



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 04:21 PM UTC

PDB ID : 1P09 / pdb_00001p09
Title : STRUCTURAL PLASTICITY AS A DETERMINANT OF ENZYME SPECIFICITY. CREATING BROADLY SPECIFIC PROTEASES
Authors : Bone, R.; Agard, D.A.
Deposited on : 1989-04-24
Resolution : 2.20 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

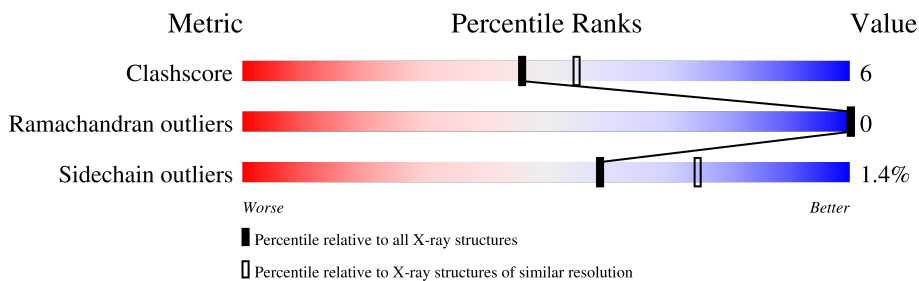
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)

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	198	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 1571 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 ALPHA-LYTIC PROTEASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	198	Total	C	N	O	S	0	2	0
			1392	847	262	275	8			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	213	ALA	MET	conflict	UNP P00778

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



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

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	168	Total 169	O 169	0	1

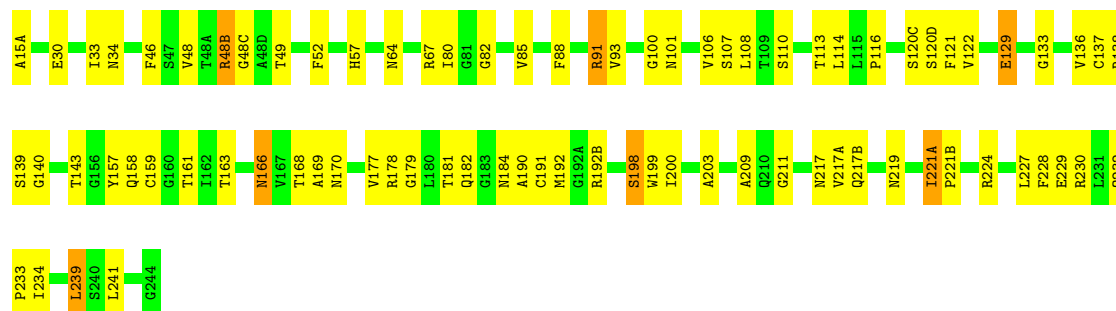
3 Residue-property plots [i](#)

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: ALPHA-LYTIC PROTEASE

Chain A: 



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	66.28Å 66.28Å 80.18Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	(Not available) – 2.20	Depositor
% Data completeness (in resolution range)	(Not available) ((Not available)-2.20)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	PROLSQ	Depositor
R, R_{free}	0.135 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	1571	wwPDB-VP
Average B, all atoms (Å ²)	13.0	wwPDB-VP

5 Model quality i

5.1 Standard geometry i

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	1.39	4/1414 (0.3%)	2.51	85/1916 (4.4%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	140	GLY	N-CA	5.40	1.51	1.45
1	A	139	SER	C-O	5.19	1.29	1.24
1	A	234	ILE	CA-CB	5.12	1.60	1.54
1	A	163	THR	CA-CB	5.03	1.61	1.54

All (85) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	91	ARG	CD-NE-CZ	16.10	146.94	124.40
1	A	48(B)	ARG	NE-CZ-NH1	-14.89	106.61	121.50
1	A	166	ASN	CA-CB-CG	-11.54	101.06	112.60
1	A	67	ARG	NE-CZ-NH1	11.51	133.01	121.50
1	A	48(B)	ARG	NH1-CZ-NH2	11.27	133.95	119.30
1	A	91	ARG	NE-CZ-NH2	11.00	129.10	119.20
1	A	120(D)	SER	N-CA-CB	-10.60	91.91	110.14
1	A	110	SER	CA-CB-OG	-9.68	91.73	111.10
1	A	67	ARG	NE-CZ-NH2	-8.68	111.39	119.20
1	A	49	THR	CA-CB-OG1	-8.63	96.65	109.60
1	A	82	GLY	CA-C-O	-8.26	110.29	119.04
1	A	192(B)[B]	ARG	NE-CZ-NH2	8.21	126.59	119.20
1	A	110	SER	O-C-N	7.96	132.74	122.39
1	A	221(A)	ILE	N-CA-CB	7.75	117.18	111.83
1	A	200	ILE	CA-C-O	-7.65	113.80	121.45
1	A	239	LEU	CA-C-O	-7.58	112.34	121.28
1	A	184	ASN	OD1-CG-ND2	-7.55	115.05	122.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	190	ALA	N-CA-C	-7.49	98.36	110.20
1	A	217(B)	GLN	CG-CD-NE2	7.44	127.56	116.40
1	A	114	LEU	CA-C-O	-7.42	112.23	120.69
1	A	168	THR	N-CA-CB	7.34	122.40	110.43
1	A	133	GLY	CA-C-O	-7.26	111.19	119.10
1	A	228	PHE	CA-CB-CG	7.12	120.92	113.80
1	A	34	ASN	OD1-CG-ND2	7.09	129.69	122.60
1	A	48(C)	GLY	N-CA-C	-7.08	99.29	111.47
1	A	230	ARG	NE-CZ-NH1	7.08	128.58	121.50
1	A	177	VAL	N-CA-CB	7.04	120.79	111.64
1	A	107	SER	O-C-N	7.00	131.20	123.22
1	A	91	ARG	NH1-CZ-NH2	-6.89	110.34	119.30
1	A	203	ALA	CA-C-O	-6.55	110.54	119.12
1	A	120(C)	SER	CA-CB-OG	-6.53	98.03	111.10
1	A	64	ASN	CA-C-O	-6.45	110.58	119.16
1	A	108	LEU	N-CA-CB	-6.38	100.60	110.56
1	A	64	ASN	N-CA-CB	6.37	123.01	113.65
1	A	178	ARG	N-CA-CB	-6.24	101.40	110.70
1	A	157	TYR	CA-C-O	-6.20	114.10	120.54
1	A	110	SER	CB-CA-C	-6.06	97.53	110.32
1	A	182	GLN	CB-CG-CD	-6.04	102.33	112.60
1	A	192(B)[B]	ARG	CA-C-O	-6.04	114.38	120.96
1	A	181	THR	CA-CB-OG1	-6.02	100.56	109.60
1	A	113	THR	O-C-N	5.93	130.61	123.13
1	A	107	SER	CA-CB-OG	-5.89	99.31	111.10
1	A	168	THR	CA-CB-OG1	-5.84	100.84	109.60
1	A	106	VAL	CA-C-O	-5.82	114.31	120.48
1	A	113	THR	CA-CB-OG1	-5.80	100.90	109.60
1	A	169	ALA	N-CA-C	-5.78	100.28	109.59
1	A	106	VAL	O-C-N	5.69	129.16	123.18
1	A	88	PHE	CA-CB-CG	-5.68	108.12	113.80
1	A	227	LEU	N-CA-CB	-5.68	101.77	110.85
1	A	170	ASN	CA-C-O	-5.60	114.32	120.70
1	A	143	THR	CA-CB-CG2	5.54	119.91	110.50
1	A	241	LEU	CA-C-O	-5.52	114.98	121.16
1	A	161	THR	CA-CB-OG1	-5.47	101.39	109.60
1	A	217(A)	VAL	CB-CA-C	5.44	118.79	111.28
1	A	48	VAL	CA-CB-CG2	5.42	119.62	110.40
1	A	57	HIS	CA-C-O	-5.42	112.00	119.05
1	A	48(B)	ARG	CA-CB-CG	-5.41	103.29	114.10
1	A	33	ILE	CB-CG1-CD1	5.39	125.13	113.80
1	A	161	THR	CA-CB-CG2	5.36	119.62	110.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	107	SER	N-CA-CB	5.36	119.00	110.16
1	A	199	TRP	N-CA-C	-5.36	99.67	108.41
1	A	209	ALA	N-CA-CB	-5.36	101.87	109.85
1	A	122	VAL	N-CA-C	-5.34	100.69	108.17
1	A	136	VAL	N-CA-CB	-5.34	101.05	112.00
1	A	224	ARG	O-C-N	5.34	129.35	123.16
1	A	158	GLN	CA-C-O	-5.34	114.86	121.11
1	A	139	SER	N-CA-CB	-5.33	101.57	110.80
1	A	48(B)	ARG	CG-CD-NE	-5.29	100.37	112.00
1	A	80	ILE	N-CA-CB	5.29	119.01	111.82
1	A	129	GLU	CG-CD-OE2	-5.22	106.40	118.40
1	A	230	ARG	NE-CZ-NH2	-5.22	114.51	119.20
1	A	116	PRO	CA-C-O	-5.14	117.51	121.31
1	A	15(A)	ALA	O-C-N	5.13	130.91	122.70
1	A	85	VAL	CB-CA-C	5.12	115.67	110.65
1	A	224	ARG	CA-C-O	-5.12	115.03	120.92
1	A	100	GLY	CA-C-O	5.11	129.46	120.57
1	A	64	ASN	O-C-N	5.10	127.23	121.87
1	A	121	PHE	CA-C-O	-5.09	115.16	121.06
1	A	30	GLU	CB-CG-CD	-5.09	103.95	112.60
1	A	239	LEU	O-C-N	5.05	128.91	123.10
1	A	64	ASN	OD1-CG-ND2	5.04	127.64	122.60
1	A	217	ASN	CB-CG-OD1	5.04	130.89	120.80
1	A	80	ILE	N-CA-C	-5.03	100.32	107.77
1	A	170	ASN	CA-CB-CG	-5.03	107.57	112.60
1	A	181	THR	CA-CB-CG2	5.01	119.02	110.50

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	1392	0	1355	16	0
2	A	10	0	0	0	0
3	A	169	0	0	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	1571	0	1355	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 6.

All (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:192[A]:MET:HE3	3:A:245[A]:HOH:O	1.38	1.20
1:A:192[A]:MET:CE	3:A:245[A]:HOH:O	1.86	1.17
1:A:48(B):ARG:HD2	3:A:364:HOH:O	2.00	0.61
1:A:48(B):ARG:HG3	1:A:239:LEU:HD23	1.88	0.55
1:A:166:ASN:ND2	1:A:179:GLY:HA2	2.21	0.55
1:A:192[A]:MET:HE1	3:A:245[A]:HOH:O	1.76	0.54
1:A:191:CYS:HA	1:A:219:ASN:OD1	2.09	0.52
1:A:166:ASN:HD22	1:A:179:GLY:HA2	1.76	0.50
1:A:129:GLU:HG3	3:A:373:HOH:O	2.12	0.48
1:A:137:CYS:HA	1:A:159:CYS:HA	1.95	0.47
1:A:138:ARG:HA	1:A:198:SER:O	2.15	0.46
1:A:232:GLN:N	1:A:233:PRO:HD2	2.31	0.46
1:A:46:PHE:O	1:A:52:PHE:HA	2.18	0.44
1:A:93:VAL:HG11	1:A:101:ASN:HD22	1.84	0.43
1:A:211:GLY:HA3	1:A:229:GLU:O	2.19	0.43
1:A:221(A):ILE:HB	1:A:221(B):PRO:HD2	2.03	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	197/198 (100%)	190 (96%)	7 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	142/141 (101%)	140 (99%)	2 (1%)	59 75

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	91	ARG
1	A	198	SER

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	101	ASN
1	A	166	ASN
1	A	182	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

2 ligands are 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	2	-	4,4,4	0.77	0	6,6,6	0.60	0
2	SO4	A	1	-	4,4,4	0.54	0	6,6,6	0.57	0

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.