



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:39 PM UTC

PDB ID : 7QNL / pdb_00007qnl
Title : LOV2-DARPIN FUSION - D4_DeltaLOV
Authors : Mittl, P.
Deposited on : 2021-12-21
Resolution : 1.71 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
Mogul : 2022.3.0, CSD as543be (2022)
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

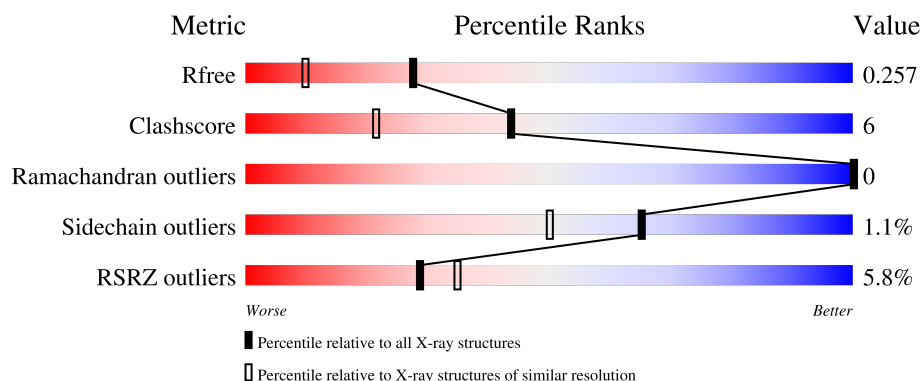
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.71 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	1039 (1.72-1.72)
Clashscore	190562	1049 (1.72-1.72)
Ramachandran outliers	187476	1041 (1.72-1.72)
Sidechain outliers	187428	1041 (1.72-1.72)
RSRZ outliers	180081	1039 (1.72-1.72)

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	AAA	240	6% 88% 12%

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
2	SO4	AAA	808	-	-	X	-

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 2066 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 DARPIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	AAA	240	Total	C	N	O	S	0	3	0
			1849	1163	320	364	2			

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	AAA	1	Total	O	S	0	0
			5	4	1		
2	AAA	1	Total	O	S	0	0
			5	4	1		
2	AAA	1	Total	O	S	0	0
			5	4	1		
2	AAA	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	AAA	1	Total	O	S	0	0
			5	4	1		
2	AAA	1	Total	O	S	0	0
			5	4	1		
2	AAA	1	Total	O	S	0	0
			5	4	1		
2	AAA	1	Total	O	S	0	0
			5	4	1		
2	AAA	1	Total	O	S	0	0
			5	4	1		

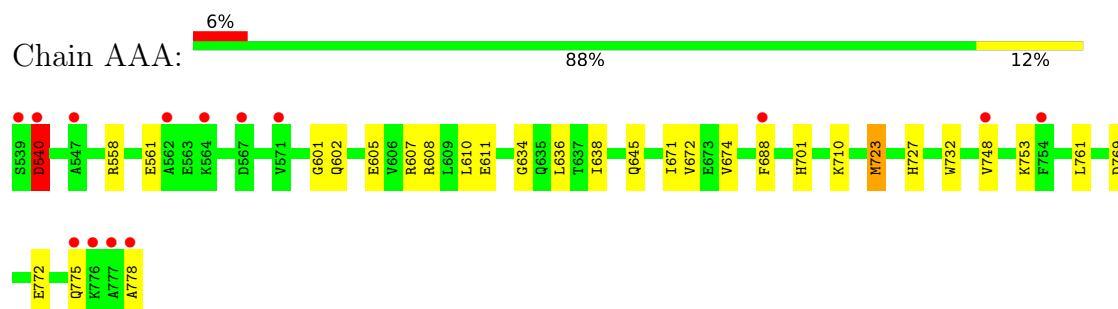
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	AAA	167	Total	O	0	0
			167	167		

3 Residue-property plots

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: DARPIN



4 Data and refinement statistics

Property	Value	Source
Space group	P 42 21 2	Depositor
Cell constants a, b, c, α , β , γ	89.74Å 89.74Å 69.30Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	63.46 – 1.71 63.46 – 1.71	Depositor EDS
% Data completeness (in resolution range)	72.4 (63.46-1.71) 72.9 (63.46-1.71)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.68 (at 1.71Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.197 , 0.264 (Not available) , 0.257	Depositor DCC
R_{free} test set	1135 reflections (4.65%)	wwPDB-VP
Wilson B-factor (Å ²)	30.5	Xtriage
Anisotropy	0.030	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 33.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	2066	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.87% 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: SO4

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	AAA	1.22	2/1873 (0.1%)	1.48	3/2536 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	AAA	701	HIS	CE1-NE2	5.69	1.38	1.32
1	AAA	672	VAL	C-O	5.31	1.30	1.24

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AAA	775	GLN	CA-C-N	7.13	130.69	120.79
1	AAA	775	GLN	C-N-CA	7.13	130.69	120.79
1	AAA	540	ASP	CA-CB-CG	5.25	117.85	112.60

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	AAA	1849	0	1862	23	0
2	AAA	50	0	0	4	0
3	AAA	167	0	0	2	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	2066	0	1862	23	1

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 (23) 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:AAA:607:ARG:NH1	2:AAA:808:SO4:O1	2.25	0.70
1:AAA:753:LYS:O	1:AAA:753:LYS:HG2	1.99	0.62
1:AAA:748:VAL:HG11	1:AAA:778:ALA:HA	1.82	0.61
1:AAA:602:GLN:OE1	3:AAA:901:HOH:O	2.17	0.60
1:AAA:688:PHE:CD2	1:AAA:688:PHE:O	2.63	0.52
1:AAA:607:ARG:NH1	2:AAA:808:SO4:S	2.87	0.47
1:AAA:723:MET:HE1	1:AAA:761:LEU:HD11	1.96	0.47
1:AAA:723:MET:HE2	1:AAA:727:HIS:HB3	1.96	0.45
1:AAA:610:LEU:HD13	1:AAA:645:GLN:HG3	1.99	0.44
1:AAA:607:ARG:O	1:AAA:611:GLU:HG3	2.17	0.44
1:AAA:732:TRP:HA	1:AAA:732:TRP:CE3	2.53	0.43
1:AAA:608[A]:ARG:HE	1:AAA:608[A]:ARG:HB3	1.36	0.43
1:AAA:710:LYS:NZ	3:AAA:909:HOH:O	2.43	0.43
1:AAA:558:ARG:HA	1:AAA:561:GLU:OE1	2.19	0.43
1:AAA:769:ASP:O	1:AAA:772:GLU:HB3	2.19	0.43
1:AAA:601:GLY:HA2	1:AAA:638:ILE:HD12	2.02	0.42
1:AAA:769:ASP:N	2:AAA:806:SO4:O4	2.44	0.42
1:AAA:605:GLU:OE2	1:AAA:608[A]:ARG:NH1	2.52	0.42
1:AAA:634:GLY:HA2	1:AAA:671:ILE:HD12	2.01	0.42
1:AAA:636:LEU:HD11	1:AAA:674:VAL:HG21	2.02	0.41
1:AAA:540:ASP:HB3	2:AAA:810:SO4:O2	2.20	0.41
1:AAA:605:GLU:OE2	1:AAA:605:GLU:HA	2.20	0.41
1:AAA:748:VAL:CG1	1:AAA:778:ALA:HA	2.49	0.40

All (1) 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)
3:AAA:1021:HOH:O	3:AAA:1040:HOH:O[8_555]	2.11	0.09

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	AAA	241/240 (100%)	239 (99%)	2 (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	AAA	189/186 (102%)	187 (99%)	2 (1%)	65	50

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	AAA	540	ASP
1	AAA	723	MET

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

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](#)

10 ligands are modelled in this entry.

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	$\# Z > 2$
2	SO4	AAA	801	-	4,4,4	0.24	0	6,6,6	0.22	0
2	SO4	AAA	802	-	4,4,4	0.32	0	6,6,6	0.12	0
2	SO4	AAA	806	-	4,4,4	0.34	0	6,6,6	0.11	0
2	SO4	AAA	808	-	4,4,4	0.25	0	6,6,6	0.12	0
2	SO4	AAA	810	-	4,4,4	0.29	0	6,6,6	0.15	0
2	SO4	AAA	803	-	4,4,4	0.27	0	6,6,6	0.14	0
2	SO4	AAA	804	-	4,4,4	0.29	0	6,6,6	0.14	0
2	SO4	AAA	807	-	4,4,4	0.32	0	6,6,6	0.14	0
2	SO4	AAA	805	-	4,4,4	0.31	0	6,6,6	0.14	0
2	SO4	AAA	809	-	4,4,4	0.34	0	6,6,6	0.07	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	AAA	806	SO4	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	AAA	808	SO4	2	0
2	AAA	810	SO4	1	0

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 [i](#)

6.1 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	AAA	240/240 (100%)	0.52	14 (5%)	29 35	14, 34, 61, 79	3 (1%)

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	AAA	688	PHE	4.1
1	AAA	539	SER	3.7
1	AAA	754	PHE	3.5
1	AAA	778	ALA	3.3
1	AAA	562	ALA	3.0
1	AAA	777	ALA	2.5
1	AAA	547	ALA	2.5
1	AAA	571	VAL	2.4
1	AAA	540	ASP	2.3
1	AAA	776	LYS	2.2
1	AAA	775	GLN	2.1
1	AAA	748	VAL	2.1
1	AAA	567	ASP	2.1
1	AAA	564	LYS	2.0

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 oligosaccharides 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. 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(Å ²)	Q<0.9
2	SO4	AAA	806	5/5	0.50	0.16	110,117,120,132	0
2	SO4	AAA	809	5/5	0.50	0.14	119,125,130,134	0
2	SO4	AAA	803	5/5	0.70	0.14	59,61,80,88	0
2	SO4	AAA	810	5/5	0.73	0.16	78,78,93,102	0
2	SO4	AAA	805	5/5	0.76	0.13	55,60,69,69	0
2	SO4	AAA	802	5/5	0.79	0.12	52,52,62,71	0
2	SO4	AAA	804	5/5	0.79	0.11	65,65,70,72	0
2	SO4	AAA	807	5/5	0.82	0.14	73,75,84,85	0
2	SO4	AAA	808	5/5	0.82	0.11	50,53,62,63	0
2	SO4	AAA	801	5/5	0.90	0.08	47,48,58,62	0

6.5 Other polymers [i](#)

There are no such residues in this entry.