



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 7, 2018 – 11:20 AM EST

PDB ID : 5N0W
Title : Crystal structure of OphA-DeltaC6 mutant R72A in complex with SAM
Authors : Song, H.; Naismith, J.H.
Deposited on : 2017-02-03
Resolution : 1.93 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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20030736
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20030736

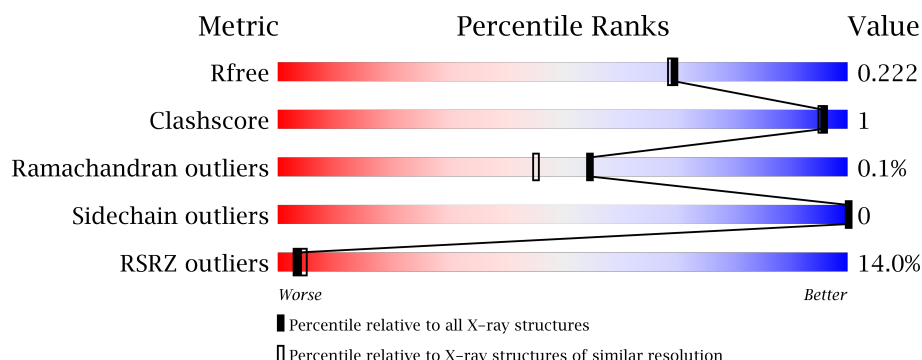
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.93 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}	100719	3233 (1.96-1.92)
Clashscore	112137	3430 (1.96-1.92)
Ramachandran outliers	110173	3395 (1.96-1.92)
Sidechain outliers	110143	3395 (1.96-1.92)
RSRZ outliers	101464	3250 (1.96-1.92)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	410	<div> <div>13%</div> <div>93%</div> <div>• •</div> </div>
1	B	410	<div> <div>14%</div> <div>93%</div> <div>• 6%</div> </div>

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	GOL	B	502	-	-	-	X



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			6	3	3		

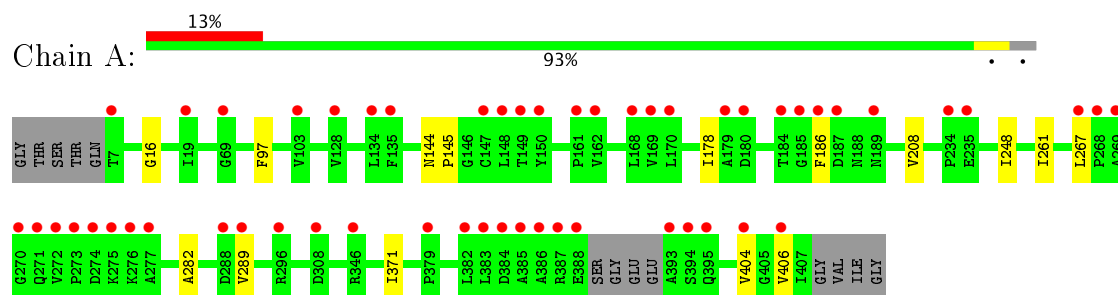
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	276	Total	O	0	0
			276	276		
4	B	258	Total	O	0	0
			258	258		

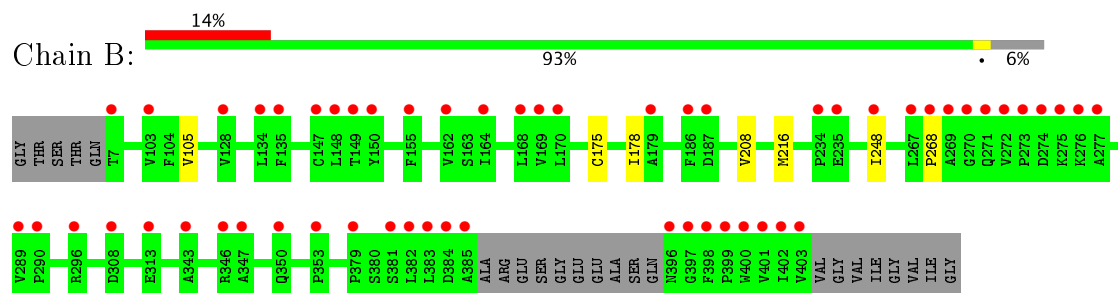
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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: Peptide N-methyltransferase



- Molecule 1: Peptide N-methyltransferase



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 ₁ 2 ₁ 2	Depositor
Cell constants a, b, c, α , β , γ	163.10 Å 92.28 Å 85.68 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.80 – 1.93 49.75 – 1.93	Depositor EDS
% Data completeness (in resolution range)	98.7 (49.80-1.93) 98.7 (49.75-1.93)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.76 (at 1.92 Å)	Xtriage
Refinement program	REFMAC 5.8.0158	Depositor
R, R_{free}	0.194 , 0.216 0.201 , 0.222	Depositor DCC
R_{free} test set	4905 reflections (5.34%)	DCC
Wilson B-factor (Å ²)	35.4	Xtriage
Anisotropy	0.126	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 51.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6707	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 49.80 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 7.0836e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² 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.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, SAM

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	0.43	0/3185	0.65	0/4335
1	B	0.43	0/3072	0.65	0/4181
All	All	0.43	0/6257	0.65	0/8516

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	3113	0	3073	10	0
1	B	3000	0	2967	8	0
2	A	27	0	22	0	0
2	B	27	0	22	0	0
3	B	6	0	8	1	0
4	A	276	0	0	0	2
4	B	258	0	0	0	0
All	All	6707	0	6092	16	2

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 1.

All (16) 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:371:ILE:HG21	1:B:216:MET:HE1	1.79	0.65
1:B:208:VAL:HG22	1:B:248:ILE:HD12	1.84	0.58
1:A:404:VAL:HG12	1:A:406:VAL:HG22	1.84	0.58
1:B:105:VAL:HG13	3:B:502:GOL:H2	1.85	0.57
1:A:261:ILE:HG21	1:A:267:LEU:HD12	1.91	0.53
1:A:178:ILE:HD13	1:A:186:PHE:HB3	1.90	0.53
1:A:371:ILE:HD13	1:B:216:MET:HE1	1.92	0.52
1:A:208:VAL:HG22	1:A:248:ILE:HD12	1.91	0.52
1:B:208:VAL:HG13	1:B:248:ILE:CD1	2.46	0.46
1:A:208:VAL:HG13	1:A:248:ILE:CD1	2.45	0.46
1:B:208:VAL:HG22	1:B:248:ILE:CD1	2.46	0.46
1:B:175:CYS:HA	1:B:178:ILE:HD12	1.99	0.45
1:A:144:ASN:HA	1:A:145:PRO:HA	1.88	0.43
1:B:208:VAL:HG13	1:B:248:ILE:HD13	2.00	0.43
1:A:282:ALA:HB1	1:A:289:VAL:HG21	2.00	0.43
1:A:16:GLY:HA2	1:A:97:PHE:O	2.20	0.42

All (2) 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)
4:A:854:HOH:O	4:A:854:HOH:O[2_755]	1.93	0.27
4:A:868:HOH:O	4:A:868:HOH:O[2_755]	2.19	0.01

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	398/410 (97%)	382 (96%)	16 (4%)	0	100	100
1	B	383/410 (93%)	366 (96%)	16 (4%)	1 (0%)	44	33

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	781/820 (95%)	748 (96%)	32 (4%)	1 (0%)	55 45

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	268	PRO

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	333/337 (99%)	333 (100%)	0	100	100
1	B	321/337 (95%)	321 (100%)	0	100	100
All	All	654/674 (97%)	654 (100%)	0	100	100

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

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

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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry

3 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	SAM	A	501	-	21,29,29	1.47	2 (9%)	17,42,42	2.16	3 (17%)
2	SAM	B	501	-	21,29,29	1.50	3 (14%)	17,42,42	2.18	2 (11%)
3	GOL	B	502	-	5,5,5	0.87	0	5,5,5	1.03	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
2	SAM	A	501	-	-	0/8/33/33	0/3/3/3
2	SAM	B	501	-	-	0/8/33/33	0/3/3/3
3	GOL	B	502	-	-	0/4/4/4	0/0/0/0

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	SAM	CG-SD	-5.27	1.69	1.80
2	A	501	SAM	CG-SD	-4.51	1.71	1.80
2	B	501	SAM	O4'-C1'	2.01	1.44	1.41
2	B	501	SAM	C5-C4	2.30	1.45	1.40
2	A	501	SAM	C5-C4	2.90	1.47	1.40

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	SAM	N3-C2-N1	-7.98	121.90	128.86
2	A	501	SAM	N3-C2-N1	-6.68	123.04	128.86
2	A	501	SAM	C4-C5-N7	-2.02	107.46	109.41
2	B	501	SAM	C2-N1-C6	2.55	123.23	118.77
2	A	501	SAM	C4'-O4'-C1'	3.46	113.46	109.77

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	502	GOL	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 ⓘ

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	397/410 (96%)	0.73	54 (13%) 3 5	25, 42, 84, 129	0
1	B	387/410 (94%)	0.77	56 (14%) 3 4	25, 41, 82, 121	0
All	All	784/820 (95%)	0.75	110 (14%) 3 4	25, 41, 83, 129	0

All (110) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	269	ALA	13.0
1	B	272	VAL	10.5
1	B	269	ALA	10.3
1	B	270	GLY	9.4
1	A	273	PRO	9.4
1	B	268	PRO	8.4
1	A	272	VAL	8.0
1	A	268	PRO	7.7
1	A	186	PHE	7.2
1	B	273	PRO	6.7
1	B	403	VAL	6.6
1	B	271	GLN	6.1
1	A	270	GLY	5.9
1	A	7	THR	5.8
1	A	382	LEU	5.7
1	B	397	GLY	5.6
1	B	400	TRP	5.5
1	A	394	SER	5.1
1	A	388	GLU	5.1
1	B	385	ALA	5.0
1	A	271	GLN	5.0
1	A	383	LEU	4.9
1	B	186	PHE	4.9
1	A	406	VAL	4.9

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Mol	Chain	Res	Type	RSRZ
1	A	387	ARG	4.9
1	A	386	ALA	4.8
1	A	393	ALA	4.8
1	B	401	VAL	4.6
1	A	267	LEU	4.5
1	A	276	LYS	4.3
1	A	395	GLN	4.0
1	A	179	ALA	3.9
1	B	276	LYS	3.8
1	B	179	ALA	3.7
1	B	343	ALA	3.7
1	B	399	PRO	3.7
1	A	277	ALA	3.7
1	B	383	LEU	3.6
1	B	402	ILE	3.6
1	A	147	CYS	3.6
1	A	148	LEU	3.6
1	A	274	ASP	3.5
1	A	69	GLY	3.5
1	B	148	LEU	3.5
1	B	277	ALA	3.4
1	A	235	GLU	3.4
1	A	162	VAL	3.4
1	A	308	ASP	3.4
1	B	346	ARG	3.3
1	A	384	ASP	3.3
1	B	187	ASP	3.3
1	A	234	PRO	3.3
1	B	162	VAL	3.2
1	B	350	GLN	3.1
1	B	275	LYS	3.1
1	B	290	PRO	3.1
1	A	103	VAL	3.0
1	A	187	ASP	3.0
1	B	353	PRO	3.0
1	B	267	LEU	3.0
1	A	135	PHE	3.0
1	B	235	GLU	2.9
1	A	346	ARG	2.9
1	A	184	THR	2.9
1	B	164	ILE	2.9
1	B	396	ASN	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	234	PRO	2.8
1	B	168	LEU	2.7
1	A	379	PRO	2.7
1	B	149	THR	2.7
1	A	189	ASN	2.7
1	B	147	CYS	2.7
1	B	347	ALA	2.6
1	B	103	VAL	2.6
1	B	398	PHE	2.6
1	B	150	TYR	2.6
1	A	404	VAL	2.6
1	A	385	ALA	2.5
1	A	161	PRO	2.5
1	B	274	ASP	2.5
1	B	169	VAL	2.4
1	B	308	ASP	2.4
1	A	180	ASP	2.3
1	A	168	LEU	2.3
1	A	169	VAL	2.3
1	B	296	ARG	2.3
1	A	275	LYS	2.3
1	B	381	SER	2.3
1	B	170	LEU	2.3
1	B	135	PHE	2.2
1	A	170	LEU	2.2
1	A	128	VAL	2.2
1	A	288	ASP	2.2
1	B	379	PRO	2.2
1	A	289	VAL	2.2
1	B	128	VAL	2.2
1	A	19	ILE	2.1
1	B	7	THR	2.1
1	B	155	PHE	2.1
1	B	384	ASP	2.1
1	B	248	ILE	2.1
1	B	313	GLU	2.1
1	A	149	THR	2.1
1	B	134	LEU	2.1
1	A	134	LEU	2.0
1	B	382	LEU	2.0
1	A	150	TYR	2.0
1	A	296	ARG	2.0

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Mol	Chain	Res	Type	RSRZ
1	B	289	VAL	2.0
1	A	185	GLY	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
3	GOL	B	502	6/6	0.76	0.27	5.41	40,43,45,45	0
2	SAM	B	501	27/27	0.95	0.15	0.04	25,28,32,36	0
2	SAM	A	501	27/27	0.94	0.15	-0.16	28,30,36,40	0

6.5 Other polymers [i](#)

There are no such residues in this entry.