



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 3, 2016 – 02:07 PM EDT

PDB ID : 5L6D
Title : Crystal structure of the human METTL3-METTL14 complex bound to SAH
Authors : Sledz, P.; Jinek, M.
Deposited on : 2016-05-29
Resolution : 1.85 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

<http://wwpdb.org/validation/2016/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	:	unknown
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20027939
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20027939

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	MG	A	602	-	-	-	X

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 7523 atoms, of which 3522 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.

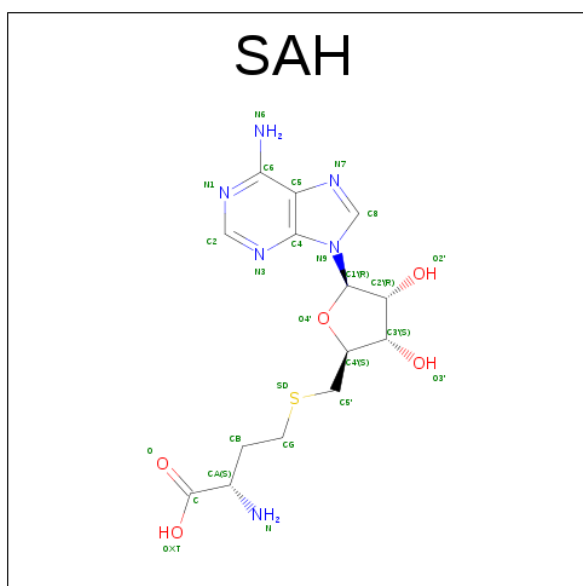
- Molecule 1 is a protein called N6-adenosine-methyltransferase 70 kDa subunit.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	199	3209	1036	1596	281	286	10	0	1	0

- Molecule 2 is a protein called N6-adenosine-methyltransferase subunit METTL14.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
2	B	238	3862	1250	1904	335	360	13	0	2	0

- Molecule 3 is S-ADENOSYL-L-HOMOCYSTEINE (three-letter code: SAH) (formula: $C_{14}H_{20}N_6O_5S$).

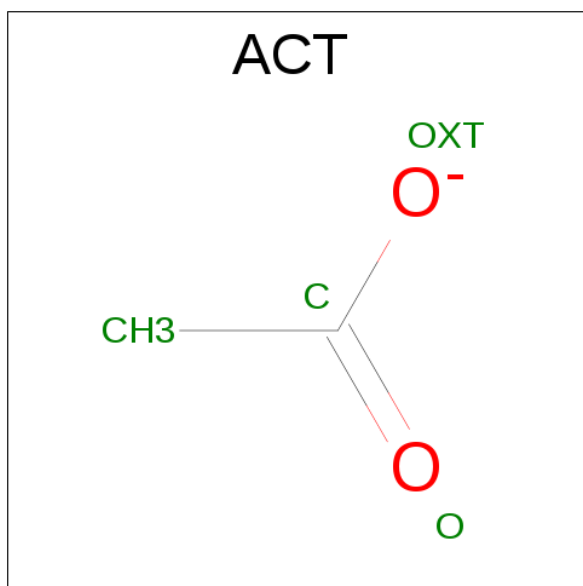


Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
			Total	C	H	N	O	S		
3	A	1	45	14	19	6	5	1	0	0

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 5 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	B	1	Total	C	H	O	0	0
			7	2	3	2		

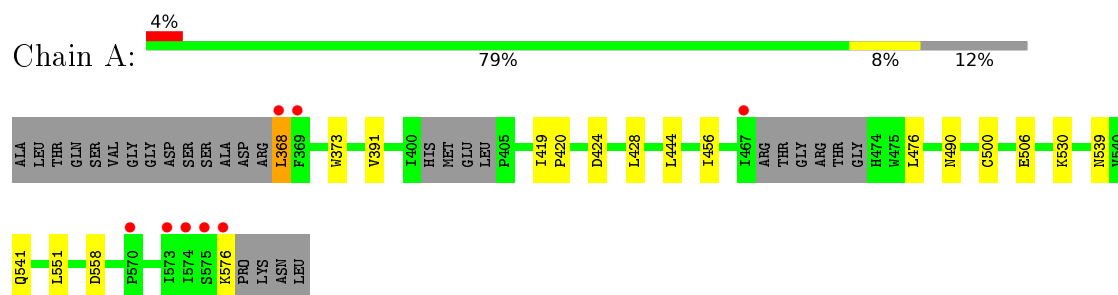
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	200	Total	O	0	0
			200	200		
6	B	199	Total	O	0	0
			199	199		

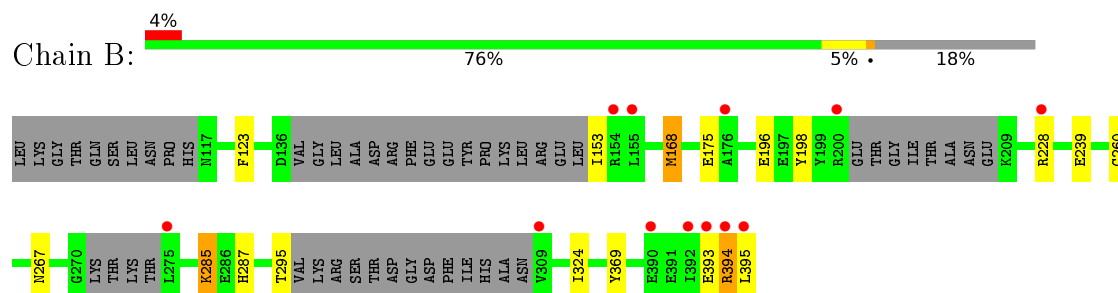
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: N6-adenosine-methyltransferase 70 kDa subunit



- Molecule 2: N6-adenosine-methyltransferase subunit METTL14



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	63.82Å 63.82Å 225.63Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.63 – 1.85 49.63 – 1.85	Depositor EDS
% Data completeness (in resolution range)	99.9 (49.63-1.85) 99.9 (49.63-1.85)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.36 (at 1.86Å)	Xtriage
Refinement program	PHENIX (1.10_2155: ???)	Depositor
R, R_{free}	0.157 , 0.192 0.160 , 0.195	Depositor DCC
R_{free} test set	2320 reflections (4.99%)	DCC
Wilson B-factor (Å ²)	24.5	Xtriage
Anisotropy	0.161	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 50.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.034 for -h,-k,l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7523	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.21% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MG, SAH, ACT

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.52	0/1657	0.70	2/2250 (0.1%)
2	B	0.52	0/2021	0.63	0/2736
All	All	0.52	0/3678	0.67	2/4986 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	368	LEU	CB-CG-CD2	-7.11	98.91	111.00
1	A	368	LEU	CA-CB-CG	-5.01	103.77	115.30

There are no chirality outliers.

There are no planarity outliers.

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	1613	1596	1596	15	0
2	B	1958	1904	1898	19	0
3	A	26	19	19	0	0
4	A	1	0	0	0	0
5	B	4	3	3	0	0
6	A	200	0	0	4	2
6	B	199	0	0	7	1

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	4001	3522	3516	32	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 (32) 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)
2:B:153:ILE:N	6:B:501:HOH:O	2.01	0.93
1:A:506:GLU:OE1	6:A:701:HOH:O	1.95	0.82
2:B:394:ARG:NH1	6:B:506:HOH:O	2.23	0.70
1:A:576:LYS:HB3	6:A:711:HOH:O	1.91	0.69
2:B:196:GLU:OE1	6:B:502:HOH:O	2.10	0.68
1:A:424:ASP:OD2	6:A:702:HOH:O	2.12	0.67
2:B:393:GLU:O	6:B:503:HOH:O	2.13	0.66
1:A:428:LEU:HD23	1:A:444:LEU:CD1	2.25	0.66
1:A:368:LEU:HD22	1:A:558:ASP:HA	1.78	0.65
1:A:500[B]:CYS:SG	2:B:123:PHE:CZ	2.95	0.59
1:A:428:LEU:HD23	1:A:444:LEU:HD12	1.85	0.58
1:A:490:ASN:HB2	6:A:738:HOH:O	2.05	0.56
1:A:428:LEU:HD23	1:A:444:LEU:HD11	1.90	0.53
2:B:260[B]:CYS:SG	2:B:287:HIS:CE1	3.03	0.51
2:B:168:MET:HE3	2:B:369:TYR:HA	1.92	0.51
2:B:239:GLU:OE2	6:B:504:HOH:O	2.20	0.50
2:B:153:ILE:CA	6:B:501:HOH:O	2.53	0.49
1:A:391:VAL:HB	1:A:530:LYS:HG2	1.94	0.49
1:A:456:ILE:HG12	2:B:285:LYS:HE2	1.96	0.47
2:B:228:ARG:NE	2:B:295:THR:HB	2.29	0.47
2:B:395:LEU:HD23	2:B:395:LEU:N	2.30	0.47
2:B:267:ASN:ND2	6:B:505:HOH:O	2.22	0.46
2:B:324:ILE:HG22	2:B:324:ILE:O	2.15	0.46
2:B:394:ARG:HG2	2:B:394:ARG:HH11	1.81	0.46
2:B:198:TYR:HB3	2:B:324:ILE:HD12	1.98	0.45
1:A:368:LEU:HD22	1:A:558:ASP:CA	2.46	0.45
1:A:373:TRP:HB2	1:A:551:LEU:HD13	1.99	0.45
2:B:175:GLU:N	2:B:175:GLU:OE2	2.42	0.45
2:B:228:ARG:CZ	2:B:295:THR:HB	2.48	0.44
1:A:419:ILE:N	1:A:420:PRO:CD	2.81	0.44
1:A:539:ASN:O	1:A:541:GLN:HG2	2.19	0.43
2:B:198:TYR:HB3	2:B:324:ILE:HG23	1.99	0.42

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)
6:A:717:HOH:O	6:A:847:HOH:O[5_445]	1.95	0.25
6:A:863:HOH:O	6:B:596:HOH:O[4_545]	2.08	0.12

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	194/227 (86%)	192 (99%)	2 (1%)	0	100	100
2	B	230/289 (80%)	227 (99%)	3 (1%)	0	100	100
All	All	424/516 (82%)	419 (99%)	5 (1%)	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	177/199 (89%)	176 (99%)	1 (1%)	90	87
2	B	216/258 (84%)	213 (99%)	3 (1%)	74	63
All	All	393/457 (86%)	389 (99%)	4 (1%)	82	76

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	476	LEU
2	B	168	MET
2	B	285	LYS
2	B	394	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 3 ligands modelled in this entry, 1 is monoatomic - leaving 2 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	$\# Z > 2$
3	SAH	A	601	-	22,28,28	0.79	1 (4%)	18,40,40	3.19	3 (16%)
5	ACT	B	401	-	0,3,3	0.00	-	0,3,3	0.00	-

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	SAH	A	601	-	-	0/7/31/31	0/3/3/3
5	ACT	B	401	-	-	0/0/0/0	0/0/0/0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	601	SAH	C2-N1	2.17	1.38	1.33

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	601	SAH	N3-C2-N1	-12.54	119.02	128.87
3	A	601	SAH	C1'-N9-C4	-2.76	123.73	126.81
3	A	601	SAH	O4'-C1'-N9	2.04	111.96	108.11

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	199/227 (87%)	-0.17	8 (4%) 42 39	16, 24, 48, 72	0
2	B	238/289 (82%)	-0.10	12 (5%) 32 31	15, 26, 58, 79	0
All	All	437/516 (84%)	-0.13	20 (4%) 36 34	15, 25, 54, 79	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	155	LEU	7.2
1	A	573	ILE	6.7
2	B	154	ARG	6.2
1	A	368	LEU	5.7
1	A	574	ILE	5.4
1	A	576	LYS	4.1
2	B	395	LEU	4.0
2	B	200	ARG	3.4
2	B	309	VAL	3.3
2	B	275	LEU	3.0
2	B	393	GLU	2.6
1	A	575	SER	2.6
2	B	390	GLU	2.6
2	B	176	ALA	2.5
2	B	228	ARG	2.5
1	A	369	PHE	2.5
1	A	467	ILE	2.5
2	B	392	ILE	2.5
1	A	570	PRO	2.2
2	B	394	ARG	2.2

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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
4	MG	A	602	1/1	0.70	0.12	2.85	43,43,43,43	0
5	ACT	B	401	4/4	0.97	0.10	0.65	21,24,29,29	0
3	SAH	A	601	26/26	0.97	0.08	-0.48	17,22,28,31	0

6.5 Other polymers [i](#)

There are no such residues in this entry.