



# Full wwPDB X-ray Structure Validation Report

Feb 26, 2014 – 04:27 PM GMT

PDB ID : 2QT3  
Title : Crystal structure of N-Isopropylammelide isopropylaminohydrolaseAtzC from Pseudomonas sp. strain ADP complexed with Zn  
Authors : Fedorov, A.A.; Fedorov, E.V.; Seffernick, J.; Wackett, L.P.; Burley, S.K.; Almo, S.C.; New York SGX Research Center for Structural Genomics (NYS-GXRC)  
Deposited on : 2007-08-01  
Resolution : 2.24 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

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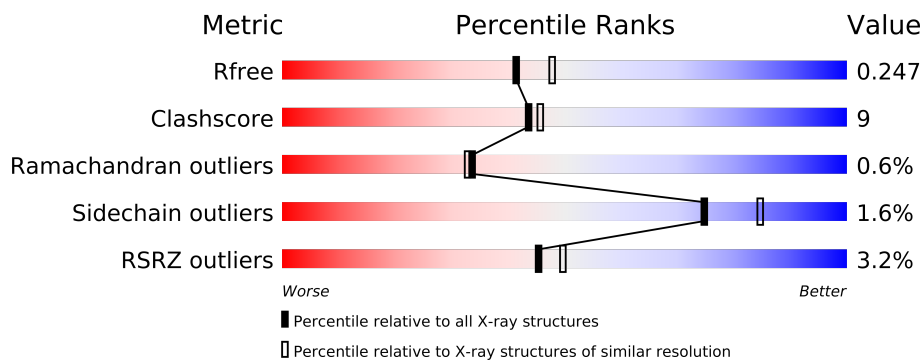
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.24 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))
$R_{free}$	66092	1112 (2.26-2.22)
Clashscore	79885	1317 (2.26-2.22)
Ramachandran outliers	78287	1282 (2.26-2.22)
Sidechain outliers	78261	1282 (2.26-2.22)
RSRZ outliers	66119	1112 (2.26-2.22)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	403	
1	B	403	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6453 atoms, of which 0 are hydrogen and 0 are deuterium.

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 N-isopropylammelide isopropyl amidohydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	401	Total	C	N	O	S	0	0	0
			3147	1998	533	603	13			
1	B	401	Total	C	N	O	S	0	0	0
			3147	1998	533	603	13			

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Zn	0	0
			1	1		
2	A	1	Total	Zn	0	0
			1	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	92	Total	O	0	0
			92	92		
3	B	65	Total	O	0	0
			65	65		



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	102.97Å 102.97Å 225.56Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	24.97 – 2.24 39.27 – 2.23	Depositor EDS
% Data completeness (in resolution range)	96.4 (24.97-2.24) 96.3 (39.27-2.23)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	6.78 (at 2.22Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.226 , 0.247 0.226 , 0.247	Depositor DCC
$R_{free}$ test set	2885 reflections (5.06%)	DCC
Wilson B-factor (Å <sup>2</sup> )	32.1	Xtriage
Anisotropy	0.146	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 35.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 60167 reflections	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	6453	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.26% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

## 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: ZN

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	0.35	0/3205	0.59	1/4335 (0.0%)
1	B	0.34	0/3205	0.58	0/4335
All	All	0.35	0/6410	0.59	1/8670 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	327	ARG	NE-CZ-NH2	-5.55	117.52	120.30

There are no chirality outliers.

There are no planarity outliers.

### 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3147	0	3150	56	0
1	B	3147	0	3150	57	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	92	0	0	2	0
3	B	65	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	6453	0	6300	109	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 9.

All (109) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:292:LEU:HD11	1:A:342:MET:HE2	1.52	0.90
1:B:292:LEU:HD11	1:B:342:MET:HE2	1.56	0.88
1:A:397:LYS:O	1:A:400:VAL:HG12	1.94	0.68
1:B:388:VAL:HB	1:B:396:VAL:HB	1.76	0.67
1:A:89:GLU:O	1:A:93:LYS:HD3	1.94	0.67
1:B:89:GLU:O	1:B:93:LYS:HD3	1.95	0.67
1:A:388:VAL:HB	1:A:396:VAL:HB	1.78	0.65
1:B:397:LYS:O	1:B:400:VAL:HG12	1.97	0.65
1:B:102:GLU:O	1:B:105:ARG:HG2	1.97	0.64
1:A:102:GLU:O	1:A:105:ARG:HG2	1.98	0.63
1:A:34:GLU:HB2	1:A:37:ILE:HG22	1.79	0.63
1:B:53:SER:HB2	1:B:54:PRO:HD2	1.81	0.62
1:A:53:SER:HB2	1:A:54:PRO:HD2	1.83	0.61
1:A:83:THR:HG23	1:A:86:ALA:HB3	1.83	0.61
1:B:83:THR:HG23	1:B:86:ALA:HB3	1.82	0.60
1:A:323:ILE:O	1:A:327:ARG:HG3	2.01	0.60
1:B:323:ILE:O	1:B:327:ARG:HG3	2.02	0.59
1:A:329:GLU:OE1	1:B:327:ARG:NH2	2.33	0.58
1:B:117:HIS:HE1	3:B:543:HOH:O	1.86	0.58
1:A:292:LEU:HD11	1:A:342:MET:CE	2.30	0.58
1:B:158:PHE:CE1	1:B:187:VAL:HG22	2.39	0.57
1:A:158:PHE:CE1	1:A:187:VAL:HG22	2.39	0.56
1:B:7:LEU:HA	1:B:43:ASP:O	2.05	0.56
1:A:116:LEU:HD21	1:A:401:ILE:HG13	1.87	0.56
1:A:288:VAL:CG1	1:A:342:MET:HE1	2.36	0.56
1:A:327:ARG:NH2	1:B:329:GLU:OE1	2.35	0.55
1:B:117:HIS:CE1	3:B:543:HOH:O	2.59	0.55
1:B:288:VAL:CG1	1:B:342:MET:HE1	2.37	0.55
1:B:292:LEU:HD11	1:B:342:MET:CE	2.33	0.55
1:B:274:LYS:HD2	1:B:350:VAL:HA	1.89	0.54
1:A:16:GLU:OE2	1:A:341:LYS:HE2	2.07	0.54
1:B:125:HIS:HE1	3:B:541:HOH:O	1.91	0.54
1:A:371:ASN:OD1	1:A:386:LEU:HD11	2.08	0.53
1:B:371:ASN:OD1	1:B:386:LEU:HD11	2.08	0.53
1:B:367:LEU:N	1:B:367:LEU:HD12	2.23	0.53

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:274:LYS:HD2	1:A:350:VAL:HA	1.89	0.53
1:B:342:MET:HA	1:B:346:GLU:HB2	1.92	0.52
1:B:116:LEU:HD21	1:B:401:ILE:HG13	1.91	0.52
1:A:367:LEU:HD12	1:A:367:LEU:N	2.24	0.51
1:A:333:ASN:ND2	1:B:75:PRO:HD2	2.26	0.51
1:B:138:GLU:HG2	1:B:179:MET:HE3	1.92	0.51
1:B:182:ASP:HB2	1:B:183:LEU:HD12	1.93	0.50
1:A:342:MET:HA	1:A:346:GLU:HB2	1.93	0.50
1:A:125:HIS:HE1	3:A:529:HOH:O	1.94	0.50
1:B:175:LYS:O	1:B:179:MET:HG3	2.11	0.50
1:B:208:LYS:HE2	1:B:239:ASN:HB3	1.93	0.50
1:A:208:LYS:HE2	1:A:239:ASN:HB3	1.93	0.49
1:A:390:LYS:HE2	3:A:510:HOH:O	2.12	0.49
1:B:402:VAL:HG12	1:B:403:ALA:N	2.28	0.49
1:B:8:ILE:CG2	1:B:37:ILE:HD11	2.43	0.49
1:A:402:VAL:HG12	1:A:403:ALA:N	2.28	0.48
1:B:248:SER:O	1:B:249:HIS:HB2	2.14	0.48
1:B:11:ASN:O	1:B:47:ALA:HB3	2.13	0.48
1:A:5:PHE:CD2	1:A:8:ILE:HD11	2.49	0.48
1:A:182:ASP:HB2	1:A:183:LEU:HD12	1.95	0.48
1:B:83:THR:HG23	1:B:86:ALA:CB	2.44	0.48
1:A:298:LEU:O	1:A:342:MET:HG3	2.14	0.47
1:A:175:LYS:O	1:A:179:MET:HG3	2.15	0.47
1:B:10:ARG:NH2	1:B:44:GLU:OE1	2.38	0.47
1:A:333:ASN:HD21	1:B:75:PRO:HD2	1.80	0.47
1:B:8:ILE:HD13	1:B:37:ILE:HD11	1.96	0.46
1:A:89:GLU:HB3	1:A:93:LYS:HE3	1.98	0.46
1:A:83:THR:HG23	1:A:86:ALA:CB	2.44	0.46
1:A:248:SER:O	1:A:249:HIS:HB2	2.15	0.46
1:A:80:ARG:HG3	1:A:80:ARG:HH11	1.81	0.46
1:A:106:HIS:HD2	1:A:109:GLU:OE1	1.99	0.45
1:A:19:SER:HB3	1:A:21:TYR:CE2	2.52	0.45
1:B:106:HIS:HD2	1:B:109:GLU:OE1	1.99	0.45
1:B:89:GLU:HB3	1:B:93:LYS:HE3	1.99	0.45
1:A:340:TRP:CG	1:A:375:PRO:HB3	2.51	0.45
1:B:340:TRP:CG	1:B:375:PRO:HB3	2.52	0.44
1:A:25:ILE:HD13	1:A:365:ALA:HB1	1.99	0.44
1:B:80:ARG:HH11	1:B:80:ARG:HG3	1.80	0.44
1:A:116:LEU:CD2	1:A:401:ILE:HG13	2.47	0.44
1:A:304:ASN:H	1:A:304:ASN:HD22	1.66	0.44
1:A:202:LEU:HD12	1:A:205:LYS:HE2	1.99	0.43
1:B:202:LEU:HD12	1:B:205:LYS:HE2	1.99	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:11:ASN:O	1:A:47:ALA:HB3	2.19	0.43
1:A:208:LYS:CE	1:A:239:ASN:HD22	2.31	0.43
1:B:220:ASP:HB3	1:B:224:VAL:HG13	2.00	0.43
1:B:7:LEU:HD21	1:B:389:ILE:HD13	2.01	0.43
1:B:374:SER:HB2	1:B:375:PRO:HD2	2.01	0.43
1:B:138:GLU:HG2	1:B:179:MET:CE	2.49	0.43
1:B:274:LYS:HG2	1:B:297:ASN:HD22	1.85	0.42
1:B:298:LEU:O	1:B:342:MET:HG3	2.19	0.42
1:B:116:LEU:CD2	1:B:401:ILE:HG13	2.49	0.42
1:B:208:LYS:CE	1:B:239:ASN:HD22	2.32	0.42
1:B:208:LYS:HE2	1:B:239:ASN:HD22	1.84	0.42
1:A:105:ARG:O	1:A:109:GLU:HG3	2.20	0.42
1:A:220:ASP:HB3	1:A:224:VAL:HG13	2.00	0.42
1:A:274:LYS:HG2	1:A:297:ASN:HD22	1.85	0.42
1:A:208:LYS:HE2	1:A:239:ASN:HD22	1.85	0.42
1:A:304:ASN:H	1:A:304:ASN:ND2	2.18	0.42
1:B:8:ILE:HA	1:B:23:ILE:O	2.20	0.42
1:B:187:VAL:HG12	1:B:188:ASP:N	2.35	0.42
1:A:7:LEU:HD21	1:A:389:ILE:HD13	2.00	0.42
1:A:179:MET:HE2	1:A:179:MET:HB3	1.76	0.41
1:B:179:MET:HB3	1:B:179:MET:HE2	1.79	0.41
1:B:374:SER:HB2	1:B:375:PRO:CD	2.50	0.41
1:B:105:ARG:O	1:B:109:GLU:HG3	2.21	0.41
1:B:187:VAL:O	1:B:188:ASP:C	2.58	0.41
1:A:187:VAL:HG12	1:A:188:ASP:N	2.36	0.41
1:A:374:SER:HB2	1:A:375:PRO:CD	2.50	0.41
1:B:8:ILE:HG23	1:B:37:ILE:HD11	2.03	0.41
1:A:374:SER:HB2	1:A:375:PRO:HD2	2.02	0.40
1:A:31:ILE:C	1:A:361:VAL:HG13	2.42	0.40
1:A:41:VAL:HG22	1:A:42:LYS:N	2.35	0.40
1:B:125:HIS:CD2	1:B:217:HIS:HE1	2.39	0.40
1:A:78:TRP:CD1	1:A:311:PRO:HB3	2.57	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	399/403 (99%)	378 (95%)	19 (5%)	2 (0%)	38	37
1	B	399/403 (99%)	376 (94%)	20 (5%)	3 (1%)	27	24
All	All	798/806 (99%)	754 (94%)	39 (5%)	5 (1%)	33	33

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	402	VAL
1	B	402	VAL
1	B	188	ASP
1	A	188	ASP
1	B	249	HIS

### 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	343/345 (99%)	337 (98%)	6 (2%)	73	83
1	B	343/345 (99%)	338 (98%)	5 (2%)	76	86
All	All	686/690 (99%)	675 (98%)	11 (2%)	75	84

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	83	THR
1	A	93	LYS
1	A	216	TYR
1	A	241	TYR
1	A	367	LEU
1	A	399	GLU
1	B	83	THR
1	B	93	LYS
1	B	216	TYR
1	B	241	TYR

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Mol	Chain	Res	Type
1	B	367	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	106	HIS
1	A	117	HIS
1	A	125	HIS
1	A	239	ASN
1	A	297	ASN
1	A	304	ASN
1	A	314	ASN
1	A	326	GLN
1	A	333	ASN
1	B	106	HIS
1	B	117	HIS
1	B	125	HIS
1	B	217	HIS
1	B	239	ASN
1	B	297	ASN
1	B	304	ASN
1	B	314	ASN
1	B	326	GLN
1	B	333	ASN

### 5.3.3 RNA ⓘ

There are no RNA chains 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 carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	401/403 (99%)	-0.23	3 (0%) 84 88	18, 33, 59, 73	0
1	B	401/403 (99%)	0.17	23 (5%) 23 26	18, 40, 58, 76	0
All	All	802/806 (99%)	-0.03	26 (3%) 45 50	18, 37, 58, 76	0

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	403	ALA	6.4
1	B	403	ALA	5.6
1	B	81	PRO	5.1
1	B	84	ARG	5.1
1	B	82	TYR	4.5
1	B	85	ASP	4.3
1	B	399	GLU	4.3
1	B	92	LEU	3.9
1	B	86	ALA	3.6
1	B	210	TYR	3.3
1	B	87	ALA	3.1
1	B	105	ARG	3.0
1	B	167	LEU	2.9
1	B	133	LYS	2.8
1	B	80	ARG	2.8
1	A	399	GLU	2.8
1	B	83	THR	2.7
1	B	95	TYR	2.7
1	B	96	LYS	2.6
1	B	175	LYS	2.3
1	A	398	ASP	2.2
1	B	176	SER	2.2
1	B	386	LEU	2.2
1	B	88	ILE	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	37	ILE	2.1
1	B	89	GLU	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

## 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands ⓘ

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	ZN	A	501	1/1	0.08	-2.14	31,31,31,31	0
2	ZN	B	502	1/1	0.06	-5.04	44,44,44,44	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.