



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:44 PM UTC

PDB ID : 5INT / pdb_00005int
Title : Crystal structure of the C-terminal Domain of Coenzyme A biosynthesis bi-functional protein CoaBC
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Deposited on : 2016-03-07
Resolution : 2.15 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

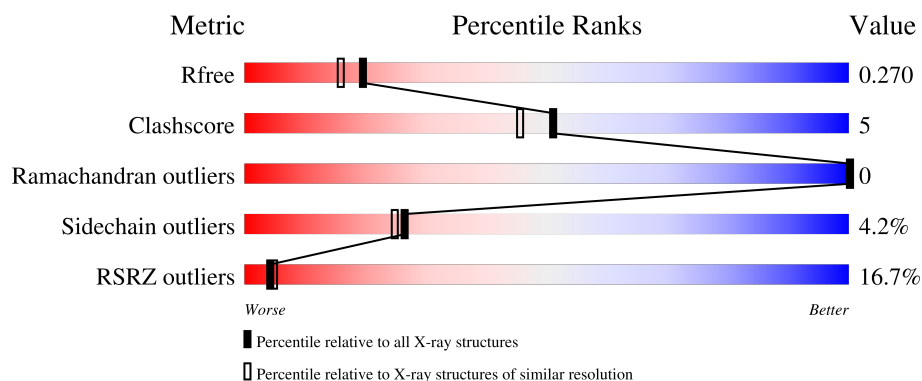
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 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	2057 (2.16-2.16)
Clashscore	190562	2159 (2.16-2.16)
Ramachandran outliers	187476	2134 (2.16-2.16)
Sidechain outliers	187428	2133 (2.16-2.16)
RSRZ outliers	180081	2059 (2.16-2.16)

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	220	13% 75% 14% 11%
1	B	220	15% 77% 13% 10%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3207 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 Phosphopantothenate--cysteine ligase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	196	Total	C	N	O	Se	0	0	0
			1494	947	252	287	8			
1	B	198	Total	C	N	O	Se	0	0	0
			1469	930	246	285	8			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	182	SER	-	expression tag	UNP Q81WG9
A	183	ASN	-	expression tag	UNP Q81WG9
A	184	ALA	-	expression tag	UNP Q81WG9
B	182	SER	-	expression tag	UNP Q81WG9
B	183	ASN	-	expression tag	UNP Q81WG9
B	184	ALA	-	expression tag	UNP Q81WG9

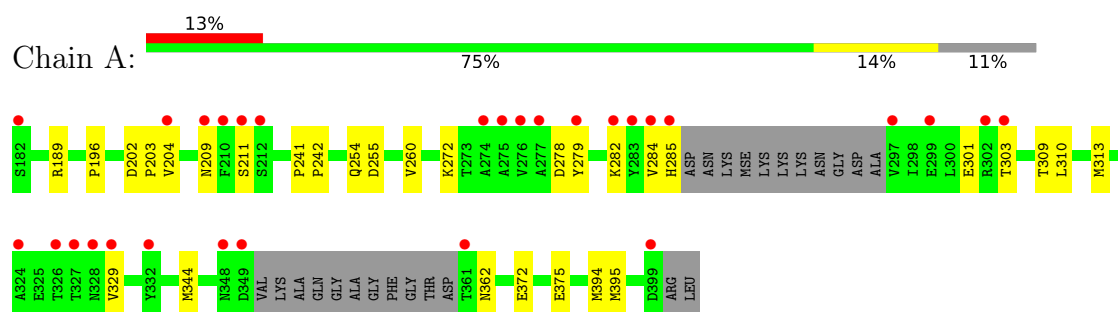
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	131	Total	O	0	1
			132	132		
2	B	112	Total	O	0	0
			112	112		

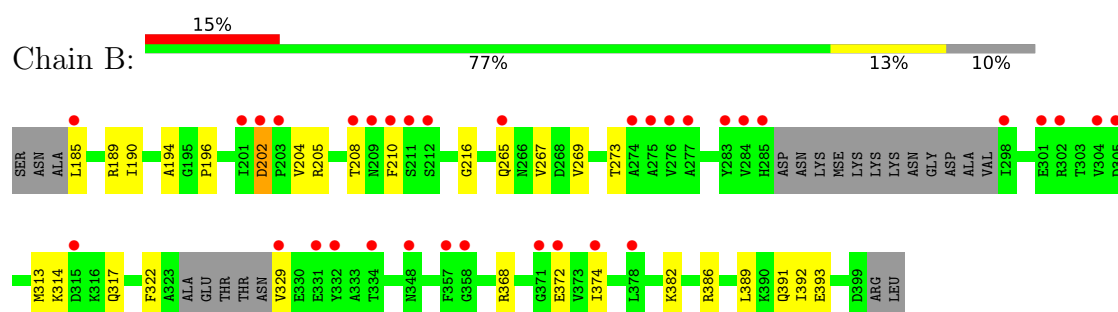
3 Residue-property plots [i](#)

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: Phosphopantothenate--cysteine ligase



- Molecule 1: Phosphopantothenate--cysteine ligase



4 Data and refinement statistics

Property	Value	Source
Space group	P 4 21 2	Depositor
Cell constants a, b, c, α , β , γ	133.17Å 133.17Å 55.77Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	32.30 – 2.15 32.30 – 2.15	Depositor EDS
% Data completeness (in resolution range)	95.2 (32.30-2.15) 95.7 (32.30-2.15)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.67 (at 2.16Å)	Xtriage
Refinement program	PHENIX dev_1888	Depositor
R, R_{free}	0.226 , 0.258 (Not available) , 0.270	Depositor DCC
R_{free} test set	1360 reflections (4.88%)	wwPDB-VP
Wilson B-factor (Å ²)	21.8	Xtriage
Anisotropy	0.037	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 50.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	3207	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.40% 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](#)

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	0.27	0/1501	0.71	0/2020
1	B	0.26	0/1476	0.71	0/1992
All	All	0.26	0/2977	0.71	0/4012

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 [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	1494	0	1524	16	0
1	B	1469	0	1464	16	0
2	A	132	0	0	3	0
2	B	112	0	0	2	0
All	All	3207	0	2988	31	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 5.

All (31) 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:282:LYS:HG2	1:A:301:GLU:HB3	1.77	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:254:GLN:HG3	1:B:313:MSE:HE3	1.78	0.66
1:A:209:ASN:HD21	1:A:279:TYR:H	1.48	0.61
1:B:329:VAL:N	2:B:506:HOH:O	2.35	0.59
1:A:344:MSE:HB2	1:A:395:MSE:HE1	1.84	0.57
1:A:278:ASP:HB2	1:A:303:THR:HG21	1.87	0.57
1:A:202:ASP:HB2	1:A:203:PRO:HD2	1.88	0.56
1:B:202:ASP:HB2	1:B:205:ARG:HB2	1.89	0.54
1:A:362:ASN:ND2	2:A:512:HOH:O	2.42	0.52
1:A:255:ASP:OD2	2:A:501:HOH:O	2.20	0.50
1:B:389:LEU:HD23	1:B:392:ILE:HD12	1.92	0.50
1:B:204:VAL:HG23	1:B:205:ARG:HG3	1.93	0.49
1:B:185:LEU:N	2:B:512:HOH:O	2.46	0.48
1:A:362:ASN:ND2	2:A:503:HOH:O	2.47	0.47
1:B:196:PRO:HB3	1:B:210:PHE:HA	1.96	0.47
1:B:368:ARG:HD3	1:B:372:GLU:HB3	1.98	0.46
1:B:194:ALA:HB2	1:B:216:GLY:HA3	1.98	0.46
1:A:209:ASN:ND2	1:A:279:TYR:H	2.12	0.46
1:B:189:ARG:HB2	1:B:267:VAL:HA	1.98	0.45
1:A:209:ASN:N	1:A:209:ASN:OD1	2.50	0.45
1:A:196:PRO:HB3	1:A:211:SER:H	1.83	0.44
1:B:368:ARG:HD2	1:B:391:GLN:NE2	2.32	0.44
1:A:260:VAL:HG11	1:A:310:LEU:HD11	1.99	0.43
1:B:374:ILE:HD12	1:B:391:GLN:NE2	2.33	0.43
1:B:314:LYS:HE3	1:B:317:GLN:O	2.20	0.42
1:A:241:PRO:HA	1:A:242:PRO:HD3	1.92	0.41
1:B:190:ILE:HG12	1:B:269:VAL:HB	2.01	0.41
1:A:309:THR:O	1:A:313:MSE:HG2	2.21	0.41
1:B:273:THR:HA	1:B:322:PHE:CD2	2.57	0.40
1:B:382:LYS:HE3	1:B:386:ARG:HH22	1.86	0.40
1:A:209:ASN:HD22	1:A:278:ASP:HA	1.86	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	190/220 (86%)	184 (97%)	6 (3%)	0	100	100
1	B	192/220 (87%)	186 (97%)	6 (3%)	0	100	100
All	All	382/440 (87%)	370 (97%)	12 (3%)	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	160/177 (90%)	151 (94%)	9 (6%)	19	15
1	B	153/177 (86%)	149 (97%)	4 (3%)	40	43
All	All	313/354 (88%)	300 (96%)	13 (4%)	26	25

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	189	ARG
1	A	204	VAL
1	A	272	LYS
1	A	284	VAL
1	A	285	HIS
1	A	329	VAL
1	A	372	GLU
1	A	375	GLU
1	A	394	MSE
1	B	202	ASP
1	B	208	THR
1	B	265	GLN
1	B	393	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	391	GLN

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	188/220 (85%)	0.81	29 (15%) 5 5	7, 28, 68, 81	0
1	B	190/220 (86%)	0.99	34 (17%) 3 4	13, 38, 66, 83	0
All	All	378/440 (85%)	0.90	63 (16%) 4 5	7, 33, 67, 83	0

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	274	ALA	7.1
1	B	275	ALA	6.4
1	A	275	ALA	6.0
1	A	284	VAL	5.5
1	B	202	ASP	5.4
1	B	298	ILE	5.2
1	A	283	TYR	5.2
1	A	276	VAL	5.1
1	B	274	ALA	5.0
1	A	277	ALA	4.9
1	B	283	TYR	4.9
1	A	326	THR	4.7
1	A	361	THR	4.5
1	B	276	VAL	4.4
1	A	297	VAL	4.4
1	A	327	THR	4.2
1	A	209	ASN	3.9
1	A	399	ASP	3.9
1	B	284	VAL	3.8
1	B	212	SER	3.8
1	A	211	SER	3.7
1	B	332	TYR	3.5
1	A	285	HIS	3.5
1	B	357	PHE	3.5

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Mol	Chain	Res	Type	RSRZ
1	A	279	TYR	3.5
1	B	210	PHE	3.4
1	A	348	ASN	3.4
1	A	212	SER	3.4
1	B	201	ILE	3.2
1	A	332	TYR	3.1
1	B	285	HIS	3.1
1	B	329	VAL	3.1
1	B	209	ASN	3.0
1	A	210	PHE	2.9
1	A	349	ASP	2.9
1	A	329	VAL	2.8
1	B	371	GLY	2.8
1	B	358	GLY	2.7
1	B	374	ILE	2.7
1	A	328	ASN	2.7
1	A	282	LYS	2.6
1	A	302	ARG	2.6
1	B	208	THR	2.6
1	B	277	ALA	2.5
1	B	315	ASP	2.4
1	B	331	GLU	2.4
1	A	324	ALA	2.4
1	A	299	GLU	2.3
1	A	204	VAL	2.3
1	B	304	VAL	2.3
1	A	303	THR	2.3
1	B	302	ARG	2.2
1	B	372	GLU	2.2
1	B	334	THR	2.2
1	B	265	GLN	2.1
1	B	348	ASN	2.1
1	B	203	PRO	2.1
1	B	301	GLU	2.1
1	B	211	SER	2.1
1	B	378	LEU	2.0
1	B	305	ASP	2.0
1	A	182	SER	2.0
1	B	185	LEU	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.