WORLD WIDE
wwPDB
PROTEIN DATA BANK

wwPDB X-ray Structure Validation Summary Report (i)

Jun 15, 2024 – 10:02 PM EDT

PDB ID	:	2AH1
Title	:	Crystal structure of aromatic amine dehydrogenase (AADH) from Alcaligenes faecalis
Authors	:	Masgrau, L.; Roujeinikova, A.; Johannissen, L.O.; Hothi, P.; Basran, J.; Ranaghan, K.E.; Mulholland, A.J.; Sutcliffe, M.J.; Scrutton, N.S.; Leys, D.
Deposited on	:	2005-07-27
Resolution	:	1.20 Å(reported)

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

|		
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
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)	:	2.37.1

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 1.20 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.
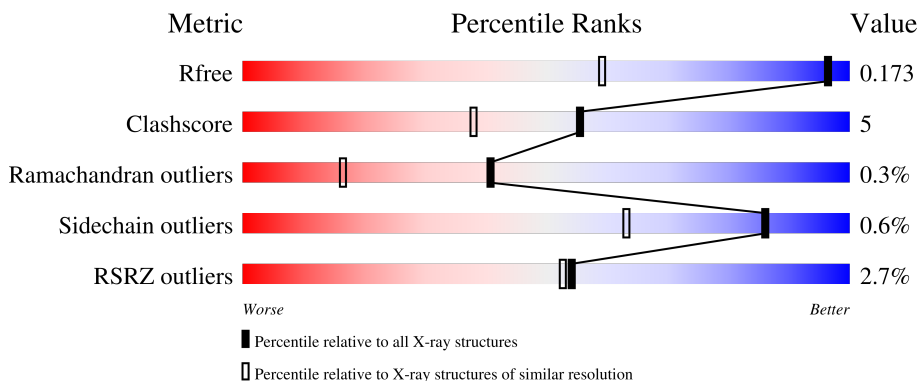
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1223 (1.22-1.18)
Clashscore	141614	1286 (1.22-1.18)
Ramachandran outliers	138981	1240 (1.22-1.18)
Sidechain outliers	138945	1239 (1.22-1.18)
RSRZ outliers	127900	1200 (1.22-1.18)

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 <=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	D	135	2% / 73% / 8% · 19%
1	H	135	3% / 70% / 10% · 19%
2	A	361	% / 89% / 10% ··
2	B	361	4% / 89% / 9% ·

2 Entry composition ⓘ

There are 3 unique types of molecules in this entry. The entry contains 9013 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 Aromatic amine dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	D	110	Total	C	N	O	S	0	0	0
			840	513	147	165	15			
1	H	110	Total	C	N	O	S	0	1	0
			845	516	148	166	15			

- Molecule 2 is a protein called Aromatic amine dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	359	Total	C	N	O	S	0	1	0
			2797	1764	486	534	13			
2	B	360	Total	C	N	O	S	0	1	0
			2787	1758	482	534	13			

- Molecule 3 is water.

| Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	193	Total	O	0	0
			193	193		
3	H	202	Total	O	0	0
			202	202		
3	A	711	Total	O	0	0
			711	711		
3	B	638	Total	O	0	0
			638	638		

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.
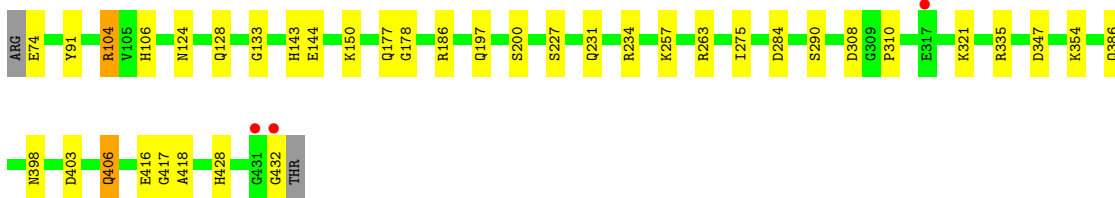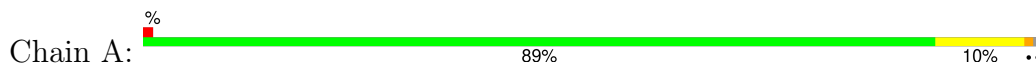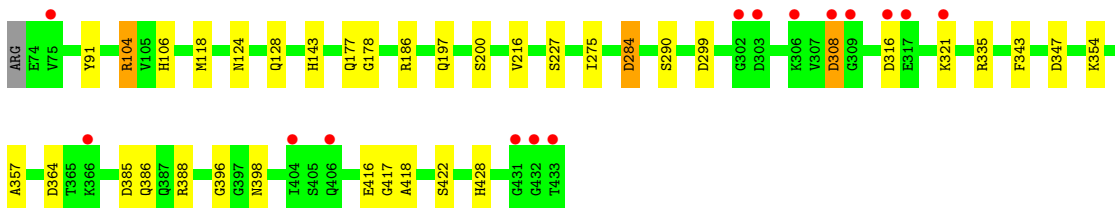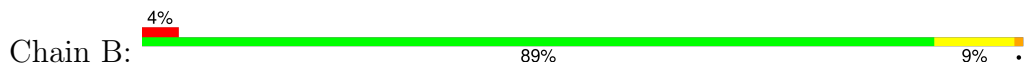
- Molecule 1: Aromatic amine dehydrogenase

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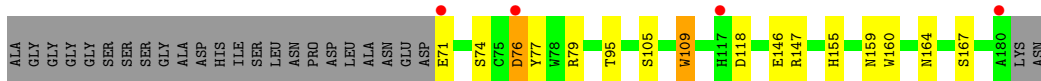
- Molecule 2: Aromatic amine dehydrogenase

- Molecule 2: Aromatic amine dehydrogenase

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	70.78Å 89.14Å 80.32Å 90.00° 90.23° 90.00°	Depositor
Resolution (Å)	12.00 – 1.20 12.00 – 1.20	Depositor EDS
% Data completeness (in resolution range)	100.0 (12.00-1.20) 98.4 (12.00-1.20)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.19 (at 1.20Å)	Xtriage
Refinement program	REFMAC 5.1.9999	Depositor
R, R_{free}	0.142 , 0.170 0.144 , 0.173	Depositor DCC
R_{free} test set	15336 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å^2)	13.6	Xtriage
Anisotropy	0.238	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 48.6	EDS
L-test for twinning[2]	$< |L| > = 0.52, < L^2 > = 0.36$	Xtriage
Estimated twinning fraction	0.010 for h,-k,-l	Xtriage
F_o,F_c correlation	0.98	EDS
Total number of atoms	9013	wwPDB-VP
Average B, all atoms (Å^2)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.16% of the height of the origin peak. No significant pseudotranslation is detected.*

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[1]Intensities estimated from amplitudes.
[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: TRQ

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	D	0.62	0/846	0.83	2/1151 (0.2%)
1	H	0.61	0/854	0.84	3/1163 (0.3%)
2	A	0.61	1/2864 (0.0%)	0.82	3/3884 (0.1%)
2	B	0.54	0/2856	0.79	7/3874 (0.2%)
All	All	0.58	1/7420 (0.0%)	0.81	15/10072 (0.1%)

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	A	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	144	GLU	CD-OE1	5.12	1.31	1.25

The worst 5 of 15 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	D	79	ARG	NE-CZ-NH2	-6.26	117.17	120.30
2	A	308	ASP	CB-CG-OD2	6.21	123.89	118.30
2	A	403	ASP	CB-CG-OD1	6.07	123.76	118.30
2	B	347	ASP	CB-CG-OD2	5.99	123.69	118.30
1	H	76	ASP	CB-CG-OD2	5.68	123.41	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	A	310	PRO	Peptide

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	D	840	0	736	10	0
1	H	845	0	744	13	0
2	A	2797	0	2731	31	0
2	B	2787	0	2701	24	0
3	A	711	0	0	12	3
3	B	638	0	0	5	2
3	D	193	0	0	3	1
3	H	202	0	0	5	0
All	All	9013	0	6912	72	3

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 72 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)
2:A:74:GLU:N	3:A:735:HOH:O	2.05	0.90
1:D:95:THR:HG22	3:D:271:HOH:O	1.80	0.80
3:H:275:HOH:O	2:B:118:MET:SD	2.40	0.80
2:A:104:ARG:HH11	2:A:106:HIS:HE1	1.30	0.77
2:A:124:ASN:HD21	2:A:178:GLY:H	1.33	0.76

All (3) 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:A:831:HOH:O	3:B:1010:HOH:O[1_655]	2.08	0.12
3:A:1093:HOH:O	3:B:705:HOH:O[1_655]	2.14	0.06
3:D:344:HOH:O	3:A:827:HOH:O[1_554]	2.19	0.01

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	D	107/135 (79%)	104 (97%)	3 (3%)	0	100	100
1	H	108/135 (80%)	106 (98%)	2 (2%)	0	100	100
2	A	358/361 (99%)	346 (97%)	11 (3%)	1 (0%)	41	16
2	B	358/361 (99%)	344 (96%)	12 (3%)	2 (1%)	25	5
All	All	931/992 (94%)	900 (97%)	28 (3%)	3 (0%)	41	16

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	216	VAL
2	A	200	SER
2	B	200	SER

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	D	94/112 (84%)	94 (100%)	0	100	100
1	H	96/112 (86%)	96 (100%)	0	100	100
2	A	303/305 (99%)	300 (99%)	3 (1%)	76	47
2	B	301/305 (99%)	298 (99%)	3 (1%)	76	47
All	All	794/834 (95%)	788 (99%)	6 (1%)	86	55

5 of 6 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	104	ARG
2	B	308	ASP
2	B	388	ARG
2	A	406[A]	GLN
2	A	104	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 42 such sidechains are listed below:

Mol	Chain	Res	Type
2	B	124	ASN
2	B	231	GLN
2	B	128	GLN
2	B	180	ASN
2	B	386	GLN

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	#$|Z| > 2$
1	TRQ	D	109	1	13,17,18	4.83	4 (30%)	13,24,26	2.79	5 (38%)
1	TRQ	H	109	1	13,17,18	5.18	6 (46%)	13,24,26	2.75	6 (46%)

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
1	TRQ	D	109	1	-	0/4/19/21	0/2/2/2
1	TRQ	H	109	1	-	0/4/19/21	0/2/2/2

The worst 5 of 10 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	H	109	TRQ	CH2-CZ2	-15.43	1.35	1.53
1	D	109	TRQ	CH2-CZ2	-14.38	1.36	1.53
1	H	109	TRQ	CE2-CZ2	-8.59	1.39	1.50
1	D	109	TRQ	CE2-CZ2	-7.85	1.40	1.50
1	H	109	TRQ	O7-CZ2	3.93	1.31	1.23

The worst 5 of 11 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	H	109	TRQ	CZ2-CE2-NE1	5.75	129.12	119.94
1	D	109	TRQ	CZ2-CE2-NE1	5.74	129.10	119.94
1	D	109	TRQ	O7-CZ2-CH2	5.44	125.04	119.01
1	H	109	TRQ	O7-CZ2-CH2	4.41	123.90	119.01
1	H	109	TRQ	CB-CG-CD1	-3.74	123.35	127.97

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	D	109	TRQ	2	0
1	H	109	TRQ	2	0

5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

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 ⓘ

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(Å²)	Q<0.9
1	D	109/135 (80%)	0.06	3 (2%)	53	51	10, 16, 22, 26	0
1	H	109/135 (80%)	0.11	4 (3%)	41	40	11, 15, 26, 34	0
2	A	359/361 (99%)	-0.12	3 (0%)	86	86	9, 13, 20, 26	0
2	B	360/361 (99%)	0.25	15 (4%)	36	36	10, 19, 28, 35	0
All	All	937/992 (94%)	0.07	25 (2%)	54	53	9, 15, 26, 35	0

The worst 5 of 25 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	433[A]	THR	4.7
2	A	431	GLY	4.5
2	B	303	ASP	4.1
2	B	306	LYS	3.9
1	H	117	HIS	3.8

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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(Å²)	Q<0.9
1	TRQ	D	109	16/17	0.95	0.08	14,15,17,23	0
1	TRQ	H	109	16/17	0.97	0.07	12,13,15,23	0

6.3 Carbohydrates ⓘ

There are no monosaccharides in this entry.

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.