



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 06:56 PM GMT

PDB ID : 1DBX
Title : Crystal structure of cysteinyl-tRNA(Pro) deacylase from H. influenzae (HI1434)
Authors : Zhang, H.; Huang, K.; Li, Z.; Herzberg, O.; Structure 2 Function Project (S2F)
Deposited on : 1999-11-03
Resolution : 1.80 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
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

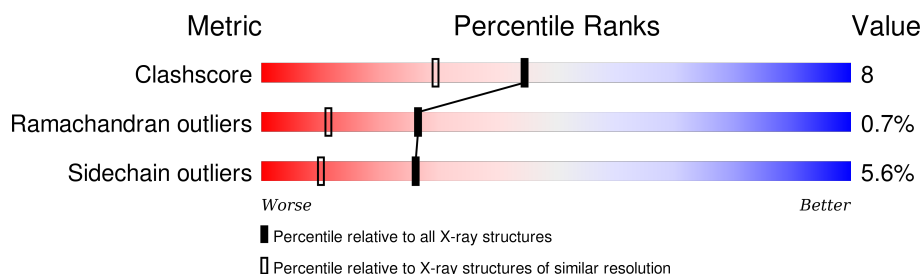
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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))
Clashscore	102246	5383 (1.80-1.80)
Ramachandran outliers	100387	5320 (1.80-1.80)
Sidechain outliers	100360	5319 (1.80-1.80)

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. 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	158	 78% 15% . . .
1	B	158	 72% 20% . . .

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2678 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 cysteinyl-tRNA(Pro) deacylase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	152	Total	C	N	O	S	Se	0	0	0
			1152	733	194	222	1	2			
1	B	152	Total	C	N	O	S	Se	0	0	0
			1152	733	194	222	1	2			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	69	MSE	MET	MODIFIED RESIDUE	UNP P45202
A	86	MSE	MET	MODIFIED RESIDUE	UNP P45202
A	117	ASN	LYS	SEE REMARK 999	UNP P45202
B	69	MSE	MET	MODIFIED RESIDUE	UNP P45202
B	86	MSE	MET	MODIFIED RESIDUE	UNP P45202
B	117	ASN	LYS	SEE REMARK 999	UNP P45202

- Molecule 2 is water.

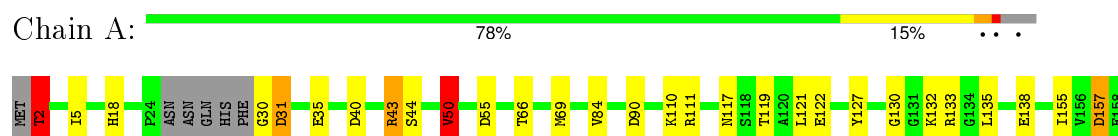
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	184	Total	O	0	0
			184	184		
2	B	190	Total	O	0	0
			190	190		

3 Residue-property plots [i](#)

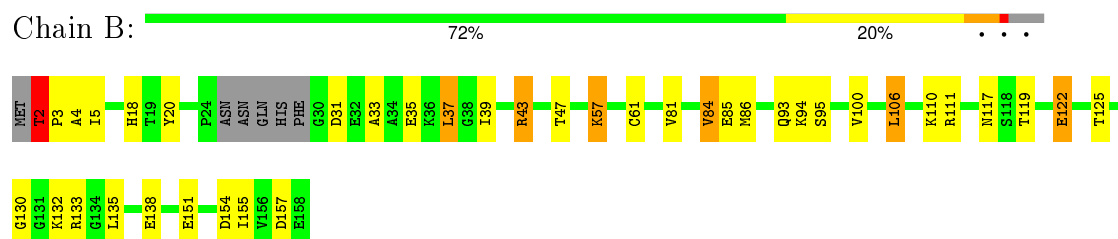
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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.

Note EDS was not executed.

- Molecule 1: cysteinyl-tRNA(Pro) deacylase



- Molecule 1: cysteinyl-tRNA(Pro) deacylase



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	32.56Å 42.21Å 54.16Å 109.07° 100.32° 90.62°	Depositor
Resolution (Å)	20.00 – 1.80	Depositor
% Data completeness (in resolution range)	97.8 (20.00-1.80)	Depositor
R_{merge}	0.03	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	REFMAC	Depositor
R, R_{free}	0.183 , 0.259	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2678	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

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	0.63	0/1164	1.28	10/1561 (0.6%)
1	B	0.67	1/1164 (0.1%)	1.33	11/1561 (0.7%)
All	All	0.65	1/2328 (0.0%)	1.31	21/3122 (0.7%)

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
1	B	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	86	MSE	SE-CE	-5.88	1.60	1.95

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	106	LEU	CA-CB-CG	9.96	138.20	115.30
1	A	111	ARG	NE-CZ-NH1	9.37	124.99	120.30
1	A	2	THR	N-CA-CB	-8.12	94.87	110.30
1	A	43	ARG	NE-CZ-NH2	-7.92	116.34	120.30
1	B	154	ASP	CB-CG-OD1	6.68	124.31	118.30
1	B	84	VAL	N-CA-CB	-6.37	97.48	111.50
1	A	50	VAL	N-CA-CB	-6.04	98.21	111.50
1	A	127	TYR	CB-CG-CD2	-5.95	117.43	121.00
1	A	50	VAL	CG1-CB-CG2	5.88	120.30	110.90
1	B	85	GLU	OE1-CD-OE2	5.81	130.28	123.30
1	B	111	ARG	NE-CZ-NH1	5.75	123.17	120.30
1	B	2	THR	N-CA-CB	-5.65	99.57	110.30
1	A	111	ARG	CD-NE-CZ	5.63	131.48	123.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	55	ASP	CB-CG-OD2	5.42	123.18	118.30
1	B	84	VAL	CA-CB-CG1	5.38	118.96	110.90
1	B	133	ARG	NE-CZ-NH2	-5.36	117.62	120.30
1	B	43	ARG	CD-NE-CZ	5.30	131.02	123.60
1	B	85	GLU	CB-CA-C	-5.29	99.83	110.40
1	B	151	GLU	CB-CG-CD	-5.21	100.14	114.20
1	A	40	ASP	CB-CG-OD2	-5.12	113.69	118.30
1	A	157	ASP	C-N-CA	5.01	134.22	121.70

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	157	ASP	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	A	1152	0	1202	15	0
1	B	1152	0	1202	22	0
2	A	184	0	0	4	0
2	B	190	0	0	3	0
All	All	2678	0	2404	37	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 8.

All (37) 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:2:THR:HG21	1:A:130:GLY:O	1.68	0.92
1:B:2:THR:HG21	1:B:130:GLY:O	1.72	0.88
1:B:93:GLN:HE21	1:B:100:VAL:H	1.23	0.84
1:B:132:LYS:HG2	1:B:135:LEU:HD12	1.77	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:31:ASP:O	1:B:35:GLU:HG3	1.97	0.65
1:B:2:THR:HG22	1:B:5:ILE:H	1.61	0.64
1:A:2:THR:HG22	1:A:5:ILE:H	1.62	0.63
1:A:132:LYS:HG2	1:A:135:LEU:HD12	1.82	0.62
1:A:117:ASN:HD22	1:A:155:ILE:H	1.48	0.60
1:A:133:ARG:HD3	2:A:263:HOH:O	2.01	0.60
1:B:18:HIS:HD2	2:B:202:HOH:O	1.85	0.59
1:A:69:MSE:SE	2:A:264:HOH:O	2.73	0.56
1:A:117:ASN:ND2	1:A:155:ILE:H	2.04	0.55
1:B:132:LYS:HG2	1:B:135:LEU:CD1	2.37	0.55
1:B:2:THR:HG23	1:B:4:ALA:H	1.72	0.55
1:A:30:GLY:O	1:A:31:ASP:HB2	2.09	0.53
1:B:119:THR:O	1:B:122:GLU:HG2	2.08	0.53
1:A:119:THR:O	1:A:122:GLU:HG2	2.10	0.52
1:B:117:ASN:ND2	1:B:155:ILE:H	2.08	0.50
1:B:93:GLN:NE2	1:B:100:VAL:H	2.03	0.49
1:A:18:HIS:HE1	1:A:138:GLU:OE1	1.95	0.49
1:B:57:LYS:HA	2:B:204:HOH:O	2.13	0.48
1:A:50:VAL:HG13	1:A:84:VAL:HB	1.96	0.48
1:A:35:GLU:HG3	2:A:284:HOH:O	2.14	0.48
1:B:43:ARG:NH2	1:B:125:THR:OG1	2.47	0.47
1:A:44:SER:O	1:A:66:THR:HG23	2.14	0.47
1:B:33:ALA:O	1:B:37:LEU:HB2	2.15	0.47
1:A:132:LYS:HG2	1:A:135:LEU:CD1	2.43	0.47
1:B:110:LYS:HD3	2:B:336:HOH:O	2.16	0.45
1:B:20:TYR:HB3	1:B:37:LEU:HD13	1.97	0.45
1:B:95:SER:O	1:B:110:LYS:HG2	2.17	0.45
1:B:37:LEU:HB3	1:B:39:ILE:HG12	1.99	0.44
1:B:47:THR:HG23	1:B:61:CYS:SG	2.58	0.43
1:A:18:HIS:HD2	2:A:209:HOH:O	2.00	0.43
1:B:18:HIS:HE1	1:B:138:GLU:OE1	2.02	0.43
1:B:117:ASN:HD22	1:B:155:ILE:H	1.69	0.41
1:B:2:THR:HA	1:B:3:PRO:HD3	1.98	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	148/158 (94%)	145 (98%)	1 (1%)	2 (1%)	14	3
1	B	148/158 (94%)	145 (98%)	3 (2%)	0	100	100
All	All	296/316 (94%)	290 (98%)	4 (1%)	2 (1%)	26	11

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	157	ASP
1	A	31	ASP

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	124/128 (97%)	118 (95%)	6 (5%)	31	14
1	B	124/128 (97%)	116 (94%)	8 (6%)	21	7
All	All	248/256 (97%)	234 (94%)	14 (6%)	26	10

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	THR
1	A	43	ARG
1	A	50	VAL
1	A	90	ASP

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Mol	Chain	Res	Type
1	A	110	LYS
1	A	121	LEU
1	B	2	THR
1	B	37	LEU
1	B	57	LYS
1	B	81	VAL
1	B	84	VAL
1	B	94	LYS
1	B	106	LEU
1	B	122	GLU

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

Mol	Chain	Res	Type
1	A	18	HIS
1	A	117	ASN
1	B	18	HIS
1	B	93	GLN
1	B	117	ASN

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

EDS was not executed - this section will therefore be empty.

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates [i](#)

EDS was not executed - this section will therefore be empty.

6.4 Ligands [i](#)

EDS was not executed - this section will therefore be empty.

6.5 Other polymers [i](#)

EDS was not executed - this section will therefore be empty.