Full wwPDB X-ray Structure Validation Report ⓘ

Oct 29, 2026 – 02:08 PM JST

PDB ID	:	25YD / pdb_000025yd
Title	:	Pseudomonas aeruginosa AchC2 (YfdX2) at pH 4.0
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Deposited on	:	2026-04-22
Resolution	:	1.94 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:

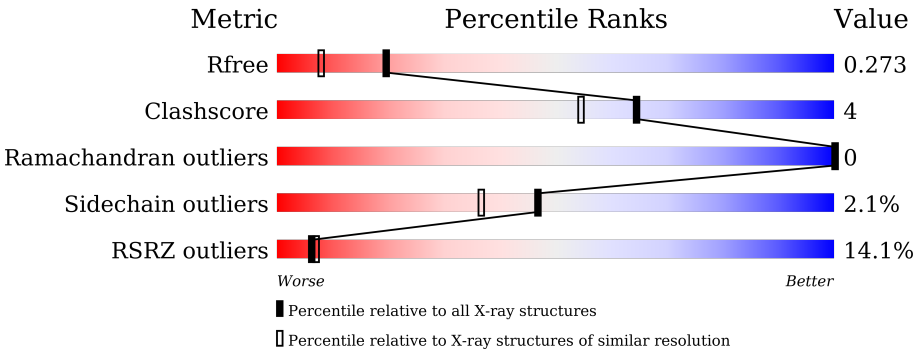
MolProbity	:	4-5-2 with Phenix2.2
Xtriage (Phenix)	:	2.2
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.51

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.94 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	1452 (1.94-1.94)
Clashscore	190562	1494 (1.94-1.94)
Ramachandran outliers	187476	1479 (1.94-1.94)
Sidechain outliers	187428	1479 (1.94-1.94)
RSRZ outliers	180081	1453 (1.94-1.94)

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	A	303	15% 76% 6% 18%
1	B	303	8% 71% 11% 18%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 3852 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 YfdX protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	248	Total	C	N	O	S	0	0	0
			1859	1166	321	370	2			
1	B	249	Total	C	N	O	S	0	0	0
			1849	1158	321	368	2			

- Molecule 2 is water.

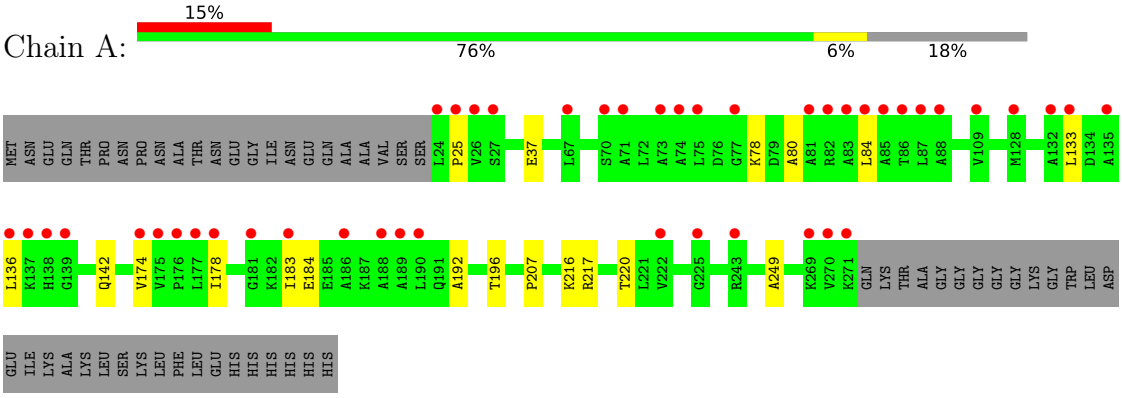
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	57	Total	O	0	0
			57	57		
2	B	87	Total	O	0	0
			87	87		

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.

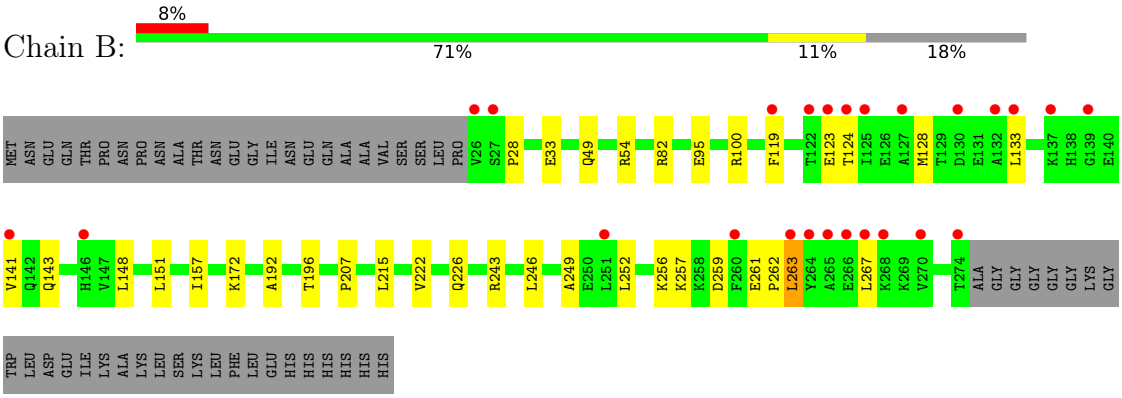
• Molecule 1: YfdX protein

Chain A:



• Molecule 1: YfdX protein

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	41.30Å 110.60Å 121.89Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.38 – 1.94 29.38 – 1.94	Depositor EDS
% Data completeness (in resolution range)	99.6 (29.38-1.94) 99.6 (29.38-1.94)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.73 (at 1.93Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.252 , 0.274 0.252 , 0.273	Depositor DCC
R_{free} test set	2107 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å^2)	38.7	Xtriage
Anisotropy	0.085	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 29.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	3852	wwPDB-VP
Average B, all atoms (Å^2)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.94% of the height of the origin peak. No significant pseudotranslation is detected.*

[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

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	0.22	0/1878	0.48	1/2555 (0.0%)
1	B	0.24	0/1867	0.49	1/2542 (0.0%)
All	All	0.23	0/3745	0.48	2/5097 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	25	PRO	N-CA-CB	10.46	110.24	103.23
1	B	28	PRO	N-CA-CB	7.69	110.16	103.31

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	1859	0	1900	11	0
1	B	1849	0	1874	22	0
2	A	57	0	0	0	0
2	B	87	0	0	6	0
All	All	3852	0	3774	29	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 4.

All (29) 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:54:ARG:NH2	2:B:401:HOH:O	2.10	0.85
1:B:49:GLN:HG2	1:B:157:ILE:HG13	1.71	0.72
1:B:172:LYS:NZ	2:B:402:HOH:O	2.21	0.72
1:B:243:ARG:NH1	2:B:404:HOH:O	2.27	0.67
1:B:95:GLU:OE2	2:B:402:HOH:O	2.14	0.64
1:B:256:LYS:N	1:B:259:ASP:OD2	2.27	0.62
1:B:100:ARG:NH1	2:B:405:HOH:O	2.33	0.57
1:A:80:ALA:HB1	1:A:178:ILE:HD11	1.86	0.57
1:A:207:PRO:HB2	1:A:249:ALA:HB2	1.89	0.55
1:B:246:LEU:HD21	1:B:263:LEU:HD23	1.90	0.54
1:B:49:GLN:CG	1:B:157:ILE:HG13	2.40	0.51
1:B:215:LEU:HD13	1:B:267:LEU:CB	2.40	0.51
1:A:136:LEU:HD21	1:B:148:LEU:HD11	1.93	0.51
1:B:192:ALA:O	1:B:196:THR:HG23	2.13	0.49
1:A:216:LYS:O	1:A:220:THR:HG23	2.14	0.48
1:A:37:GLU:OE2	1:A:217:ARG:NH1	2.48	0.47
1:A:192:ALA:O	1:A:196:THR:HG23	2.16	0.46
1:A:136:LEU:HB3	1:B:133:LEU:HG	1.99	0.45
1:A:133:LEU:HD23	1:B:133:LEU:HD23	2.00	0.44
1:B:207:PRO:HB2	1:B:249:ALA:HB2	1.99	0.43
1:A:142:GLN:HG2	1:B:252:LEU:O	2.20	0.42
1:B:222:VAL:O	1:B:226:GLN:HG2	2.18	0.42
1:B:261:GLU:N	1:B:262:PRO:HD2	2.34	0.42
1:B:151:LEU:HD23	1:B:151:LEU:HA	1.85	0.41
1:A:84:LEU:HG	1:A:178:ILE:HD13	2.01	0.41
1:A:174:VAL:O	1:A:178:ILE:HG22	2.20	0.41
1:B:123:GLU:HG2	1:B:124:THR:N	2.36	0.41
1:B:119:PHE:HZ	1:B:128:MET:HE2	1.85	0.41
1:B:82:ARG:NH1	2:B:411:HOH:O	2.52	0.40

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	246/303 (81%)	240 (98%)	6 (2%)	0	100	100
1	B	247/303 (82%)	243 (98%)	4 (2%)	0	100	100
All	All	493/606 (81%)	483 (98%)	10 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	A	194/245 (79%)	191 (98%)	3 (2%)	57	49
1	B	190/245 (78%)	185 (97%)	5 (3%)	40	28
All	All	384/490 (78%)	376 (98%)	8 (2%)	47	37

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	78	LYS
1	A	183	ILE
1	A	184	GLU
1	B	33	GLU
1	B	141	VAL
1	B	143	GLN
1	B	257	LYS
1	B	263	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	121	ASN
1	A	143	GLN

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Mol	Chain	Res	Type
1	B	49	GLN
1	B	116	HIS
1	B	244	GLN

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

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, 95^{th} 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	248/303 (81%)	1.08	45 (18%)	3	3	31, 46, 75, 82	0
1	B	249/303 (82%)	0.69	25 (10%)	12	14	24, 38, 72, 81	0
All	All	497/606 (82%)	0.88	70 (14%)	6	7	24, 43, 73, 82	0

All (70) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	265	ALA	5.7
1	A	24	LEU	4.8
1	A	270	VAL	4.7
1	A	74	ALA	4.3
1	B	26	VAL	4.0
1	B	267	LEU	3.9
1	A	183	ILE	3.8
1	A	71	ALA	3.8
1	A	133	LEU	3.7
1	B	263	LEU	3.6
1	B	274	THR	3.5
1	A	136	LEU	3.4
1	A	75	LEU	3.4
1	A	81	ALA	3.3
1	B	270	VAL	3.2
1	A	73	ALA	3.2
1	A	26	VAL	3.2
1	A	84	LEU	3.1
1	A	82	ARG	3.1
1	A	88	ALA	3.1
1	B	133	LEU	3.0
1	A	85	ALA	3.0
1	B	251	LEU	3.0
1	A	67	LEU	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	266	GLU	2.9
1	A	181	GLY	2.9
1	B	264	TYR	2.8
1	A	178	ILE	2.8
1	A	128	MET	2.8
1	B	124	THR	2.7
1	A	135	ALA	2.7
1	A	87	LEU	2.7
1	B	139	GLY	2.6
1	A	271	LYS	2.6
1	A	83	ALA	2.6
1	A	174	VAL	2.6
1	A	177	LEU	2.6
1	A	77	GLY	2.6
1	A	269	LYS	2.6
1	B	125	ILE	2.5
1	A	25	PRO	2.5
1	A	138	HIS	2.5
1	B	137	LYS	2.5
1	B	27	SER	2.5
1	B	260	PHE	2.4
1	A	86	THR	2.4
1	A	132	ALA	2.4
1	A	186	ALA	2.4
1	B	130	ASP	2.4
1	A	175	VAL	2.4
1	A	189	ALA	2.3
1	B	123	GLU	2.3
1	B	132	ALA	2.3
1	A	176	PRO	2.3
1	A	225	GLY	2.2
1	B	146	HIS	2.2
1	B	268	LYS	2.2
1	A	109	VAL	2.2
1	A	222	VAL	2.2
1	B	122	THR	2.2
1	A	243	ARG	2.2
1	B	141	VAL	2.2
1	A	190	LEU	2.1
1	B	119	PHE	2.1
1	A	27	SER	2.1
1	A	137	LYS	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	188	ALA	2.1
1	B	127	ALA	2.1
1	A	70	SER	2.0
1	A	139	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 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.