



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 22, 2026 – 06:08 PM JST

PDB ID : 9VAD / pdb_00009vad
Title : Crystal structure of the PDZ tandem of syntenin
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Deposited on : 2025-06-03
Resolution : 2.05 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 ⓘ 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
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

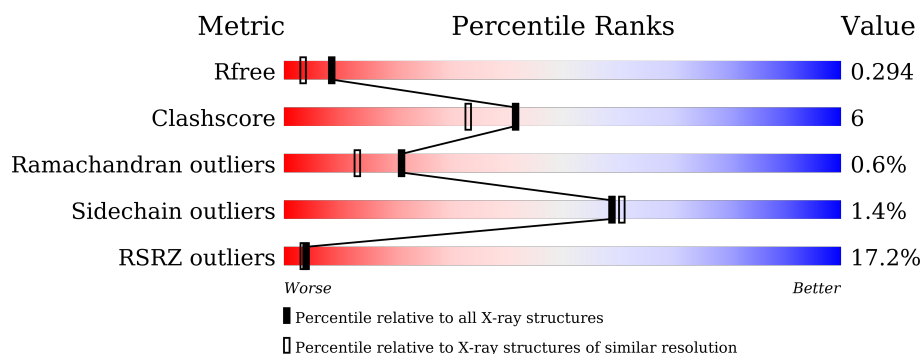
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.05 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	2260 (2.04-2.04)
Clashscore	190562	2333 (2.04-2.04)
Ramachandran outliers	187476	2318 (2.04-2.04)
Sidechain outliers	187428	2318 (2.04-2.04)
RSRZ outliers	180081	2260 (2.04-2.04)

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 $\leq 5\%$. 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	166	 6% 80% 20%
1	B	166	 28% 89% 11%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2621 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 Syntenin-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	166	Total	C	N	O	S	0	0	0
			1270	797	225	242	6			
1	B	166	Total	C	N	O	S	0	0	0
			1270	797	225	242	6			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	275	TYR	PHE	engineered mutation	UNP O00560
B	275	TYR	PHE	engineered mutation	UNP O00560

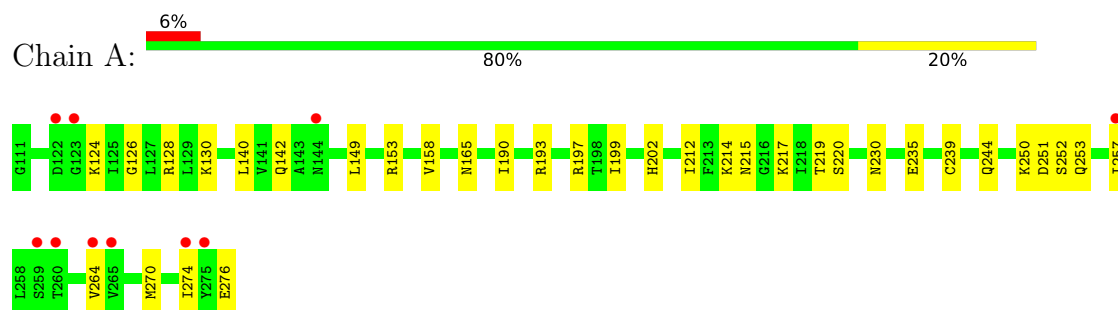
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	45	Total	O	0	0
			45	45		
2	B	36	Total	O	0	0
			36	36		

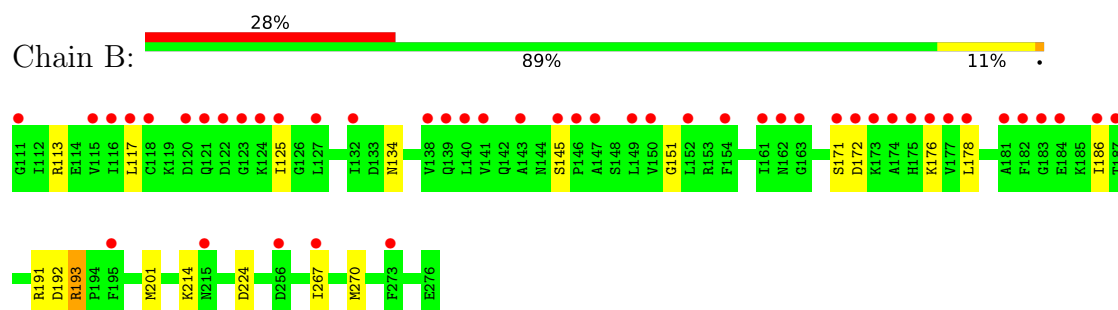
3 Residue-property plots [i](#)

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: Syntenin-1



• Molecule 1: Syntenin-1



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	57.02Å 57.02Å 154.75Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	47.04 – 2.05 47.04 – 2.05	Depositor EDS
% Data completeness (in resolution range)	100.0 (47.04-2.05) 92.6 (47.04-2.05)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.80 (at 2.05Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.243 , 0.294 0.243 , 0.294	Depositor DCC
R_{free} test set	949 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	34.5	Xtriage
Anisotropy	0.420	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 38.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.035 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	2621	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.01% 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

5.1 Standard geometry

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	0.10	0/1286	0.28	0/1730
1	B	0.09	0/1286	0.29	0/1730
All	All	0.10	0/2572	0.28	0/3460

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

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	1270	0	1300	25	0
1	B	1270	0	1300	10	0
2	A	45	0	0	0	0
2	B	36	0	0	1	0
All	All	2621	0	2600	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 6.

All (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:B:193:ARG:HD2	1:B:270:MET:HE1	1.58	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:239:CYS:SG	1:A:270:MET:HE2	2.32	0.69
1:A:197:ARG:NH2	1:A:199:ILE:HD11	2.09	0.68
1:A:244:GLN:N	1:A:244:GLN:OE1	2.30	0.63
1:A:126:GLY:HA2	1:A:142:GLN:HG3	1.87	0.56
1:A:165:ASN:HD22	1:B:224:ASP:HA	1.72	0.55
1:B:117:LEU:HD23	1:B:125:ILE:HD11	1.89	0.53
1:A:128:ARG:HG3	1:A:140:LEU:HD22	1.93	0.51
1:A:257:ILE:HD12	1:A:257:ILE:H	1.75	0.50
1:A:158:VAL:HA	1:A:190:ILE:HD13	1.94	0.49
1:B:178:LEU:HD23	1:B:186:ILE:HG21	1.94	0.49
1:B:214:LYS:NZ	2:B:302:HOH:O	2.27	0.49
1:A:250:LYS:H	1:A:253:GLN:NE2	2.11	0.49
1:B:201:MET:HE1	1:B:267:ILE:HD12	1.94	0.49
1:A:214:LYS:HZ3	1:A:219:THR:HG21	1.79	0.47
1:B:172:ASP:OD2	1:B:176:LYS:NZ	2.45	0.47
1:A:126:GLY:CA	1:A:142:GLN:HG3	2.44	0.47
1:A:130:LYS:NZ	1:A:274:ILE:O	2.34	0.47
1:A:153:ARG:HE	1:A:153:ARG:HB2	1.57	0.46
1:A:257:ILE:HD12	1:A:257:ILE:N	2.30	0.46
1:A:214:LYS:HG2	1:A:219:THR:HG21	1.97	0.46
1:A:217:LYS:HE3	1:A:235:GLU:HA	1.99	0.45
1:A:215:ASN:N	1:A:251:ASP:OD2	2.38	0.44
1:A:128:ARG:HB2	1:A:140:LEU:HB3	1.99	0.44
1:A:250:LYS:H	1:A:253:GLN:HE21	1.65	0.44
1:A:230:ASN:O	1:B:191:ARG:NH1	2.42	0.44
1:A:130:LYS:NZ	1:A:276:GLU:HB2	2.33	0.43
1:A:202:HIS:ND1	1:A:264:VAL:HG22	2.34	0.42
1:B:113:ARG:NH1	1:B:192:ASP:OD1	2.53	0.41
1:A:124:LYS:HB2	1:A:124:LYS:HE2	1.90	0.41
1:A:212:ILE:HD12	1:A:220:SER:OG	2.20	0.41
1:A:149:LEU:HD23	1:A:149:LEU:HA	1.87	0.40
1:B:113:ARG:HH22	1:B:151:GLY:C	2.29	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	164/166 (99%)	160 (98%)	3 (2%)	1 (1%)	21	13
1	B	164/166 (99%)	152 (93%)	11 (7%)	1 (1%)	21	13
All	All	328/332 (99%)	312 (95%)	14 (4%)	2 (1%)	21	13

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	193	ARG
1	B	193	ARG

5.3.2 Protein sidechains ⓘ

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	141/141 (100%)	140 (99%)	1 (1%)	76	80
1	B	141/141 (100%)	138 (98%)	3 (2%)	47	46
All	All	282/282 (100%)	278 (99%)	4 (1%)	59	60

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

Mol	Chain	Res	Type
1	A	252	SER
1	B	134	ASN
1	B	145	SER
1	B	171	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	253	GLN

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Mol	Chain	Res	Type
1	B	134	ASN
1	B	237	ASN

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 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.

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	166/166 (100%)	0.87	10 (6%)	27 28	26, 43, 62, 68	0
1	B	166/166 (100%)	1.28	47 (28%)	1 1	28, 47, 98, 110	0
All	All	332/332 (100%)	1.08	57 (17%)	4 3	26, 44, 84, 110	0

All (57) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	182	PHE	5.4
1	A	257	ILE	4.1
1	A	275	TYR	4.1
1	B	115	VAL	3.9
1	B	147	ALA	3.9
1	B	181	ALA	3.8
1	B	117	LEU	3.8
1	B	125	ILE	3.8
1	B	124	LYS	3.6
1	B	183	GLY	3.5
1	B	178	LEU	3.5
1	B	177	VAL	3.4
1	B	186	ILE	3.4
1	B	161	ILE	3.3
1	B	141	VAL	3.2
1	B	118	CYS	3.2
1	B	143	ALA	3.1
1	B	123	GLY	3.0
1	B	146	PRO	3.0
1	B	162	ASN	3.0
1	B	172	ASP	3.0
1	B	175	HIS	3.0
1	B	149	LEU	2.9
1	A	260	THR	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	127	LEU	2.8
1	B	139	GLN	2.7
1	B	122	ASP	2.7
1	B	111	GLY	2.6
1	B	121	GLN	2.5
1	A	264	VAL	2.5
1	B	140	LEU	2.5
1	A	265	VAL	2.5
1	B	150	VAL	2.5
1	A	274	ILE	2.5
1	B	171	SER	2.4
1	B	154	PHE	2.4
1	B	256	ASP	2.4
1	B	195	PHE	2.4
1	B	174	ALA	2.3
1	B	145	SER	2.3
1	B	132	ILE	2.3
1	B	116	ILE	2.2
1	A	144	ASN	2.2
1	B	173	LYS	2.2
1	A	122	ASP	2.2
1	B	120	ASP	2.2
1	B	176	LYS	2.1
1	B	215	ASN	2.1
1	B	187	THR	2.1
1	B	273	PHE	2.1
1	B	152	LEU	2.1
1	B	163	GLY	2.1
1	B	267	ILE	2.1
1	B	138	VAL	2.1
1	A	259	SER	2.1
1	A	123	GLY	2.0
1	B	184	GLU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.