



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 9, 2026 – 06:04 AM UTC

PDB ID : 5LNN / pdb_00005lnn
Title : Crystal structure of D1050A mutant of the receiver domain of the histidine kinase CKI1 from Arabidopsis thaliana
Authors : Otrusanova, O.; Demo, G.; Kaderavek, P.; Jansen, S.; Jasenakova, Z.; Pekarova, B.; Janda, L.; Wimmerova, M.; Hejatk, J.; Zidek, L.
Deposited on : 2016-08-05
Resolution : 1.60 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

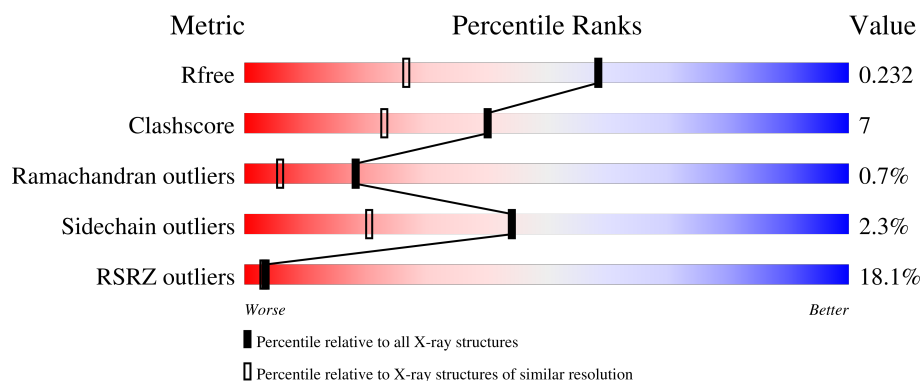
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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 $\leq 5\%$. 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	206	13% 55% 15% • 28%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 1346 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 Histidine kinase CKI1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	149	Total	C	N	O	S	0	6	0
			1209	749	211	242	7			

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	925	GLY	-	expression tag	UNP O22267
A	926	SER	-	expression tag	UNP O22267
A	927	SER	-	expression tag	UNP O22267
A	928	HIS	-	expression tag	UNP O22267
A	929	HIS	-	expression tag	UNP O22267
A	930	HIS	-	expression tag	UNP O22267
A	931	HIS	-	expression tag	UNP O22267
A	932	HIS	-	expression tag	UNP O22267
A	933	HIS	-	expression tag	UNP O22267
A	934	SER	-	expression tag	UNP O22267
A	935	SER	-	expression tag	UNP O22267
A	936	GLY	-	expression tag	UNP O22267
A	937	LEU	-	expression tag	UNP O22267
A	938	VAL	-	expression tag	UNP O22267
A	939	PRO	-	expression tag	UNP O22267
A	940	ARG	-	expression tag	UNP O22267
A	941	GLY	-	expression tag	UNP O22267
A	942	SER	-	expression tag	UNP O22267
A	943	HIS	-	expression tag	UNP O22267
A	1050	ALA	ASP	engineered mutation	UNP O22267
A	1123	LEU	-	expression tag	UNP O22267
A	1124	GLU	-	expression tag	UNP O22267
A	1125	HIS	-	expression tag	UNP O22267
A	1126	HIS	-	expression tag	UNP O22267
A	1127	HIS	-	expression tag	UNP O22267
A	1128	HIS	-	expression tag	UNP O22267
A	1129	HIS	-	expression tag	UNP O22267

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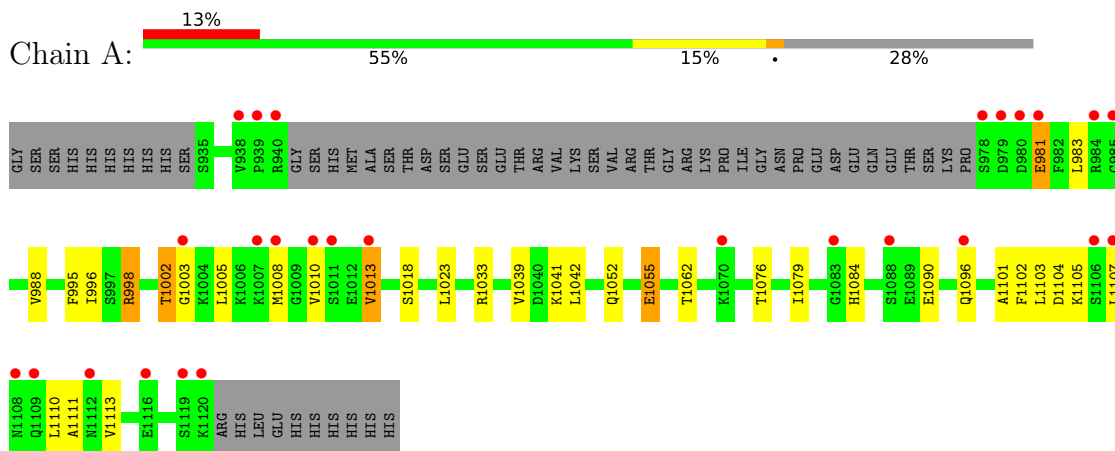
Chain	Residue	Modelled	Actual	Comment	Reference
A	1130	HIS	-	expression tag	UNP O22267

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	137	Total 137	O 137	0	0

i

- Molecule 1: Histidine kinase CKI1



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	53.18Å 101.00Å 79.46Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.51 – 1.60 42.51 – 1.60	Depositor EDS
% Data completeness (in resolution range)	98.4 (42.51-1.60) 98.4 (42.51-1.60)	Depositor EDS
R_{merge}	0.02	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.00 (at 1.60Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.192 , 0.228 0.204 , 0.232	Depositor DCC
R_{free} test set	1423 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	24.4	Xtriage
Anisotropy	0.014	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 41.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	1346	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.23% 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	1.70	9/1224 (0.7%)	1.48	12/1639 (0.7%)

The worst 5 of 9 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	995	PHE	N-CA	-9.24	1.35	1.46
1	A	1101	ALA	CA-C	-6.24	1.45	1.52
1	A	1023	LEU	CA-C	-6.05	1.44	1.52
1	A	1062	THR	CA-C	-5.92	1.45	1.52
1	A	1042	LEU	CA-C	-5.87	1.45	1.52

The worst 5 of 12 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	998	ARG	NE-CZ-NH2	8.25	126.63	119.20
1	A	1013	VAL	CB-CA-C	-6.81	101.10	110.77
1	A	983	LEU	O-C-N	-5.99	114.89	122.26
1	A	1003	GLY	CA-C-N	5.79	128.03	120.28
1	A	1003	GLY	C-N-CA	5.79	128.03	120.28

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	1209	0	1209	16	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	137	0	0	2	0
All	All	1346	0	1209	16	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 16 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:998:ARG:O	1:A:1002:THR:HG23	1.98	0.62
1:A:1010[A]:VAL:HG11	1:A:1013:VAL:HG22	1.84	0.58
1:A:1084:HIS:HD2	2:A:1323:HOH:O	1.90	0.53
1:A:1084:HIS:HE1	1:A:1090:GLU:OE2	1.93	0.51
1:A:1010[A]:VAL:HG11	1:A:1013:VAL:CG2	2.40	0.51

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	151/206 (73%)	150 (99%)	0	1 (1%)	18 6

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1055	GLU

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	134/180 (74%)	131 (98%)	3 (2%)	45 22

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	981	GLU
1	A	996	ILE
1	A	1096	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1084	HIS
1	A	1096	GLN

5.3.3 RNA ⓘ

There are no RNA molecules 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 oligosaccharides 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 [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	149/206 (72%)	0.86	27 (18%) 3 3	11, 29, 50, 70	9 (6%)

The worst 5 of 27 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1007	LYS	4.3
1	A	940	ARG	4.2
1	A	1108	ASN	3.9
1	A	1003	GLY	3.7
1	A	1008	MET	2.9

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.