



wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 7, 2020 – 03:33 AM BST

PDB ID : 6RXJ
Title : Crystal structure of CobB wt in complex with H4K16-Acetyl peptide
Authors : Spinck, M.; Gasper, R.; Neumann, H.
Deposited on : 2019-06-08
Resolution : 1.60 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.13.1
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.13.1

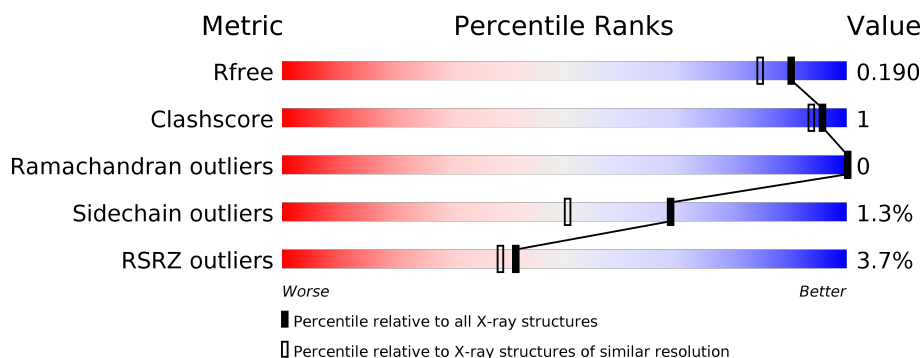
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.60 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}	130704	3398 (1.60-1.60)
Clashscore	141614	3665 (1.60-1.60)
Ramachandran outliers	138981	3564 (1.60-1.60)
Sidechain outliers	138945	3563 (1.60-1.60)
RSRZ outliers	127900	3321 (1.60-1.60)

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 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	254	<div> <div>3%</div> <div> <div></div> <div>89%</div> <div>7%</div> </div> </div>
1	B	254	<div> <div>4%</div> <div> <div></div> <div>89%</div> <div>7%</div> </div> </div>
2	C	10	<div> <div></div> <div> <div>70%</div> <div>30%</div> </div> </div>
2	D	10	<div> <div>10%</div> <div> <div></div> <div>70%</div> <div>10%</div> <div>20%</div> </div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7800 atoms, of which 3710 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 NAD-dependent protein deacylase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	235	Total	C	H	N	O	S	0	0	0
			3624	1162	1789	325	339	9			
1	B	235	Total	C	H	N	O	S	0	1	0
			3634	1165	1794	326	340	9			

There are 78 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-14	MET	-	initiating methionine	UNP P75960
A	-13	GLY	-	expression tag	UNP P75960
A	-12	SER	-	expression tag	UNP P75960
A	-11	SER	-	expression tag	UNP P75960
A	-10	HIS	-	expression tag	UNP P75960
A	-9	HIS	-	expression tag	UNP P75960
A	-8	HIS	-	expression tag	UNP P75960
A	-7	HIS	-	expression tag	UNP P75960
A	-6	HIS	-	expression tag	UNP P75960
A	-5	HIS	-	expression tag	UNP P75960
A	-4	SER	-	expression tag	UNP P75960
A	-3	GLN	-	expression tag	UNP P75960
A	-2	ASP	-	expression tag	UNP P75960
A	-1	PRO	-	expression tag	UNP P75960
A	255	TYR	-	expression tag	UNP P75960
A	256	GLY	-	expression tag	UNP P75960
A	257	PRO	-	expression tag	UNP P75960
A	258	ALA	-	expression tag	UNP P75960
A	259	SER	-	expression tag	UNP P75960
A	260	GLN	-	expression tag	UNP P75960
A	261	VAL	-	expression tag	UNP P75960
A	262	VAL	-	expression tag	UNP P75960
A	263	PRO	-	expression tag	UNP P75960
A	264	GLU	-	expression tag	UNP P75960
A	265	PHE	-	expression tag	UNP P75960

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Chain	Residue	Modelled	Actual	Comment	Reference
A	266	VAL	-	expression tag	UNP P75960
A	267	GLU	-	expression tag	UNP P75960
A	268	LYS	-	expression tag	UNP P75960
A	269	LEU	-	expression tag	UNP P75960
A	270	LEU	-	expression tag	UNP P75960
A	271	LYS	-	expression tag	UNP P75960
A	272	GLY	-	expression tag	UNP P75960
A	273	LEU	-	expression tag	UNP P75960
A	274	LYS	-	expression tag	UNP P75960
A	275	ALA	-	expression tag	UNP P75960
A	276	GLY	-	expression tag	UNP P75960
A	277	SER	-	expression tag	UNP P75960
A	278	ILE	-	expression tag	UNP P75960
A	279	ALA	-	expression tag	UNP P75960
B	-14	MET	-	initiating methionine	UNP P75960
B	-13	GLY	-	expression tag	UNP P75960
B	-12	SER	-	expression tag	UNP P75960
B	-11	SER	-	expression tag	UNP P75960
B	-10	HIS	-	expression tag	UNP P75960
B	-9	HIS	-	expression tag	UNP P75960
B	-8	HIS	-	expression tag	UNP P75960
B	-7	HIS	-	expression tag	UNP P75960
B	-6	HIS	-	expression tag	UNP P75960
B	-5	HIS	-	expression tag	UNP P75960
B	-4	SER	-	expression tag	UNP P75960
B	-3	GLN	-	expression tag	UNP P75960
B	-2	ASP	-	expression tag	UNP P75960
B	-1	PRO	-	expression tag	UNP P75960
B	255	TYR	-	expression tag	UNP P75960
B	256	GLY	-	expression tag	UNP P75960
B	257	PRO	-	expression tag	UNP P75960
B	258	ALA	-	expression tag	UNP P75960
B	259	SER	-	expression tag	UNP P75960
B	260	GLN	-	expression tag	UNP P75960
B	261	VAL	-	expression tag	UNP P75960
B	262	VAL	-	expression tag	UNP P75960
B	263	PRO	-	expression tag	UNP P75960
B	264	GLU	-	expression tag	UNP P75960
B	265	PHE	-	expression tag	UNP P75960
B	266	VAL	-	expression tag	UNP P75960
B	267	GLU	-	expression tag	UNP P75960
B	268	LYS	-	expression tag	UNP P75960

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Chain	Residue	Modelled	Actual	Comment	Reference
B	269	LEU	-	expression tag	UNP P75960
B	270	LEU	-	expression tag	UNP P75960
B	271	LYS	-	expression tag	UNP P75960
B	272	GLY	-	expression tag	UNP P75960
B	273	LEU	-	expression tag	UNP P75960
B	274	LYS	-	expression tag	UNP P75960
B	275	ALA	-	expression tag	UNP P75960
B	276	GLY	-	expression tag	UNP P75960
B	277	SER	-	expression tag	UNP P75960
B	278	ILE	-	expression tag	UNP P75960
B	279	ALA	-	expression tag	UNP P75960

- Molecule 2 is a protein called Histone H4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	7	Total	C	H	N	O	0	0	0
			114	33	57	16	8			
2	D	8	Total	C	H	N	O	0	0	0
			136	39	70	18	9			

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Zn	0	0
			1	1		
3	A	1	Total	Zn	0	0
			1	1		

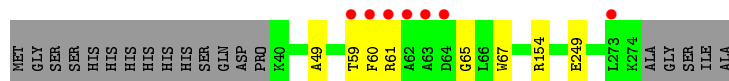
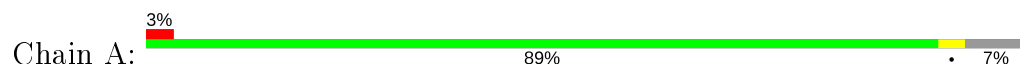
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	134	Total	O	0	0
			134	134		
4	B	149	Total	O	0	0
			149	149		
4	C	4	Total	O	0	0
			4	4		
4	D	3	Total	O	0	0
			3	3		

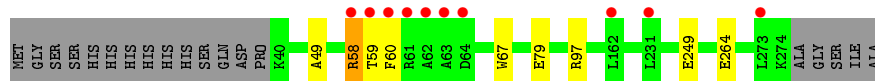
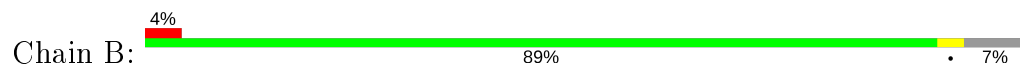
3 Residue-property plots [i](#)

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: NAD-dependent protein deacylase



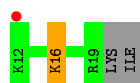
- Molecule 1: NAD-dependent protein deacylase



- Molecule 2: Histone H4



- Molecule 2: Histone H4



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	131.10Å 131.26Å 58.54Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.38 – 1.60 46.38 – 1.60	Depositor EDS
% Data completeness (in resolution range)	99.9 (46.38-1.60) 99.5 (46.38-1.60)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.36 (at 1.60Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.173 , 0.188 0.175 , 0.190	Depositor DCC
R_{free} test set	2013 reflections (3.01%)	wwPDB-VP
Wilson B-factor (Å ²)	27.8	Xtriage
Anisotropy	0.222	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 40.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.439 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	7800	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.01% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.41	0/1880	0.55	0/2553
1	B	0.40	0/1888	0.54	0/2564
2	C	0.32	0/44	0.64	0/54
2	D	0.33	0/53	0.81	0/65
All	All	0.40	0/3865	0.55	0/5236

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
2	D	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	D	16	ALY	Mainchain

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	A	1835	1789	1788	3	0
1	B	1840	1794	1794	5	0
2	C	57	57	57	0	0
2	D	66	70	70	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	134	0	0	0	0
4	B	149	0	0	3	0
4	C	4	0	0	0	0
4	D	3	0	0	0	0
All	All	4090	3710	3709	8	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 1.

The worst 5 of 8 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:79:GLU:OE2	4:B:401:HOH:O	1.80	0.99
1:B:264:GLU:OE2	4:B:402:HOH:O	2.09	0.69
1:B:60:PHE:HB3	1:B:67:TRP:HB2	1.90	0.53
1:A:60:PHE:HB3	1:A:67:TRP:HB2	1.91	0.51
1:A:61:ARG:O	1:A:65:GLY:N	2.45	0.49

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 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	233/254 (92%)	232 (100%)	1 (0%)	0	100	100
1	B	234/254 (92%)	233 (100%)	1 (0%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	C	4/10 (40%)	4 (100%)	0	0	100	100
2	D	5/10 (50%)	5 (100%)	0	0	100	100
All	All	476/528 (90%)	474 (100%)	2 (0%)	0	100	100

There are no Ramachandran outliers to report.

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	193/208 (93%)	191 (99%)	2 (1%)	76	61
1	B	194/208 (93%)	191 (98%)	3 (2%)	65	44
2	C	3/6 (50%)	3 (100%)	0	100	100
2	D	4/6 (67%)	4 (100%)	0	100	100
All	All	394/428 (92%)	389 (99%)	5 (1%)	69	50

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	154	ARG
1	A	249	GLU
1	B	58	ARG
1	B	97	ARG
1	B	249	GLU

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 ⓘ

2 non-standard protein/DNA/RNA residues 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	ALY	C	16	2	10,11,12	0.81	0	7,12,14	0.73	0
2	ALY	D	16	2	10,11,12	1.43	1 (10%)	7,12,14	0.72	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	ALY	C	16	2	-	0/9/10/12	-
2	ALY	D	16	2	-	0/9/10/12	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	16	ALY	O-C	3.80	1.35	1.19

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

5.6 Ligand geometry

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

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.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	A	235/254 (92%)	0.06	7 (2%) 50 48	21, 32, 59, 104	0
1	B	235/254 (92%)	0.04	10 (4%) 35 32	21, 31, 61, 99	0
2	C	6/10 (60%)	0.21	0 100 100	27, 33, 48, 60	0
2	D	7/10 (70%)	0.59	1 (14%) 2 2	26, 34, 63, 64	0
All	All	483/528 (91%)	0.06	18 (3%) 41 39	21, 32, 61, 104	0

The worst 5 of 18 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	63	ALA	8.8
1	A	62	ALA	7.6
1	B	64	ASP	6.4
1	B	62	ALA	6.1
1	A	61	ARG	5.7

6.2 Non-standard residues in protein, DNA, RNA chains [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	ALY	C	16	12/13	0.94	0.08	21,27,37,37	0
2	ALY	D	16	12/13	0.97	0.08	21,26,37,39	0

6.3 Carbohydrates [i](#)

There are no monosaccharides 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(\AA^2)	Q<0.9
3	ZN	A	301	1/1	0.98	0.11	27,27,27,27	1
3	ZN	B	301	1/1	0.99	0.07	32,32,32,32	0

6.5 Other polymers [i](#)

There are no such residues in this entry.