



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 03:28 AM UTC

PDB ID : 1KGG / pdb_00001kgg
Title : STRUCTURE OF BETA-LACTAMASE GLU166GLN:ASN170ASP MUTANT
Authors : Chen, C.C.H.; Herzberg, O.
Deposited on : 1999-05-20
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
Mogul	:	2022.3.0, CSD as543be (2022)
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

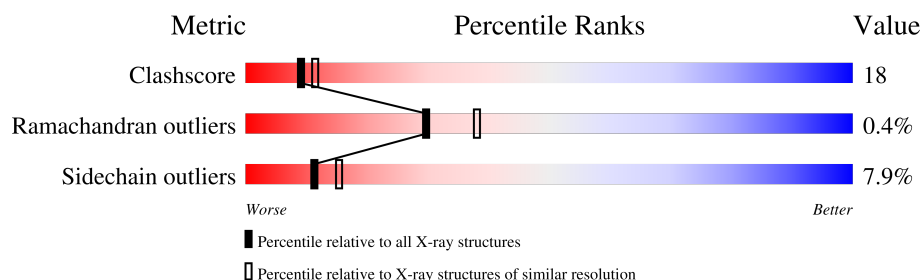
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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)

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	258	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 2101 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 PROTEIN (BETA-LACTAMASE).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	257	2029	1293	341	392	3	42	0	0

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	30	MET	ALA	SEE REMARK 999	UNP P00807
A	166	GLN	GLU	engineered mutation	UNP P00807
A	170	ASP	ASN	engineered mutation	UNP P00807

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	S		
2	A	1	5	4	1	0	0

- Molecule 3 is water.

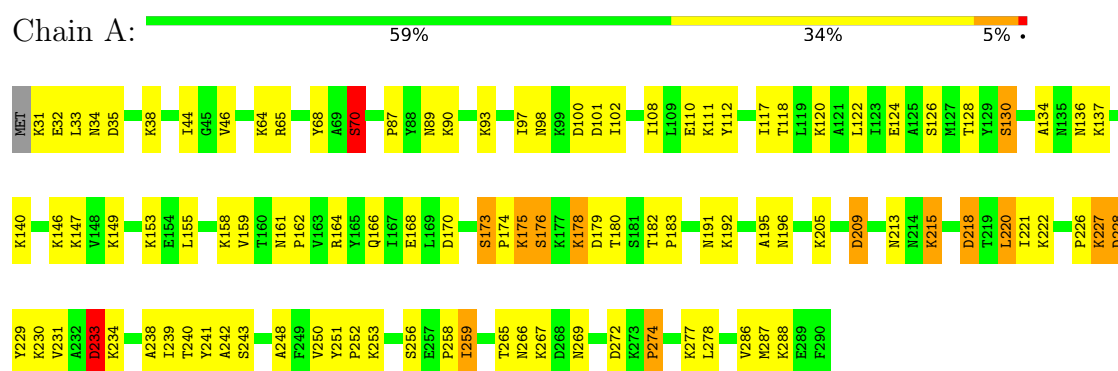
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	67	Total	O	0	0
			67	67		

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. 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: PROTEIN (BETA-LACTAMASE)



4 Data and refinement statistics

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

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	54.00Å 94.80Å 138.60Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 2.30	Depositor
% Data completeness (in resolution range)	82.0 (8.00-2.30)	Depositor
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.7	Depositor
R, R_{free}	0.178 , 0.263	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2101	wwPDB-VP
Average B, all atoms (Å ²)	29.0	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: SO4

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	0.98	1/2059 (0.0%)	1.22	13/2764 (0.5%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	117	ILE	CA-CB	5.40	1.62	1.54

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	173	SER	CA-C-N	10.11	129.91	119.19
1	A	173	SER	C-N-CA	10.11	129.91	119.19
1	A	218	ASP	N-CA-C	6.75	119.48	111.71
1	A	102	ILE	N-CA-C	6.26	117.10	108.84
1	A	233	ASP	N-CA-C	6.24	118.90	109.41
1	A	259	ILE	N-CA-C	-6.10	99.64	108.36
1	A	220	LEU	N-CA-C	5.87	114.39	108.75
1	A	192	LYS	N-CA-C	5.78	118.52	111.82
1	A	176	SER	N-CA-C	5.50	118.70	109.95
1	A	108	ILE	N-CA-C	5.43	116.83	111.67
1	A	128	THR	N-CA-C	5.40	119.97	112.90
1	A	100	ASP	N-CA-C	-5.23	106.92	113.15
1	A	248	ALA	N-CA-C	5.01	116.42	108.96

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	2029	0	2115	74	0
2	A	5	0	0	0	0
3	A	67	0	0	4	0
All	All	2101	0	2115	74	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 18.

All (74) 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:SER:OG	1:A:175:LYS:HG2	1.73	0.89
1:A:146:LYS:HD2	1:A:146:LYS:H	1.39	0.88
1:A:227:LYS:HE3	1:A:227:LYS:HA	1.60	0.83
1:A:149:LYS:HD2	1:A:153:LYS:HE2	1.65	0.79
1:A:251:TYR:CE1	1:A:258:PRO:HB3	2.25	0.72
1:A:44:ILE:HD11	1:A:278:LEU:HD11	1.73	0.71
1:A:238:ALA:HB3	1:A:243:SER:HB2	1.73	0.70
1:A:31:LYS:HD2	1:A:35:ASP:HB3	1.74	0.69
1:A:259:ILE:HD13	1:A:287:MET:HE1	1.75	0.68
1:A:136:ASN:O	1:A:140:LYS:HG2	1.93	0.67
1:A:159:VAL:HG11	1:A:182:THR:HG23	1.78	0.66
1:A:32:GLU:HB3	1:A:35:ASP:HB2	1.79	0.64
1:A:120:LYS:O	1:A:124:GLU:HG3	2.00	0.62
1:A:240:THR:HB	1:A:241:TYR:CD1	2.34	0.62
1:A:222:LYS:HE3	1:A:233:ASP:OD1	2.00	0.61
1:A:168:GLU:HG3	1:A:168:GLU:O	2.01	0.61
1:A:228:ASP:HB3	1:A:253:LYS:NZ	2.16	0.60
1:A:46:VAL:HG11	1:A:286:VAL:HG21	1.84	0.59
1:A:252:PRO:HG3	1:A:259:ILE:HD12	1.84	0.58
1:A:196:ASN:HA	3:A:369:HOH:O	2.02	0.58
1:A:228:ASP:HB2	1:A:253:LYS:HB2	1.87	0.57
1:A:153:LYS:HD3	1:A:158:LYS:HD3	1.86	0.57
1:A:252:PRO:HG3	1:A:259:ILE:CD1	2.35	0.57
1:A:32:GLU:HG3	1:A:34:ASN:H	1.71	0.55

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:175:LYS:HG3	1:A:176:SER:N	2.22	0.55
1:A:222:LYS:HG2	1:A:231:VAL:HB	1.89	0.55
1:A:149:LYS:HD2	1:A:153:LYS:CE	2.37	0.54
1:A:159:VAL:CG1	1:A:182:THR:HG23	2.38	0.53
1:A:170:ASP:O	1:A:238:ALA:HB1	2.09	0.52
1:A:87:PRO:HB2	1:A:90:LYS:HD3	1.91	0.51
1:A:242:ALA:HB1	1:A:274:PRO:HB3	1.94	0.50
1:A:230:LYS:HB3	1:A:251:TYR:HB2	1.95	0.49
1:A:228:ASP:HB3	1:A:253:LYS:HZ3	1.76	0.49
1:A:98:ASN:O	1:A:101:ASP:HB2	2.12	0.48
1:A:110:GLU:HG3	1:A:111:LYS:N	2.26	0.48
1:A:250:VAL:HG11	1:A:287:MET:HE2	1.95	0.48
1:A:140:LYS:HE3	3:A:328:HOH:O	2.14	0.48
1:A:161:ASN:ND2	1:A:178:LYS:HA	2.28	0.47
1:A:164:ARG:HA	3:A:362:HOH:O	2.14	0.47
1:A:70:SER:O	1:A:234:LYS:HE3	2.14	0.47
1:A:178:LYS:HD3	1:A:178:LYS:H	1.79	0.47
1:A:68:TYR:OH	1:A:162:PRO:HA	2.15	0.46
1:A:137:LYS:O	1:A:140:LYS:HB2	2.15	0.46
1:A:241:TYR:HA	1:A:269:ASN:O	2.16	0.46
1:A:226:PRO:HG2	1:A:229:TYR:CD1	2.51	0.46
1:A:64:LYS:O	1:A:183:PRO:HD3	2.15	0.46
1:A:32:GLU:HG3	1:A:34:ASN:N	2.30	0.46
1:A:126:SER:O	1:A:130:SER:HA	2.15	0.46
1:A:65:ARG:HD3	1:A:180:THR:OG1	2.16	0.45
1:A:222:LYS:HG3	3:A:305:HOH:O	2.16	0.45
1:A:97:ILE:HB	1:A:112:TYR:O	2.16	0.45
1:A:44:ILE:HG22	1:A:46:VAL:HG23	1.99	0.44
1:A:122:LEU:HD22	1:A:134:ALA:HA	2.00	0.44
1:A:266:ASN:OD1	1:A:266:ASN:C	2.60	0.43
1:A:173:SER:HA	1:A:174:PRO:HD3	1.68	0.43
1:A:251:TYR:CZ	1:A:258:PRO:HB3	2.54	0.43
1:A:227:LYS:HA	1:A:227:LYS:CE	2.42	0.43
1:A:215:LYS:HA	1:A:215:LYS:CE	2.49	0.42
1:A:229:TYR:HE2	1:A:287:MET:CE	2.32	0.42
1:A:267:LYS:HD2	1:A:272:ASP:HB3	2.02	0.42
1:A:209:ASP:O	1:A:213:ASN:ND2	2.52	0.42
1:A:215:LYS:HD3	1:A:218:ASP:OD2	2.19	0.42
1:A:153:LYS:HD3	1:A:153:LYS:HA	1.79	0.41
1:A:87:PRO:HB2	1:A:89:ASN:OD1	2.21	0.41
1:A:68:TYR:OH	1:A:179:ASP:HA	2.20	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:136:ASN:HD22	1:A:136:ASN:HA	1.71	0.41
1:A:155:LEU:HD23	1:A:155:LEU:HA	1.94	0.41
1:A:93:LYS:O	1:A:118:THR:HA	2.21	0.41
1:A:228:ASP:HB3	1:A:253:LYS:HZ2	1.83	0.41
1:A:239:ILE:HD12	1:A:239:ILE:HA	1.80	0.41
1:A:265:THR:O	1:A:266:ASN:HB3	2.20	0.41
1:A:191:ASN:ND2	1:A:195:ALA:HB3	2.36	0.40
1:A:220:LEU:O	1:A:221:ILE:C	2.63	0.40
1:A:229:TYR:HE2	1:A:287:MET:HE3	1.86	0.40

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	255/258 (99%)	240 (94%)	14 (6%)	1 (0%)	30	38

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	70	SER

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	227/228 (100%)	209 (92%)	18 (8%)	11	15

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	33	LEU
1	A	38	LYS
1	A	70	SER
1	A	130	SER
1	A	147	LYS
1	A	166	GLN
1	A	175	LYS
1	A	178	LYS
1	A	205	LYS
1	A	209	ASP
1	A	215	LYS
1	A	227	LYS
1	A	228	ASP
1	A	233	ASP
1	A	256	SER
1	A	274	PRO
1	A	277	LYS
1	A	288	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	34	ASN
1	A	41	ASN
1	A	98	ASN
1	A	132	ASN
1	A	136	ASN
1	A	161	ASN
1	A	191	ASN
1	A	213	ASN

5.3.3 RNA ⓘ

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](#)

1 ligand is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SO4	A	292	-	4,4,4	0.40	0	6,6,6	1.18	1 (16%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	A	292	SO4	O4-S-O1	2.18	120.94	109.56

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 ⓘ

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.