



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 10, 2026 – 11:45 AM UTC

PDB ID : 9VG0 / pdb_00009vg0
Title : SIRT2 structure in complex with H3K18myr peptide
Authors : Zhang, N.; Hao, Q.
Deposited on : 2025-06-12
Resolution : 1.61 Å(reported)

This is a wwPDB X-ray Structure 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/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 : ?? (???), CSD ??CSD?? (????)
Xtriage (Phenix) : 2.0
EDS : 3.0
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

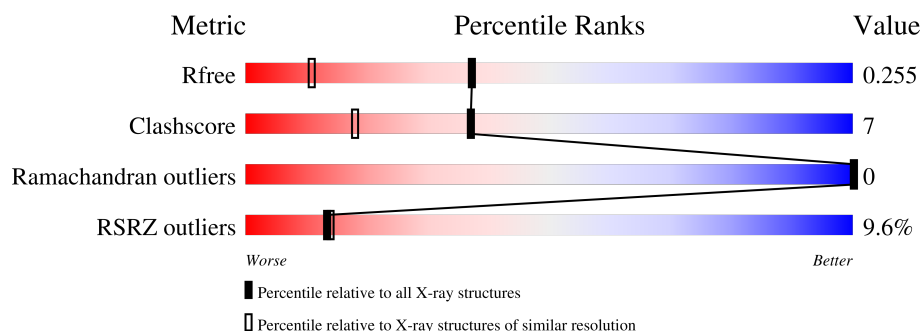
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.61 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	6728 (1.64-1.60)
Clashscore	190562	7023 (1.64-1.60)
Ramachandran outliers	187476	6898 (1.64-1.60)
RSRZ outliers	180081	6727 (1.64-1.60)

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	EDO	A	404	-	-	X	-
3	EDO	A	405	-	-	X	-

ENTRY-COMPOSITION INFOmissingINFO

SEQUENCE-PLOTS INFOmissingINFO

2 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	36.15Å 73.50Å 55.73Å 90.00° 94.09° 90.00°	Depositor
Resolution (Å)	27.81 – 1.61 27.81 – 1.61	Depositor EDS
% Data completeness (in resolution range)	91.7 (27.81-1.61) 91.7 (27.81-1.61)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.86 (at 1.61Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.188 , 0.231 (Not available) , 0.255	Depositor DCC
R_{free} test set	1746 reflections (4.63%)	wwPDB-VP
Wilson B-factor (Å ²)	21.9	Xtriage
Anisotropy	0.108	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 49.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	2525	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.16% 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.

3 Model quality

3.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MYR, ZN, EDO

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	1.19	4/2320 (0.2%)	1.31	5/3137 (0.2%)
2	C	1.34	0/43	1.35	0/55
All	All	1.19	4/2363 (0.2%)	1.31	5/3192 (0.2%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	225	GLN	C-N	7.68	1.44	1.33
1	A	350	ALA	C-O	-6.81	1.16	1.24
1	A	341	LEU	C-O	5.85	1.30	1.24
1	A	230	PRO	C-O	-5.71	1.17	1.23

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	65	GLU	CB-CG-CD	5.93	122.68	112.60
1	A	68	ALA	N-CA-C	-5.77	105.07	111.36
1	A	124	PHE	CA-C-N	5.28	127.36	120.28
1	A	124	PHE	C-N-CA	5.28	127.36	120.28
1	A	227	LEU	CA-C-O	-5.06	116.18	121.55

There are no chirality outliers.

There are no planarity outliers.

3.2 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	2271	0	2220	30	0
2	C	44	0	38	2	0
3	A	20	0	29	14	0
4	A	1	0	0	0	0
5	C	15	0	27	1	0
6	A	170	0	0	3	0
6	C	4	0	0	1	0
All	All	2525	0	2314	33	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 7.

The worst 5 of 33 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:170:ASP:OD2	3:A:404:EDO:H21	1.79	0.81
1:A:170:ASP:OD2	3:A:404:EDO:C2	2.39	0.70
1:A:238:SER:HG	2:C:6:ALA:N	1.92	0.68
1:A:148[B]:CYS:SG	1:A:324:CYS:HB3	2.35	0.67
1:A:125:LYS:O	3:A:405:EDO:H12	1.94	0.66

There are no symmetry-related clashes.

3.3 Torsion angles [i](#)

3.3.1 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	287/307 (94%)	279 (97%)	8 (3%)	0	100	100
2	C	4/7 (57%)	3 (75%)	1 (25%)	0	100	100
All	All	291/314 (93%)	282 (97%)	9 (3%)	0	100	100

There are no Ramachandran outliers to report.

3.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

3.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

3.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

3.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 2 could not be matched to an existing wwPDB Chemical Component Dictionary definition at this stage - leaving 5 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	EDO	A	403	-	3,3,3	0.48	0	2,2,2	0.52	0
3	EDO	A	405	-	3,3,3	0.31	0	2,2,2	0.42	0
3	EDO	A	401	-	3,3,3	0.94	0	2,2,2	0.32	0
3	EDO	A	402	-	3,3,3	0.12	0	2,2,2	0.47	0
3	EDO	A	404	-	3,3,3	0.52	0	2,2,2	0.24	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
3	EDO	A	403	-	-	1/1/1/1	-
3	EDO	A	405	-	-	0/1/1/1	-
3	EDO	A	401	-	-	1/1/1/1	-
3	EDO	A	402	-	-	1/1/1/1	-
3	EDO	A	404	-	-	1/1/1/1	-
5	MYR	C	101	2	-	6/12/12/13	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

5 of 10 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	C	101	MYR	O1-C1-C2-C3
5	C	101	MYR	C9-C10-C11-C12
5	C	101	MYR	C10-C11-C12-C13
3	A	404	EDO	O1-C1-C2-O2
5	C	101	MYR	C7-C8-C9-C10

There are no ring outliers.

4 monomers are involved in 15 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	403	EDO	3	0
3	A	405	EDO	4	0
3	A	404	EDO	7	0
5	C	101	MYR	1	0

3.7 Other polymers [i](#)

There are no such residues in this entry.

3.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

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

4.1 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	287/307 (93%)	0.46	25 (8%) 16 17	8, 25, 43, 68	6 (2%)
2	C	6/7 (85%)	1.96	3 (50%) 0 0	35, 40, 50, 56	0
All	All	293/314 (93%)	0.49	28 (9%) 13 14	8, 26, 46, 68	6 (2%)

The worst 5 of 28 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	110	TYR	4.7
1	A	297	LEU	4.5
1	A	114	TYR	3.8
1	A	296	PHE	3.6
1	A	217	VAL	3.5

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

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

4.3 Carbohydrates [i](#)

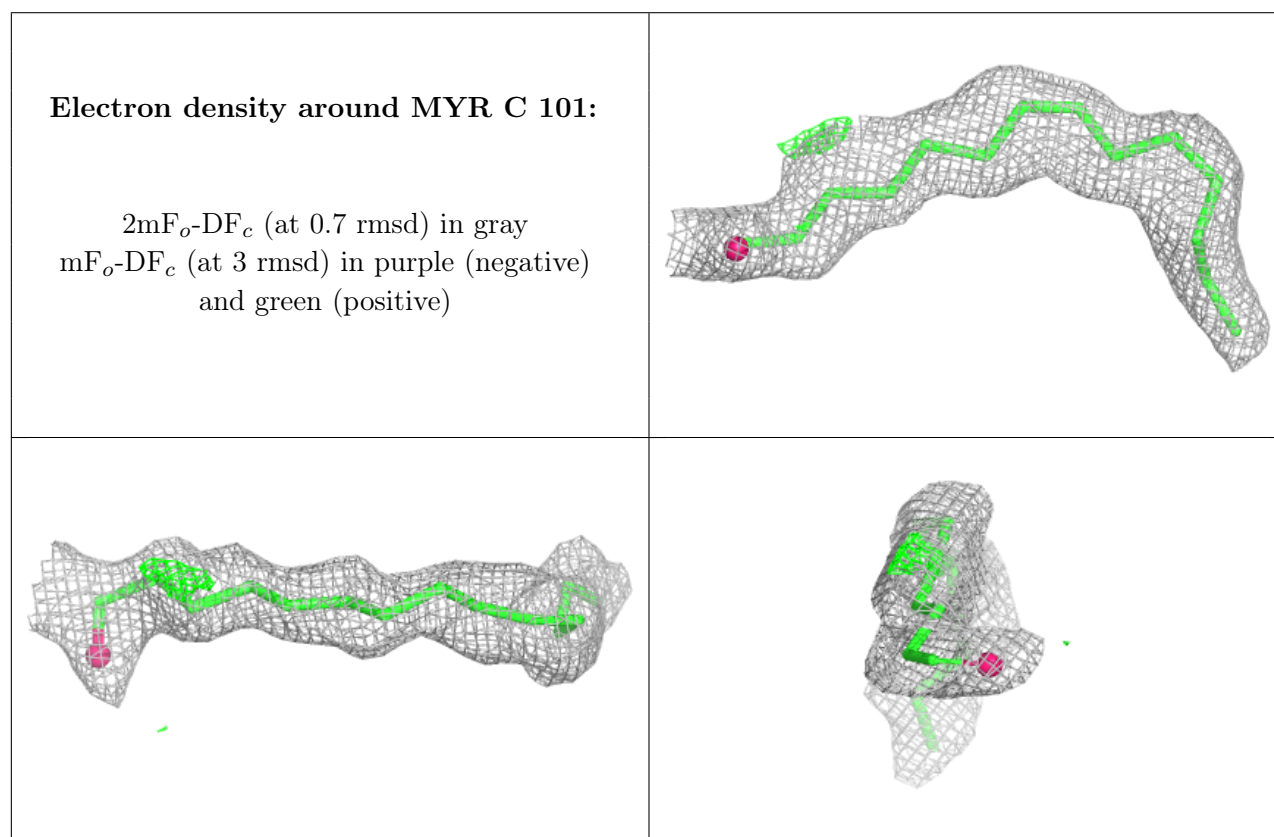
There are no oligosaccharides in this entry.

4.4 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($\text{\AA}^2$)	Q<0.9
3	EDO	A	405	4/4	0.79	0.14	43,45,45,46	0
3	EDO	A	403	4/4	0.83	0.15	37,37,37,38	0
3	EDO	A	401	4/4	0.85	0.16	28,34,35,38	0
3	EDO	A	404	4/4	0.85	0.23	30,34,38,43	0
3	EDO	A	402	4/4	0.85	0.12	30,35,37,37	0
5	MYR	C	101	15/16	0.86	0.11	29,34,40,40	0
4	ZN	A	406	1/1	0.99	0.02	21,21,21,21	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.



4.5 Other polymers ⓘ

There are no such residues in this entry.