



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 2, 2023 – 03:12 PM EDT

PDB ID : 6NM4  
Title : Crystal structure of SAM-bound PRDM9 in complex with MRK-740 inhibitor  
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Deposited on : 2019-01-10  
Resolution : 2.58 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](mailto: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>.

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

MolProbity : **FAILED**  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : **FAILED**  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.35.1

## 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 2.58 Å.

There are no overall percentile quality scores available for this entry.

MolProbity and EDS failed to run properly - the sequence quality summary graphics cannot be shown.

## 2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 3073 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.

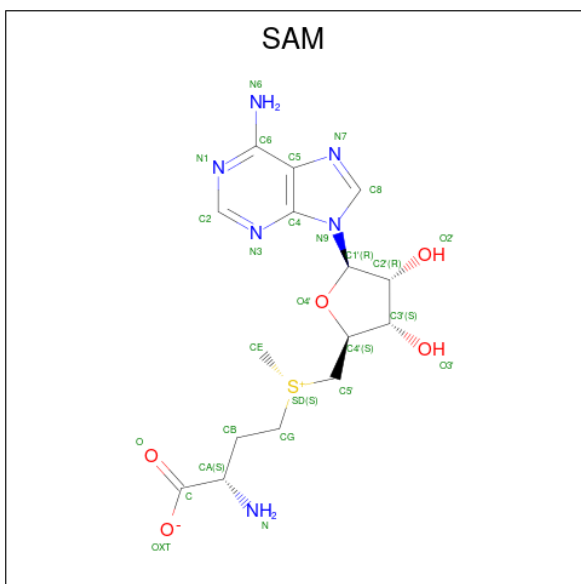
- Molecule 1 is a protein called Histone-lysine N-methyltransferase PRDM9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	185	Total 1462	C 930	N 247	O 275	S 10	0	0	0
1	B	183	Total 1461	C 928	N 252	O 271	S 10	0	0	0

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

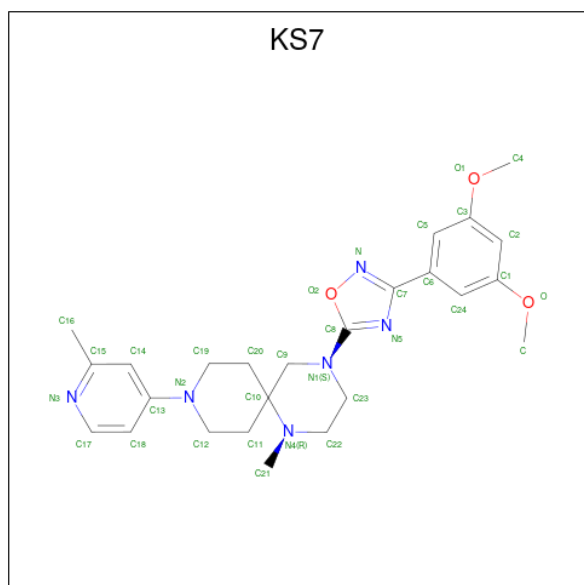
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Zn		
2	A	1	Total 1	Zn 1	0	0
2	B	1	Total 1	Zn 1	0	0

- Molecule 3 is S-ADENOSYLMETHIONINE (three-letter code: SAM) (formula: C<sub>15</sub>H<sub>22</sub>N<sub>6</sub>O<sub>5</sub>S).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
3	A	1	Total	C	N	O	S	0	0
			27	15	6	5	1		
3	B	1	Total	C	N	O	S	0	0
			27	15	6	5	1		

- Molecule 4 is 4-[3-(3,5-dimethoxyphenyl)-1,2,4-oxadiazol-5-yl]-1-methyl-9-(2-methylpyridin-4-yl)-1,4,9-triazaspiro[5.5]undecane (three-letter code: KS7) (formula: C<sub>25</sub>H<sub>32</sub>N<sub>6</sub>O<sub>3</sub>).



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<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>		<b>ZeroOcc</b>	<b>AltConf</b>
6	B	8	Total	O	0	0
			8	8		

MolProbity and EDS failed to run properly - this section is therefore empty.

### 3 Data and refinement statistics

EDS failed to run properly - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	38.03Å 74.80Å 141.44Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.88 – 2.58	Depositor
% Data completeness (in resolution range)	99.3 (39.88-2.58)	Depositor
$R_{merge}$	0.13	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.47 (at 2.58Å)	Xtrriage
Refinement program	REFMAC 5.8.0218	Depositor
R, $R_{free}$	0.207 , 0.258	Depositor
Wilson B-factor (Å <sup>2</sup> )	61.6	Xtrriage
Anisotropy	0.259	Xtrriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.35$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
Total number of atoms	3073	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	63.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.03% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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.

## 4 Model quality [i](#)

### 4.1 Standard geometry [i](#)

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### 4.2 Too-close contacts [i](#)

MolProbity failed to run properly - this section is therefore empty.

### 4.3 Torsion angles [i](#)

#### 4.3.1 Protein backbone [i](#)

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#### 4.3.2 Protein sidechains [i](#)

MolProbity failed to run properly - this section is therefore empty.

#### 4.3.3 RNA [i](#)

MolProbity failed to run properly - this section is therefore empty.

### 4.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

### 4.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

### 4.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 2 are monoatomic and 3 are unknown - leaving 4 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	SAM	A	402	-	24,29,29	1.00	1 (4%)	23,42,42	1.66	3 (13%)
4	KS7	B	403	-	33,38,38	0.59	1 (3%)	38,55,55	1.35	5 (13%)
4	KS7	A	403	-	33,38,38	0.87	2 (6%)	38,55,55	1.00	2 (5%)
3	SAM	B	402	-	24,29,29	1.07	1 (4%)	23,42,42	1.67	5 (21%)

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	SAM	A	402	-	-	2/12/33/33	0/3/3/3
4	KS7	B	403	-	-	0/12/44/44	1/5/5/5
4	KS7	A	403	-	-	0/12/44/44	0/5/5/5
3	SAM	B	402	-	-	2/12/33/33	0/3/3/3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	403	KS7	C8-N5	-3.10	1.30	1.35
4	A	403	KS7	C9-C10	3.00	1.57	1.54
4	B	403	KS7	C8-N5	-2.34	1.31	1.35
3	A	402	SAM	C5-C4	2.33	1.47	1.40
3	B	402	SAM	C5-C4	2.11	1.46	1.40

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	403	KS7	C20-C10-C9	-4.53	100.53	109.14
3	A	402	SAM	CG-SD-C5'	4.36	114.53	103.40
3	A	402	SAM	N3-C2-N1	-4.34	121.89	128.68
3	B	402	SAM	CG-SD-C5'	3.56	112.50	103.40
3	B	402	SAM	C3'-C2'-C1'	3.51	106.26	100.98
3	B	402	SAM	N3-C2-N1	-3.20	123.67	128.68
4	B	403	KS7	C6-C7-N	3.17	123.24	119.12
3	B	402	SAM	C4-C5-N7	-3.08	106.19	109.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	403	KS7	C6-C7-N	3.01	123.04	119.12
4	B	403	KS7	C23-N1-C8	2.76	128.72	119.10
4	B	403	KS7	O1-C3-C5	-2.40	108.46	119.94
3	A	402	SAM	C2-N1-C6	2.25	122.59	118.75
3	B	402	SAM	O2'-C2'-C3'	-2.21	104.66	111.82
4	A	403	KS7	C20-C10-C9	-2.15	105.05	109.14
4	B	403	KS7	O1-C3-C2	2.14	130.20	119.94

There are no chirality outliers.

All (4) torsion outliers are listed below:

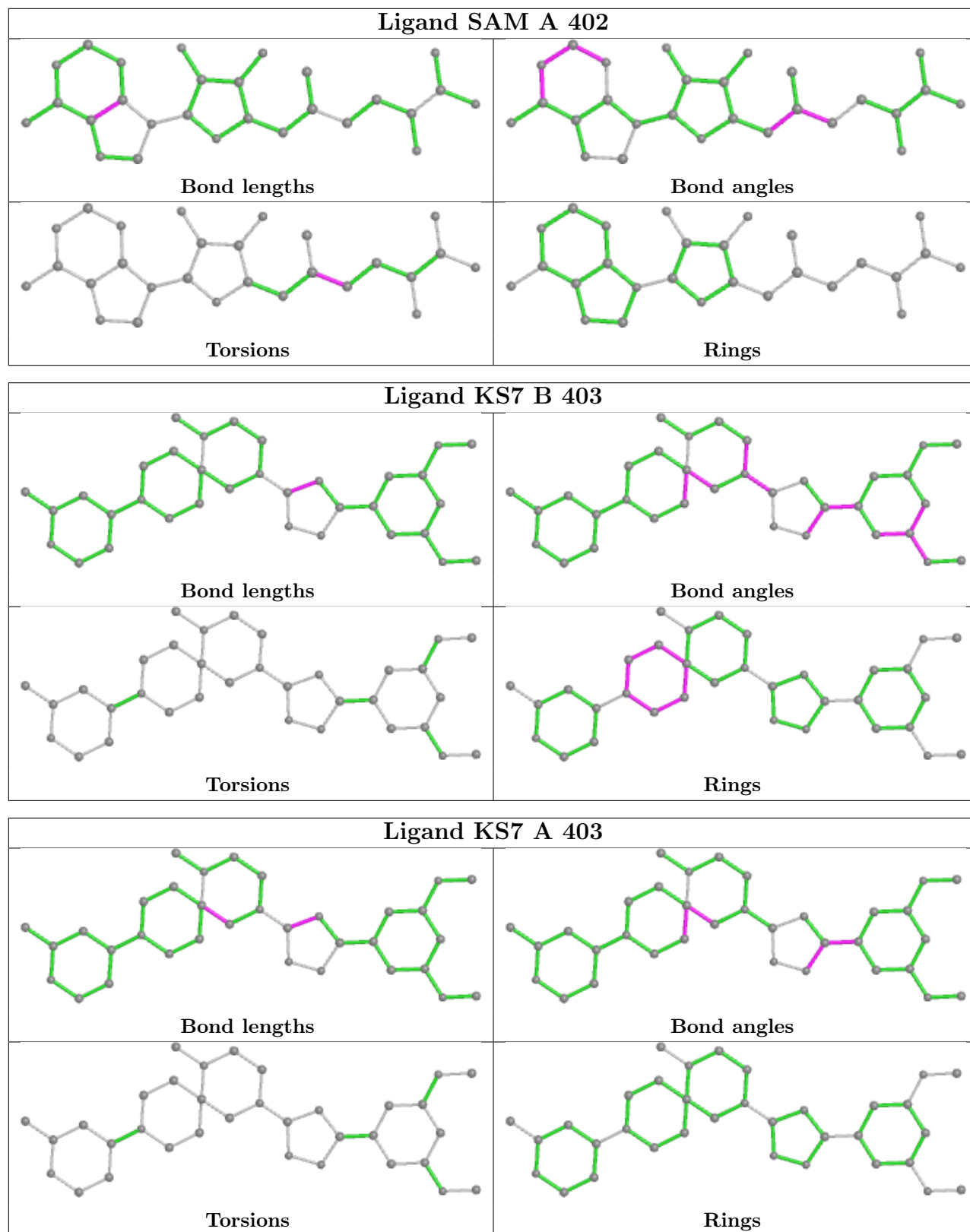
Mol	Chain	Res	Type	Atoms
3	A	402	SAM	CB-CG-SD-CE
3	A	402	SAM	CB-CG-SD-C5'
3	B	402	SAM	CB-CG-SD-CE
3	B	402	SAM	CB-CG-SD-C5'

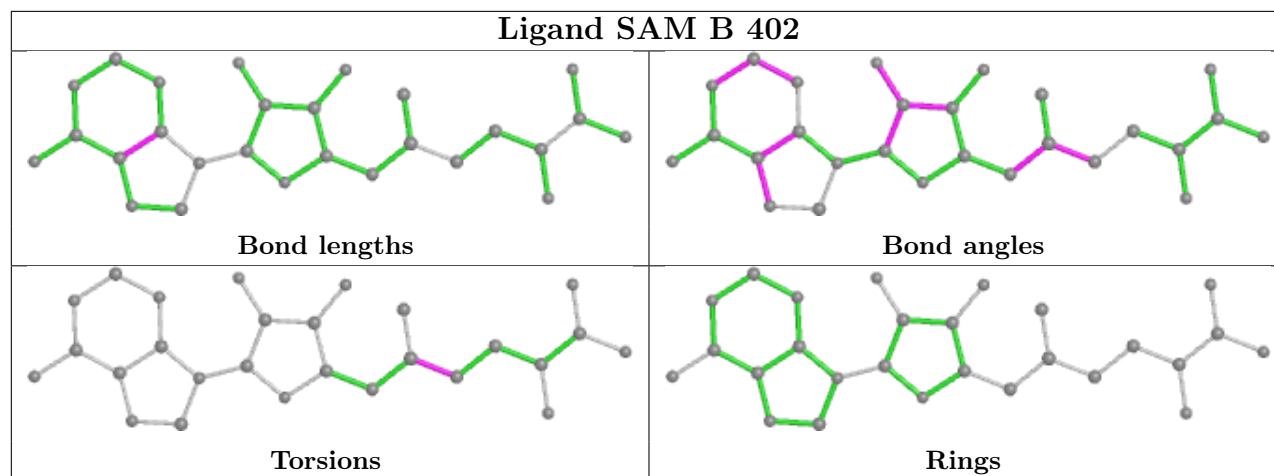
All (1) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	403	KS7	C10-C11-C12-C19-C20-N2

No monomer is involved in short contacts.

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.





#### 4.7 Other polymers [i](#)

There are no such residues in this entry.

#### 4.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 5 Fit of model and data [i](#)

### 5.1 Protein, DNA and RNA chains [i](#)

EDS failed to run properly - this section is therefore empty.

### 5.2 Non-standard residues in protein, DNA, RNA chains [i](#)

EDS failed to run properly - this section is therefore empty.

### 5.3 Carbohydrates [i](#)

EDS failed to run properly - this section is therefore empty.

### 5.4 Ligands [i](#)

EDS failed to run properly - this section is therefore empty.

### 5.5 Other polymers [i](#)

EDS failed to run properly - this section is therefore empty.