



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:48 PM UTC

PDB ID : 3CEG / pdb_00003ceg
Title : Crystal structure of the UBC domain of baculoviral IAP repeat-containing protein 6
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Deposited on : 2008-02-29
Resolution : 2.01 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)	:	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

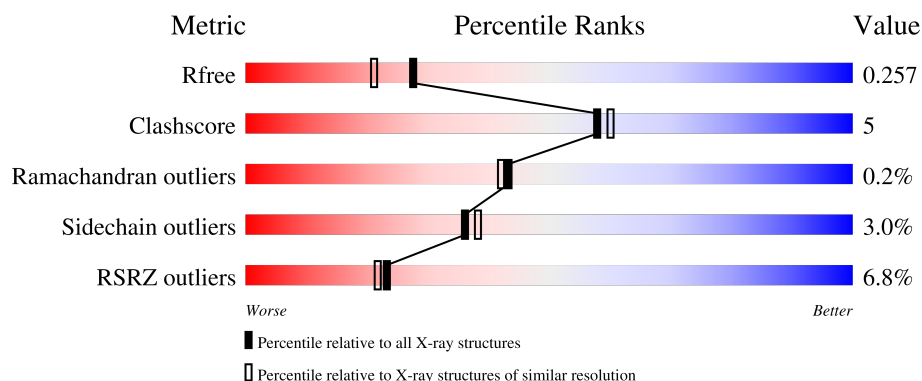
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.01 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (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 $\leq 5\%$. 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	323	5% 81% 9% • 10%
1	B	323	7% 79% 11% • 10%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4968 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 Baculoviral IAP repeat-containing protein 6.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	292	Total	C	N	O	S	Se	0	2	0
			2314	1467	390	443	6	8			
1	B	292	Total	C	N	O	S	Se	0	5	0
			2326	1471	396	446	6	7			

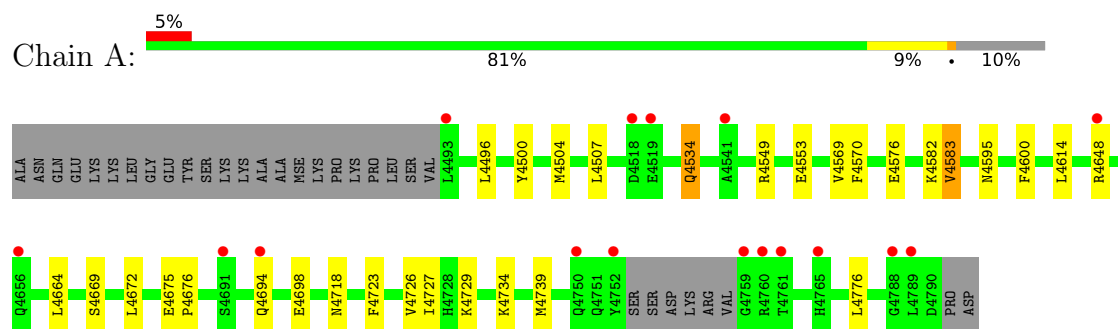
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	182	Total	O	0	8
			190	190		
2	B	133	Total	O	0	5
			138	138		

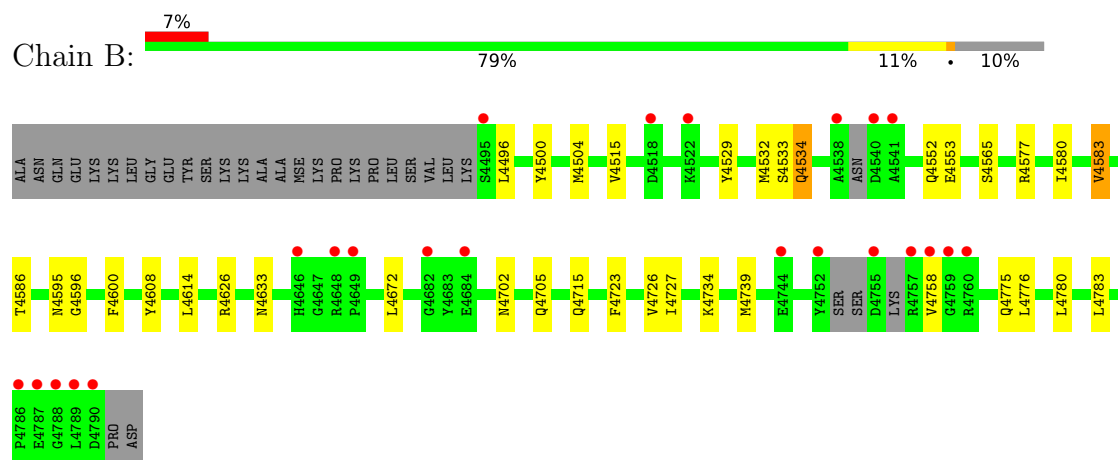
3 Residue-property plots [i](#)

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 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: Baculoviral IAP repeat-containing protein 6



- Molecule 1: Baculoviral IAP repeat-containing protein 6



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	78.20Å 97.81Å 129.89Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.50 – 2.01 44.50 – 2.01	Depositor EDS
% Data completeness (in resolution range)	92.9 (44.50-2.01) 92.9 (44.50-2.01)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.09 (at 2.01Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.211 , 0.252 0.212 , 0.257	Depositor DCC
R_{free} test set	3127 reflections (4.66%)	wwPDB-VP
Wilson B-factor (Å ²)	37.1	Xtriage
Anisotropy	0.410	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 36.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4968	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.93% 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	0.68	0/2362	0.89	1/3194 (0.0%)
1	B	0.65	0/2373	0.88	0/3208
All	All	0.67	0/4735	0.88	1/6402 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	4669	SER	N-CA-C	5.43	117.20	111.28

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	2314	0	2216	17	0
1	B	2326	0	2206	25	0
2	A	190	0	0	2	0
2	B	138	0	0	2	0
All	All	4968	0	4422	41	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 5.

All (41) 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:B:4739:MSE:HE3	1:B:4780:LEU:HD12	1.56	0.85
1:A:4595:ASN:OD1	1:A:4734:LYS:HE2	1.77	0.85
1:B:4739:MSE:CE	1:B:4780:LEU:HD12	2.12	0.80
1:B:4595:ASN:OD1	1:B:4734:LYS:HE2	1.84	0.77
1:B:4534:GLN:HE21	1:B:4534:GLN:H	1.36	0.71
1:B:4583:VAL:HG13	1:B:4600:PHE:HB2	1.75	0.68
1:A:4595:ASN:OD1	1:A:4734:LYS:CE	2.49	0.60
1:A:4739:MSE:HA	1:A:4739:MSE:HE2	1.84	0.59
1:A:4694:GLN:O	1:A:4698:GLU:HG3	2.04	0.58
1:B:4739:MSE:CE	1:B:4780:LEU:CD1	2.84	0.54
2:A:104:HOH:O	1:B:4775:GLN:HG2	2.10	0.52
1:A:4507:LEU:HD11	1:A:4729:LYS:HD2	1.92	0.50
1:A:4648:ARG:HA	1:B:4532:MSE:HE1	1.95	0.49
1:A:4534:GLN:H	1:A:4534:GLN:HE21	1.59	0.49
1:B:4702:ASN:O	1:B:4705:GLN:HG3	2.13	0.49
1:B:4723:PHE:O	1:B:4727[A]:ILE:HG12	2.14	0.48
1:B:4739:MSE:HE3	1:B:4780:LEU:CD1	2.37	0.48
1:A:4739:MSE:HE1	1:A:4776:LEU:HD23	1.95	0.47
1:A:4576:GLU:HG2	2:A:101:HOH:O	2.15	0.47
1:B:4529:TYR:O	1:B:4532:MSE:HG3	2.15	0.47
1:A:4504:MSE:CE	1:A:4726:VAL:HG13	2.45	0.46
1:A:4569:VAL:HG11	1:A:4664:LEU:HD13	1.96	0.46
1:B:4626:ARG:HB2	1:B:4633:ASN:OD1	2.17	0.45
1:B:4739:MSE:HE2	1:B:4780:LEU:CD1	2.46	0.45
1:B:4739:MSE:CE	1:B:4776:LEU:HG	2.47	0.45
1:A:4500:TYR:CE1	1:A:4734:LYS:HD2	2.51	0.45
1:B:4715:GLN:HA	1:B:4715:GLN:OE1	2.17	0.44
1:A:4549:ARG:NH1	1:A:4553:GLU:OE2	2.51	0.44
1:A:4723:PHE:O	1:A:4727[A]:ILE:HG12	2.17	0.44
1:A:4504:MSE:HE3	1:A:4570:PHE:CE1	2.52	0.44
1:B:4726:VAL:HG23	2:B:303:HOH:O	2.18	0.44
1:B:4534:GLN:HE21	1:B:4534:GLN:N	2.11	0.43
1:B:4586:THR:HA	1:B:4596:GLY:O	2.19	0.43
1:B:4577:ARG:HD3	1:B:4580:ILE:HD12	2.01	0.43
1:B:4500:TYR:CZ	1:B:4504:MSE:HG3	2.55	0.42
1:B:4553:GLU:OE2	1:B:4608:TYR:OH	2.35	0.42
1:B:4565:SER:O	1:B:4734:LYS:NZ	2.51	0.41
1:B:4552:GLN:NE2	2:B:293:HOH:O	2.48	0.41
1:A:4583:VAL:HG22	1:A:4600:PHE:HB2	2.03	0.40
1:A:4675:GLU:N	1:A:4676:PRO:CD	2.84	0.40
1:B:4739:MSE:HE1	1:B:4776:LEU:HG	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	290/323 (90%)	286 (99%)	4 (1%)	0	100	100
1	B	290/323 (90%)	283 (98%)	6 (2%)	1 (0%)	36	35
All	All	580/646 (90%)	569 (98%)	10 (2%)	1 (0%)	43	42

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	4758	VAL

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	250/278 (90%)	243 (97%)	7 (3%)	38	41
1	B	250/278 (90%)	242 (97%)	8 (3%)	34	35
All	All	500/556 (90%)	485 (97%)	15 (3%)	36	38

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4496	LEU
1	A	4534	GLN
1	A	4582	LYS
1	A	4583	VAL
1	A	4614	LEU

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Mol	Chain	Res	Type
1	A	4672	LEU
1	A	4718	ASN
1	B	4496	LEU
1	B	4515	VAL
1	B	4533	SER
1	B	4534	GLN
1	B	4583	VAL
1	B	4614	LEU
1	B	4672	LEU
1	B	4783	LEU

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	4534	GLN
1	A	4616	ASN
1	A	4750	GLN
1	A	4775	GLN
1	B	4534	GLN
1	B	4537	ASN
1	B	4552	GLN
1	B	4668	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 [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

6.1 Protein, DNA and RNA chains

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	285/323 (88%)	0.25	16 (5%) 30 29	15, 37, 65, 84	1 (0%)
1	B	285/323 (88%)	0.40	23 (8%) 18 17	13, 38, 76, 96	5 (1%)
All	All	570/646 (88%)	0.32	39 (6%) 23 22	13, 38, 72, 96	6 (1%)

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	4518	ASP	6.0
1	B	4538	ALA	4.9
1	B	4789	LEU	4.6
1	B	4758	VAL	4.4
1	A	4759	GLY	4.1
1	B	4755	ASP	3.9
1	B	4757	ARG	3.5
1	A	4519	GLU	3.2
1	B	4541	ALA	3.1
1	A	4761	THR	3.1
1	B	4649	PRO	3.0
1	B	4684	GLU	2.9
1	A	4493	LEU	2.7
1	A	4541	ALA	2.7
1	B	4540	ASP	2.7
1	A	4788	GLY	2.7
1	A	4752	TYR	2.7
1	B	4646	HIS	2.6
1	B	4752	TYR	2.6
1	A	4760	ARG	2.6
1	A	4789	LEU	2.5
1	B	4790	ASP	2.4
1	B	4522	LYS	2.4
1	B	4682	GLY	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	4518	ASP	2.3
1	B	4744	GLU	2.3
1	A	4691	SER	2.2
1	A	4750	GLN	2.2
1	B	4760	ARG	2.2
1	A	4648	ARG	2.2
1	B	4648	ARG	2.2
1	B	4788	GLY	2.2
1	A	4765	HIS	2.2
1	B	4787	GLU	2.1
1	B	4786	PRO	2.1
1	A	4656	GLN	2.1
1	B	4495	SER	2.1
1	B	4759	GLY	2.0
1	A	4694	GLN	2.0

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.