Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 07:25 am BST

PDB ID	:	4C7I
Title	:	Leismania major N-myristoyltransferase in complex with a peptidomimetic (-OH) molecule
Authors	:	Olaleye, T.O.; Brannigan, J.A.; Goncalves, V.; Roberts, S.M.; Leatherbarrow, R.J.; Wilkinson, A.J.; Tate, E.W.
Deposited on	:	2013-09-20
Resolution	:	1.30 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:

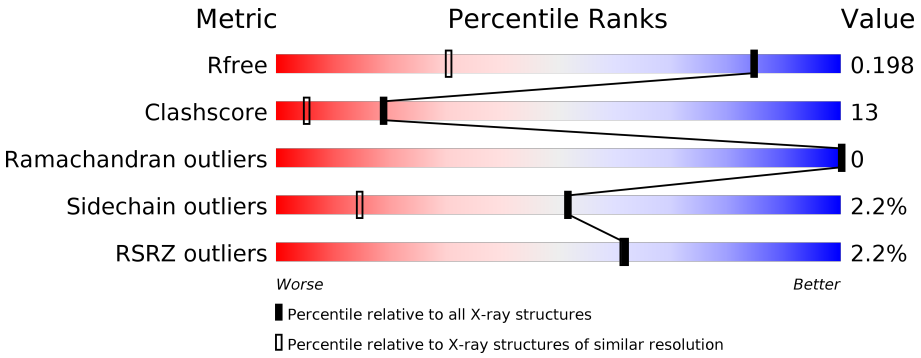
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

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 1.30 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	1058 (1.30-1.30)
Clashscore	141614	1101 (1.30-1.30)
Ramachandran outliers	138981	1058 (1.30-1.30)
Sidechain outliers	138945	1058 (1.30-1.30)
RSRZ outliers	127900	1029 (1.30-1.30)

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 on 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	411	2% 77% 19%

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
5	DMS	A	1425[A]	-	-	X	-
5	DMS	A	1425[B]	-	X	X	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	DMS	A	1426[B]	-	-	X	-

2 Entry composition

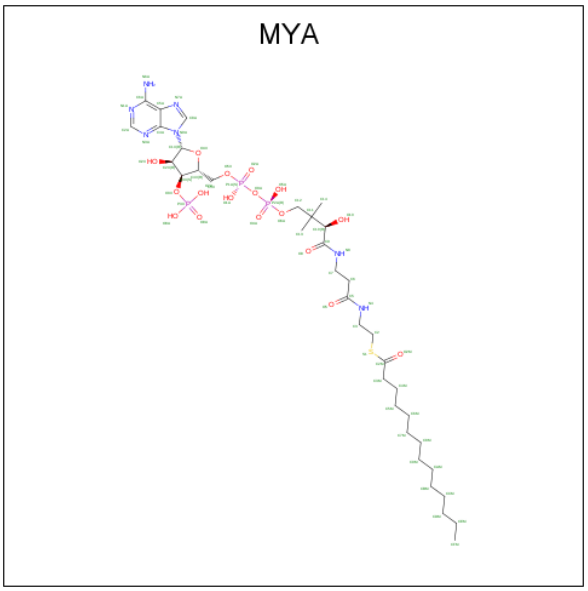
There are 6 unique types of molecules in this entry. The entry contains 4348 atoms, of which 0 are hydrogens and 0 are deuteriums.

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.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	411	Total	C	N	O	S	0	51	0
			3650	2377	602	653	18			

- Molecule 2 is TETRADECANOYL-COA (three-letter code: MYA) (formula: $C_{35}H_{62}N_7O_{17}P_3S$).



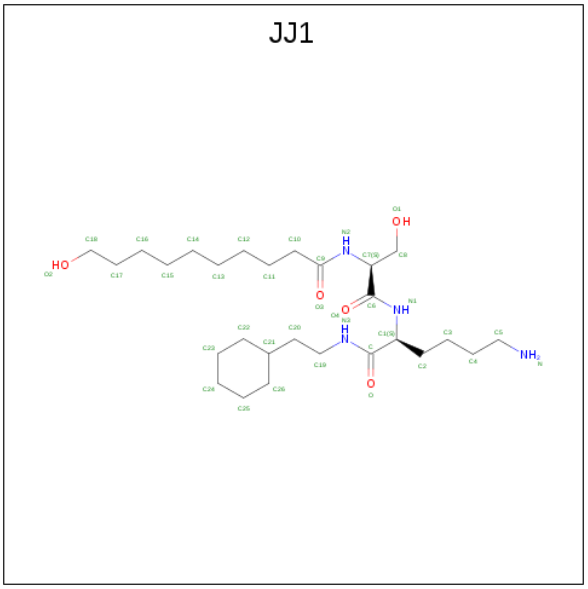
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	S	0	0
			63	35	7	17	3	1		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

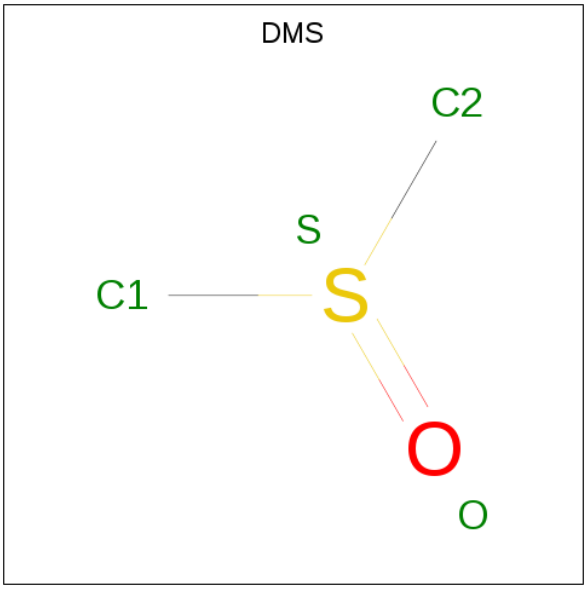
- Molecule 4 is N-[(1S)-1-{[(1S)-5-AMINO-1-[(2-CYCLOHEXYLETHYL)CARBAMO

YL]PENTYL]CARBAMOYL}-2-HYDROXYETHYL]-10-HYDROXYDECANAMIDE (three-letter code: JJ1) (formula: $C_{27}H_{52}N_4O_5$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			36	27	4	5		

- Molecule 5 is DIMETHYL SULFOXIDE (three-letter code: DMS) (formula: C_2H_6OS).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	O	S	0	1
			8	4	2	2		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	O	S	0	1
			8	4	2	2		

- Molecule 6 is water.

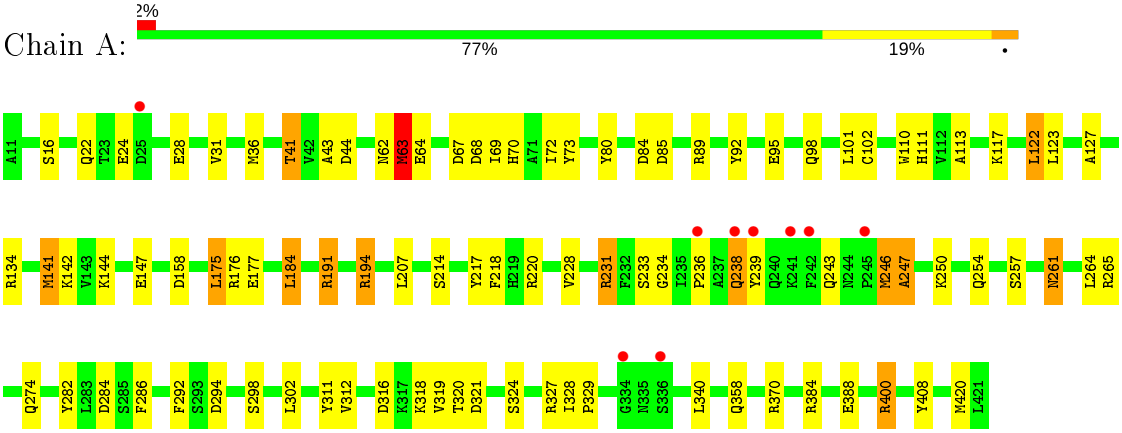
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	579	Total	O	0	3
			582	582		

3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA 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

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	48.64Å 91.06Å 53.60Å 90.00° 113.98° 90.00°	Depositor
Resolution (Å)	45.57 – 1.30 45.53 – 1.30	Depositor EDS
% Data completeness (in resolution range)	96.1 (45.57-1.30) 96.1 (45.53-1.30)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.41 (at 1.30Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.168 , 0.199 0.167 , 0.198	Depositor DCC
R_{free} test set	5043 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å^2)	11.2	Xtriage
Anisotropy	0.392	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 39.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	0.022 for h,-k,-h-l	Xtriage
F_o,F_c correlation	0.97	EDS
Total number of atoms	4348	wwPDB-VP
Average B, all atoms (Å^2)	16.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.33% of the height of the origin peak. No significant pseudotranslation is detected.*

[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: MYA, MG, DMS, JJ1

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	1.51	23/3888 (0.6%)	1.48	64/5274 (1.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	1	1

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	191	ARG	CZ-NH1	9.61	1.45	1.33
1	A	110	TRP	CG-CD1	8.09	1.48	1.36
1	A	388[A]	GLU	CD-OE2	7.00	1.33	1.25
1	A	388[B]	GLU	CD-OE2	7.00	1.33	1.25
1	A	147	GLU	CD-OE2	6.83	1.33	1.25
1	A	388[A]	GLU	CG-CD	6.34	1.61	1.51
1	A	388[B]	GLU	CG-CD	6.34	1.61	1.51
1	A	265	ARG	CZ-NH2	-6.30	1.24	1.33
1	A	239	TYR	CG-CD2	5.99	1.47	1.39
1	A	239	TYR	CB-CG	5.94	1.60	1.51
1	A	384	ARG	CZ-NH2	5.67	1.40	1.33
1	A	408	TYR	CE1-CZ	-5.66	1.31	1.38
1	A	320	THR	CB-OG1	-5.61	1.32	1.43
1	A	257	SER	CA-CB	5.60	1.61	1.52
1	A	234	GLY	CA-C	-5.49	1.43	1.51
1	A	110	TRP	CD2-CE2	5.35	1.47	1.41
1	A	89	ARG	CZ-NH1	5.35	1.40	1.33
1	A	142	LYS	C-O	5.35	1.33	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	95	GLU	CD-OE1	5.22	1.31	1.25
1	A	324	SER	CA-CB	-5.21	1.45	1.52
1	A	282	TYR	CG-CD1	5.13	1.45	1.39
1	A	67[A]	ASP	CB-CG	5.00	1.62	1.51
1	A	67[B]	ASP	CB-CG	5.00	1.62	1.51

All (64) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	191	ARG	NE-CZ-NH2	-10.70	114.95	120.30
1	A	286	PHE	CB-CG-CD1	9.26	127.28	120.80
1	A	134	ARG	NE-CZ-NH2	-8.62	115.99	120.30
1	A	370	ARG	NE-CZ-NH2	-8.54	116.03	120.30
1	A	141	MET	CG-SD-CE	-8.26	86.98	100.20
1	A	84[A]	ASP	CB-CG-OD1	8.21	125.69	118.30
1	A	84[B]	ASP	CB-CG-OD1	8.21	125.69	118.30
1	A	95	GLU	OE1-CD-OE2	-8.17	113.50	123.30
1	A	408	TYR	CB-CG-CD2	8.04	125.82	121.00
1	A	247	ALA	N-CA-CB	8.03	121.34	110.10
1	A	89	ARG	NE-CZ-NH1	7.79	124.19	120.30
1	A	384	ARG	NE-CZ-NH2	-7.59	116.51	120.30
1	A	85	ASP	CB-CG-OD2	7.54	125.08	118.30
1	A	294	ASP	CB-CG-OD1	7.36	124.92	118.30
1	A	294	ASP	CB-CG-OD2	-7.16	111.85	118.30
1	A	36	MET	CA-CB-CG	7.15	125.46	113.30
1	A	89	ARG	NE-CZ-NH2	-7.11	116.74	120.30
1	A	44	ASP	CB-CG-OD1	7.01	124.61	118.30
1	A	191	ARG	NE-CZ-NH1	6.94	123.77	120.30
1	A	265	ARG	NE-CZ-NH2	-6.94	116.83	120.30
1	A	311	TYR	CB-CG-CD1	6.79	125.07	121.00
1	A	265	ARG	NE-CZ-NH1	-6.67	116.97	120.30
1	A	246[A]	MET	C-N-CA	-6.66	105.06	121.70
1	A	246[B]	MET	C-N-CA	-6.66	105.06	121.70
1	A	80	TYR	CB-CG-CD2	-6.65	117.01	121.00
1	A	184[A]	LEU	CB-CG-CD2	6.64	122.28	111.00
1	A	184[B]	LEU	CB-CG-CD2	6.64	122.28	111.00
1	A	63[A]	MET	N-CA-CB	-6.50	98.91	110.60
1	A	63[B]	MET	N-CA-CB	-6.50	98.91	110.60
1	A	220	ARG	NE-CZ-NH2	-6.48	117.06	120.30
1	A	68	ASP	CB-CG-OD2	-6.37	112.56	118.30
1	A	286	PHE	CB-CG-CD2	-6.35	116.35	120.80
1	A	247	ALA	O-C-N	6.10	132.46	122.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	95	GLU	CG-CD-OE1	6.10	130.50	118.30
1	A	265	ARG	NH1-CZ-NH2	6.04	126.04	119.40
1	A	321	ASP	CB-CG-OD1	-6.01	112.89	118.30
1	A	122	LEU	CB-CG-CD1	6.01	121.21	111.00
1	A	246[A]	MET	N-CA-CB	5.96	121.33	110.60
1	A	246[B]	MET	N-CA-CB	5.96	121.33	110.60
1	A	194[A]	ARG	NE-CZ-NH2	-5.89	117.36	120.30
1	A	194[B]	ARG	NE-CZ-NH2	-5.89	117.36	120.30
1	A	158	ASP	CB-CG-OD1	5.86	123.57	118.30
1	A	420[A]	MET	CG-SD-CE	-5.80	90.92	100.20
1	A	420[B]	MET	CG-SD-CE	-5.80	90.92	100.20
1	A	246[A]	MET	CG-SD-CE	-5.72	91.05	100.20
1	A	246[B]	MET	CG-SD-CE	-5.72	91.05	100.20
1	A	44	ASP	CB-CG-OD2	-5.71	113.16	118.30
1	A	92	TYR	CB-CG-CD1	5.70	124.42	121.00
1	A	340	LEU	CB-CG-CD2	-5.55	101.57	111.00
1	A	400[A]	ARG	NE-CZ-NH2	-5.47	117.56	120.30
1	A	400[B]	ARG	NE-CZ-NH2	-5.47	117.56	120.30
1	A	63[A]	MET	CA-C-N	-5.47	105.17	117.20
1	A	63[B]	MET	CA-C-N	-5.47	105.17	117.20
1	A	218	PHE	CB-CG-CD2	-5.40	117.02	120.80
1	A	316	ASP	CB-CG-OD2	-5.39	113.45	118.30
1	A	284	ASP	CB-CG-OD1	5.32	123.09	118.30
1	A	73	TYR	CB-CG-CD2	5.31	124.19	121.00
1	A	84[A]	ASP	CB-CG-OD2	-5.24	113.58	118.30
1	A	84[B]	ASP	CB-CG-OD2	-5.24	113.58	118.30
1	A	292	PHE	CB-CG-CD2	-5.15	117.19	120.80
1	A	175[A]	LEU	CA-CB-CG	-5.15	103.46	115.30
1	A	175[B]	LEU	CA-CB-CG	-5.15	103.46	115.30
1	A	63[A]	MET	CB-CA-C	-5.12	100.17	110.40
1	A	63[B]	MET	CB-CA-C	-5.12	100.17	110.40

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	247	ALA	CA

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	261	ASN	Peptide

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	3650	0	3721	84	0
2	A	63	0	58	0	0
3	A	1	0	0	0	0
4	A	36	0	52	9	0
5	A	16	0	24	17	0
6	A	582	0	0	17	0
All	All	4348	0	3855	99	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 13.

All (99) 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:63[B]:MET:HE1	1:A:72[B]:ILE:CD1	1.54	1.37
1:A:63[B]:MET:CE	1:A:72[B]:ILE:HD12	1.64	1.27
1:A:63[B]:MET:CE	1:A:72[B]:ILE:CD1	2.15	1.25
1:A:63[B]:MET:HE1	1:A:72[B]:ILE:HD11	1.27	1.09
1:A:194[B]:ARG:CG	1:A:194[B]:ARG:HH11	1.67	1.07
1:A:63[B]:MET:HG3	1:A:69[B]:ILE:HD11	1.39	1.05
1:A:24[B]:GLU:OE2	6:A:2033:HOH:O	1.75	1.01
1:A:250:LYS:O	1:A:254[B]:GLN:HG3	1.62	1.00
1:A:194[B]:ARG:HG3	1:A:194[B]:ARG:HH11	0.81	0.97
1:A:194[B]:ARG:NH1	1:A:194[B]:ARG:HG3	1.60	0.95
1:A:72[A]:ILE:HD13	1:A:122:LEU:CD2	2.02	0.90
1:A:318[B]:LYS:NZ	6:A:2483:HOH:O	2.05	0.88
1:A:141:MET:SD	1:A:144[B]:LYS:HD2	2.17	0.84
5:A:1426[A]:DMS:H11	6:A:2508:HOH:O	1.78	0.83
1:A:63[B]:MET:HE2	1:A:72[B]:ILE:CD1	2.08	0.83
1:A:231[A]:ARG:HD3	6:A:2376:HOH:O	1.79	0.81
1:A:62[B]:ASN:HD21	1:A:64[B]:GLU:HG2	1.47	0.80
1:A:72[A]:ILE:CD1	1:A:122:LEU:HD21	2.16	0.76
1:A:63[B]:MET:HE3	1:A:69[B]:ILE:HD13	1.68	0.75
1:A:72[A]:ILE:HD13	1:A:122:LEU:HD22	1.67	0.74
1:A:72[A]:ILE:HD13	1:A:122:LEU:HD21	1.71	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:231[A]:ARG:CD	6:A:2376:HOH:O	2.38	0.72
1:A:117[A]:LYS:HE2	6:A:2111:HOH:O	1.90	0.70
1:A:63[B]:MET:HG3	1:A:69[B]:ILE:CD1	2.19	0.70
1:A:28[B]:GLU:OE1	1:A:28[B]:GLU:HA	1.91	0.70
4:A:1424:JJ1:H2	5:A:1425[B]:DMS:H12	1.56	0.69
1:A:63[A]:MET:HG3	1:A:111[A]:HIS:CE1	2.28	0.68
1:A:250:LYS:O	1:A:254[B]:GLN:CG	2.38	0.67
1:A:63[A]:MET:HB3	1:A:102:CYS:SG	2.35	0.67
1:A:63[B]:MET:HE1	1:A:72[B]:ILE:HD12	1.33	0.64
1:A:63[B]:MET:HE2	1:A:72[B]:ILE:HD12	1.66	0.63
1:A:117[A]:LYS:CE	6:A:2111:HOH:O	2.46	0.62
1:A:101[A]:LEU:HD23	6:A:2220[A]:HOH:O	1.99	0.62
1:A:16:SER:HA	1:A:22:GLN:HE22	1.65	0.62
1:A:274:GLN:NE2	1:A:319:VAL:H	1.99	0.60
1:A:62[B]:ASN:ND2	1:A:64[B]:GLU:HG2	2.15	0.57
1:A:63[B]:MET:CE	1:A:69[B]:ILE:HD13	2.34	0.57
4:A:1424:JJ1:H2	5:A:1425[B]:DMS:C1	2.17	0.56
4:A:1424:JJ1:H2	5:A:1425[A]:DMS:H12	1.71	0.55
1:A:327:ARG:NH1	6:A:2206:HOH:O	2.38	0.55
1:A:274:GLN:HE22	1:A:319:VAL:H	1.55	0.54
1:A:72[A]:ILE:CD1	1:A:122:LEU:CD2	2.75	0.54
1:A:63[A]:MET:CG	1:A:111[A]:HIS:ND1	2.71	0.54
1:A:228:VAL:O	1:A:231[A]:ARG:NE	2.40	0.54
1:A:123:LEU:HD13	1:A:175[B]:LEU:HD21	1.90	0.53
1:A:41:THR:HG23	1:A:43:ALA:H	1.72	0.53
1:A:31:VAL:HA	1:A:141:MET:CE	2.39	0.53
1:A:214[B]:SER:OG	1:A:400[B]:ARG:HG2	2.09	0.52
5:A:1425[B]:DMS:H13	6:A:2508:HOH:O	2.08	0.52
1:A:63[A]:MET:HG2	1:A:111[A]:HIS:ND1	2.25	0.52
1:A:63[B]:MET:HB3	1:A:102:CYS:SG	2.48	0.52
4:A:1424:JJ1:H101	5:A:1425[A]:DMS:H12	1.92	0.52
1:A:194[B]:ARG:NH1	1:A:194[B]:ARG:CG	2.39	0.51
1:A:72[A]:ILE:HD12	1:A:122:LEU:HD21	1.93	0.51
1:A:98[B]:GLN:OE1	6:A:2216:HOH:O	2.19	0.51
1:A:31:VAL:HA	1:A:141:MET:HE2	1.92	0.51
1:A:70:HIS:HD2	6:A:2140:HOH:O	1.94	0.51
4:A:1424:JJ1:H2	5:A:1425[A]:DMS:C1	2.21	0.50
4:A:1424:JJ1:N2	5:A:1425[B]:DMS:H12	2.24	0.50
1:A:261:ASN:HD21	1:A:358:GLN:HE21	1.58	0.50
4:A:1424:JJ1:H101	5:A:1425[A]:DMS:C1	2.42	0.50
1:A:261:ASN:ND2	1:A:358:GLN:HE21	2.09	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:28[B]:GLU:OE1	1:A:28[B]:GLU:CA	2.58	0.50
1:A:176:ARG:O	1:A:177[B]:GLU:HB2	2.11	0.50
1:A:16:SER:HA	1:A:22:GLN:NE2	2.26	0.49
1:A:236:PRO:CB	1:A:238:GLN:OE1	2.62	0.48
1:A:63[B]:MET:HE2	1:A:72[B]:ILE:HD13	1.91	0.48
1:A:63[A]:MET:CG	1:A:111[A]:HIS:CE1	2.96	0.47
1:A:63[A]:MET:CE	1:A:111[A]:HIS:ND1	2.78	0.47
4:A:1424:JJ1:H101	5:A:1425[B]:DMS:H12	1.97	0.46
1:A:207[B]:LEU:HD22	1:A:400[B]:ARG:NE	2.30	0.46
1:A:194[B]:ARG:HA	1:A:194[B]:ARG:HD2	1.34	0.46
1:A:62[A]:ASN:C	1:A:62[A]:ASN:OD1	2.54	0.45
1:A:70:HIS:HE1	6:A:2149:HOH:O	2.00	0.45
1:A:264:LEU:HA	1:A:312:VAL:O	2.16	0.45
1:A:123:LEU:HB3	1:A:175[B]:LEU:HD22	1.98	0.45
1:A:113:ALA:HB1	1:A:122:LEU:CD1	2.47	0.45
5:A:1425[A]:DMS:H13	6:A:2508:HOH:O	2.15	0.45
5:A:1426[B]:DMS:C1	6:A:2574:HOH:O	2.65	0.45
1:A:101[A]:LEU:HD22	1:A:127:ALA:HB1	1.99	0.44
1:A:31:VAL:CB	1:A:141:MET:HE1	2.48	0.43
1:A:62[A]:ASN:OD1	1:A:63[A]:MET:N	2.50	0.43
1:A:246[B]:MET:HE3	1:A:250:LYS:HG3	2.01	0.43
1:A:238:GLN:HE21	1:A:238:GLN:HB3	1.68	0.43
1:A:236:PRO:HB2	1:A:238:GLN:OE1	2.19	0.42
4:A:1424:JJ1:N2	5:A:1425[A]:DMS:H12	2.34	0.42
1:A:207[B]:LEU:HA	1:A:207[B]:LEU:HD13	1.83	0.42
1:A:328:ILE:HD11	5:A:1426[B]:DMS:H21	2.02	0.42
1:A:231[A]:ARG:HA	1:A:231[A]:ARG:HD3	1.81	0.41
1:A:217:TYR:CE2	5:A:1426[A]:DMS:H12	2.56	0.41
1:A:111[B]:HIS:HD2	1:A:127:ALA:HA	1.85	0.41
1:A:298[B]:SER:OG	1:A:302:LEU:HD12	2.19	0.41
1:A:247:ALA:HB1	6:A:2400:HOH:O	2.20	0.41
1:A:274:GLN:HE22	1:A:318[B]:LYS:HA	1.86	0.41
5:A:1426[B]:DMS:H12	6:A:2574:HOH:O	2.20	0.41
1:A:236:PRO:HB3	1:A:238:GLN:OE1	2.21	0.40

There are no symmetry-related clashes.

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	461/411 (112%)	452 (98%)	9 (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	413/362 (114%)	401 (97%)	12 (3%)	42 7

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	41	THR
1	A	63[A]	MET
1	A	63[B]	MET
1	A	184[A]	LEU
1	A	184[B]	LEU
1	A	231[A]	ARG
1	A	231[B]	ARG
1	A	233[A]	SER
1	A	233[B]	SER
1	A	238	GLN
1	A	243	GLN
1	A	329	PRO

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	22	GLN
1	A	70	HIS
1	A	91	ASN
1	A	243	GLN
1	A	252	ASN
1	A	261	ASN
1	A	274	GLN
1	A	339	ASN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no carbohydrates in this entry.

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	$\#\|Z\| > 2$
5	DMS	A	1425[A]	-	3,3,3	1.26	0	3,3,3	1.31	1 (33%)
2	MYA	A	1422	3	54,65,65	1.18	4 (7%)	67,91,91	1.37	5 (7%)
5	DMS	A	1426[A]	-	3,3,3	0.53	0	3,3,3	1.32	1 (33%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
4	JJ1	A	1424	-	36,36,36	1.03	1 (2%)	42,42,42	1.00	2 (4%)
5	DMS	A	1426[B]	-	3,3,3	0.54	0	3,3,3	1.32	1 (33%)
5	DMS	A	1425[B]	-	3,3,3	0.96	0	3,3,3	3.46	3 (100%)

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	JJ1	A	1424	-	-	2/39/47/47	0/1/1/1
2	MYA	A	1422	3	-	2/59/80/80	0/3/3/3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1422	MYA	P3X-O3X	3.74	1.66	1.59
4	A	1424	JJ1	C8-C7	3.40	1.60	1.53
2	A	1422	MYA	O4X-C4X	2.90	1.51	1.45
2	A	1422	MYA	C13-C11	2.50	1.59	1.53
2	A	1422	MYA	P1A-O2A	-2.05	1.43	1.50

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	1422	MYA	O2M-C2M-C3M	6.54	120.67	109.02
5	A	1425[B]	DMS	O-S-C2	-4.35	84.37	106.54
2	A	1422	MYA	O6A-C12-C11	-2.99	105.75	110.55
5	A	1425[B]	DMS	O-S-C1	-2.93	91.61	106.54
5	A	1425[B]	DMS	C2-S-C1	2.91	113.43	98.44
4	A	1424	JJ1	C14-C13-C12	2.88	129.06	114.42
2	A	1422	MYA	C4M-C3M-C2M	-2.77	106.16	113.80
2	A	1422	MYA	C13-C11-C10	2.57	113.27	108.82
2	A	1422	MYA	C5A-C6A-N6A	2.19	123.69	120.35
5	A	1426[B]	DMS	C2-S-C1	2.14	109.46	98.44
5	A	1426[A]	DMS	C2-S-C1	2.14	109.44	98.44
5	A	1425[A]	DMS	C2-S-C1	2.05	108.98	98.44
4	A	1424	JJ1	C17-C16-C15	-2.00	104.25	114.42

There are no chirality outliers.

All (4) torsion outliers are listed below:

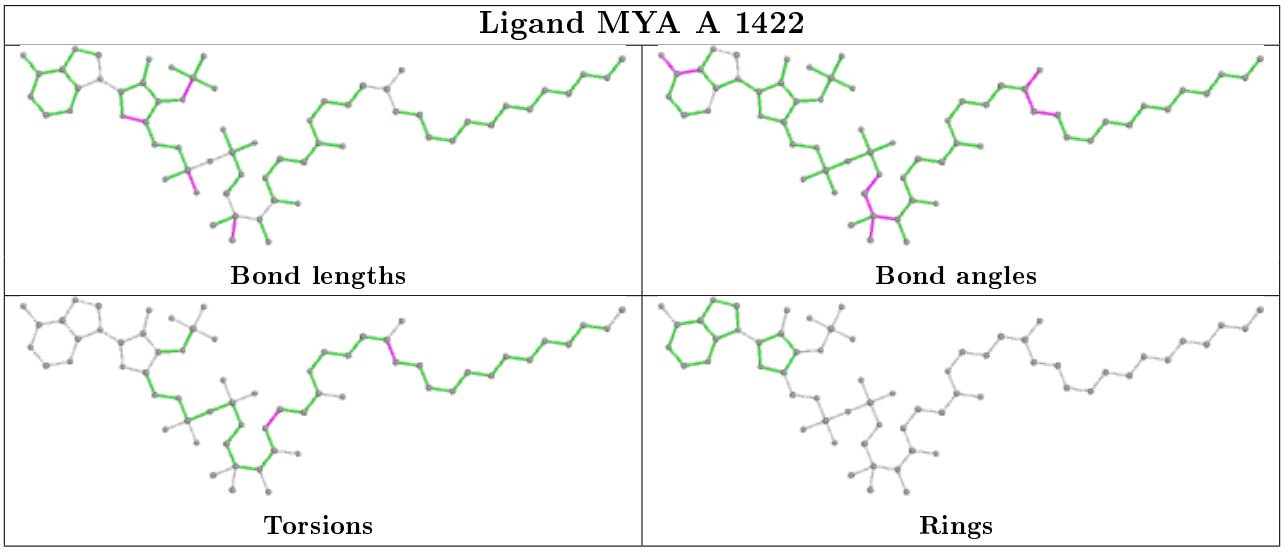
Mol	Chain	Res	Type	Atoms
2	A	1422	MYA	S1-C2M-C3M-C4M
2	A	1422	MYA	C6-C7-N8-C9
4	A	1424	JJ1	O-C-C1-N1
4	A	1424	JJ1	N3-C-C1-N1

There are no ring outliers.

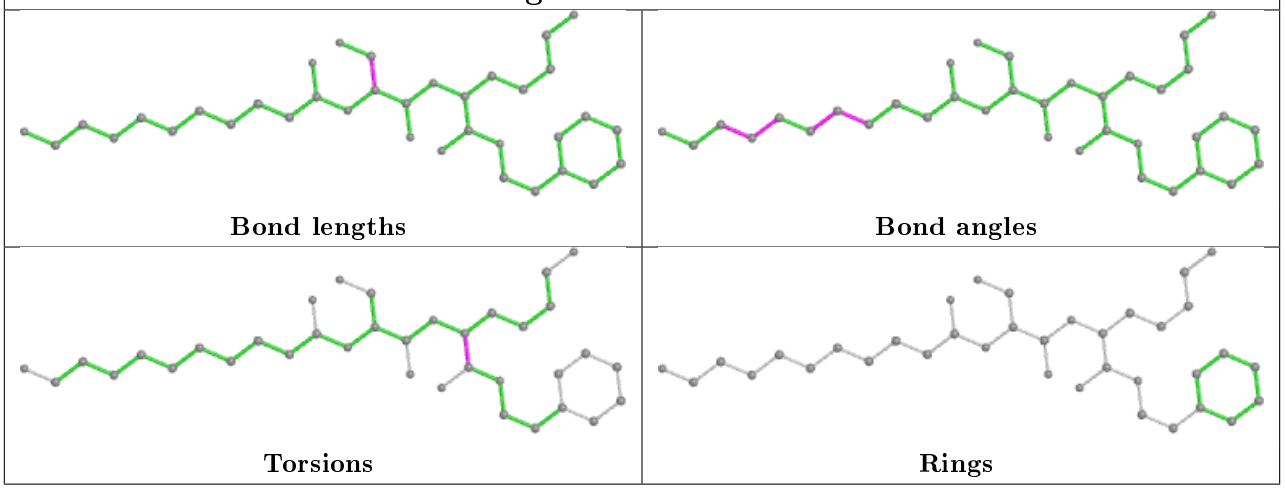
5 monomers are involved in 17 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	1425[A]	DMS	6	0
5	A	1426[A]	DMS	2	0
4	A	1424	JJ1	9	0
5	A	1426[B]	DMS	4	0
5	A	1425[B]	DMS	5	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 JJ1 A 1424



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, 95^{th} 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	411/411 (100%)	0.05	9 (2%) 62 61	7, 12, 28, 42	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	242	PHE	5.4
1	A	245	PRO	3.4
1	A	334	GLY	3.2
1	A	336	SER	3.1
1	A	241	LYS	2.8
1	A	238	GLN	2.7
1	A	25[A]	ASP	2.5
1	A	239	TYR	2.2
1	A	236	PRO	2.1

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 carbohydrates in this entry.

6.4 Ligands

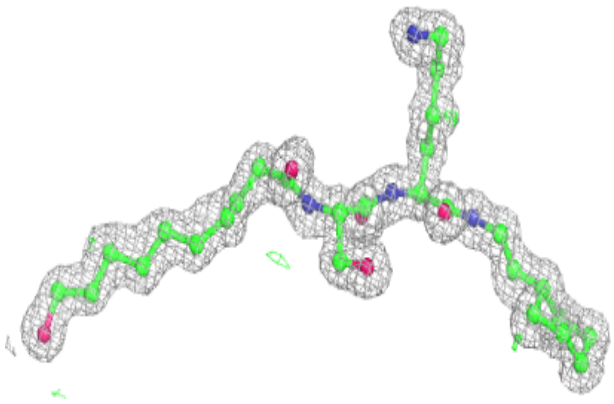
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9'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

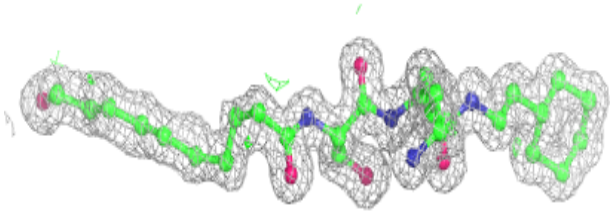
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
5	DMS	A	1426[A]	4/4	0.91	0.15	9,15,17,22	4
5	DMS	A	1426[B]	4/4	0.91	0.15	10,10,12,13	4
5	DMS	A	1425[A]	4/4	0.93	0.16	13,15,17,19	4
5	DMS	A	1425[B]	4/4	0.93	0.16	14,15,20,20	4
4	JJ1	A	1424	36/36	0.97	0.07	8,12,21,21	0
2	MYA	A	1422	63/63	0.98	0.06	7,10,12,14	0
3	MG	A	1423	1/1	0.99	0.04	19,19,19,19	0

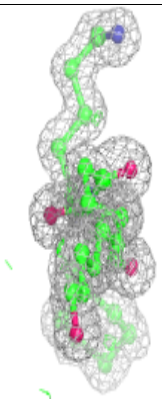
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around JJ1 A 1424:

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

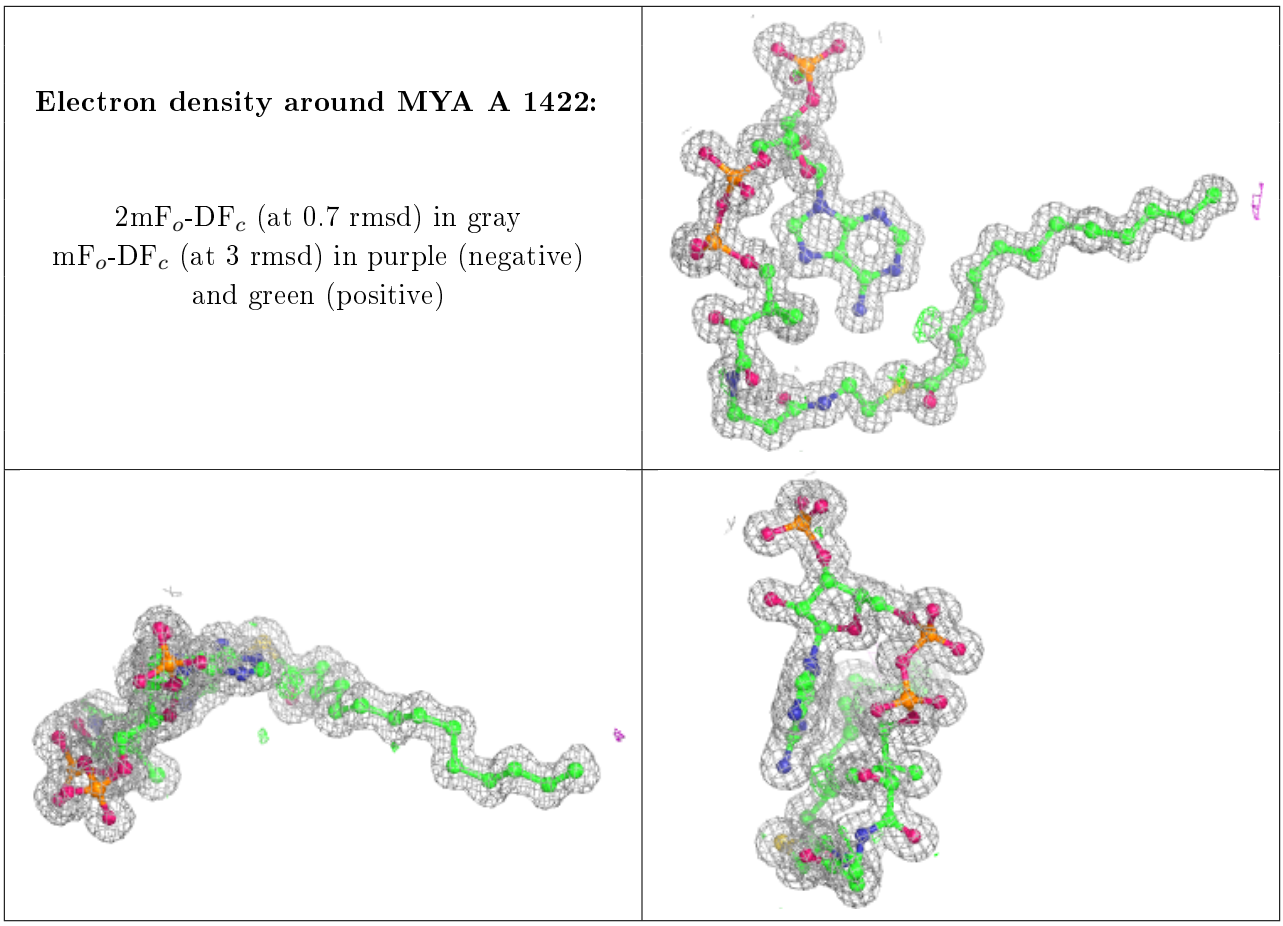






Electron density around MYA A 1422:

$2mF_o$-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.