



wwPDB X-ray Structure Validation Summary Report ⓘ

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PDB ID : 4EMY
Title : Crystal structure of aminotransferase from anaerococcus prevotii dsm 20548.
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Deposited on : 2012-04-12
Resolution : 2.86 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 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	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.21
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.004 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.41.4

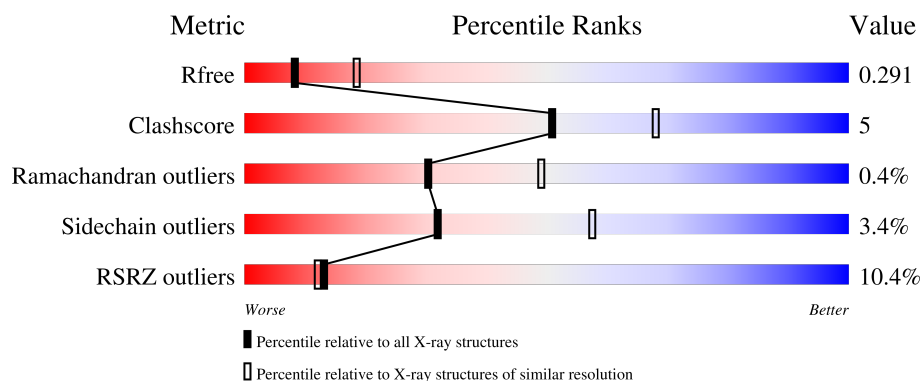
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION





The reported resolution of this entry is 2.86 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}	164625	1268 (2.88-2.84)
Clashscore	180529	1351 (2.88-2.84)
Ramachandran outliers	177936	1318 (2.88-2.84)
Sidechain outliers	177891	1319 (2.88-2.84)
RSRZ outliers	164620	1269 (2.88-2.84)

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	413	
1	B	413	
1	C	413	
1	D	413	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 13055 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 Aminotransferase class I and II.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	409	Total	C	N	O	S	Se	0	0	0
			3249	2080	538	621	4	6			
1	B	409	Total	C	N	O	S	Se	0	0	0
			3249	2080	538	621	4	6			
1	C	407	Total	C	N	O	S	Se	0	0	0
			3230	2065	536	619	4	6			
1	D	407	Total	C	N	O	S	Se	0	0	0
			3235	2073	536	616	4	6			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	UNP C7REB0
A	-1	ASN	-	expression tag	UNP C7REB0
A	0	ALA	-	expression tag	UNP C7REB0
B	-2	SER	-	expression tag	UNP C7REB0
B	-1	ASN	-	expression tag	UNP C7REB0
B	0	ALA	-	expression tag	UNP C7REB0
C	-2	SER	-	expression tag	UNP C7REB0
C	-1	ASN	-	expression tag	UNP C7REB0
C	0	ALA	-	expression tag	UNP C7REB0
D	-2	SER	-	expression tag	UNP C7REB0
D	-1	ASN	-	expression tag	UNP C7REB0
D	0	ALA	-	expression tag	UNP C7REB0

- Molecule 2 is PYRIDOXAL-5'-PHOSPHATE (three-letter code: PLP) (formula: C₈H₁₀NO₆P).



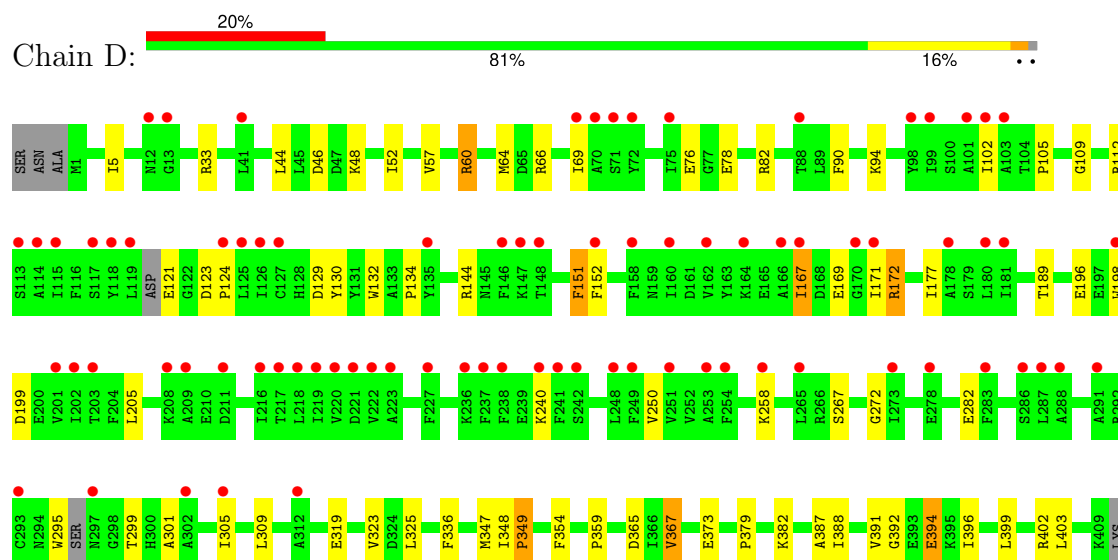
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			16	8	1	6	1		
2	B	1	Total	C	N	O	P	0	0
			16	8	1	6	1		
2	C	1	Total	C	N	O	P	0	0
			16	8	1	6	1		
2	D	1	Total	C	N	O	P	0	0
			16	8	1	6	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	7	Total	O	0	0
			7	7		
3	B	10	Total	O	0	0
			10	10		
3	C	4	Total	O	0	0
			4	4		
3	D	7	Total	O	0	0
			7	7		



● Molecule 1: Aminotransferase class I and II



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	64.30Å 131.18Å 239.91Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.86 50.00 – 2.86	Depositor EDS
% Data completeness (in resolution range)	99.4 (50.00-2.86) 99.8 (50.00-2.86)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.04 (at 2.86Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.245 , 0.290 0.246 , 0.291	Depositor DCC
R_{free} test set	2418 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	50.4	Xtriage
Anisotropy	0.960	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 37.3	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	0.92	EDS
Total number of atoms	13055	wwPDB-VP
Average B, all atoms (Å ²)	75.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.05% 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: PLP

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	0.42	0/3315	0.55	0/4470
1	B	0.40	0/3315	0.54	0/4470
1	C	0.41	1/3293 (0.0%)	0.51	0/4437
1	D	0.41	1/3299 (0.0%)	0.51	0/4445
All	All	0.41	2/13222 (0.0%)	0.53	0/17822

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	10	TRP	CD2-CE2	5.11	1.47	1.41
1	D	198	TRP	CD2-CE2	5.05	1.47	1.41

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	3249	0	3200	46	0
1	B	3249	0	3200	32	0
1	C	3230	0	3178	21	0
1	D	3235	0	3189	40	0
2	A	16	0	7	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	16	0	8	0	0
2	C	16	0	8	0	0
2	D	16	0	8	2	0
3	A	7	0	0	1	0
3	B	10	0	0	1	0
3	C	4	0	0	0	0
3	D	7	0	0	0	0
All	All	13055	0	12798	132	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.

The worst 5 of 132 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:C:228:ALA:HA	1:C:352:GLY:HA3	1.46	0.94
1:A:112:ARG:HD2	1:B:294:ASN:HA	1.64	0.78
1:D:394:GLU:CD	1:D:394:GLU:H	1.89	0.75
1:A:111:ILE:HD13	1:A:135:TYR:HE1	1.57	0.70
1:B:105:PRO:HD2	1:B:109:GLY:HA3	1.74	0.69

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	407/413 (98%)	386 (95%)	20 (5%)	1 (0%)	44 63
1	B	407/413 (98%)	391 (96%)	14 (3%)	2 (0%)	25 43
1	C	401/413 (97%)	370 (92%)	29 (7%)	2 (0%)	25 43
1	D	401/413 (97%)	375 (94%)	24 (6%)	2 (0%)	25 43

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	1616/1652 (98%)	1522 (94%)	87 (5%)	7 (0%)	30	49

5 of 7 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	33	ARG
1	C	381	ALA
1	B	296	SER
1	D	172	ARG
1	D	387	ALA

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	344/341 (101%)	334 (97%)	10 (3%)	37	63
1	B	344/341 (101%)	332 (96%)	12 (4%)	31	57
1	C	342/341 (100%)	331 (97%)	11 (3%)	34	60
1	D	342/341 (100%)	329 (96%)	13 (4%)	28	54
All	All	1372/1364 (101%)	1326 (97%)	46 (3%)	32	57

5 of 46 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	278	GLU
1	D	151	PHE
1	C	296	SER
1	C	346	THR
1	D	196	GLU

Sometimes 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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

4 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	PLP	A	501	-	16,16,16	1.30	1 (6%)	20,23,23	1.29	3 (15%)
2	PLP	C	501	-	16,16,16	1.28	2 (12%)	20,23,23	1.24	3 (15%)
2	PLP	D	501	-	16,16,16	1.26	2 (12%)	20,23,23	1.39	2 (10%)
2	PLP	B	501	-	16,16,16	1.30	1 (6%)	20,23,23	1.44	4 (20%)

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	PLP	A	501	-	-	5/8/8/8	0/1/1/1
2	PLP	C	501	-	-	2/8/8/8	0/1/1/1
2	PLP	D	501	-	-	2/8/8/8	0/1/1/1
2	PLP	B	501	-	-	4/8/8/8	0/1/1/1

The worst 5 of 6 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	501	PLP	C2-N1	3.14	1.39	1.33
2	C	501	PLP	C2-N1	3.06	1.39	1.33
2	B	501	PLP	C2-N1	2.97	1.39	1.33
2	A	501	PLP	C2-N1	2.84	1.38	1.33
2	D	501	PLP	C6-N1	2.33	1.39	1.34

The worst 5 of 12 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	PLP	C5-C6-N1	-2.80	119.27	123.83
2	D	501	PLP	C3-C4-C5	2.80	120.52	118.28
2	D	501	PLP	C5-C6-N1	-2.76	119.34	123.83
2	A	501	PLP	C3-C4-C4A	-2.65	116.20	119.84
2	C	501	PLP	C3-C4-C5	2.43	120.23	118.28

There are no chirality outliers.

5 of 13 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	PLP	C3-C4-C4A-O4A
2	A	501	PLP	C5A-O4P-P-O1P
2	A	501	PLP	C5A-O4P-P-O2P
2	B	501	PLP	C3-C4-C4A-O4A
2	B	501	PLP	C5A-O4P-P-O3P

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	PLP	1	0
2	D	501	PLP	2	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	A	403/413 (97%)	-0.36	0 100 100	34, 48, 64, 86	0
1	B	403/413 (97%)	-0.35	2 (0%) 87 86	33, 46, 70, 96	0
1	C	401/413 (97%)	1.27	84 (20%) 3 3	61, 108, 144, 233	0
1	D	401/413 (97%)	1.17	81 (20%) 3 4	50, 95, 130, 220	0
All	All	1608/1652 (97%)	0.43	167 (10%) 13 12	33, 67, 136, 233	0

The worst 5 of 167 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	384	ILE	7.2
1	D	70	ALA	7.2
1	C	17	ALA	5.4
1	D	241	PHE	5.0
1	C	363	ALA	4.4

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 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
2	PLP	C	501	16/16	0.79	0.14	59,68,74,79	0
2	PLP	D	501	16/16	0.88	0.10	38,46,57,57	0
2	PLP	A	501	16/16	0.96	0.09	23,28,33,42	0
2	PLP	B	501	16/16	0.96	0.09	30,37,43,50	0

6.5 Other polymers [i](#)

There are no such residues in this entry.