



Full wwPDB X-ray Structure Validation Report

Mar 1, 2014 – 01:05 AM GMT

PDB ID : 3A1K
Title : Crystal structure of Rhodococcus sp. N771 Amidase
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Deposited on : 2009-04-09
Resolution : 2.17 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

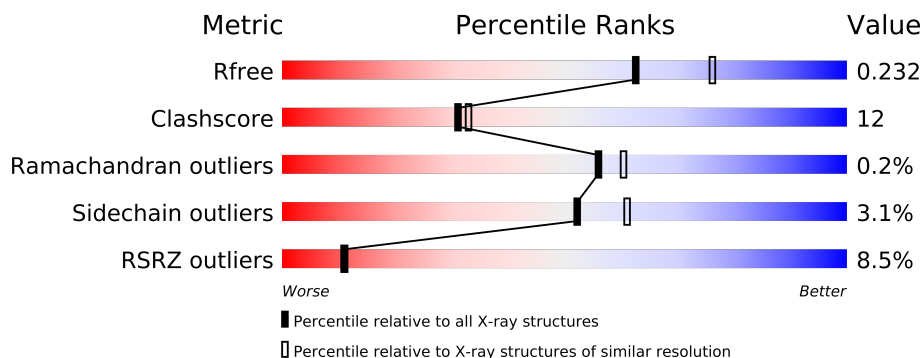
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.17 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	3841 (2.20-2.16)
Clashscore	79885	4835 (2.20-2.16)
Ramachandran outliers	78287	4740 (2.20-2.16)
Sidechain outliers	78261	4741 (2.20-2.16)
RSRZ outliers	66119	3842 (2.20-2.16)

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. 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	521	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3929 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 Amidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	508	Total	C	N	O	S	0	0	0
			3759	2354	666	725	14			

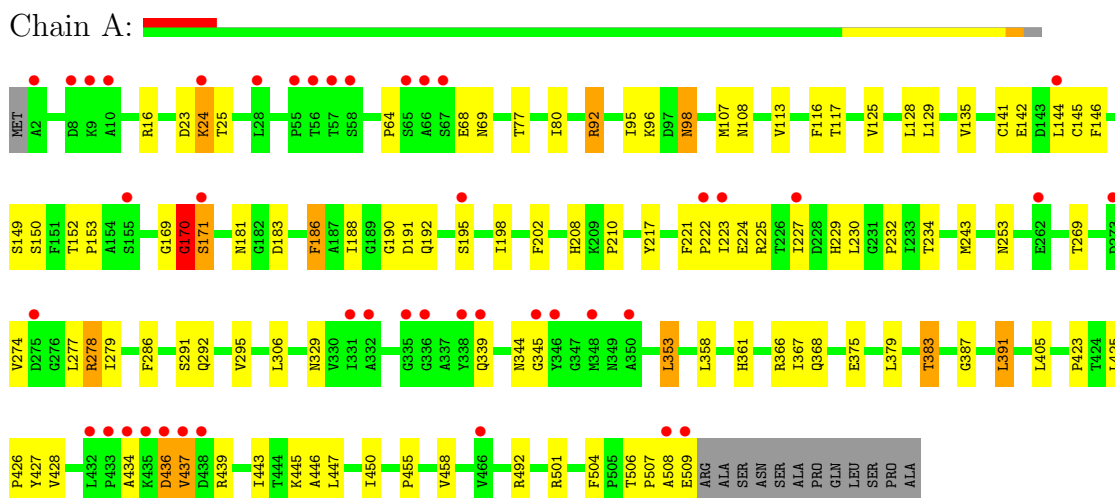
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	170	Total	O	0	0
			170	170		

3 Residue-property plots

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.

• Molecule 1: Amidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	92.67Å 92.67Å 162.31Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.17 44.56 – 2.17	Depositor EDS
% Data completeness (in resolution range)	(Not available) (50.00-2.17) 90.3 (44.56-2.17)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.56 (at 2.16Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.201 , 0.232 0.200 , 0.232	Depositor DCC
R_{free} test set	1722 reflections (4.99%)	DCC
Wilson B-factor (Å ²)	43.8	Xtriage
Anisotropy	0.327	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 46.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 37144 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3929	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.30% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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	0.42	3/3844 (0.1%)	0.63	2/5249 (0.0%)

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	A	0	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	171	SER	N-CA	12.00	1.70	1.46
1	A	170	GLY	C-O	-8.16	1.10	1.23
1	A	170	GLY	CA-C	7.60	1.64	1.51

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	170	GLY	CA-C-O	-10.19	102.25	120.60
1	A	170	GLY	CA-C-N	6.99	132.58	117.20

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	170	GLY	Mainchain

5.2 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 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	3759	0	3677	87	0
2	A	170	0	0	2	0
All	All	3929	0	3677	87	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 12.

All (87) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:171:SER:N	1:A:171:SER:CA	1.70	1.49
1:A:506:THR:HG22	1:A:508:ALA:H	1.37	0.88
1:A:186:PHE:HB3	1:A:234:THR:HG22	1.56	0.87
1:A:225:ARG:H	1:A:339:GLN:HE22	1.21	0.84
1:A:367:ILE:HD12	1:A:368:GLN:N	1.93	0.82
1:A:279:ILE:HD13	1:A:306:LEU:HD13	1.66	0.78
1:A:142:GLU:HG2	1:A:171:SER:HA	1.66	0.77
1:A:379:LEU:O	1:A:383:THR:HG23	1.89	0.71
1:A:405:LEU:HG	2:A:568:HOH:O	1.91	0.69
1:A:95:ILE:HD13	1:A:125:VAL:HG22	1.73	0.68
1:A:24:LYS:HD3	1:A:24:LYS:H	1.59	0.68
1:A:221:PHE:HE1	1:A:358:LEU:HD21	1.60	0.66
1:A:367:ILE:HD12	1:A:368:GLN:H	1.61	0.65
1:A:279:ILE:CD1	1:A:306:LEU:HD13	2.28	0.63
1:A:224:GLU:HB3	1:A:227:ILE:HG12	1.81	0.62
1:A:358:LEU:O	1:A:358:LEU:HD23	2.01	0.61
1:A:152:THR:N	1:A:153:PRO:CD	2.64	0.61
1:A:292:GLN:HE22	1:A:428:VAL:HG22	1.66	0.60
1:A:190:GLY:O	1:A:195:SER:HB2	2.02	0.59
1:A:116:PHE:HB2	1:A:361:HIS:CD2	2.37	0.59
1:A:149:SER:O	1:A:170:GLY:HA3	2.02	0.59
1:A:253:ASN:HB3	1:A:353:LEU:HB2	1.84	0.59
1:A:117:THR:O	1:A:117:THR:HG23	2.03	0.58
1:A:152:THR:N	1:A:153:PRO:HD3	2.18	0.58
1:A:142:GLU:OE2	1:A:150:SER:HA	2.04	0.58

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:434:ALA:O	1:A:437:VAL:HG22	2.03	0.58
1:A:375:GLU:OE2	1:A:439:ARG:HD3	2.05	0.57
1:A:506:THR:HG23	1:A:507:PRO:HD2	1.87	0.57
1:A:64:PRO:HB2	1:A:69:ASN:HB2	1.87	0.56
1:A:202:PHE:CE2	1:A:425:LEU:HD13	2.40	0.56
1:A:358:LEU:C	1:A:358:LEU:HD23	2.26	0.56
1:A:64:PRO:HG3	1:A:77:THR:HG23	1.89	0.55
1:A:221:PHE:CE1	1:A:358:LEU:HD21	2.40	0.54
1:A:434:ALA:O	1:A:436:ASP:N	2.41	0.54
1:A:24:LYS:N	1:A:24:LYS:HD3	2.22	0.54
1:A:198:ILE:HG23	1:A:425:LEU:HD21	1.90	0.54
1:A:269:THR:O	1:A:492:ARG:HD3	2.07	0.54
1:A:455:PRO:O	1:A:458:VAL:HG12	2.08	0.53
1:A:225:ARG:H	1:A:339:GLN:NE2	1.98	0.52
1:A:145:CYS:HB2	1:A:223:ILE:HD11	1.90	0.52
1:A:232:PRO:HB3	1:A:243:MET:SD	2.50	0.52
1:A:16:ARG:HH11	1:A:16:ARG:HG2	1.73	0.52
1:A:129:LEU:HD21	1:A:135:VAL:HG23	1.92	0.52
1:A:171:SER:N	1:A:171:SER:CB	2.63	0.51
1:A:227:ILE:O	1:A:229:HIS:HD2	1.93	0.51
1:A:24:LYS:CD	1:A:24:LYS:H	2.17	0.50
1:A:152:THR:H	1:A:153:PRO:HD3	1.76	0.50
1:A:92:ARG:HH11	1:A:92:ARG:HG3	1.76	0.50
1:A:191:ASP:HA	1:A:195:SER:HB2	1.91	0.50
1:A:170:GLY:C	1:A:171:SER:CA	2.71	0.50
1:A:98:ASN:HD22	1:A:98:ASN:H	1.60	0.50
1:A:198:ILE:HD11	1:A:450:ILE:HG22	1.94	0.49
1:A:192:GLN:HG2	1:A:223:ILE:HD12	1.94	0.49
1:A:291:SER:CB	1:A:426:PRO:HB2	2.43	0.49
1:A:387:GLY:HA2	1:A:391:LEU:HB2	1.94	0.48
1:A:278:ARG:HH21	1:A:278:ARG:HB3	1.79	0.48
1:A:222:PRO:HB2	1:A:345:GLY:O	2.14	0.47
1:A:379:LEU:HD22	1:A:443:ILE:HD13	1.97	0.47
1:A:23:ASP:OD1	1:A:25:THR:HG22	2.15	0.46
1:A:96:LYS:HG2	1:A:98:ASN:HD22	1.80	0.46
1:A:181:ASN:HB3	1:A:183:ASP:OD1	2.16	0.46
1:A:274:VAL:HB	1:A:277:LEU:HD12	1.97	0.46
1:A:169:GLY:CA	1:A:202:PHE:HD1	2.29	0.46
1:A:210:PRO:HB2	2:A:525:HOH:O	2.17	0.45
1:A:507:PRO:C	1:A:509:GLU:H	2.19	0.44
1:A:141:CYS:HB2	1:A:171:SER:OG	2.17	0.44
1:A:286:PHE:CZ	1:A:423:PRO:HD3	2.52	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:95:ILE:CD1	1:A:125:VAL:HG22	2.45	0.44
1:A:217:TYR:CD1	1:A:230:LEU:HG	2.53	0.44
1:A:80:ILE:HB	1:A:135:VAL:HB	1.99	0.43
1:A:445:LYS:HD2	1:A:445:LYS:HA	1.78	0.43
1:A:95:ILE:HD13	1:A:125:VAL:CG2	2.46	0.43
1:A:64:PRO:HB3	1:A:68:GLU:HG3	2.01	0.43
1:A:198:ILE:CD1	1:A:450:ILE:HG22	2.49	0.42
1:A:501:ARG:CZ	1:A:504:PHE:HB2	2.51	0.41
1:A:379:LEU:O	1:A:383:THR:CG2	2.64	0.41
1:A:367:ILE:H	1:A:367:ILE:HG13	1.66	0.41
1:A:344:ASN:HD21	1:A:366:ARG:HH21	1.69	0.41
1:A:295:VAL:HG21	1:A:427:TYR:HA	2.01	0.41
1:A:208:HIS:O	1:A:210:PRO:HD3	2.20	0.41
1:A:446:ALA:O	1:A:447:LEU:HD12	2.21	0.41
1:A:107:MET:HA	1:A:113:VAL:HG12	2.03	0.41
1:A:188:ILE:HD13	1:A:188:ILE:HA	1.91	0.40
1:A:150:SER:HB3	1:A:169:GLY:O	2.21	0.40
1:A:23:ASP:OD1	1:A:24:LYS:N	2.55	0.40
1:A:108:ASN:ND2	1:A:144:LEU:HA	2.37	0.40
1:A:279:ILE:HD13	1:A:306:LEU:CD1	2.45	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	506/521 (97%)	489 (97%)	16 (3%)	1 (0%)	56	60

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	436	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	389/399 (98%)	377 (97%)	12 (3%)	52 61

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	LYS
1	A	92	ARG
1	A	98	ASN
1	A	128	LEU
1	A	146	PHE
1	A	186	PHE
1	A	278	ARG
1	A	329	ASN
1	A	353	LEU
1	A	383	THR
1	A	391	LEU
1	A	437	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	98	ASN
1	A	108	ASN
1	A	229	HIS
1	A	292	GLN
1	A	339	GLN
1	A	344	ASN
1	A	401	ASN
1	A	416	GLN

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 ⓘ

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(Å ²)	Q<0.9
1	A	508/521 (97%)	0.39	43 (8%) 11 11	33, 49, 71, 92	2 (0%)

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	435	LYS	4.3
1	A	508	ALA	3.8
1	A	2	ALA	3.5
1	A	28	LEU	3.5
1	A	66	ALA	3.5
1	A	273	ASP	3.4
1	A	56	THR	3.2
1	A	437	VAL	3.2
1	A	432	LEU	3.2
1	A	24	LYS	3.1
1	A	227	ILE	3.1
1	A	223	ILE	3.0
1	A	65	SER	3.0
1	A	9	LYS	3.0
1	A	58	SER	2.9
1	A	8	ASP	2.8
1	A	436	ASP	2.8
1	A	338	TYR	2.8
1	A	350	ALA	2.8
1	A	336	GLY	2.8
1	A	331	ILE	2.7
1	A	195	SER	2.7
1	A	438	ASP	2.6
1	A	339	GLN	2.5
1	A	335	GLY	2.5
1	A	67	SER	2.4
1	A	509	GLU	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	332	ALA	2.4
1	A	55	PRO	2.4
1	A	348	MET	2.3
1	A	345	GLY	2.3
1	A	275	ASP	2.3
1	A	222	PRO	2.2
1	A	57	THR	2.2
1	A	434	ALA	2.2
1	A	346	TYR	2.2
1	A	10	ALA	2.2
1	A	144	LEU	2.1
1	A	155	SER	2.1
1	A	466	VAL	2.1
1	A	433	PRO	2.0
1	A	171	SER	2.0
1	A	262	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 ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.