Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 12:28 AM UTC

PDB ID	:	6NP1 / pdb_00006np1
Title	:	Product state mimicry leads to aminoglycoside discrimination in an antibiotic acetyltransferase
Authors	:	Kumar, P.; Cuneo, M.J.
Deposited on	:	2019-01-17
Resolution	:	1.60 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:

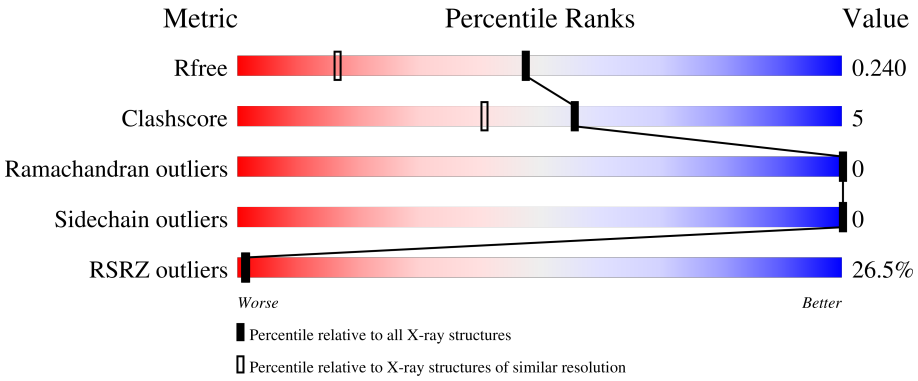
MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

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 1.60 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	4673 (1.60-1.60)
Clashscore	190562	4931 (1.60-1.60)
Ramachandran outliers	187476	4831 (1.60-1.60)
Sidechain outliers	187428	4830 (1.60-1.60)
RSRZ outliers	180081	4672 (1.60-1.60)

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	268	26% 93% 6%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 2268 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 Aminoglycoside N(3)-acetyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	268	Total	C	N	O	S	0	4	0
			2059	1298	372	383	6			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	132	SER	THR	variant	UNP Q47030

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Mg	0	0
			2	2		

- Molecule 3 is water.

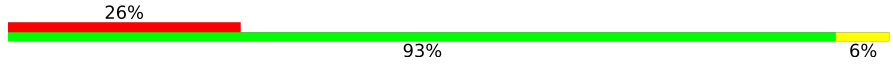
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	207	Total	O	0	0
			207	207		

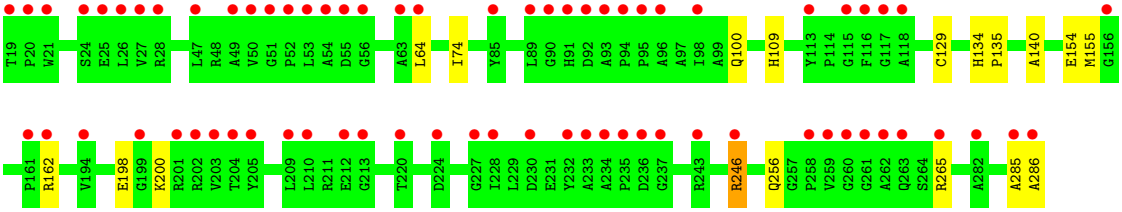
3 Residue-property plots (i)

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: Aminoglycoside N(3)-acetyltransferase

Chain A:





4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	85.35Å 86.58Å 50.35Å 90.00° 118.92° 90.00°	Depositor
Resolution (Å)	43.29 – 1.60 43.29 – 1.60	Depositor EDS
% Data completeness (in resolution range)	96.9 (43.29-1.60) 93.0 (43.29-1.60)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	5.33 (at 1.60Å)	Xtriage
Refinement program	PHENIX 1.11.1_2575	Depositor
R, R_{free}	0.207 , 0.238 0.209 , 0.240	Depositor DCC
R_{free} test set	2000 reflections (4.87%)	wwPDB-VP
Wilson B-factor (Å^2)	24.4	Xtriage
Anisotropy	0.841	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 37.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	2268	wwPDB-VP
Average B, all atoms (Å^2)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.62% 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

Bond lengths and bond angles in the following residue types are not validated in this section: MG

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	0.54	0/2117	0.71	2/2894 (0.1%)

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	246	ARG	NE-CZ-NH2	5.71	124.33	119.20
1	A	246	ARG	NE-CZ-NH1	-5.31	116.19	121.50

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	2059	0	1993	19	3
2	A	2	0	0	0	0
3	A	207	0	0	7	1
All	All	2268	0	1993	19	4

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.

All (19) 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:265:ARG:NH2	3:A:401:HOH:O	1.90	1.02
1:A:162:ARG:NH1	3:A:402:HOH:O	1.97	0.91
1:A:154[A]:GLU:OE1	3:A:403:HOH:O	1.99	0.80
1:A:135:PRO:HD2	1:A:155:MET:HE3	1.80	0.62
1:A:135:PRO:HD2	1:A:155:MET:CE	2.31	0.61
1:A:64:LEU:HD22	1:A:74:ILE:HD11	1.87	0.57
1:A:64:LEU:HD22	1:A:74:ILE:HG12	1.91	0.53
1:A:64:LEU:HD22	1:A:74:ILE:CD1	2.40	0.52
1:A:64:LEU:HD22	1:A:74:ILE:CG1	2.41	0.51
1:A:100[B]:GLN:NE2	3:A:411:HOH:O	2.45	0.48
1:A:256:GLN:HG2	1:A:265:ARG:HG3	1.94	0.48
1:A:198:GLU:HG2	1:A:285:ALA:O	2.15	0.46
1:A:200:LYS:NZ	3:A:415:HOH:O	2.48	0.45
1:A:135:PRO:CD	1:A:155:MET:HE3	2.47	0.45
1:A:129:CYS:SG	1:A:140:ALA:HB1	2.59	0.43
1:A:134:HIS:CD2	1:A:155:MET:HE2	2.54	0.42
1:A:286:ALA:HB2	3:A:476:HOH:O	2.19	0.42
1:A:246:ARG:HH11	1:A:246:ARG:HD3	1.58	0.41
1:A:100[A]:GLN:NE2	3:A:423:HOH:O	2.54	0.40

All (4) 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)
1:A:109:HIS:CD2	1:A:246:ARG:NH1[4_455]	1.09	1.11
1:A:109:HIS:NE2	1:A:246:ARG:NH1[4_455]	1.23	0.97
1:A:109:HIS:NE2	1:A:246:ARG:CZ[4_455]	1.95	0.25
3:A:509:HOH:O	3:A:565:HOH:O[2_556]	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	A	270/268 (101%)	262 (97%)	8 (3%)	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	205/201 (102%)	205 (100%)	0	100 100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	226	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 oligosaccharides in this entry.

5.6 Ligand geometry (i)

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.

No monomer is involved in short contacts.

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

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	268/268 (100%)	1.45	71 (26%) 1 1	13, 34, 56, 88	4 (1%)

All (71) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	234	ALA	5.8
1	A	19	THR	5.6
1	A	21	TRP	5.0
1	A	286	ALA	4.4
1	A	115	GLY	4.4
1	A	285	ALA	4.2
1	A	50	VAL	4.2
1	A	54	ALA	4.2
1	A	64	LEU	4.1
1	A	209	LEU	4.0
1	A	233	ALA	3.9
1	A	265	ARG	3.8
1	A	47	LEU	3.8
1	A	85	TYR	3.8
1	A	237	GLY	3.6
1	A	282	ALA	3.6
1	A	235	PRO	3.5
1	A	260	GLY	3.4
1	A	53	LEU	3.4
1	A	55	ASP	3.3
1	A	259	VAL	3.3
1	A	236	ASP	3.2
1	A	116	PHE	3.1
1	A	232	TYR	3.1
1	A	20	PRO	3.0
1	A	52	PRO	3.0
1	A	227	GLY	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	261	GLY	2.9
1	A	56	GLY	2.9
1	A	63	ALA	2.8
1	A	26	LEU	2.8
1	A	89	LEU	2.8
1	A	90	GLY	2.7
1	A	156	GLY	2.7
1	A	24	SER	2.7
1	A	93	ALA	2.7
1	A	203	VAL	2.7
1	A	199	GLY	2.7
1	A	27	VAL	2.6
1	A	49	ALA	2.6
1	A	92	ASP	2.6
1	A	202	ARG	2.5
1	A	212	GLU	2.5
1	A	25	GLU	2.5
1	A	51	GLY	2.5
1	A	117	GLY	2.5
1	A	224	ASP	2.5
1	A	28	ARG	2.5
1	A	162	ARG	2.4
1	A	213	GLY	2.4
1	A	246	ARG	2.4
1	A	118	ALA	2.4
1	A	204	THR	2.4
1	A	228	ILE	2.4
1	A	161	PRO	2.3
1	A	96	ALA	2.3
1	A	194	VAL	2.3
1	A	262	ALA	2.3
1	A	243	ARG	2.3
1	A	98	ILE	2.3
1	A	210	LEU	2.2
1	A	91	HIS	2.2
1	A	220	THR	2.2
1	A	230	ASP	2.2
1	A	201	ARG	2.2
1	A	95	PRO	2.2
1	A	205	TYR	2.1
1	A	94	PRO	2.1
1	A	113	TYR	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	258	PRO	2.0
1	A	263	GLN	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)

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, 95^{th} 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(Å^2)	Q<0.9
2	MG	A	302	1/1	0.92	0.21	43,43,43,43	0
2	MG	A	301	1/1	0.97	0.06	24,24,24,24	1

6.5 Other polymers (i)

There are no such residues in this entry.