



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 9, 2026 – 01:30 AM UTC

PDB ID : 4J8N / pdb_00004j8n
Title : Aurora A Kinase Apo
Authors : Meyerowitz, J.G.; Gustafson, W.C.; Shokat, K.M.; Weiss, W.A.
Deposited on : 2013-02-14
Resolution : 3.13 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

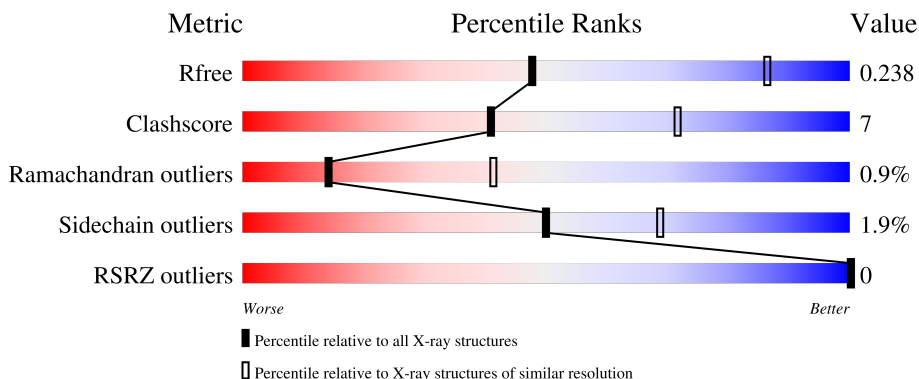
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

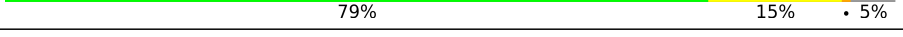
The reported resolution of this entry is 3.13 Å.

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	2351 (3.18-3.10)
Clashscore	190562	2452 (3.18-3.10)
Ramachandran outliers	187476	2324 (3.18-3.10)
Sidechain outliers	187428	2324 (3.18-3.10)
RSRZ outliers	180081	2351 (3.18-3.10)

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 $\leq 5\%$. 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.

Mol	Chain	Length	Quality of chain
1	A	279	
1	B	279	
1	C	279	
1	D	279	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 17546 atoms, of which 8807 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 Aurora kinase A.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	266	Total	C	H	N	O	S	0	0	0
			4390	1403	2205	387	389	6			
1	B	266	Total	C	H	N	O	S	0	0	0
			4390	1403	2205	387	389	6			
1	C	266	Total	C	H	N	O	S	0	0	0
			4390	1403	2205	387	389	6			
1	D	265	Total	C	H	N	O	S	0	0	0
			4368	1397	2192	385	388	6			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	287	ASP	THR	conflict	UNP O14965
B	287	ASP	THR	conflict	UNP O14965
C	287	ASP	THR	conflict	UNP O14965
D	287	ASP	THR	conflict	UNP O14965

- Molecule 2 is water.

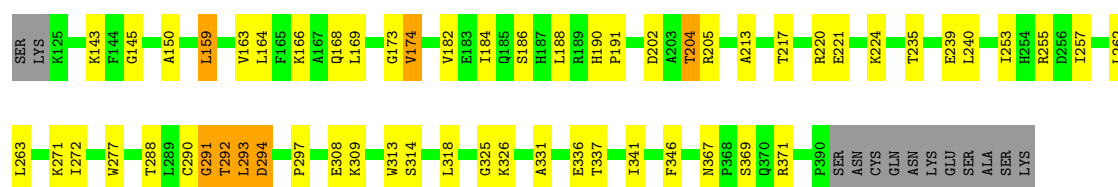
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	4	Total	O	0	0
			4	4		
2	C	2	Total	O	0	0
			2	2		
2	D	2	Total	O	0	0
			2	2		

3 Residue-property plots

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.

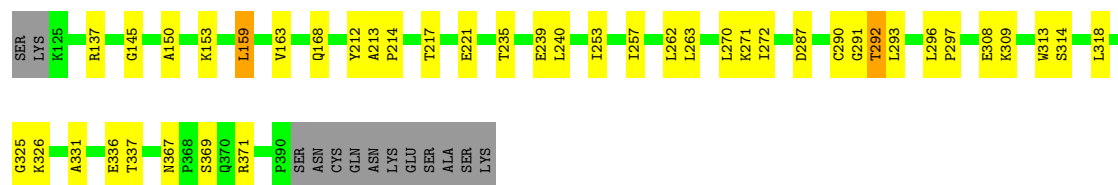
• Molecule 1: Aurora kinase A

Chain A: 



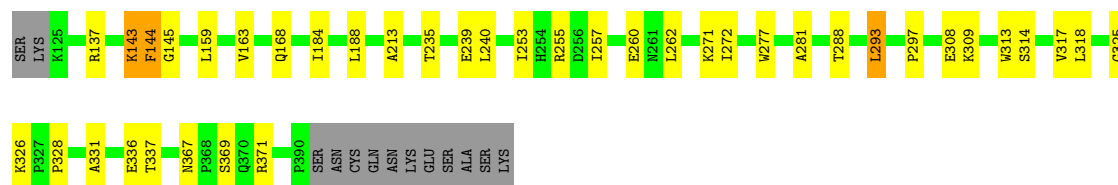
• Molecule 1: Aurora kinase A

Chain B: 



• Molecule 1: Aurora kinase A

Chain C: 



• Molecule 1: Aurora kinase A

Chain D: 



E308	K309	W313	S314	L318	A331	E336	T337	F346	D350	N367	P368	S369	Q370	R371	H380	P390	SER	ASN	CYS	GLN	ASN	LYS	GLU	SER	ALA	SER	LYS
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4 Data and refinement statistics

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, α , β , γ	83.78Å 83.78Å 171.78Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	44.95 – 3.13 44.95 – 3.13	Depositor EDS
% Data completeness (in resolution range)	99.8 (44.95-3.13) 97.1 (44.95-3.13)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.28 (at 3.12Å)	Xtriage
Refinement program	PHENIX 1.8_1069	Depositor
R, R_{free}	0.184 , 0.234 0.195 , 0.238	Depositor DCC
R_{free} test set	2033 reflections (4.30%)	wwPDB-VP
Wilson B-factor (Å ²)	93.6	Xtriage
Anisotropy	0.012	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 78.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtriage
Estimated twinning fraction	0.480 for -h,-k,l 0.487 for h,-h-k,-l 0.488 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	17546	wwPDB-VP
Average B, all atoms (Å ²)	114.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.32% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

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	0.38	0/2240	0.78	3/3027 (0.1%)
1	B	0.34	0/2240	0.73	2/3027 (0.1%)
1	C	0.34	0/2240	0.74	2/3027 (0.1%)
1	D	0.35	0/2231	0.76	4/3016 (0.1%)
All	All	0.35	0/8951	0.75	11/12097 (0.1%)

There are no bond length outliers.

The worst 5 of 11 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	213	ALA	CA-C-N	5.79	126.19	119.47
1	D	213	ALA	C-N-CA	5.79	126.19	119.47
1	A	294	ASP	N-CA-C	-5.71	106.44	113.41
1	C	213	ALA	CA-C-N	5.67	126.04	119.47
1	C	213	ALA	C-N-CA	5.67	126.04	119.47

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-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 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	2185	2205	2195	40	0
1	B	2185	2205	2195	25	0
1	C	2185	2205	2195	25	0
1	D	2176	2192	2182	31	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	4	0	0	0	0
2	C	2	0	0	2	0
2	D	2	0	0	0	0
All	All	8739	8807	8767	114	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 7.

The worst 5 of 114 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:173:GLY:O	1:A:174:VAL:HG22	1.53	1.09
1:C:367:ASN:OD1	1:C:369:SER:OG	1.99	0.80
1:B:367:ASN:OD1	1:B:369:SER:OG	2.01	0.79
1:A:173:GLY:O	1:A:174:VAL:CG2	2.33	0.75
1:D:367:ASN:OD1	1:D:369:SER:OG	2.04	0.74

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	264/279 (95%)	245 (93%)	14 (5%)	5 (2%)	6	24
1	B	264/279 (95%)	248 (94%)	14 (5%)	2 (1%)	16	44
1	C	264/279 (95%)	248 (94%)	14 (5%)	2 (1%)	16	44
1	D	263/279 (94%)	248 (94%)	14 (5%)	1 (0%)	30	59
All	All	1055/1116 (94%)	989 (94%)	56 (5%)	10 (1%)	14	41

5 of 10 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	174	VAL
1	B	292	THR
1	A	291	GLY
1	A	292	THR
1	B	287	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	235/247 (95%)	229 (97%)	6 (3%)	40	64
1	B	235/247 (95%)	232 (99%)	3 (1%)	61	73
1	C	235/247 (95%)	230 (98%)	5 (2%)	47	67
1	D	234/247 (95%)	230 (98%)	4 (2%)	53	70
All	All	939/988 (95%)	921 (98%)	18 (2%)	50	69

5 of 18 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	159	LEU
1	D	271	LYS
1	D	260	GLU
1	B	271	LYS
1	C	293	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	366	HIS
1	B	223	GLN
1	C	254	HIS

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](#)

There are no ligands in this entry.

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 [i](#)

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

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	266/279 (95%)	-0.91	0 100 100	75, 106, 167, 185	0
1	B	266/279 (95%)	-0.93	0 100 100	75, 107, 167, 196	0
1	C	266/279 (95%)	-0.91	0 100 100	75, 107, 167, 199	0
1	D	265/279 (94%)	-0.91	0 100 100	75, 106, 165, 186	0
All	All	1063/1116 (95%)	-0.92	0 100 100	75, 107, 167, 199	0

There are no RSRZ outliers to report.

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.