



wwPDB EM Validation Summary Report ⓘ

Jun 7, 2026 – 12:15 AM JST

PDB ID : 9WSQ / pdb_00009wsq
EMDB ID : EMD-66203
Title : Structure of dimeric mouse NLRP14-KDM2A-SKP1 complex
Authors : Liu, S.; Jiao, H.; Yan, L.; Qi, Q.; Chi, P.; Lu, Y.; Li, J.H.; Li, J.L.; Ju, S.; Wang, X.; Hu, H.; Deng, D.
Deposited on : 2025-09-15
Resolution : 3.35 Å(reported)

This is a wwPDB EM Validation Summary 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/EMValidationReportHelp>
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:

EMDB validation analysis : **FAILED**
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : **NOT EXECUTED**
MapQ : **FAILED**
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

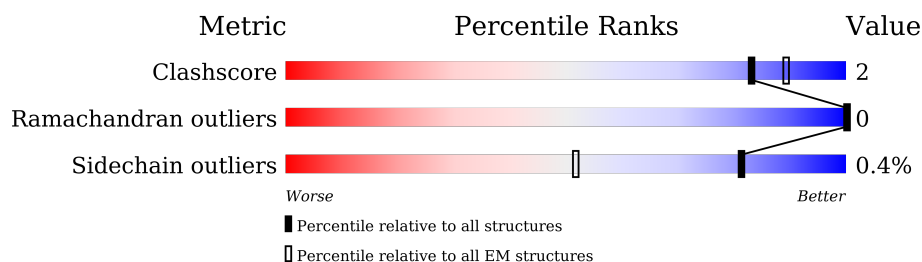
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.35 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)	EM structures (#Entries)
Clashscore	229148	23984
Ramachandran outliers	224038	23583
Sidechain outliers	223484	23102

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the 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 $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	B	358	70% 6% 24%
1	E	358	71% 6% 24%
2	C	174	86% 8% 6%
2	F	174	86% 8% 6%
3	A	1012	71% . 25%
3	D	1012	71% . 25%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 18805 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, 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 Lysine-specific demethylase 2A.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	B	273	Total	C	N	O	S	0	0
			2184	1379	393	394	18		
1	E	273	Total	C	N	O	S	0	0
			2187	1381	393	394	19		

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	804	MET	-	initiating methionine	UNP P59997
B	805	TRP	-	expression tag	UNP P59997
B	806	SER	-	expression tag	UNP P59997
B	807	HIS	-	expression tag	UNP P59997
B	808	PRO	-	expression tag	UNP P59997
B	809	GLN	-	expression tag	UNP P59997
B	810	PHE	-	expression tag	UNP P59997
B	811	GLU	-	expression tag	UNP P59997
B	812	LYS	-	expression tag	UNP P59997
B	813	GLY	-	expression tag	UNP P59997
B	814	THR	-	expression tag	UNP P59997
E	804	MET	-	initiating methionine	UNP P59997
E	805	TRP	-	expression tag	UNP P59997
E	806	SER	-	expression tag	UNP P59997
E	807	HIS	-	expression tag	UNP P59997
E	808	PRO	-	expression tag	UNP P59997
E	809	GLN	-	expression tag	UNP P59997
E	810	PHE	-	expression tag	UNP P59997
E	811	GLU	-	expression tag	UNP P59997
E	812	LYS	-	expression tag	UNP P59997
E	813	GLY	-	expression tag	UNP P59997
E	814	THR	-	expression tag	UNP P59997

- Molecule 2 is a protein called S-phase kinase-associated protein 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	C	163	Total	C	N	O	S	0	0
			1274	797	207	266	4		
2	F	163	Total	C	N	O	S	0	0
			1274	797	207	266	4		

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-10	MET	-	initiating methionine	UNP Q9WTX5
C	-9	TRP	-	expression tag	UNP Q9WTX5
C	-8	SER	-	expression tag	UNP Q9WTX5
C	-7	HIS	-	expression tag	UNP Q9WTX5
C	-6	PRO	-	expression tag	UNP Q9WTX5
C	-5	GLN	-	expression tag	UNP Q9WTX5
C	-4	PHE	-	expression tag	UNP Q9WTX5
C	-3	GLU	-	expression tag	UNP Q9WTX5
C	-2	LYS	-	expression tag	UNP Q9WTX5
C	-1	GLY	-	expression tag	UNP Q9WTX5
C	0	THR	-	expression tag	UNP Q9WTX5
F	-10	MET	-	initiating methionine	UNP Q9WTX5
F	-9	TRP	-	expression tag	UNP Q9WTX5
F	-8	SER	-	expression tag	UNP Q9WTX5
F	-7	HIS	-	expression tag	UNP Q9WTX5
F	-6	PRO	-	expression tag	UNP Q9WTX5
F	-5	GLN	-	expression tag	UNP Q9WTX5
F	-4	PHE	-	expression tag	UNP Q9WTX5
F	-3	GLU	-	expression tag	UNP Q9WTX5
F	-2	LYS	-	expression tag	UNP Q9WTX5
F	-1	GLY	-	expression tag	UNP Q9WTX5
F	0	THR	-	expression tag	UNP Q9WTX5

- Molecule 3 is a protein called NACHT, LRR and PYD domains-containing protein 14.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	A	756	Total	C	N	O	S	0	0
			5942	3755	1038	1092	57		
3	D	756	Total	C	N	O	S	0	0
			5944	3757	1038	1092	57		

There are 38 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-18	ASP	-	expression tag	UNP Q6B966

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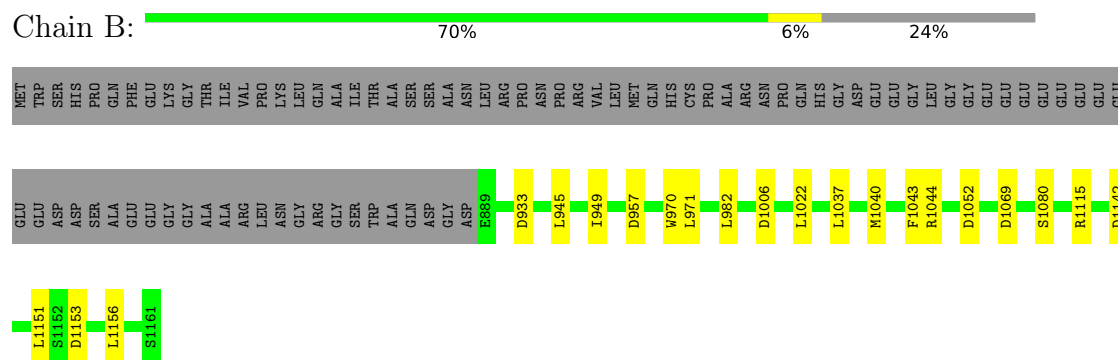
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Chain	Residue	Modelled	Actual	Comment	Reference
A	-17	TYR	-	expression tag	UNP Q6B966
A	-16	LYS	-	expression tag	UNP Q6B966
A	-15	ASP	-	expression tag	UNP Q6B966
A	-14	ASP	-	expression tag	UNP Q6B966
A	-13	ASP	-	expression tag	UNP Q6B966
A	-12	ASP	-	expression tag	UNP Q6B966
A	-11	LYS	-	expression tag	UNP Q6B966
A	-10	GLY	-	expression tag	UNP Q6B966
A	-9	ASP	-	expression tag	UNP Q6B966
A	-8	TYR	-	expression tag	UNP Q6B966
A	-7	LYS	-	expression tag	UNP Q6B966
A	-6	ASP	-	expression tag	UNP Q6B966
A	-5	ASP	-	expression tag	UNP Q6B966
A	-4	ASP	-	expression tag	UNP Q6B966
A	-3	ASP	-	expression tag	UNP Q6B966
A	-2	LYS	-	expression tag	UNP Q6B966
A	-1	GLY	-	expression tag	UNP Q6B966
A	0	SER	-	expression tag	UNP Q6B966
D	-18	ASP	-	expression tag	UNP Q6B966
D	-17	TYR	-	expression tag	UNP Q6B966
D	-16	LYS	-	expression tag	UNP Q6B966
D	-15	ASP	-	expression tag	UNP Q6B966
D	-14	ASP	-	expression tag	UNP Q6B966
D	-13	ASP	-	expression tag	UNP Q6B966
D	-12	ASP	-	expression tag	UNP Q6B966
D	-11	LYS	-	expression tag	UNP Q6B966
D	-10	GLY	-	expression tag	UNP Q6B966
D	-9	ASP	-	expression tag	UNP Q6B966
D	-8	TYR	-	expression tag	UNP Q6B966
D	-7	LYS	-	expression tag	UNP Q6B966
D	-6	ASP	-	expression tag	UNP Q6B966
D	-5	ASP	-	expression tag	UNP Q6B966
D	-4	ASP	-	expression tag	UNP Q6B966
D	-3	ASP	-	expression tag	UNP Q6B966
D	-2	LYS	-	expression tag	UNP Q6B966
D	-1	GLY	-	expression tag	UNP Q6B966
D	0	SER	-	expression tag	UNP Q6B966

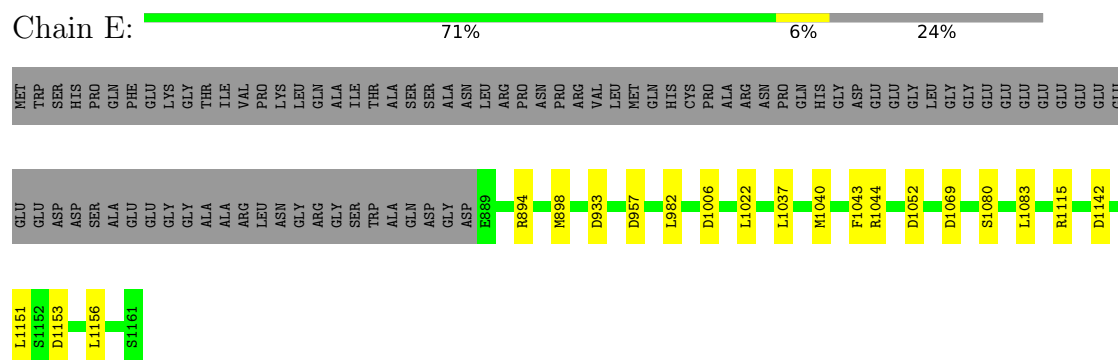
3 Residue-property plots [i](#)

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. 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. 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.

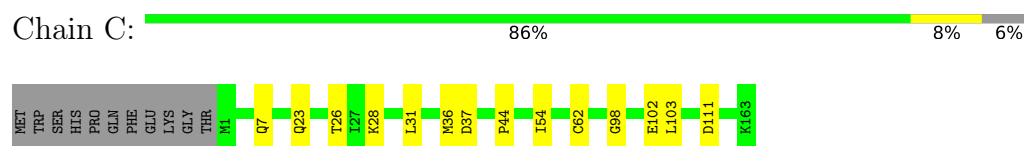
- Molecule 1: Lysine-specific demethylase 2A



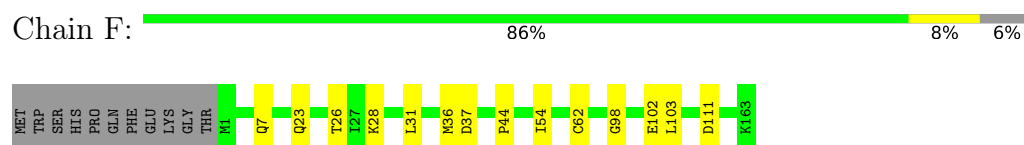
- Molecule 1: Lysine-specific demethylase 2A

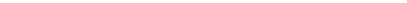


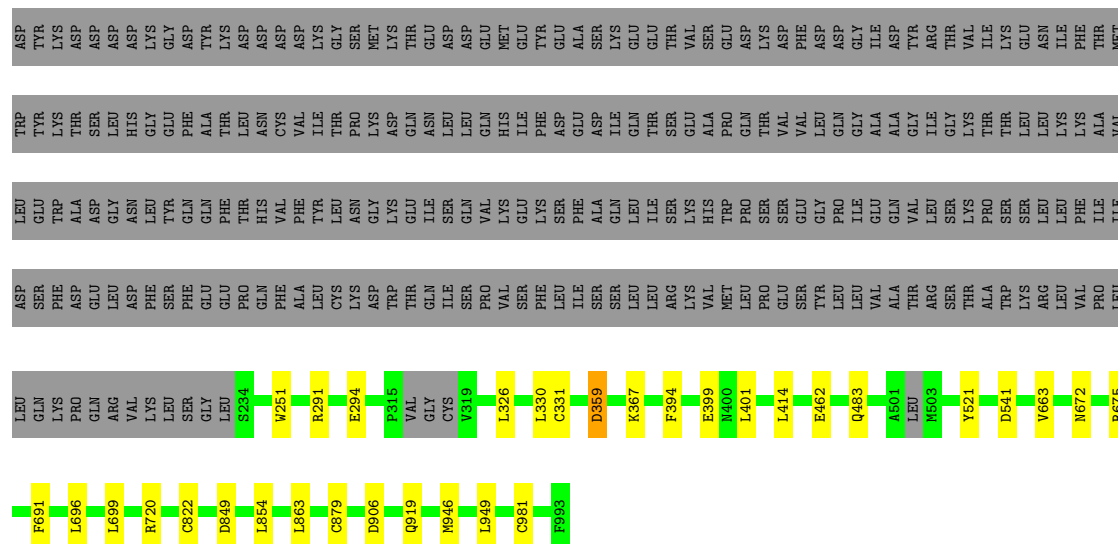
- Molecule 2: S-phase kinase-associated protein 1



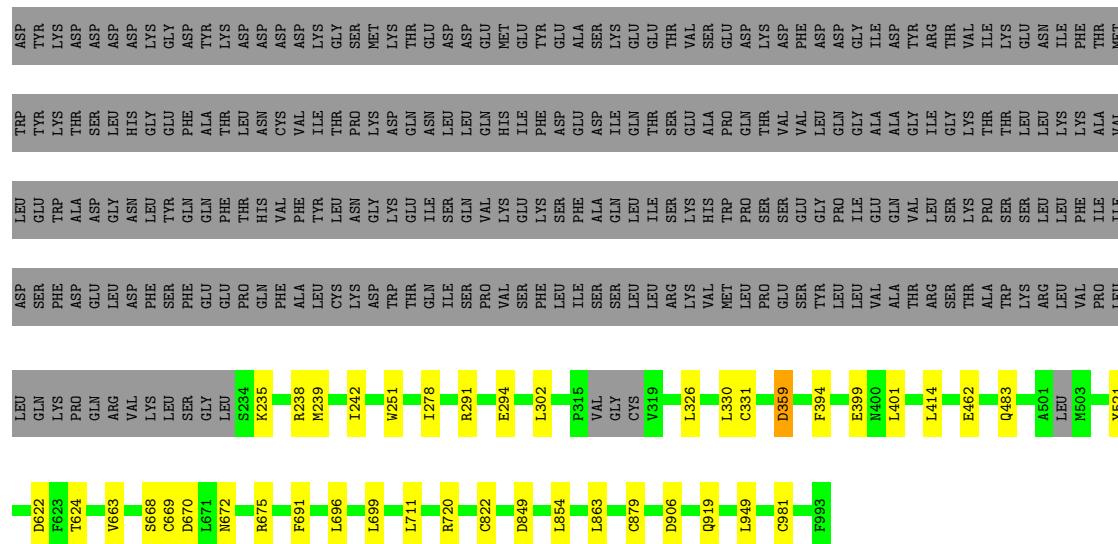
- Molecule 2: S-phase kinase-associated protein 1



- Chain A:  71% 25%



- Chain D:  71% . 25%



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	49652	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	53.633	Depositor
Minimum defocus (nm)	1200	Depositor
Maximum defocus (nm)	1800	Depositor
Magnification	105000	Depositor
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	B	0.08	0/2223	0.24	0/3006
1	E	0.07	0/2226	0.24	0/3009
2	C	0.09	0/1297	0.23	0/1765
2	F	0.09	0/1297	0.23	0/1765
3	A	0.10	0/6042	0.25	0/8163
3	D	0.09	0/6044	0.25	0/8166
All	All	0.09	0/19129	0.24	0/25874

There are no bond length outliers.

There are no bond angle outliers.

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	B	2184	0	2262	14	0
1	E	2187	0	2269	11	0
2	C	1274	0	1203	8	0
2	F	1274	0	1203	8	0
3	A	5942	0	5958	24	0
3	D	5944	0	5965	28	0
All	All	18805	0	18860	92	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 2.

The worst 5 of 92 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)
3:D:668:SER:HB2	3:D:696:LEU:HD23	1.29	1.14
3:A:414:LEU:HD12	3:A:462:GLU:CD	1.83	1.03
3:D:414:LEU:HD12	3:D:462:GLU:CD	1.83	1.03
3:A:414:LEU:HD12	3:A:462:GLU:OE2	1.79	0.82
3:D:699:LEU:HD12	3:D:699:LEU:O	1.79	0.82

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 PDB entries followed by that with respect to all EM entries.

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	B	271/358 (76%)	268 (99%)	3 (1%)	0	100	100
1	E	271/358 (76%)	268 (99%)	3 (1%)	0	100	100
2	C	161/174 (92%)	158 (98%)	3 (2%)	0	100	100
2	F	161/174 (92%)	158 (98%)	3 (2%)	0	100	100
3	A	750/1012 (74%)	739 (98%)	11 (2%)	0	100	100
3	D	750/1012 (74%)	739 (98%)	11 (2%)	0	100	100
All	All	2364/3088 (77%)	2330 (99%)	34 (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 PDB entries followed by that with respect to all EM entries.

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	B	251/319 (79%)	250 (100%)	1 (0%)	84	82
1	E	252/319 (79%)	251 (100%)	1 (0%)	84	82
2	C	140/160 (88%)	140 (100%)	0	100	100
2	F	140/160 (88%)	140 (100%)	0	100	100
3	A	668/926 (72%)	665 (100%)	3 (0%)	84	82
3	D	669/926 (72%)	666 (100%)	3 (0%)	84	82
All	All	2120/2810 (75%)	2112 (100%)	8 (0%)	81	82

5 of 8 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	D	949	LEU
3	D	359	ASP
1	E	1043	PHE
3	A	696	LEU
3	D	330	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 9 such sidechains are listed below:

Mol	Chain	Res	Type
3	D	236	ASN
3	D	559	GLN
3	A	559	GLN
3	A	799	HIS
1	E	1033	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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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.