	MYA	C5A-C6A	-4.45	1.34	1.43
4	C	501[A]	MYA	C5A-C6A	-4.44	1.34	1.43
4	B	501	MYA	C5A-C6A	-4.12	1.34	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	501	MYA	C6A-N1A	-2.80	1.34	1.37
4	C	501[B]	MYA	P2A-O5A	-2.79	1.48	1.56
4	C	501[B]	MYA	C5A-N7A	-2.78	1.33	1.39
4	C	501[A]	MYA	C5A-N7A	-2.76	1.33	1.39
4	C	501[A]	MYA	C6A-N1A	-2.75	1.34	1.37
4	C	501[B]	MYA	C6A-N1A	-2.73	1.34	1.37
4	C	501[A]	MYA	P2A-O5A	-2.71	1.48	1.56
4	B	501	MYA	C2A-N1A	2.68	1.39	1.33
4	B	501	MYA	P2A-O5A	-2.57	1.49	1.56
4	C	501[B]	MYA	C2A-N1A	2.53	1.39	1.33
4	C	501[A]	MYA	C2A-N1A	2.52	1.39	1.33
4	B	501	MYA	C5A-N7A	-2.44	1.34	1.39
4	C	501[A]	MYA	C2A-N3A	2.04	1.40	1.34

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	501[B]	MYA	N3A-C2A-N1A	-5.50	118.42	128.82
4	C	501[A]	MYA	N3A-C2A-N1A	-5.44	118.52	128.82
4	B	501	MYA	N3A-C2A-N1A	-5.36	118.67	128.82
4	C	501[A]	MYA	C5A-C4A-N3A	-3.60	120.66	127.41
4	C	501[A]	MYA	O5A-P2A-O4A	3.57	119.76	108.50
4	C	501[B]	MYA	C5A-C4A-N3A	-3.54	120.77	127.41
4	C	501[B]	MYA	O5A-P2A-O4A	3.54	119.66	108.50
4	B	501	MYA	C5A-C4A-N3A	-3.45	120.94	127.41
4	B	501	MYA	O5A-P2A-O4A	3.33	119.01	108.50
4	B	501	MYA	C2X-C3X-C4X	-2.94	98.01	103.22
4	C	501[B]	MYA	N6A-C6A-N1A	2.88	121.03	117.01
4	C	501[A]	MYA	N6A-C6A-N1A	2.72	120.81	117.01
4	C	501[B]	MYA	N9A-C8A-N7A	-2.70	108.31	113.39
4	C	501[A]	MYA	N9A-C8A-N7A	-2.62	108.45	113.39
4	B	501	MYA	N9A-C8A-N7A	-2.53	108.63	113.39
4	C	501[B]	MYA	O4X-C1X-C2X	-2.48	101.23	106.64
4	C	501[B]	MYA	O7A-P3X-O8A	2.36	115.70	105.60
4	B	501	MYA	N6A-C6A-N1A	2.35	120.30	117.01
4	C	501[A]	MYA	C2X-C3X-C4X	-2.27	99.20	103.22
4	C	501[B]	MYA	C5A-C6A-N6A	-2.27	117.27	121.61
4	C	501[A]	MYA	C5A-C6A-N6A	-2.16	117.46	121.61
4	C	501[A]	MYA	O4X-C1X-C2X	-2.16	101.92	106.64
4	B	501	MYA	C8A-N7A-C5A	2.16	108.15	104.24
4	C	501[B]	MYA	C5M-C4M-C3M	-2.11	108.53	112.88
4	C	501[B]	MYA	C8A-N7A-C5A	2.10	108.05	104.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	C	501[A]	MYA	C8A-N7A-C5A	2.07	107.99	104.24
4	C	501[B]	MYA	C2X-C3X-C4X	-2.06	99.56	103.22
4	C	501[B]	MYA	O4X-C4X-C5X	-2.06	102.58	109.37
4	C	501[A]	MYA	P1A-O5X-C5X	2.01	122.40	118.86

There are no chirality outliers.
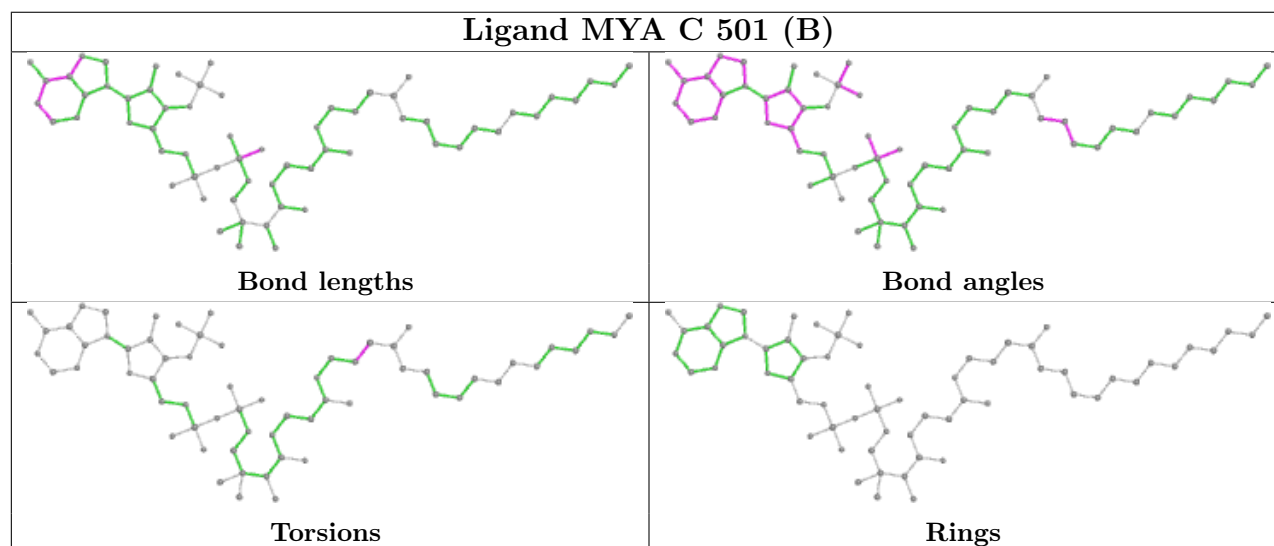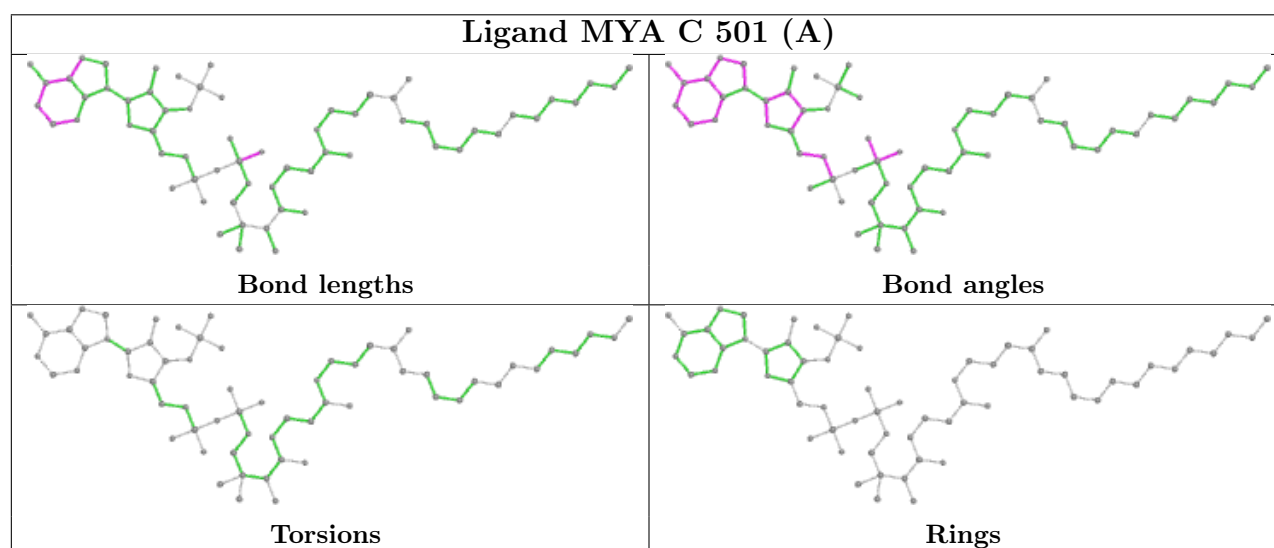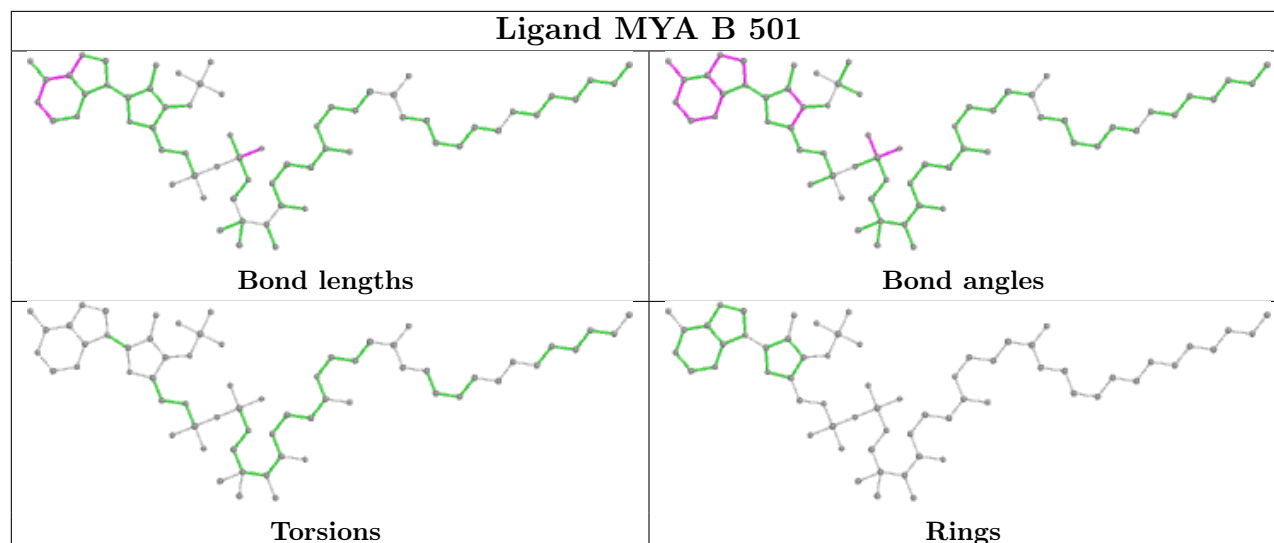
All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	504[A]	GOL	O1-C1-C2-C3
3	B	504[B]	GOL	O1-C1-C2-C3
3	B	504[A]	GOL	O1-C1-C2-O2
3	B	504[A]	GOL	O2-C2-C3-O3
4	C	501[B]	MYA	C3-C2-S1-C2M
3	B	504[B]	GOL	O1-C1-C2-O2

There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	504[A]	GOL	2	0
4	B	501	MYA	1	0
4	C	501[A]	MYA	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 then 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.

Ligand MYA B 501

Bond lengths

Bond angles

Torsions

Rings

Ligand MYA C 501 (A)

Bond lengths

Bond angles

Torsions

Rings

Ligand MYA C 501 (B)

Bond lengths

Bond angles

Torsions

Rings

5.7 Other polymers ⓘ

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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å²)	Q<0.9
1	A	392/402 (97%)	0.37	32 (8%)	11	12	11, 20, 53, 75	0
1	B	387/402 (96%)	0.49	40 (10%)	6	6	10, 21, 52, 71	0
2	C	8/8 (100%)	4.19	6 (75%)	0	0	32, 41, 50, 57	0
All	All	787/812 (96%)	0.47	78 (9%)	7	7	10, 21, 53, 75	0

All (78) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	C	2[A]	GLY	8.6
2	C	8	LEU	7.3
1	B	114	LYS	7.1
1	A	116	SER	7.0
1	B	314	LEU	6.1
1	B	109	MET	6.0
1	B	108	THR	5.9
1	B	185	ASP	5.8
1	A	105	PRO	5.7
1	A	315	SER	5.7
1	B	408	MET	4.7
1	A	114	LYS	4.7
1	B	315	SER	4.6
1	A	318	MET	4.6
1	B	107	LYS	4.5
1	B	112	ALA	4.5
1	B	181	VAL	4.4
1	B	186	ASN	4.4
2	C	9	ARG	4.3
1	A	133	ASN	4.3
1	A	316	ARG	4.2
1	B	192	TYR	4.1
1	A	113	SER	4.0

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Mol	Chain	Res	Type	RSRZ
1	A	313	HIS	3.9
2	C	4	GLN	3.9
1	B	316	ARG	3.9
1	A	317	ASN	3.8
1	A	106	ALA	3.7
1	B	106	ALA	3.7
1	A	112	ALA	3.5
1	B	110	GLU	3.4
1	B	313	HIS	3.4
1	B	187	MET	3.4
1	A	107	LYS	3.3
1	A	321	GLN	3.3
1	B	188	PHE	3.3
1	B	407	ILE	3.2
1	B	183	ASP	3.2
1	A	320	MET	3.2
1	B	307	ILE	3.2
1	A	134	THR	3.1
1	A	314	LEU	3.1
1	A	322	ARG	3.1
1	A	258	ARG	2.9
2	C	5	ASN	2.9
1	A	131	VAL	2.9
2	C	6	SER	2.9
1	A	324	MET	2.8
1	A	184	ASP	2.7
1	B	256	SER	2.7
1	B	312	SER	2.7
1	B	115	ARG	2.6
1	A	109	MET	2.6
1	A	130	GLU	2.6
1	A	108	THR	2.6
1	A	319	THR	2.6
1	A	332	THR	2.6
1	B	182	GLU	2.6
1	B	414	LYS	2.5
1	B	184	ASP	2.5
1	A	129	GLY	2.5
1	B	113	SER	2.4
1	B	309	VAL	2.4
1	B	311	PHE	2.4
1	B	322	ARG	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	284	GLY	2.3
1	B	306	LEU	2.2
1	B	416	LEU	2.2
1	B	319	THR	2.2
1	B	409	ASN	2.2
1	B	326	LEU	2.2
1	A	307	ILE	2.2
1	B	116	SER	2.2
1	A	304	ARG	2.1
1	B	304	ARG	2.1
1	B	320	MET	2.1
1	A	325	LYS	2.1
1	A	118	GLN	2.0

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

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

6.3   Carbohydrates ⓘ
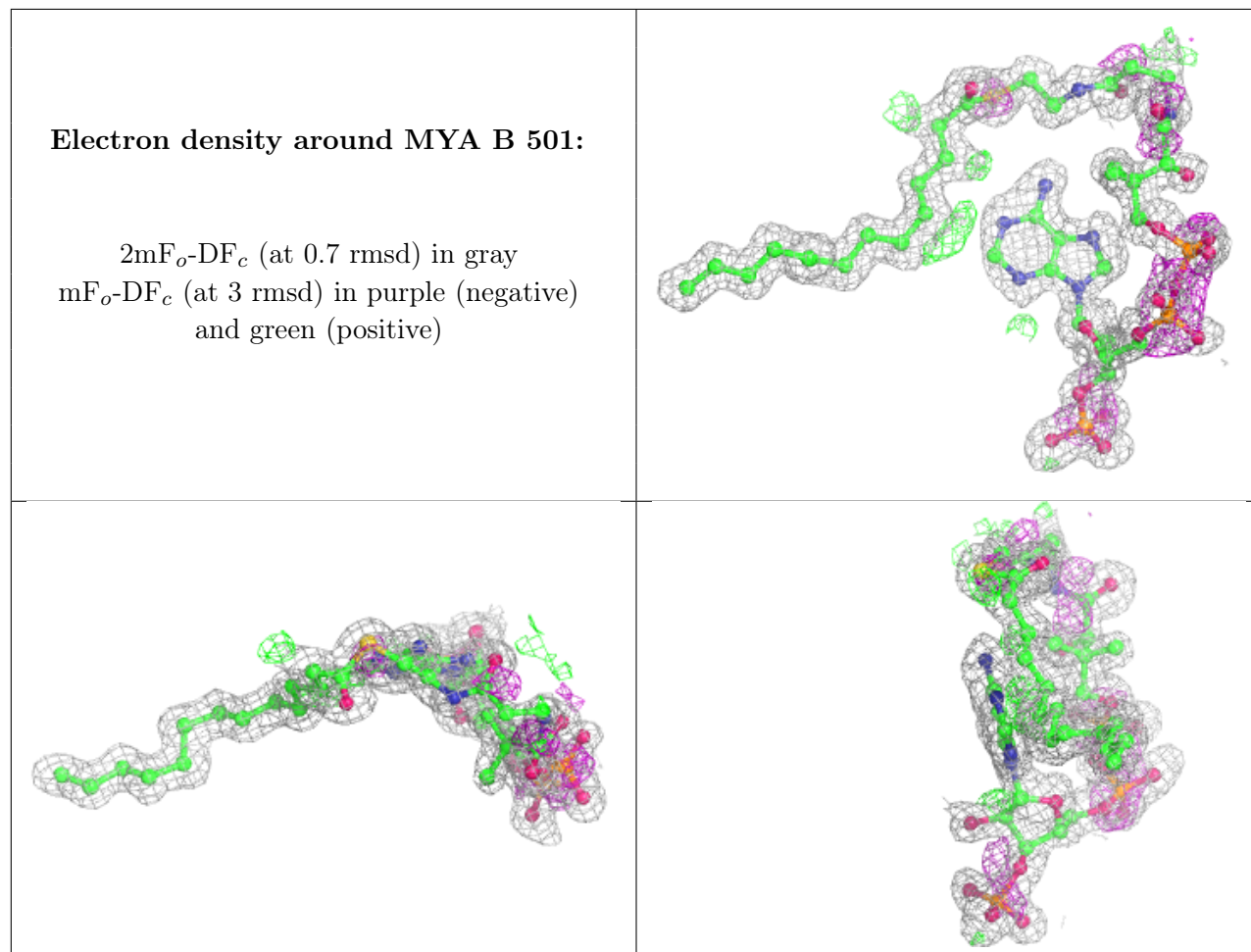
There are no monosaccharides in this entry.

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.

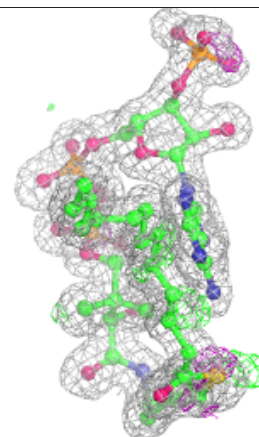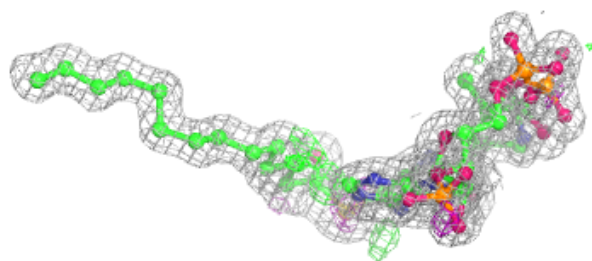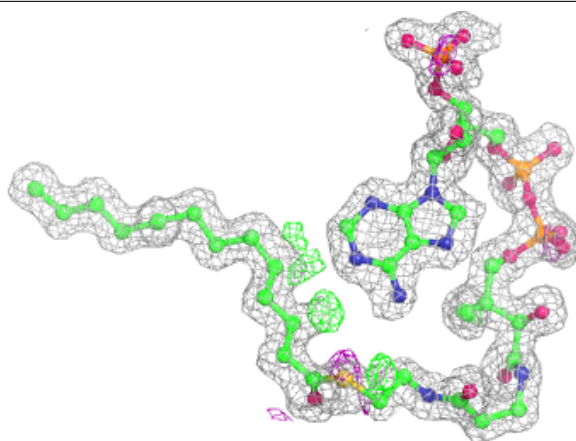| Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	GOL	B	504[A]	6/6	0.84	0.23	19,28,31,36	6
3	GOL	B	504[B]	6/6	0.84	0.23	19,25,26,27	6
4	MYA	B	501	63/63	0.85	0.17	12,32,40,41	0
4	MYA	C	501[A]	63/63	0.90	0.12	12,24,34,39	63
4	MYA	C	501[B]	63/63	0.90	0.12	13,22,27,31	63
3	GOL	B	503	6/6	0.92	0.12	40,45,47,48	0
3	GOL	A	601	6/6	0.98	0.08	17,21,22,22	0
3	GOL	B	502	6/6	0.98	0.08	14,17,20,22	0

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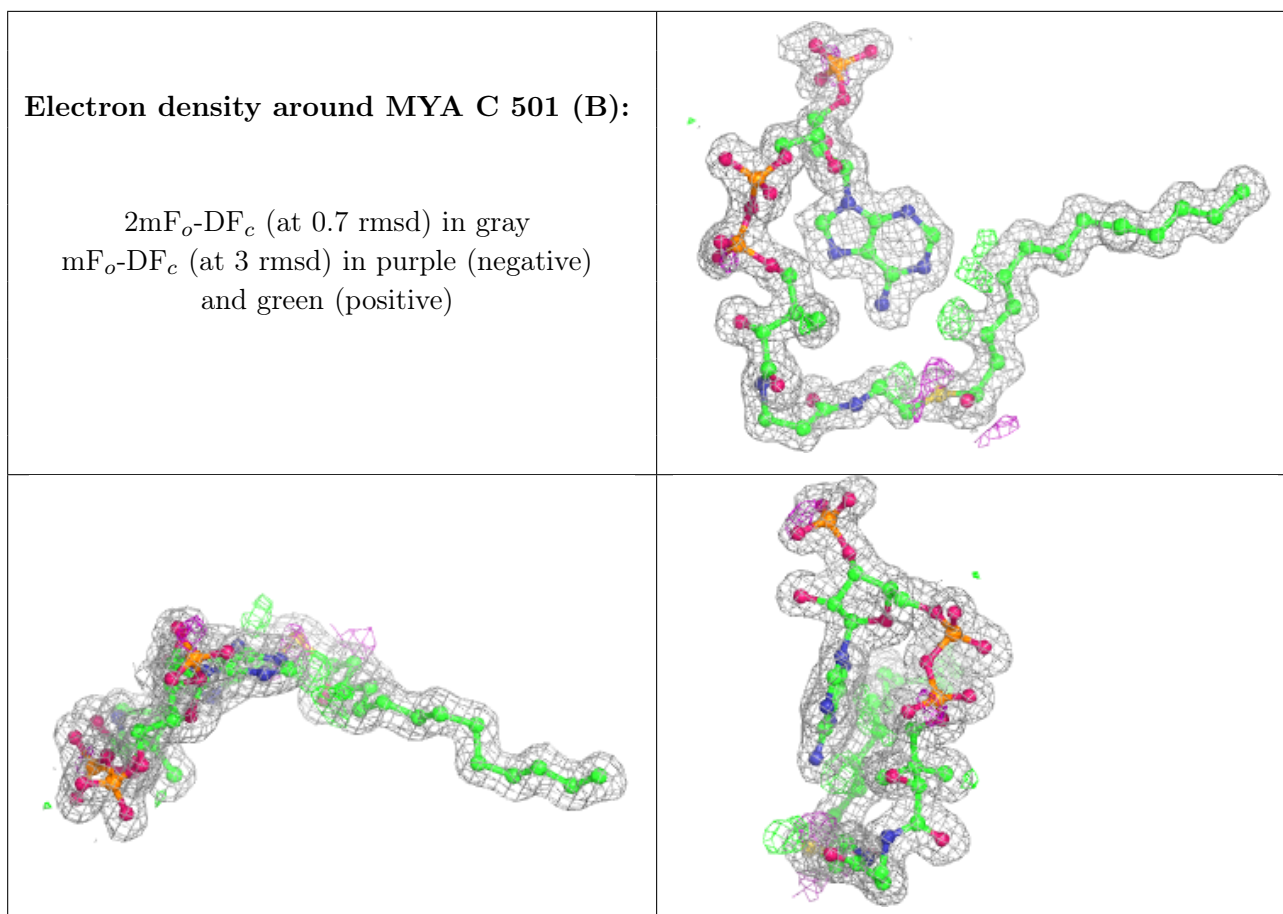
instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around MYA B 501:

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

Electron density around MYA C 501 (A):

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

Electron density around MYA C 501 (B):

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

6.5 Other polymers ⓘ

There are no such residues in this entry.