



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 03:35 AM UTC

PDB ID : 1AKL / pdb_00001akl
Title : ALKALINE PROTEASE FROM PSEUDOMONAS AERUGINOSA IFO3080
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Deposited on : 1995-09-16
Resolution : 2.00 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) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

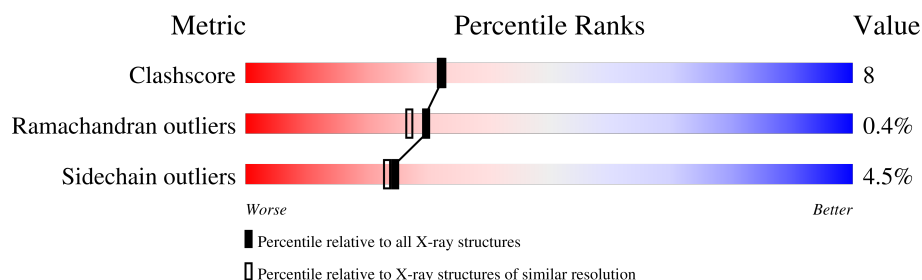
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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))
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	470	 82% 16% •

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3845 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 ALKALINE PROTEASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	470	Total	C	N	O	S	0	0	0
			3505	2195	583	726	1			

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

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

- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	8	Total	Ca	0	0
			8	8		

- Molecule 4 is water.

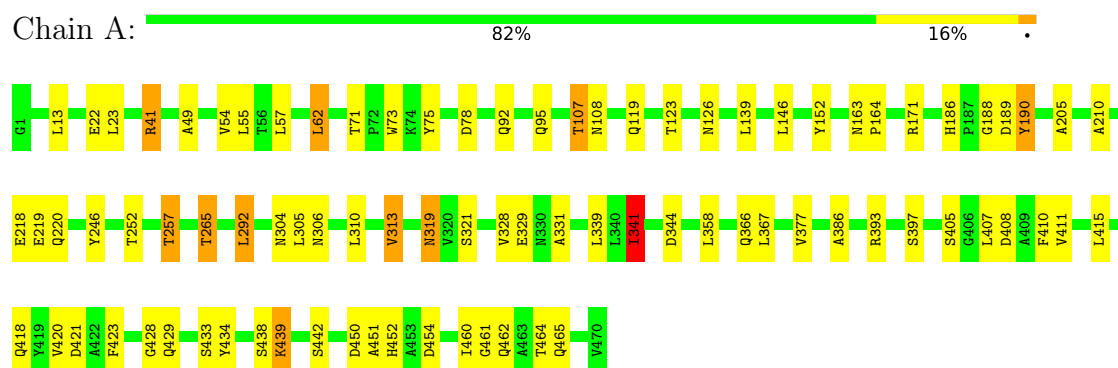
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	331	Total	O	0	0
			331	331		

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. 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. 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.

Note EDS was not executed.

• Molecule 1: ALKALINE PROTEASE



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	77.16Å 176.69Å 51.12Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 2.00	Depositor
% Data completeness (in resolution range)	75.1 (8.00-2.00)	Depositor
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.197 , 0.225	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3845	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CA, 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	0.53	1/3581 (0.0%)	0.94	17/4871 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	341	ILE	CA-CB	5.21	1.60	1.54

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	78	ASP	N-CA-C	6.45	120.15	112.93
1	A	461	GLY	N-CA-C	-6.40	103.35	110.96
1	A	205	ALA	N-CA-C	6.39	118.24	111.28
1	A	186	HIS	N-CA-C	-6.36	101.31	110.08
1	A	252	THR	N-CA-C	5.92	118.38	108.96
1	A	41	ARG	N-CA-C	5.84	119.23	111.75
1	A	439	LYS	N-CA-C	-5.56	105.80	112.58
1	A	265	THR	N-CA-C	-5.51	98.77	108.20
1	A	313	VAL	N-CA-C	5.50	116.50	108.53
1	A	344	ASP	N-CA-C	5.47	118.17	111.82
1	A	454	ASP	N-CA-C	-5.38	107.26	113.88
1	A	405	SER	N-CA-C	5.36	119.31	112.34
1	A	49	ALA	CA-C-N	5.34	126.18	119.98
1	A	49	ALA	C-N-CA	5.34	126.18	119.98
1	A	329	GLU	N-CA-C	5.21	119.69	112.45
1	A	386	ALA	N-CA-C	-5.19	105.80	111.82
1	A	377	VAL	N-CA-C	5.05	117.10	109.12

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	3505	0	3264	54	0
2	A	1	0	0	0	0
3	A	8	0	0	0	0
4	A	331	0	0	1	0
All	All	3845	0	3264	54	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 8.

All (54) 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:304:ASN:HD21	1:A:306:ASN:HD22	1.24	0.83
1:A:423:PHE:HB3	1:A:451:ALA:HA	1.69	0.75
1:A:292:LEU:HD13	1:A:328:VAL:HG11	1.70	0.73
1:A:366:GLN:HE21	1:A:393:ARG:HH11	1.44	0.64
1:A:310:LEU:HD22	1:A:321:SER:HB3	1.80	0.63
1:A:420:VAL:HG22	1:A:421:ASP:H	1.64	0.62
1:A:439:LYS:NZ	1:A:460:ILE:HG22	2.15	0.62
1:A:319:ASN:ND2	1:A:319:ASN:H	2.03	0.57
1:A:366:GLN:NE2	1:A:393:ARG:HH11	2.03	0.57
1:A:397:SER:HB2	1:A:462:GLN:O	2.04	0.57
1:A:305:LEU:HD21	1:A:331:ALA:HB2	1.87	0.56
1:A:319:ASN:H	1:A:319:ASN:HD22	1.52	0.55
1:A:464:THR:HG22	4:A:1314:HOH:O	2.05	0.55
1:A:313:VAL:H	1:A:319:ASN:HD21	1.54	0.54
1:A:304:ASN:ND2	1:A:306:ASN:HD22	1.99	0.54
1:A:62:LEU:HD13	1:A:123:THR:HB	1.89	0.53
1:A:438:SER:O	1:A:439:LYS:HB3	2.09	0.52
1:A:439:LYS:HZ2	1:A:460:ILE:HG22	1.74	0.52
1:A:415:LEU:HD21	1:A:428:GLY:HA2	1.90	0.52
1:A:304:ASN:C	1:A:304:ASN:HD22	2.17	0.52
1:A:418:GLN:O	1:A:429:GLN:HA	2.12	0.50
1:A:415:LEU:HD21	1:A:428:GLY:CA	2.41	0.50
1:A:95:GLN:NE2	1:A:171:ARG:HD3	2.26	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:439:LYS:O	1:A:439:LYS:HD3	2.12	0.49
1:A:218:GLU:HB3	1:A:220:GLN:OE1	2.13	0.49
1:A:393:ARG:HD2	1:A:460:ILE:HD12	1.95	0.49
1:A:420:VAL:HG22	1:A:421:ASP:N	2.28	0.49
1:A:210:ALA:HB2	1:A:219:GLU:HB2	1.95	0.48
1:A:304:ASN:HD21	1:A:306:ASN:ND2	2.04	0.48
1:A:304:ASN:ND2	1:A:306:ASN:H	2.11	0.48
1:A:423:PHE:HB3	1:A:451:ALA:CA	2.42	0.48
1:A:54:VAL:HG23	1:A:108:ASN:HB3	1.95	0.48
1:A:92:GLN:HE22	1:A:126:ASN:HD21	1.61	0.48
1:A:75:TYR:HB3	1:A:146:LEU:HD11	1.96	0.47
1:A:407:LEU:HB2	1:A:410:PHE:HD2	1.81	0.46
1:A:119:GLN:OE1	1:A:119:GLN:N	2.46	0.45
1:A:304:ASN:HD22	1:A:306:ASN:H	1.65	0.45
1:A:107:THR:HG21	1:A:246:TYR:HB2	1.98	0.44
1:A:407:LEU:HD12	1:A:410:PHE:HE2	1.82	0.44
1:A:257:THR:CG2	1:A:265:THR:HG22	2.47	0.44
1:A:13:LEU:HD21	1:A:341:ILE:CD1	2.48	0.43
1:A:152:TYR:CD1	1:A:152:TYR:N	2.87	0.43
1:A:95:GLN:HE21	1:A:171:ARG:HD3	1.84	0.42
1:A:107:THR:CG2	1:A:246:TYR:HB2	2.48	0.42
1:A:433:SER:HB2	1:A:442:SER:HB3	2.00	0.42
1:A:450:ASP:OD1	1:A:452:HIS:HB2	2.18	0.42
1:A:464:THR:HG22	1:A:465:GLN:H	1.83	0.42
1:A:41:ARG:HD2	1:A:190:TYR:CE2	2.55	0.41
1:A:71:THR:HG23	1:A:73:TRP:CE2	2.55	0.41
1:A:71:THR:O	1:A:71:THR:HG22	2.20	0.41
1:A:188:GLY:O	1:A:189:ASP:HB2	2.19	0.41
1:A:95:GLN:HG3	1:A:164:PRO:HB2	2.02	0.41
1:A:163:ASN:CG	1:A:163:ASN:O	2.64	0.41
1:A:13:LEU:HD21	1:A:341:ILE:HD13	2.01	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	468/470 (100%)	442 (94%)	24 (5%)	2 (0%)	30	27

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	190	TYR
1	A	434	TYR

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	358/358 (100%)	342 (96%)	16 (4%)	24	23

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	GLU
1	A	23	LEU
1	A	55	LEU
1	A	57	LEU
1	A	62	LEU
1	A	107	THR
1	A	139	LEU
1	A	257	THR
1	A	292	LEU
1	A	319	ASN
1	A	339	LEU
1	A	341	ILE
1	A	358	LEU
1	A	367	LEU
1	A	408	ASP
1	A	411	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	43	GLN
1	A	70	ASN
1	A	95	GLN
1	A	108	ASN
1	A	116	GLN
1	A	126	ASN
1	A	166	ASN
1	A	243	GLN
1	A	262	ASN
1	A	304	ASN
1	A	306	ASN
1	A	319	ASN
1	A	366	GLN
1	A	399	GLN
1	A	452	HIS
1	A	462	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 ⓘ

Of 9 ligands modelled in this entry, 9 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

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

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.