



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 06:58 PM UTC

PDB ID : 1KRA / pdb_00001kra
Title : CRYSTAL STRUCTURE OF KLEBSIELLA AEROGENES UREASE, ITS
APOENZYME AND TWO ACTIVE SITE MUTANTS
Authors : Jabri, E.; Karplus, P.A.
Deposited on : 1995-06-20
Resolution : 2.30 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

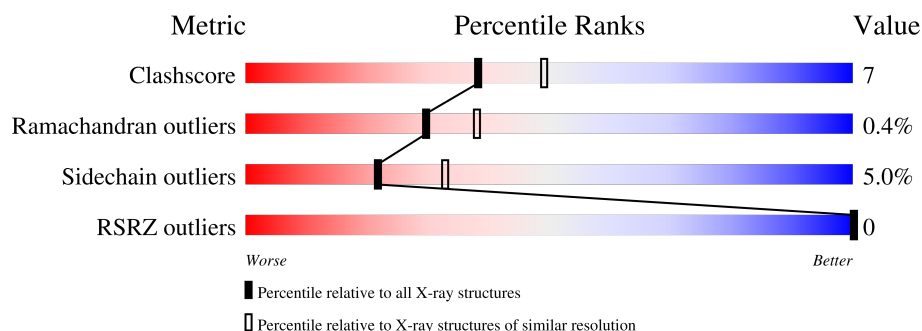
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	100	
2	B	106	
3	C	567	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5963 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 UREASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	100	Total	C	N	O	S	0	0	0
			775	491	134	145	5			

- Molecule 2 is a protein called UREASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	101	Total	C	N	O	S	0	0	0
			784	496	150	135	3			

- Molecule 3 is a protein called UREASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	566	Total	C	N	O	S	0	0	0
			4225	2651	741	810	23			

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	16	Total	O	0	0
			16	16		
4	B	12	Total	O	0	0
			12	12		
4	C	151	Total	O	0	0
			151	151		

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: UREASE

Chain A:  89% 9% .



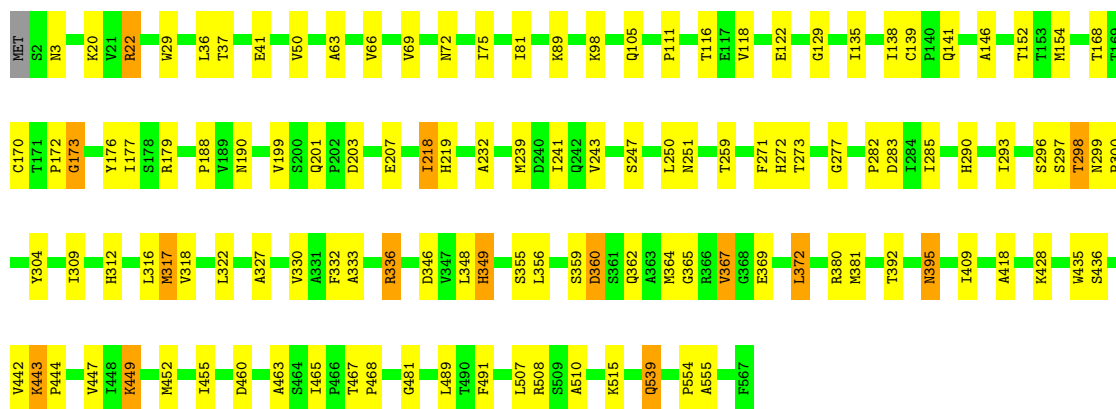
• Molecule 2: UREASE

Chain B:  77% 17% 5% .



• Molecule 3: UREASE

Chain C:  78% 19% .



4 Data and refinement statistics

Property	Value	Source
Space group	I 21 3	Depositor
Cell constants a, b, c, α , β , γ	170.80Å 170.80Å 170.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.30 10.00 – 2.30	Depositor EDS
% Data completeness (in resolution range)	95.0 (10.00-2.30) 95.7 (10.00-2.30)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.38 (at 2.31Å)	Xtriage
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.190 , (Not available) 0.183 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	12.5	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 22.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.40$, $\langle L^2 \rangle = 0.22$	Xtriage
Estimated twinning fraction	0.056 for -l,-k,-h	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5963	wwPDB-VP
Average B, all atoms (Å ²)	8.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.82% 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

5.1 Standard geometry

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	0.55	0/786	0.89	0/1061
2	B	0.49	0/804	0.96	3/1087 (0.3%)
3	C	0.53	0/4310	1.04	20/5874 (0.3%)
All	All	0.53	0/5900	1.01	23/8022 (0.3%)

There are no bond length outliers.

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	312	HIS	N-CA-C	11.34	123.72	111.36
3	C	436	SER	N-CA-C	-8.87	98.94	109.93
3	C	247	SER	N-CA-C	8.60	120.85	110.19
3	C	201	GLN	CA-C-N	7.14	127.46	119.32
3	C	201	GLN	C-N-CA	7.14	127.46	119.32
3	C	298	THR	N-CA-C	-7.05	100.37	110.59
3	C	299	ASN	N-CA-C	6.70	124.61	109.81
2	B	38	SER	N-CA-C	6.05	118.67	111.71
3	C	173	GLY	N-CA-C	6.04	124.66	112.34
3	C	409	ILE	N-CA-C	5.92	116.69	110.72
3	C	367	VAL	N-CA-C	5.75	119.11	111.17
3	C	283	ASP	N-CA-C	5.62	119.34	111.30
3	C	116	THR	N-CA-C	5.51	117.73	108.96
3	C	50	VAL	N-CA-C	5.36	116.84	111.00
3	C	443	LYS	CA-C-N	5.32	125.78	120.14
3	C	443	LYS	C-N-CA	5.32	125.78	120.14
3	C	173	GLY	CA-C-N	5.29	125.35	119.32
3	C	173	GLY	C-N-CA	5.29	125.35	119.32
3	C	309	ILE	N-CA-C	5.29	115.50	110.53
3	C	239	MET	N-CA-C	5.23	119.64	113.16
2	B	46	ASN	CA-C-N	5.22	124.94	119.82
2	B	46	ASN	C-N-CA	5.22	124.94	119.82
3	C	318	VAL	N-CA-C	5.21	115.98	110.72

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	775	0	807	5	0
2	B	784	0	775	13	0
3	C	4225	0	4178	65	0
4	A	16	0	0	0	0
4	B	12	0	0	1	0
4	C	151	0	0	3	1
All	All	5963	0	5760	80	1

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.

All (80) 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)
2:B:62:ASN:HD22	2:B:82:VAL:HB	1.43	0.84
3:C:317:MET:HE3	3:C:333:ALA:HB2	1.68	0.76
2:B:100:PRO:O	2:B:101:LEU:HB2	1.91	0.70
2:B:90:VAL:HB	2:B:97:VAL:HG11	1.75	0.68
2:B:88:ARG:HH11	2:B:88:ARG:HB3	1.60	0.64
3:C:300:PRO:HD3	3:C:365:GLY:HA2	1.79	0.63
3:C:296:SER:HB3	3:C:356:LEU:HB2	1.80	0.63
3:C:327:ALA:O	3:C:330:VAL:HG22	2.03	0.58
3:C:218:ILE:HD11	3:C:232:ALA:CB	2.33	0.58
3:C:296:SER:CB	3:C:356:LEU:HB2	2.34	0.58
2:B:4:GLY:HA2	3:C:22:ARG:O	2.03	0.58
3:C:460:ASP:HB3	3:C:463:ALA:HB2	1.86	0.58
3:C:355:SER:HB2	3:C:356:LEU:HD12	1.87	0.57
2:B:88:ARG:HB3	2:B:88:ARG:NH1	2.21	0.56
3:C:317:MET:HE3	3:C:333:ALA:CB	2.35	0.56
3:C:442:VAL:HG12	3:C:443:LYS:HG3	1.88	0.55
3:C:435:TRP:CZ3	3:C:444:PRO:HG3	2.42	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:317:MET:HE2	3:C:322:LEU:HB2	1.90	0.53
3:C:395:ASN:C	3:C:395:ASN:HD22	2.16	0.53
3:C:72:ASN:HB2	3:C:122:GLU:HB3	1.91	0.53
3:C:346:ASP:HB3	3:C:381:MET:HE3	1.90	0.52
3:C:380:ARG:NE	3:C:555:ALA:HB2	2.25	0.52
3:C:218:ILE:HD11	3:C:232:ALA:HB3	1.92	0.51
2:B:90:VAL:HG13	2:B:93:PHE:HE2	1.75	0.51
3:C:290:HIS:HB2	3:C:293:ILE:HG12	1.91	0.51
3:C:304:TYR:CE2	3:C:554:PRO:HD3	2.46	0.50
3:C:355:SER:C	3:C:356:LEU:HD12	2.36	0.50
2:B:90:VAL:HB	2:B:97:VAL:CG1	2.40	0.50
3:C:369:GLU:HG2	3:C:372:LEU:HD13	1.93	0.50
3:C:173:GLY:O	3:C:177:ILE:HG13	2.12	0.50
3:C:139:CYS:HB2	3:C:141:GLN:OE1	2.13	0.49
3:C:170:CYS:O	3:C:172:PRO:HD3	2.13	0.49
3:C:297:SER:OG	3:C:349:HIS:HE1	1.95	0.49
3:C:135:ILE:HG13	3:C:154:MET:HB3	1.95	0.48
3:C:447:VAL:HB	3:C:455:ILE:HG22	1.94	0.48
3:C:172:PRO:O	3:C:176:TYR:HB2	2.13	0.48
3:C:359:SER:O	3:C:365:GLY:HA3	2.14	0.48
3:C:418:ALA:O	3:C:428:LYS:HE2	2.13	0.48
2:B:7:HIS:HB3	3:C:20:LYS:HB2	1.96	0.47
1:A:75:GLU:HG3	2:B:2:ILE:HG23	1.95	0.47
1:A:65:THR:HB	1:A:67:GLU:OE2	2.15	0.47
3:C:539:GLN:HE21	3:C:539:GLN:H	1.62	0.47
3:C:300:PRO:HG2	4:C:687:HOH:O	2.15	0.46
3:C:63:ALA:O	3:C:89:LYS:NZ	2.49	0.46
2:B:97:VAL:HG13	2:B:98:MET:H	1.80	0.46
3:C:515:LYS:HB2	4:C:664:HOH:O	2.16	0.45
3:C:138:ILE:HD12	3:C:362:GLN:HB2	1.97	0.45
3:C:489:LEU:HD23	3:C:510:ALA:HB3	1.99	0.45
1:A:29:LYS:HA	1:A:68:GLN:O	2.16	0.45
3:C:218:ILE:HD12	3:C:243:VAL:HG13	1.99	0.45
2:B:50:LYS:HB2	4:B:115:HOH:O	2.16	0.45
3:C:203:ASP:O	3:C:207:GLU:HG3	2.17	0.44
3:C:41:GLU:O	3:C:98:LYS:NZ	2.49	0.44
3:C:75:ILE:O	3:C:81:ILE:HA	2.18	0.43
3:C:69:VAL:CG1	3:C:118:VAL:HG12	2.49	0.43
3:C:298:THR:HG22	3:C:360:ASP:HB2	1.99	0.43
3:C:129:GLY:HA3	3:C:152:THR:OG1	2.19	0.43
3:C:188:PRO:O	3:C:449:LYS:HD2	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:41:GLU:OE1	3:C:105:GLN:HG2	2.19	0.42
3:C:297:SER:OG	3:C:349:HIS:CE1	2.72	0.42
3:C:467:THR:N	3:C:468:PRO:CD	2.83	0.42
3:C:168:THR:HG22	3:C:219:HIS:CG	2.54	0.42
1:A:10:LYS:HD2	1:A:44:MET:SD	2.58	0.42
2:B:90:VAL:HG13	2:B:93:PHE:CE2	2.55	0.42
3:C:66:VAL:HG22	3:C:111:PRO:O	2.19	0.42
3:C:146:ALA:HB2	3:C:367:VAL:HG22	2.01	0.42
3:C:37:THR:HB	4:C:668:HOH:O	2.19	0.41
3:C:22:ARG:NH1	3:C:29:TRP:CZ2	2.89	0.41
3:C:36:LEU:N	3:C:36:LEU:HD12	2.36	0.41
3:C:298:THR:CG2	3:C:360:ASP:HB2	2.50	0.41
3:C:443:LYS:HA	3:C:444:PRO:HD3	1.87	0.41
1:A:43:ILE:HD11	1:A:95:VAL:HG21	2.03	0.41
3:C:190:ASN:HB3	3:C:491:PHE:CE2	2.56	0.41
3:C:273:THR:O	3:C:285:ILE:HA	2.21	0.41
3:C:285:ILE:HD11	3:C:348:LEU:CD1	2.50	0.41
3:C:332:PHE:C	3:C:332:PHE:CD1	2.95	0.41
3:C:277:GLY:O	3:C:336:ARG:NH2	2.54	0.40
3:C:250:LEU:O	3:C:251:ASN:HB2	2.21	0.40
3:C:271:PHE:O	3:C:272:HIS:C	2.63	0.40
3:C:359:SER:O	3:C:360:ASP:HB3	2.21	0.40

All (1) 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)
4:C:718:HOH:O	4:C:718:HOH:O[15_556]	1.02	1.18

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	98/100 (98%)	96 (98%)	2 (2%)	0	100	100
2	B	99/106 (93%)	92 (93%)	7 (7%)	0	100	100
3	C	564/567 (100%)	528 (94%)	33 (6%)	3 (0%)	24	31
All	All	761/773 (98%)	716 (94%)	42 (6%)	3 (0%)	30	38

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	C	360	ASP
3	C	481	GLY
3	C	364	MET

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	85/85 (100%)	81 (95%)	4 (5%)	23	35
2	B	78/83 (94%)	73 (94%)	5 (6%)	16	23
3	C	443/444 (100%)	422 (95%)	21 (5%)	23	35
All	All	606/612 (99%)	576 (95%)	30 (5%)	22	33

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	58	GLU
1	A	67	GLU
1	A	75	GLU
1	A	100	ILE
2	B	13	ILE
2	B	54	GLN
2	B	55	GLN
2	B	81	LEU
2	B	97	VAL
3	C	3	ASN

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Mol	Chain	Res	Type
3	C	22	ARG
3	C	179	ARG
3	C	199	VAL
3	C	218	ILE
3	C	241	ILE
3	C	259	THR
3	C	282	PRO
3	C	316	LEU
3	C	317	MET
3	C	336	ARG
3	C	349	HIS
3	C	372	LEU
3	C	392	THR
3	C	395	ASN
3	C	449	LYS
3	C	452	MET
3	C	465	ILE
3	C	507	LEU
3	C	508	ARG
3	C	539	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	12	GLN
2	B	29	HIS
2	B	55	GLN
2	B	62	ASN
3	C	107	ASN
3	C	142	GLN
3	C	349	HIS
3	C	362	GLN
3	C	395	ASN
3	C	419	HIS
3	C	472	HIS
3	C	486	HIS
3	C	521	GLN
3	C	539	GLN

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	100/100 (100%)	-1.89	0 100 100	3, 7, 12, 14	0
2	B	101/106 (95%)	-1.81	0 100 100	6, 10, 14, 18	0
3	C	566/567 (99%)	-1.84	0 100 100	2, 7, 17, 29	0
All	All	767/773 (99%)	-1.84	0 100 100	2, 7, 16, 29	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.