Full wwPDB X-ray Structure Validation Report (i)

Dec 1, 2022 – 12:50 pm GMT

PDB ID	:	6SK2
Title	:	HsNMT1 in complex with both MyrCoA and Acetylated-GKSFSKPR peptide reveals N-terminal Lysine Myristoylation
Authors	:	Dian, C.; Riviere, F.B.; Asensio, T.; Giglione, C.; Meinnel, T.
Deposited on	:	2019-08-14
Resolution	:	1.90 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 (i) 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 (i)) were used in the production of this report:

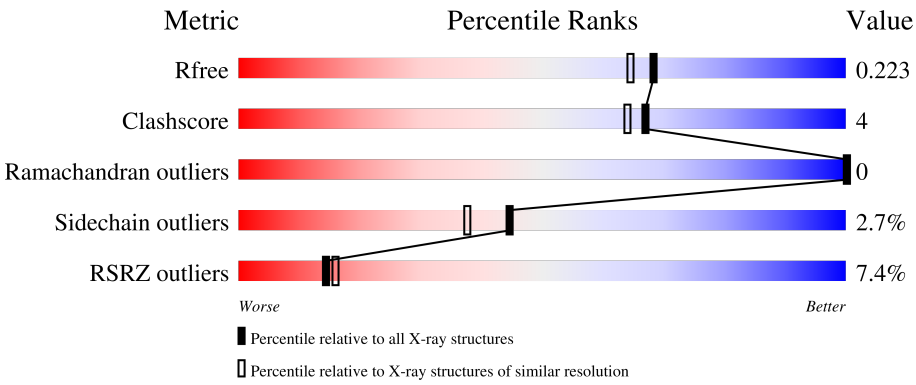
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.31.3
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0267
CCP4	:	7.1.010 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
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 1.90 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	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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 <=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	5%, 88%, 9%, ·
1	B	402	7%, 89%, 8%, ·
2	D	9	33%, 67%, 22%, 11%
2	F	9	89%, 44%, 22%, 33%

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
3	GOL	A	502	-	-	X	-

2 Entry composition (i)

There are 7 unique types of molecules in this entry. The entry contains 7411 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
			Total	C	N	O	S			
1	A	392	3223	2087	546	574	16	0	7	0
1	B	390	3199	2069	535	578	17	0	7	0

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 Apoptosis-inducing factor 3.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	D	9	61	39	10	12	0	0	0
2	F	9	57	36	9	12	0	0	0

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	101	ACE	-	acetylation	UNP Q96NN9
D	2	GLY	-	engineered mutation	UNP Q96NN9
D	3	LYS	-	engineered mutation	UNP Q96NN9
D	4	SER	-	engineered mutation	UNP Q96NN9

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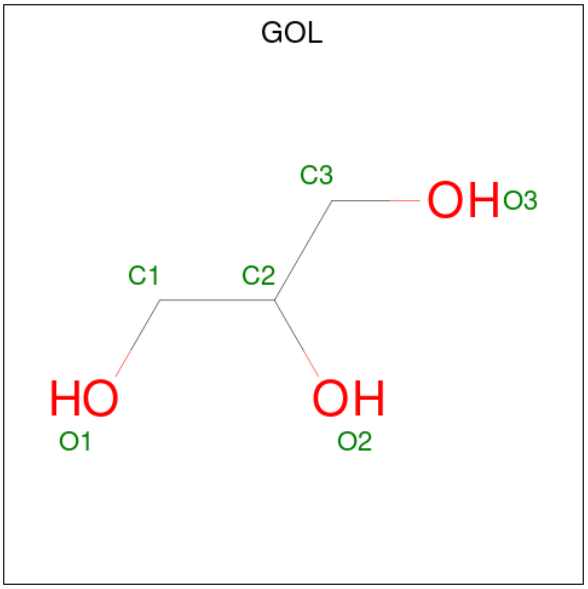
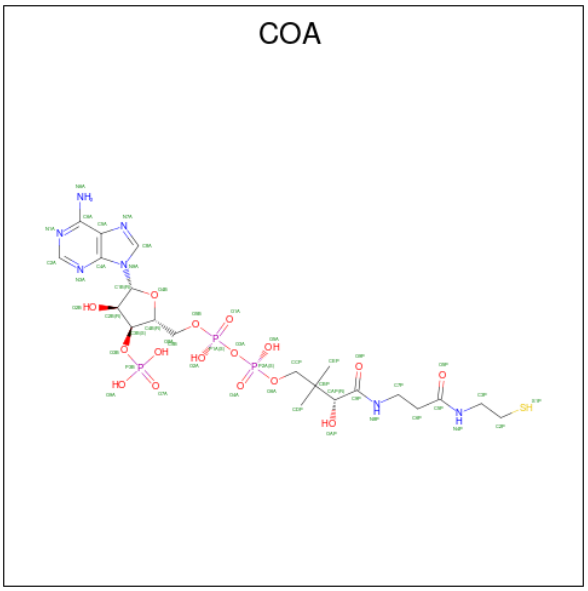
Chain	Residue	Modelled	Actual	Comment	Reference
D	9	ARG	-	engineered mutation	UNP Q96NN9
F	101	ACE	-	acetylation	UNP Q96NN9
F	2	GLY	-	engineered mutation	UNP Q96NN9
F	3	LYS	-	engineered mutation	UNP Q96NN9
F	4	SER	-	engineered mutation	UNP Q96NN9
F	9	ARG	-	engineered mutation	UNP Q96NN9

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



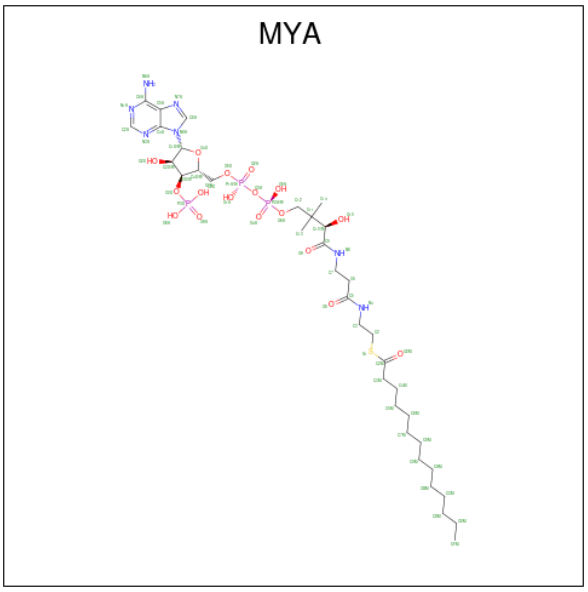
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
3	A	1	6	3	3	0	0
			Total	C	O		
3	A	1	6	3	3	0	0
			Total	C	O		
3	B	1	6	3	3	0	0

- Molecule 4 is COENZYME A (three-letter code: COA) (formula: $C_{21}H_{36}N_7O_{16}P_3S$).



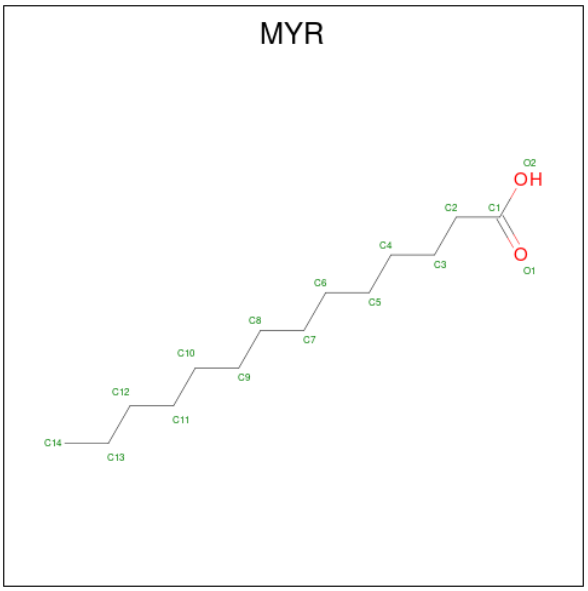
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
			Total	C	N	O	P	S		
4	A	1	48	21	7	16	3	1	0	0

- Molecule 5 is TETRADECANOYL-COA (three-letter code: MYA) (formula: $C_{35}H_{62}N_7O_{17}P_3S$) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 6 is MYRISTIC ACID (three-letter code: MYR) (formula: $C_{14}H_{28}O_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	D	1	Total C O 15 14 1	0	0

- Molecule 7 is water.

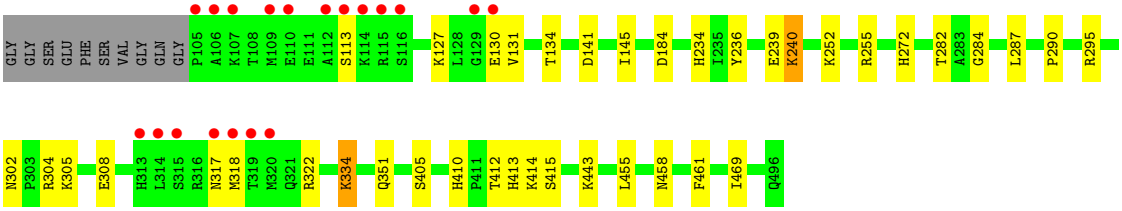
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	370	Total O 370 370	0	2
7	B	354	Total O 354 354	0	0
7	D	2	Total O 2 2	0	0
7	F	1	Total O 1 1	0	0

3 Residue-property plots (i)

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.

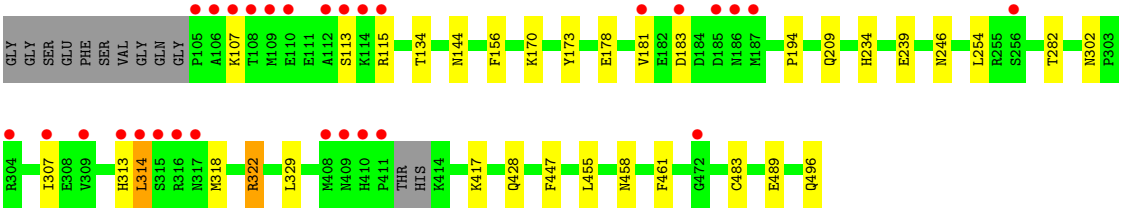
• Molecule 1: Glycylpeptide N-tetradecanoyltransferase 1

Chain A:



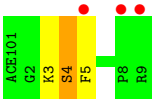
• Molecule 1: Glycylpeptide N-tetradecanoyltransferase 1

Chain B:



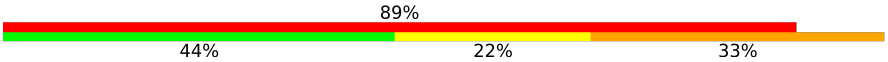
• Molecule 2: Apoptosis-inducing factor 3

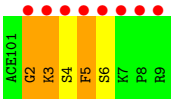
Chain D:



• Molecule 2: Apoptosis-inducing factor 3

Chain F:





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α, β, γ	78.82Å 178.44Å 58.27Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.79 – 1.90 48.79 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.9 (48.79-1.90) 99.9 (48.79-1.90)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.02 (at 1.90Å)	Xtriage
Refinement program	PHENIX 1.10.1_2155, PHENIX 1.10.1_2155	Depositor
R, R_{free}	0.172 , 0.220 0.175 , 0.223	Depositor DCC
R_{free} test set	3256 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å^2)	27.8	Xtriage
Anisotropy	0.742	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	(Not available) , (Not available)	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	7411	wwPDB-VP
Average B, all atoms (Å^2)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 16.72% 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 (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, ACE, COA, MYA, MYR

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	0.36	0/3314	0.52	0/4503
1	B	0.36	0/3288	0.52	0/4473
2	D	0.51	0/60	0.72	0/77
2	F	0.52	0/56	0.65	0/73
All	All	0.36	0/6718	0.53	0/9126

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
2	F	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	F	2	GLY	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	3223	0	3174	27	0
1	B	3199	0	3124	22	0
2	D	61	0	58	3	0
2	F	57	0	49	8	0
3	A	12	0	16	5	1
3	B	6	0	8	0	0
4	A	48	0	32	1	0
5	B	63	0	58	2	0
6	D	15	0	27	0	0
7	A	370	0	0	5	0
7	B	354	0	0	5	1
7	D	2	0	0	0	0
7	F	1	0	0	0	0
All	All	7411	0	6546	52	2

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 4.

All (52) 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)
4:A:503:COA:O4B	4:A:503:COA:C1B	1.67	1.15
1:A:405:SER:HG	2:D:5:PHE:HE1	1.06	0.98
1:A:302:ASN:OD1	1:A:443:LYS:NZ	2.14	0.79
1:B:282:THR:OG1	2:F:3:LYS:HG2	1.84	0.77
1:A:414:LYS:H	3:A:502:GOL:H32	1.52	0.75
5:B:501:MYA:S1	2:F:3:LYS:HB3	2.35	0.67
1:A:141:ASP:OD2	7:A:601:HOH:O	2.13	0.67
1:B:173[B]:TYR:CE2	1:B:194:PRO:HG3	2.30	0.66
1:B:170:LYS:HA	1:B:173[B]:TYR:CE2	2.30	0.66
1:A:410:HIS:CD2	1:A:413:HIS:H	2.16	0.64
1:A:304:ARG:O	1:A:308:GLU:HG3	1.98	0.64
1:A:239:GLU:OE2	7:A:602[A]:HOH:O	2.15	0.64
1:A:351:GLN:NE2	1:B:144[B]:ASN:OD1	2.32	0.62
1:B:115:ARG:NH2	5:B:501:MYA:O8A	2.35	0.59
1:B:417:LYS:HG2	1:B:447:PHE:HA	1.84	0.58
1:A:318:MET:HB2	1:A:322:ARG:HG2	1.86	0.57
1:B:183:ASP:HA	2:F:5:PHE:HB3	1.87	0.57
1:A:240:LYS:NZ	7:A:604:HOH:O	2.34	0.55
1:B:458:ASN:HA	1:B:461:PHE:CE2	2.42	0.55
1:A:410:HIS:HD2	1:A:412:THR:H	1.56	0.54
1:B:178:GLU:OE2	7:B:602:HOH:O	2.19	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:318:MET:HG2	1:B:322:ARG:HG2	1.91	0.53
1:B:307:ILE:HD13	1:B:314:LEU:HD12	1.90	0.52
1:B:489:GLU:OE1	7:B:601:HOH:O	2.18	0.52
2:F:2:GLY:C	2:F:4:SER:H	2.12	0.52
1:A:414:LYS:H	3:A:502:GOL:C3	2.23	0.51
1:A:458:ASN:HA	1:A:461:PHE:CE2	2.47	0.49
1:B:156:PHE:CE1	1:B:254:LEU:HD11	2.48	0.49
1:B:496:GLN:OXT	2:F:3:LYS:HE2	2.14	0.48
1:A:252:LYS:HA	1:A:255:ARG:HG3	1.96	0.48
1:A:305:LYS:NZ	3:A:502:GOL:H11	2.31	0.46
1:A:334:LYS:NZ	7:A:612:HOH:O	2.44	0.46
2:F:5:PHE:HD2	2:F:6:SER:N	2.13	0.46
1:B:134:THR:O	1:B:483:CYS:HA	2.17	0.44
1:B:170:LYS:HA	1:B:173[B]:TYR:CD2	2.51	0.44
1:A:234[A]:HIS:NE2	1:A:236:TYR:O	2.51	0.44
1:B:181:VAL:HG12	2:F:5:PHE:HB2	2.00	0.44
1:B:246:ASN:OD1	2:F:3:LYS:HD2	2.17	0.44
1:B:428:GLN:NE2	7:B:617:HOH:O	2.48	0.43
1:A:127:LYS:H	1:A:127:LYS:HG3	1.62	0.43
1:A:234[B]:HIS:ND1	7:A:606[B]:HOH:O	2.36	0.43
1:B:209:GLN:HG3	7:B:736:HOH:O	2.19	0.43
1:A:295:ARG:HE	1:A:469:ILE:HD11	1.84	0.43
1:A:305:LYS:HZ2	3:A:502:GOL:H11	1.83	0.42
1:A:284:GLY:HA3	2:D:4:SER:HB2	2.01	0.42
1:B:234:HIS:ND1	1:B:239[B]:GLU:HG3	2.35	0.41
1:A:145:ILE:HD12	1:A:272:HIS:HB3	2.03	0.41
1:A:410:HIS:HD2	1:A:413:HIS:H	1.66	0.41
1:A:415:SER:H	3:A:502:GOL:H2	1.86	0.41
1:A:282:THR:O	2:D:3:LYS:HG3	2.21	0.40
1:A:287:LEU:O	1:A:290:PRO:HD3	2.22	0.40
1:B:329:LEU:O	7:B:603:HOH:O	2.22	0.40

All (2) 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)
3:A:502:GOL:O3	3:A:502:GOL:O3[2_555]	2.14	0.06
7:B:740:HOH:O	7:B:896:HOH:O[4_457]	2.18	0.02

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	397/402 (99%)	389 (98%)	8 (2%)	0	100	100
1	B	393/402 (98%)	385 (98%)	8 (2%)	0	100	100
2	D	7/9 (78%)	6 (86%)	1 (14%)	0	100	100
2	F	7/9 (78%)	6 (86%)	1 (14%)	0	100	100
All	All	804/822 (98%)	786 (98%)	18 (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	349/362 (96%)	340 (97%)	9 (3%)	46	39
1	B	348/362 (96%)	341 (98%)	7 (2%)	55	51
2	D	6/7 (86%)	5 (83%)	1 (17%)	2	0
2	F	5/7 (71%)	3 (60%)	2 (40%)	0	0
All	All	708/738 (96%)	689 (97%)	19 (3%)	44	38

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	113	SER
1	A	130	GLU

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Mol	Chain	Res	Type
1	A	131	VAL
1	A	134	THR
1	A	184	ASP
1	A	240	LYS
1	A	317	ASN
1	A	334	LYS
1	A	455	LEU
1	B	107	LYS
1	B	113	SER
1	B	302	ASN
1	B	313	HIS
1	B	314	LEU
1	B	322	ARG
1	B	455	LEU
2	D	4	SER
2	F	3	LYS
2	F	5	PHE

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	351	GLN
1	A	391	ASN
1	A	410	HIS
1	B	302	ASN
1	B	351	GLN

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

6 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	$\#\|Z\| > 2$
5	MYA	B	501	-	50,65,65	1.29	5 (10%)	65,91,91	1.44	9 (13%)
3	GOL	B	502	-	5,5,5	0.41	0	5,5,5	0.33	0
4	COA	A	503	-	41,50,50	4.33	13 (31%)	52,75,75	1.99	8 (15%)
3	GOL	A	502	-	5,5,5	0.36	0	5,5,5	0.50	0
6	MYR	D	201	2	14,14,15	0.46	0	13,13,15	0.94	1 (7%)
3	GOL	A	501	-	5,5,5	0.35	0	5,5,5	0.25	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
5	MYA	B	501	-	-	0/44/80/80	0/3/3/3
3	GOL	B	502	-	-	0/4/4/4	-
4	COA	A	503	-	-	9/44/64/64	0/3/3/3
3	GOL	A	502	-	-	0/4/4/4	-
6	MYR	D	201	2	-	0/11/12/13	-
3	GOL	A	501	-	-	0/4/4/4	-

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	503	COA	O4B-C1B	18.74	1.67	1.41
4	A	503	COA	C2B-C1B	-13.81	1.32	1.53
4	A	503	COA	C9P-N8P	6.55	1.47	1.33
4	A	503	COA	C5P-N4P	6.25	1.47	1.33
4	A	503	COA	O4B-C4B	-6.13	1.31	1.45
4	A	503	COA	C6A-N6A	4.81	1.51	1.34
4	A	503	COA	P3B-O3B	4.41	1.67	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	501	MYA	C5A-C6A	-4.29	1.34	1.43
5	B	501	MYA	C6A-N1A	-2.98	1.33	1.37
5	B	501	MYA	C2A-N1A	2.78	1.40	1.33
4	A	503	COA	C2A-N3A	2.71	1.36	1.32
4	A	503	COA	O3B-C3B	-2.65	1.34	1.44
5	B	501	MYA	P2A-O5A	-2.65	1.48	1.56
5	B	501	MYA	C5A-N7A	-2.56	1.34	1.39
4	A	503	COA	P2A-O6A	2.28	1.68	1.59
4	A	503	COA	C3B-C4B	2.27	1.59	1.52
4	A	503	COA	O9P-C9P	-2.21	1.19	1.23
4	A	503	COA	C7P-N8P	2.05	1.50	1.46

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	A	503	COA	C5A-C6A-N6A	8.64	133.48	120.35
4	A	503	COA	N6A-C6A-N1A	-5.66	106.83	118.57
4	A	503	COA	N3A-C2A-N1A	-5.28	120.42	128.68
5	B	501	MYA	N3A-C2A-N1A	-5.25	118.88	128.82
4	A	503	COA	P2A-O3A-P1A	-3.85	119.63	132.83
5	B	501	MYA	C5A-C4A-N3A	-3.50	120.85	127.41
4	A	503	COA	C3P-N4P-C5P	-3.45	116.43	122.84
5	B	501	MYA	O5A-P2A-O4A	3.37	119.14	108.50
5	B	501	MYA	C2X-C3X-C4X	-3.05	97.81	103.22
4	A	503	COA	O6A-CCP-CBP	-2.91	105.88	110.55
5	B	501	MYA	N9A-C8A-N7A	-2.73	108.25	113.39
5	B	501	MYA	N6A-C6A-N1A	2.71	120.80	117.01
4	A	503	COA	C6P-C5P-N4P	2.29	120.27	116.42
5	B	501	MYA	C8A-N7A-C5A	2.20	108.23	104.24
4	A	503	COA	O5P-C5P-N4P	-2.18	118.90	123.01
6	D	201	MYR	C3-C2-C1	-2.09	104.98	114.26
5	B	501	MYA	O7A-P3X-O8A	2.03	114.26	105.60
5	B	501	MYA	C5A-C6A-N6A	-2.00	117.78	121.61

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	503	COA	C5B-O5B-P1A-O1A
4	A	503	COA	C5B-O5B-P1A-O2A
4	A	503	COA	C5B-O5B-P1A-O3A
4	A	503	COA	C3B-C4B-C5B-O5B

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Mol	Chain	Res	Type	Atoms
4	A	503	COA	O4B-C4B-C5B-O5B
4	A	503	COA	P1A-O3A-P2A-O4A
4	A	503	COA	C3B-O3B-P3B-O8A
4	A	503	COA	P1A-O3A-P2A-O5A
4	A	503	COA	O5P-C5P-C6P-C7P

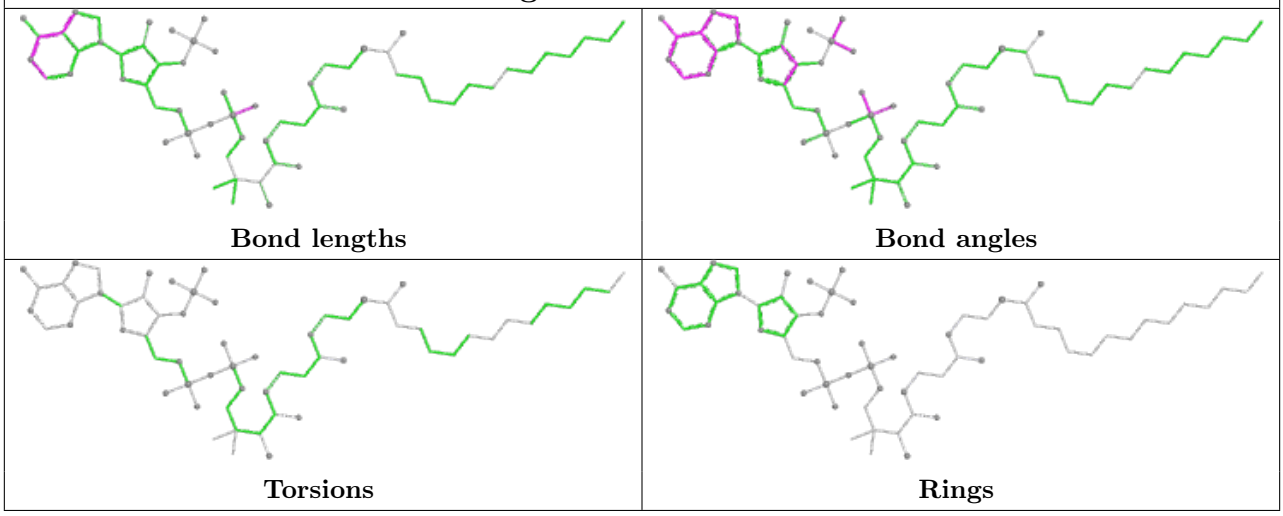
There are no ring outliers.

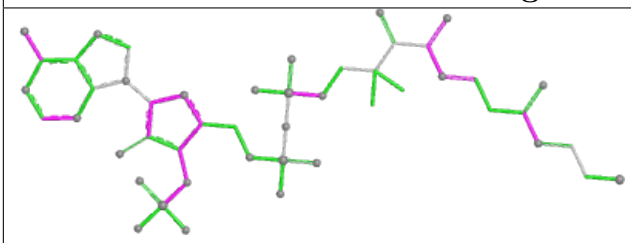
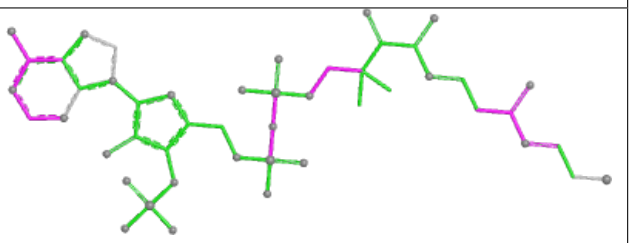
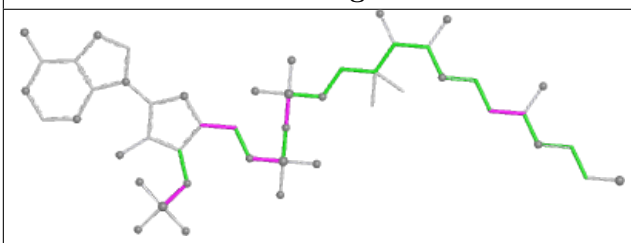
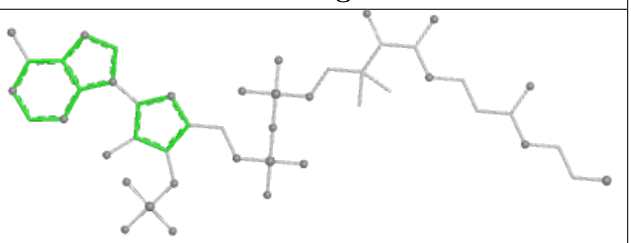
3 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	501	MYA	2	0
4	A	503	COA	1	0
3	A	502	GOL	5	1

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



Ligand COA A 503	
 Bond lengths	 Bond angles
 Torsions	 Rings

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	392/402 (97%)	0.04	19 (4%)	30	33	18, 27, 60, 82	0
1	B	390/402 (97%)	0.20	29 (7%)	14	16	16, 26, 61, 87	0
2	D	8/9 (88%)	2.35	3 (37%)	0	0	41, 46, 51, 63	0
2	F	8/9 (88%)	5.09	8 (100%)	0	0	53, 58, 63, 71	0
All	All	798/822 (97%)	0.19	59 (7%)	14	16	16, 27, 62, 87	0

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	F	5	PHE	9.6
1	B	106	ALA	8.1
1	B	108	THR	7.7
1	B	109	MET	6.2
1	A	114	LYS	5.9
2	F	8	PRO	5.9
2	F	4	SER	5.1
2	F	3	LYS	5.1
2	F	6	SER	4.8
1	B	410	HIS	4.8
2	D	5	PHE	4.7
1	B	110	GLU	4.5
2	F	2	GLY	4.4
1	B	107	LYS	4.4
1	B	105	PRO	4.2
1	A	317	ASN	3.8
1	A	110	GLU	3.7
1	A	112	ALA	3.7
2	D	8	PRO	3.7
2	F	9	ARG	3.7
1	B	408	MET	3.6

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Mol	Chain	Res	Type	RSRZ
1	A	318	MET	3.6
1	A	106	ALA	3.6
1	A	116	SER	3.5
1	B	316	ARG	3.4
1	B	187	MET	3.4
1	B	315	SER	3.3
1	A	129	GLY	3.2
1	B	115	ARG	3.2
1	B	472	GLY	3.2
1	B	186	ASN	3.1
1	A	313	HIS	3.1
1	B	409	ASN	3.1
1	B	314	LEU	3.1
1	B	411	PRO	3.0
2	D	9	ARG	3.0
1	A	115	ARG	3.0
1	B	304	ARG	2.6
1	A	105	PRO	2.6
1	B	181	VAL	2.5
1	B	185	ASP	2.5
1	B	317	ASN	2.5
1	A	113	SER	2.4
1	B	309	VAL	2.4
1	A	130	GLU	2.4
1	A	315	SER	2.3
1	B	112	ALA	2.3
1	A	319	THR	2.2
1	B	307	ILE	2.1
1	B	183	ASP	2.1
1	B	114	LYS	2.1
1	B	113	SER	2.1
1	A	109	MET	2.1
1	A	314	LEU	2.1
1	B	256	SER	2.1
2	F	7	LYS	2.1
1	A	320	MET	2.0
1	A	107	LYS	2.0
1	B	313	HIS	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)

There are no monosaccharides in this entry.

6.4 Ligands (i)

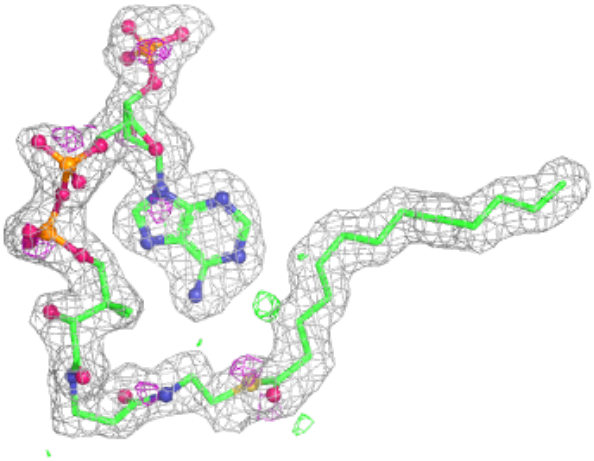
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

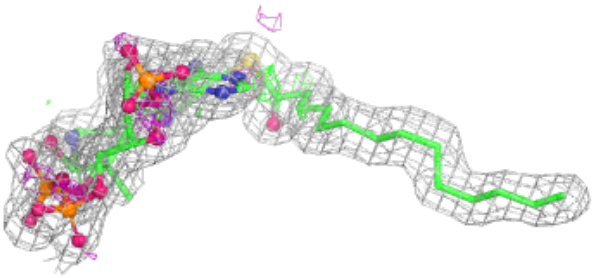
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	GOL	A	502	6/6	0.79	0.29	40,51,52,58	0
5	MYA	B	501	63/63	0.86	0.18	11,38,46,51	0
4	COA	A	503	48/48	0.90	0.15	26,41,53,72	0
6	MYR	D	201	15/16	0.95	0.14	15,20,27,35	0
3	GOL	B	502	6/6	0.96	0.13	21,24,32,34	0
3	GOL	A	501	6/6	0.98	0.07	23,30,33,34	0

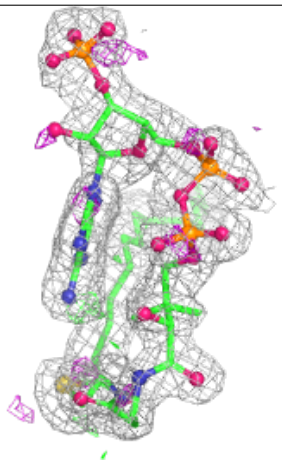
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around MYA B 501:

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

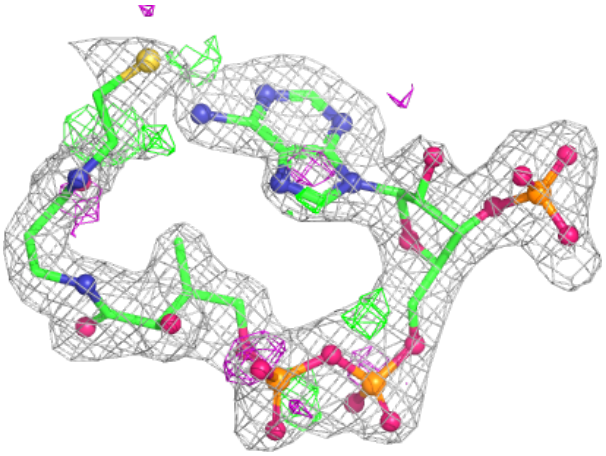


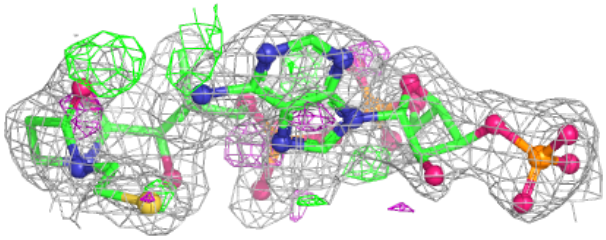


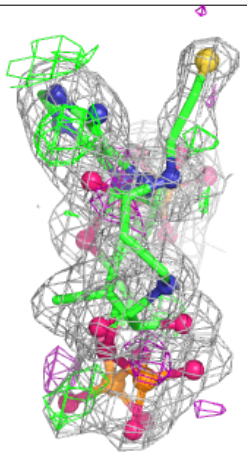


Electron density around COA A 503:

$2mF_o$-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.