



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 03:52 PM UTC

PDB ID : 4C7I / pdb_00004c7i
Title : Leishmania 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

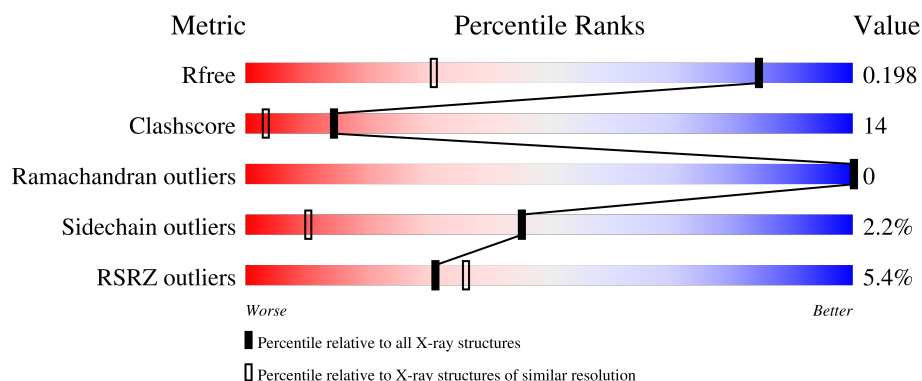
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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}	180053	1553 (1.30-1.30)
Clashscore	190562	1595 (1.30-1.30)
Ramachandran outliers	187476	1551 (1.30-1.30)
Sidechain outliers	187428	1551 (1.30-1.30)
RSRZ outliers	180081	1549 (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 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	411	5% 72% 24% .

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	-
5	DMS	A	1426[A]	-	-	X	-
5	DMS	A	1426[B]	-	-	X	-

2 Entry composition [i](#)

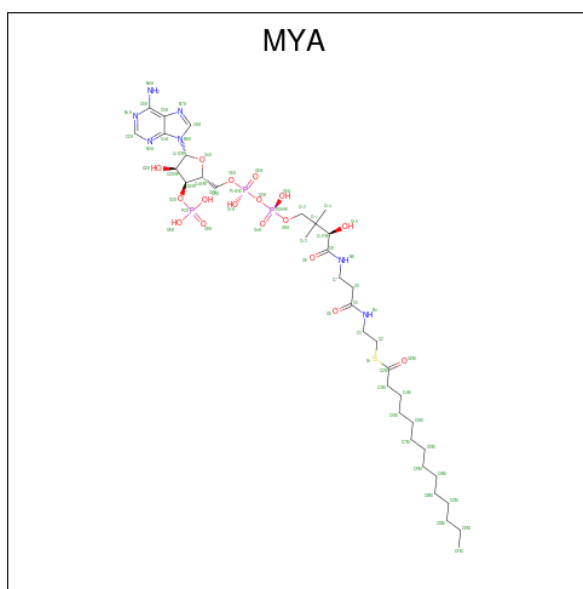
There are 6 unique types of molecules in this entry. The entry contains 4348 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 (CCD ID: MYA) (formula: $C_{35}H_{62}N_7O_{17}P_3S$).



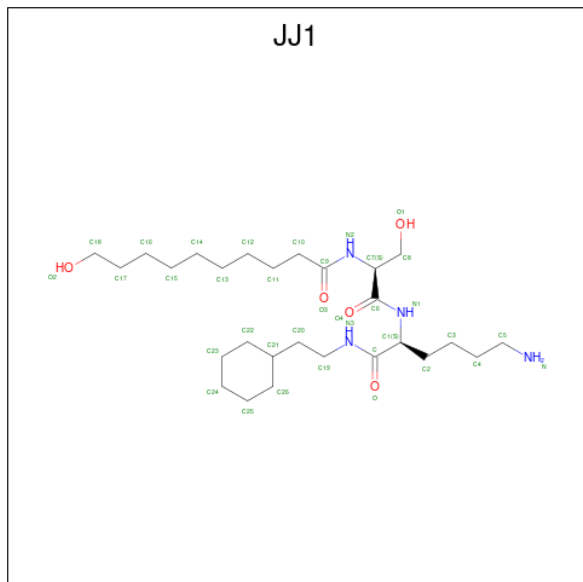
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P S	0	0
			63	35	7	17	3 1		

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

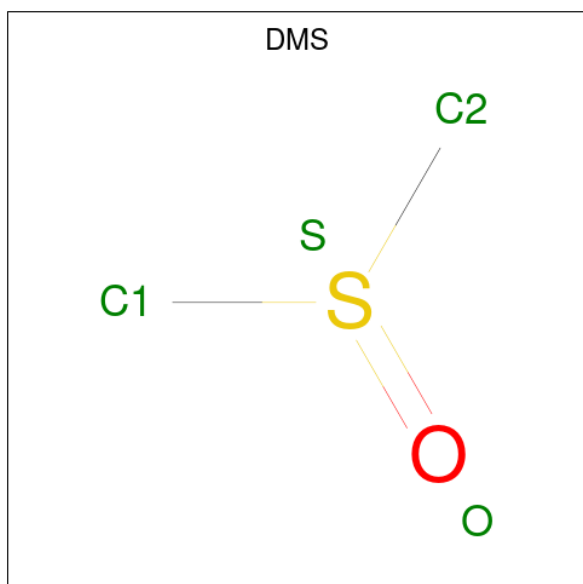
- Molecule 4 is N-[(1S)-1-{[(1S)-5-AMINO-1-[(2-CYCLOHEXYLETHYL)CARBAMOYL]PE

NTYL]CARBAMOYL}-2-HYDROXYETHYL]-10-HYDROXYDECANAMIDE (CCD ID: 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 (CCD ID: 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		

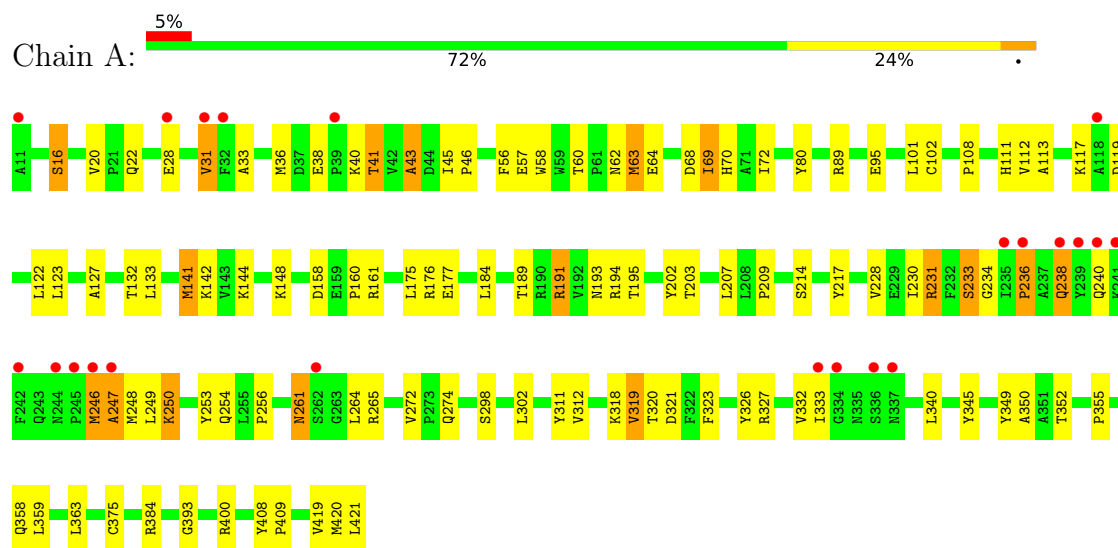
- Molecule 6 is water.

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

3 Residue-property plots

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



4 Data and refinement statistics

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 45.57 – 1.30	Depositor EDS
% Data completeness (in resolution range)	96.1 (45.57-1.30) 96.1 (45.57-1.30)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ 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 , 35.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	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.*

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: DMS, JJ1, MYA, MG

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	1.84	62/3888 (1.6%)	1.49	29/5274 (0.5%)

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 (62) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	234	GLY	CA-C	-10.40	1.43	1.52
1	A	333	ILE	CA-C	9.39	1.62	1.53
1	A	191	ARG	CZ-NH1	9.14	1.45	1.32
1	A	350	ALA	N-CA	8.35	1.55	1.46
1	A	234	GLY	N-CA	7.70	1.52	1.44
1	A	142	LYS	C-O	7.49	1.33	1.24
1	A	320	THR	CB-OG1	-7.33	1.32	1.43
1	A	230	ILE	C-O	7.24	1.32	1.24
1	A	40	LYS	CA-C	-6.92	1.44	1.52
1	A	355	PRO	CA-CB	6.68	1.62	1.53
1	A	265	ARG	CZ-NH2	-6.61	1.24	1.33
1	A	20	VAL	C-O	6.61	1.28	1.24
1	A	56	PHE	CA-C	-6.57	1.44	1.52
1	A	141	MET	C-O	6.43	1.32	1.24
1	A	46	PRO	CA-CB	6.41	1.61	1.53
1	A	311	TYR	CA-C	-6.40	1.44	1.52
1	A	38	GLU	CA-CB	6.32	1.62	1.53
1	A	332	VAL	C-N	-6.26	1.27	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	393	GLY	N-CA	6.16	1.51	1.45
1	A	133	LEU	C-O	6.12	1.31	1.24
1	A	333	ILE	C-N	6.11	1.40	1.33
1	A	359	LEU	CA-CB	-6.04	1.43	1.53
1	A	160	PRO	N-CA	6.03	1.54	1.47
1	A	80	TYR	N-CA	5.92	1.52	1.46
1	A	384	ARG	N-CA	5.91	1.53	1.46
1	A	60	THR	CA-C	-5.89	1.45	1.52
1	A	236	PRO	C-O	5.82	1.30	1.23
1	A	69[A]	ILE	N-CA	5.81	1.53	1.46
1	A	69[B]	ILE	N-CA	5.81	1.53	1.46
1	A	352	THR	C-O	5.79	1.31	1.24
1	A	112	VAL	C-O	5.74	1.30	1.24
1	A	256	PRO	C-O	5.71	1.30	1.23
1	A	349	TYR	C-N	5.68	1.40	1.33
1	A	33	ALA	C-O	5.66	1.31	1.23
1	A	323	PHE	N-CA	5.53	1.53	1.45
1	A	58	TRP	CA-C	-5.50	1.45	1.52
1	A	112	VAL	N-CA	5.49	1.52	1.46
1	A	375	CYS	N-CA	-5.47	1.39	1.46
1	A	384	ARG	CD-NE	5.44	1.53	1.46
1	A	409	PRO	C-O	5.41	1.30	1.23
1	A	359	LEU	N-CA	5.39	1.52	1.46
1	A	202	TYR	CA-C	-5.38	1.46	1.52
1	A	408	TYR	CA-C	5.37	1.59	1.52
1	A	247	ALA	C-O	5.36	1.31	1.24
1	A	384	ARG	CZ-NH2	5.36	1.40	1.33
1	A	119	ASP	N-CA	5.31	1.53	1.46
1	A	240	GLN	N-CA	5.31	1.53	1.46
1	A	189	THR	C-O	5.29	1.30	1.24
1	A	253	TYR	C-N	5.27	1.40	1.33
1	A	319	VAL	N-CA	5.25	1.52	1.46
1	A	57	GLU	C-O	5.21	1.30	1.23
1	A	89	ARG	CZ-NH1	5.18	1.40	1.32
1	A	43	ALA	CA-C	5.18	1.59	1.52
1	A	68	ASP	CA-C	5.17	1.59	1.52
1	A	132	THR	CA-C	-5.11	1.46	1.52
1	A	45	ILE	CA-C	-5.10	1.47	1.52
1	A	161	ARG	C-O	5.06	1.30	1.23
1	A	108	PRO	N-CD	5.05	1.54	1.47
1	A	195	THR	C-O	5.03	1.30	1.24
1	A	31	VAL	C-O	5.01	1.30	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	419	VAL	N-CA	5.01	1.52	1.46
1	A	38	GLU	C-N	5.01	1.39	1.33

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	246[A]	MET	CA-C-N	-9.43	105.06	121.66
1	A	246[A]	MET	C-N-CA	-9.43	105.06	121.66
1	A	246[B]	MET	CA-C-N	-9.43	105.06	121.66
1	A	246[B]	MET	C-N-CA	-9.43	105.06	121.66
1	A	247	ALA	O-C-N	7.30	132.46	122.46
1	A	247	ALA	N-CA-C	7.22	122.19	113.23
1	A	247	ALA	N-CA-CB	6.99	121.34	110.44
1	A	20	VAL	O-C-N	-6.56	117.56	121.69
1	A	141	MET	CG-SD-CE	-6.33	86.98	100.90
1	A	248	MET	N-CA-C	-6.22	104.50	111.28
1	A	333	ILE	N-CA-CB	-6.14	105.27	111.90
1	A	249	LEU	N-CA-C	-6.11	104.70	111.36
1	A	193	ASN	CB-CG-ND2	6.05	125.47	116.40
1	A	272	VAL	O-C-N	-6.05	116.55	120.42
1	A	233[A]	SER	CA-C-N	-6.04	116.75	121.82
1	A	233[A]	SER	C-N-CA	-6.04	116.75	121.82
1	A	233[B]	SER	CA-C-N	-6.04	116.75	121.82
1	A	233[B]	SER	C-N-CA	-6.04	116.75	121.82
1	A	193	ASN	CA-C-O	-5.76	113.59	120.10
1	A	250	LYS	N-CA-C	5.75	117.55	111.28
1	A	36	MET	N-CA-C	-5.73	105.96	113.12
1	A	36	MET	CA-CB-CG	5.68	125.46	114.10
1	A	209	PRO	CA-C-N	5.57	131.73	121.70
1	A	209	PRO	C-N-CA	5.57	131.73	121.70
1	A	321	ASP	O-C-N	-5.53	116.95	123.25
1	A	16	SER	CA-C-O	-5.53	112.81	119.49
1	A	158	ASP	N-CA-C	5.46	118.15	111.82
1	A	95	GLU	CG-CD-OE1	5.26	130.50	118.40
1	A	265	ARG	NH1-CZ-NH2	5.18	126.04	119.30

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 ⓘ

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	93	0
2	A	63	0	58	0	0
3	A	1	0	0	0	0
4	A	36	0	52	8	0
5	A	16	0	24	21	0
6	A	582	0	0	15	0
All	All	4348	0	3855	107	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 14.

All (107) 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	1.47	1.40
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	1.23	1.10
1:A:194[B]:ARG:HH11	1:A:194[B]:ARG:CG	1.67	1.07
1:A:63[B]:MET:HG3	1:A:69[B]:ILE:HD11	1.39	1.05
1:A:250:LYS:O	1:A:254[B]:GLN:HG3	1.62	1.00
1:A:194[B]:ARG:HH11	1:A:194[B]:ARG:HG3	0.81	0.97
1:A:194[B]:ARG:HG3	1:A:194[B]:ARG:NH1	1.60	0.95
5:A:1426[B]:DMS:H12	6:A:2574:HOH:O	1.70	0.92
1:A:72[A]:ILE:HD13	1:A:122:LEU:CD2	2.02	0.90
5:A:1425[B]:DMS:H11	6:A:2508:HOH:O	1.72	0.88
1:A:31:VAL:HA	1:A:141:MET:HE2	1.56	0.86
1:A:141:MET:SD	1:A:144[B]:LYS:HD2	2.17	0.84
1:A:318[B]:LYS:NZ	6:A:2483:HOH:O	2.05	0.84
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:72[A]:ILE:CD1	1:A:122:LEU:HD21	2.16	0.76
1:A:217:TYR:OH	5:A:1426[B]:DMS:H13	1.85	0.76
1:A:217:TYR:OH	5:A:1426[A]:DMS:H13	1.87	0.75
1:A:63[B]:MET:HE1	1:A:72[B]:ILE:HD12	1.32	0.74
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
1:A:63[B]:MET:HE2	1:A:72[B]:ILE:CD1	2.17	0.72
1:A:231[A]:ARG:CD	6:A:2376:HOH:O	2.38	0.72
4:A:1424:JJ1:H101	5:A:1425[B]:DMS:H12	1.70	0.71
1:A:117[A]:LYS:HE2	6:A:2111:HOH:O	1.90	0.70
4:A:1424:JJ1:H2	5:A:1425[B]:DMS:H13	1.55	0.70
1:A:63[B]:MET:HG3	1:A:69[B]:ILE:CD1	2.19	0.70
1:A:28[B]:GLU:HA	1:A:28[B]:GLU:OE1	1.91	0.70
5:A:1425[B]:DMS:H21	6:A:2507:HOH:O	1.92	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:HE3	1:A:69[B]:ILE:HD13	1.80	0.64
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:421:LEU:OXT	5:A:1426[B]:DMS:H22	2.00	0.62
1:A:16:SER:HA	1:A:22:GLN:HE22	1.65	0.62
1:A:63[B]:MET:HE2	1:A:72[B]:ILE:HD12	1.70	0.62
4:A:1424:JJ1:H101	5:A:1425[A]:DMS:H13	1.82	0.60
1:A:274:GLN:NE2	1:A:319:VAL:H	1.99	0.60
1:A:203:THR:HG23	1:A:420[A]:MET:HE3	1.83	0.60
1:A:31:VAL:CA	1:A:141:MET:HE2	2.31	0.58
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
1:A:345:TYR:HE2	5:A:1426[A]:DMS:H12	1.70	0.56
4:A:1424:JJ1:H2	5:A:1425[B]:DMS:C1	2.17	0.56
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:345:TYR:CE2	5:A:1426[A]:DMS:H12	2.43	0.53
1:A:214[B]:SER:OG	1:A:400[B]:ARG:HG2	2.09	0.52
1:A:63[A]:MET:HG2	1:A:111[A]:HIS:ND1	2.25	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:63[B]:MET:HB3	1:A:102:CYS:SG	2.48	0.52
4:A:1424:JJ1:H2	5:A:1425[A]:DMS:H13	1.75	0.52
1:A:194[B]:ARG:CG	1:A:194[B]:ARG:NH1	2.39	0.51
1:A:63[B]:MET:HE3	1:A:72[B]:ILE:HD12	1.81	0.51
1:A:72[A]:ILE:HD12	1:A:122:LEU:HD21	1.93	0.51
4:A:1424:JJ1:H2	5:A:1425[A]:DMS:C1	2.21	0.51
1:A:70:HIS:HD2	6:A:2140:HOH:O	1.94	0.51
1:A:62[A]:ASN:OD1	1:A:62[A]:ASN:C	2.54	0.50
1:A:261:ASN:HD21	1:A:358:GLN:HE21	1.58	0.50
1:A:28[B]:GLU:OE1	1:A:28[B]:GLU:CA	2.58	0.50
1:A:261:ASN:ND2	1:A:358:GLN:HE21	2.09	0.50
1:A:176:ARG:O	1:A:177[B]:GLU:HB2	2.11	0.50
5:A:1426[A]:DMS:H11	6:A:2508:HOH:O	2.11	0.49
1:A:16:SER:HA	1:A:22:GLN:NE2	2.26	0.49
5:A:1426[A]:DMS:C1	6:A:2508:HOH:O	2.61	0.48
1:A:236:PRO:CB	1:A:238:GLN:OE1	2.62	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
1:A:207[B]:LEU:HD22	1:A:400[B]:ARG:NE	2.30	0.46
1:A:31:VAL:HA	1:A:141:MET:CE	2.39	0.46
1:A:194[B]:ARG:HD2	1:A:194[B]:ARG:HA	1.34	0.46
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
1:A:101[A]:LEU:HD22	1:A:127:ALA:HB1	1.99	0.44
1:A:63[A]:MET:HE3	1:A:111[A]:HIS:CG	2.53	0.44
1:A:217:TYR:CZ	5:A:1426[A]:DMS:H13	2.53	0.44
4:A:1424:JJ1:H101	5:A:1425[A]:DMS:C1	2.42	0.44
1:A:246[A]:MET:HB2	1:A:246[A]:MET:HE2	1.83	0.43
1:A:63[A]:MET:HE3	1:A:111[A]:HIS:ND1	2.33	0.43
1:A:62[A]:ASN:OD1	1:A:63[A]:MET:N	2.50	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
1:A:141:MET:HE3	1:A:141:MET:HB3	1.54	0.42
1:A:326:TYR:OH	5:A:1426[B]:DMS:H23	2.19	0.42
1:A:207[B]:LEU:HD13	1:A:207[B]:LEU:HA	1.83	0.42
1:A:231[A]:ARG:HD3	1:A:231[A]:ARG:HA	1.81	0.41
1:A:31:VAL:HB	1:A:141:MET:HE1	2.02	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
1:A:236:PRO:HB3	1:A:238:GLN:OE1	2.21	0.40
4:A:1424:JJ1:C10	5:A:1425[B]:DMS:H12	2.45	0.40
1:A:363:LEU:C	1:A:363:LEU:HD23	2.46	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	461/411 (112%)	452 (98%)	9 (2%)	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	413/362 (114%)	401 (97%)	12 (3%)	37	5

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

Mol	Chain	Res	Type
1	A	41	THR

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Mol	Chain	Res	Type
1	A	63[A]	MET
1	A	63[B]	MET
1	A	148	LYS
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	340	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	22	GLN
1	A	70	HIS
1	A	91	ASN
1	A	146	GLN
1	A	243	GLN
1	A	252	ASN
1	A	261	ASN
1	A	274	GLN
1	A	315	ASN
1	A	339	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 oligosaccharides in this entry.

5.6 Ligand geometry

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	1426[A]	-	3,3,3	0.57	0	3,3,3	1.33	1 (33%)
5	DMS	A	1426[B]	-	3,3,3	0.57	0	3,3,3	1.33	1 (33%)
5	DMS	A	1425[A]	-	3,3,3	1.30	0	3,3,3	1.33	1 (33%)
2	MYA	A	1422	3	63,65,65	1.42	7 (11%)	85,91,91	1.17	8 (9%)
4	JJ1	A	1424	-	36,36,36	0.98	2 (5%)	41,42,42	0.99	2 (4%)
5	DMS	A	1425[B]	-	3,3,3	1.01	0	3,3,3	3.51	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
2	MYA	A	1422	3	-	1/64/80/80	0/3/3/3
4	JJ1	A	1424	-	-	2/39/47/47	0/1/1/1

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1422	MYA	P1A-O3A	5.26	1.65	1.59
2	A	1422	MYA	P3X-O3X	3.91	1.66	1.59
2	A	1422	MYA	C8A-N7A	3.20	1.37	1.31
2	A	1422	MYA	O4X-C4X	2.92	1.51	1.45
2	A	1422	MYA	C3M-C2M	2.68	1.53	1.50
2	A	1422	MYA	C13-C11	2.67	1.59	1.53
4	A	1424	JJ1	C8-C7	2.54	1.60	1.53
2	A	1422	MYA	P1A-O2A	-2.08	1.43	1.50
4	A	1424	JJ1	C1-N1	-2.01	1.41	1.45

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	1425[B]	DMS	O-S-C2	-4.42	84.37	106.49
2	A	1422	MYA	C4A-N9A-C8A	3.17	109.07	105.74
2	A	1422	MYA	C3M-C2M-S1	3.13	117.13	113.40
2	A	1422	MYA	O6A-C12-C11	-2.99	105.75	110.55
5	A	1425[B]	DMS	O-S-C1	-2.97	91.61	106.49
5	A	1425[B]	DMS	C2-S-C1	2.93	113.43	98.42
4	A	1424	JJ1	C14-C13-C12	2.91	129.06	114.37
2	A	1422	MYA	O2M-C2M-C3M	-2.87	120.67	123.98
2	A	1422	MYA	C4M-C3M-C2M	-2.77	106.16	112.27
2	A	1422	MYA	C13-C11-C10	2.64	113.27	108.77
2	A	1422	MYA	N9A-C8A-N7A	-2.63	110.21	113.94
5	A	1426[B]	DMS	C2-S-C1	2.15	109.46	98.42
5	A	1426[A]	DMS	C2-S-C1	2.15	109.44	98.42
5	A	1425[A]	DMS	C2-S-C1	2.06	108.98	98.42
2	A	1422	MYA	O9-C9-N8	-2.02	118.71	122.98
4	A	1424	JJ1	C17-C16-C15	-2.00	104.25	114.37

There are no chirality outliers.

All (3) torsion outliers are listed below:

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

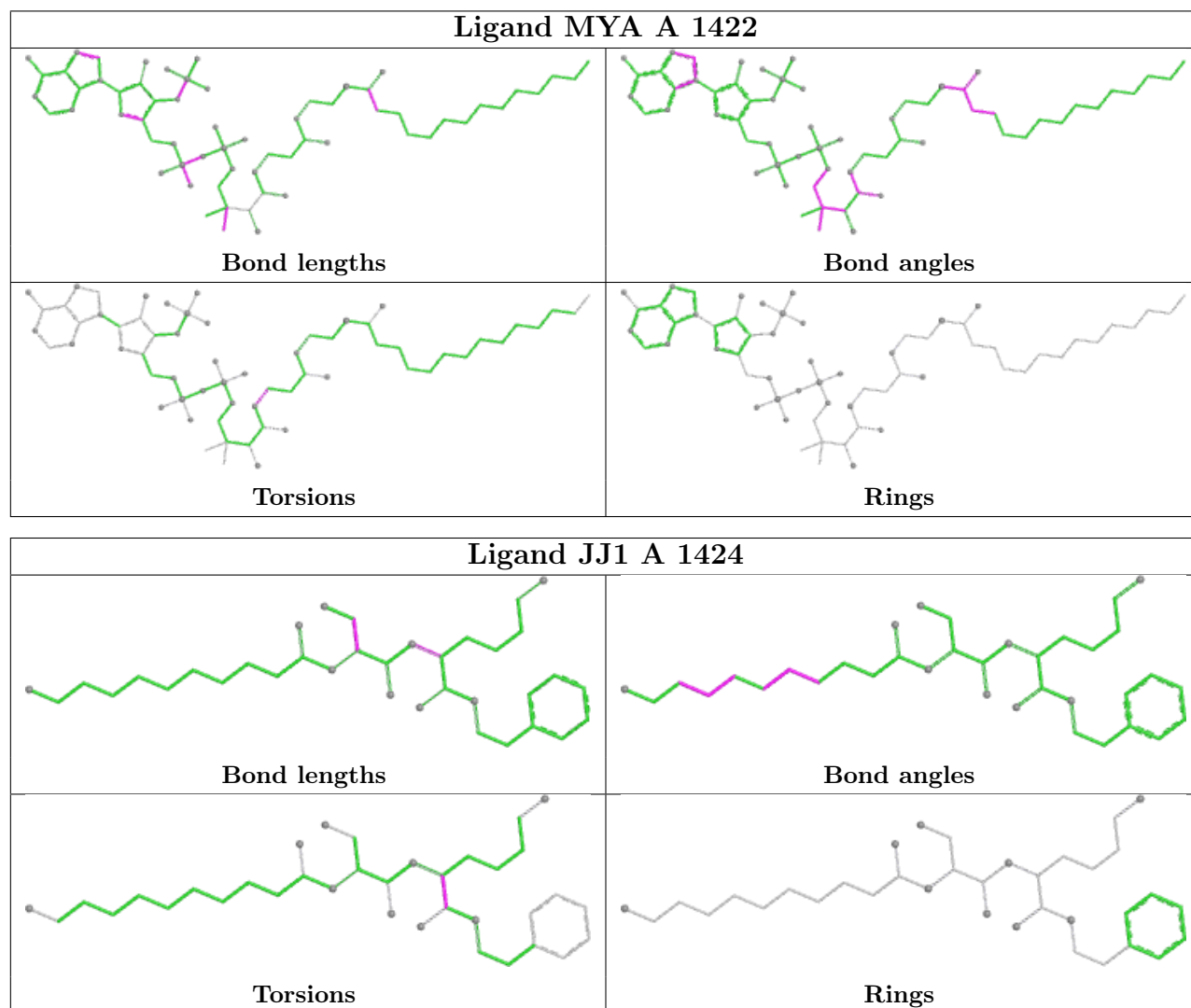
There are no ring outliers.

5 monomers are involved in 21 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	1426[A]	DMS	6	0
5	A	1426[B]	DMS	5	0
5	A	1425[A]	DMS	4	0
4	A	1424	JJ1	8	0
5	A	1425[B]	DMS	6	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

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(Å ²)	Q < 0.9
1	A	411/411 (100%)	0.09	22 (5%)	31 36	5, 12, 27, 42	51 (12%)

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	242	PHE	5.0
1	A	334	GLY	4.1
1	A	245	PRO	3.9
1	A	11	ALA	3.9
1	A	336	SER	3.4
1	A	238	GLN	3.2
1	A	239	TYR	3.1
1	A	333	ILE	3.1
1	A	236	PRO	3.0
1	A	235	ILE	3.0
1	A	241	LYS	2.9
1	A	247	ALA	2.8
1	A	118	ALA	2.8
1	A	244	ASN	2.7
1	A	262[A]	SER	2.6
1	A	246[A]	MET	2.6
1	A	31	VAL	2.5
1	A	28[A]	GLU	2.4
1	A	240	GLN	2.3
1	A	337	ASN	2.2
1	A	32	PHE	2.1
1	A	39	PRO	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 oligosaccharides in this entry.

6.4 Ligands

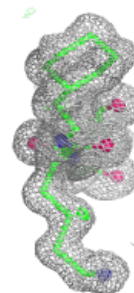
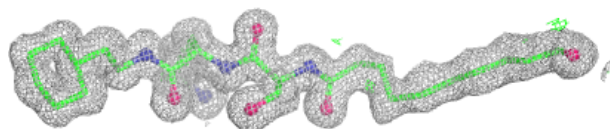
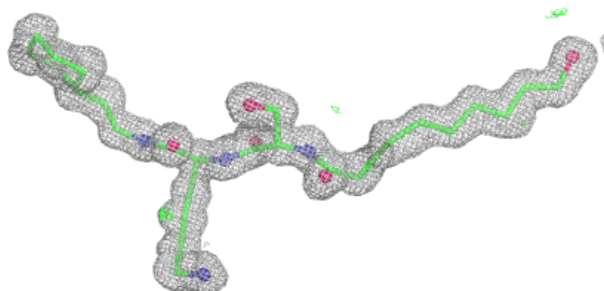
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	DMS	A	1425[A]	4/4	0.86	0.14	13,15,17,19	4
5	DMS	A	1425[B]	4/4	0.86	0.14	14,15,20,20	4
5	DMS	A	1426[A]	4/4	0.88	0.14	9,15,17,22	4
5	DMS	A	1426[B]	4/4	0.88	0.14	10,10,12,13	4
4	JJ1	A	1424	36/36	0.97	0.06	8,12,21,21	0
3	MG	A	1423	1/1	0.98	0.05	19,19,19,19	0
2	MYA	A	1422	63/63	0.99	0.04	7,10,12,14	0

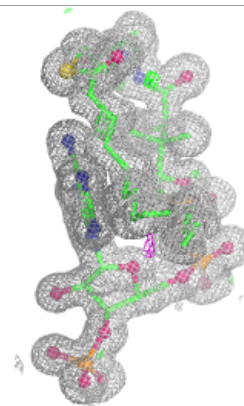
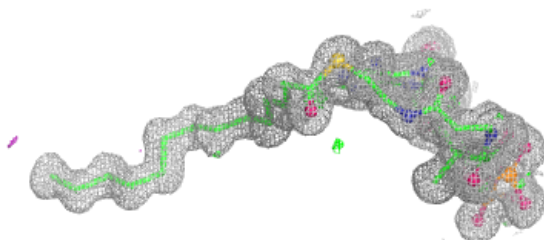
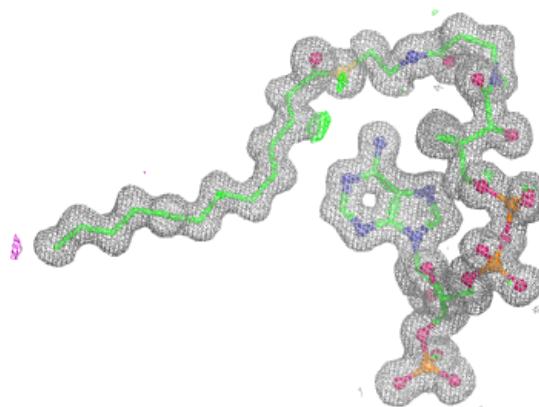
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

There are no such residues in this entry.