



wwPDB X-ray Structure Validation Summary Report ⓘ

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PDB ID : 1C4Z
Title : STRUCTURE OF E6AP: INSIGHTS INTO UBIQUITINATION PATHWAY
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Deposited on : 1999-10-14
Resolution : 2.60 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtrriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

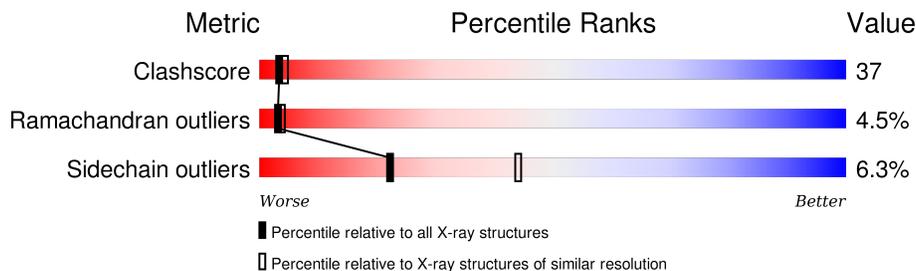
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	102246	2679 (2.60-2.60)
Ramachandran outliers	100387	2635 (2.60-2.60)
Sidechain outliers	100360	2635 (2.60-2.60)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	358	
1	B	358	
1	C	358	
2	D	154	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10040 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 UBIQUITIN-PROTEIN LIGASE E3A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	350	2861	1839	467	541	14	0	0	0
1	B	350	2861	1839	467	541	14	0	0	0
1	C	350	2861	1839	467	541	14	0	0	0

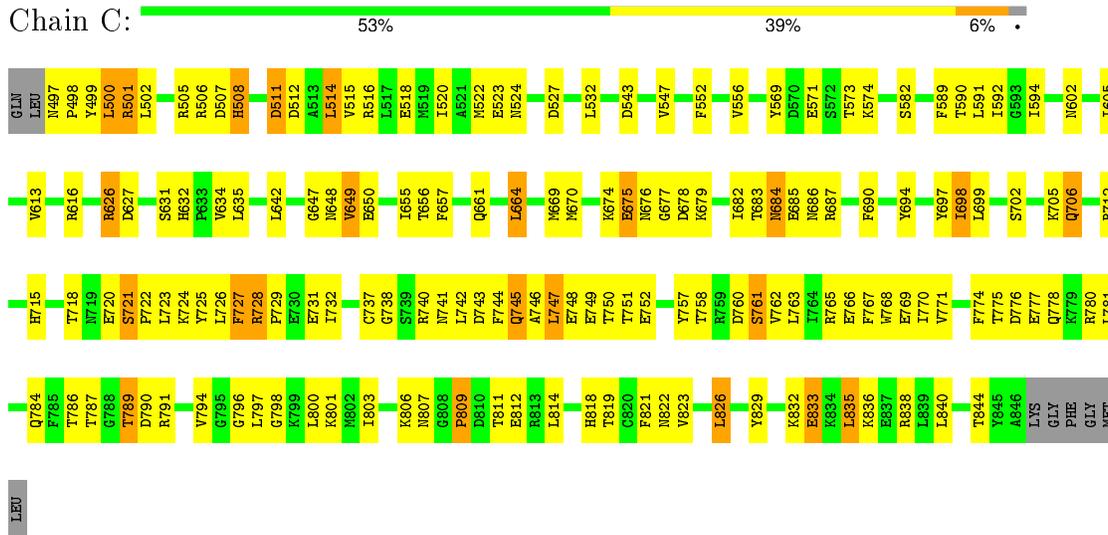
- Molecule 2 is a protein called UBIQUITIN CONJUGATING ENZYME E2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	D	144	1098	715	189	190	4	0	0	0

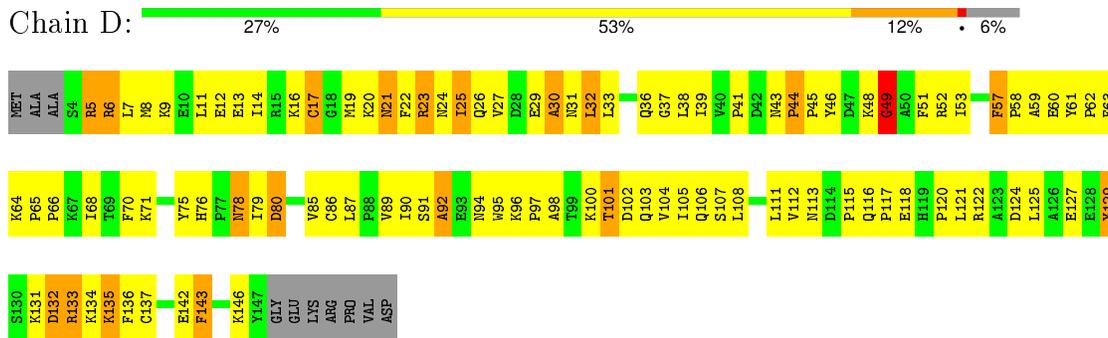
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	117	Total 117	O 117	0	0
3	B	99	Total 99	O 99	0	0
3	C	113	Total 113	O 113	0	0
3	D	30	Total 30	O 30	0	0

- Molecule 1: UBIQUITIN-PROTEIN LIGASE E3A



- Molecule 2: UBIQUITIN CONJUGATING ENZYME E2



4 Data and refinement statistics

Xtrriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	103.40Å 112.70Å 123.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 2.60	Depositor
% Data completeness (in resolution range)	90.0 (15.00-2.60)	Depositor
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.237 , 0.290	Depositor
Estimated twinning fraction	No twinning to report.	Xtrriage
Total number of atoms	10040	wwPDB-VP
Average B, all atoms (Å ²)	65.0	wwPDB-VP

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	0/2923	0.81	5/3943 (0.1%)
1	B	0.52	0/2923	0.80	2/3943 (0.1%)
1	C	0.56	1/2923 (0.0%)	0.83	2/3943 (0.1%)
2	D	0.44	0/1127	0.81	2/1536 (0.1%)
All	All	0.53	1/9896 (0.0%)	0.81	11/13365 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	626	ARG	CD-NE	-5.69	1.36	1.46

The worst 5 of 11 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	745	GLN	N-CA-C	-8.83	87.16	111.00
1	B	745	GLN	N-CA-C	-8.48	88.10	111.00
1	A	745	GLN	N-CA-C	-8.19	88.88	111.00
1	C	626	ARG	NE-CZ-NH1	7.19	123.89	120.30
1	A	663	ASP	CA-C-N	-7.18	101.40	117.20

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	671	TYR	Sidechain

5.2 Too-close contacts [i](#)

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	2861	0	2827	224	9
1	B	2861	0	2827	206	22
1	C	2861	0	2827	179	30
2	D	1098	0	1063	155	0
3	A	117	0	0	28	0
3	B	99	0	0	18	0
3	C	113	0	0	19	1
3	D	30	0	0	7	0
All	All	10040	0	9544	712	31

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 37.

The worst 5 of 712 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:626:ARG:HD2	3:A:60:HOH:O	1.38	1.19
2:D:80:ASP:OD1	2:D:121:LEU:HG	1.44	1.14
1:A:531:GLN:NE2	1:B:626:ARG:NE	1.99	1.09
1:A:531:GLN:NE2	1:B:626:ARG:CZ	2.16	1.09
1:A:531:GLN:NE2	1:B:626:ARG:CD	2.17	1.06

The worst 5 of 31 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)
1:B:813:ARG:NH1	1:C:762:VAL:CG1[3_656]	0.85	1.35
1:B:660:SER:OG	1:C:523:GLU:OE1[4_456]	0.89	1.31
1:B:813:ARG:NH2	1:C:762:VAL:CG2[3_656]	0.98	1.22
1:B:813:ARG:NH1	1:C:762:VAL:CB[3_656]	1.10	1.10

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:813:ARG:CZ	1:C:762:VAL:CB[3_656]	1.15	1.05

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	348/358 (97%)	306 (88%)	32 (9%)	10 (3%)	6	9
1	B	348/358 (97%)	305 (88%)	32 (9%)	11 (3%)	5	8
1	C	348/358 (97%)	309 (89%)	26 (8%)	13 (4%)	4	5
2	D	142/154 (92%)	91 (64%)	32 (22%)	19 (13%)	0	0
All	All	1186/1228 (97%)	1011 (85%)	122 (10%)	53 (4%)	3	4

5 of 53 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	508	HIS
1	A	698	ILE
1	A	776	ASP
1	B	508	HIS
1	B	675	GLU

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	317/323 (98%)	299 (94%)	18 (6%)	25	49
1	B	317/323 (98%)	295 (93%)	22 (7%)	19	38
1	C	317/323 (98%)	296 (93%)	21 (7%)	21	40
2	D	108/138 (78%)	102 (94%)	6 (6%)	26	50
All	All	1059/1107 (96%)	992 (94%)	67 (6%)	22	44

5 of 67 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	684	ASN
1	B	743	ASP
2	D	23	ARG
1	B	686	ASN
1	B	706	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 20 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	531	GLN
1	C	602	ASN
2	D	26	GLN
1	B	772	HIS
1	C	497	ASN

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 carbohydrates 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

6.1 Protein, DNA and RNA chains

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates

EDS was not executed - this section will therefore be empty.

6.4 Ligands

EDS was not executed - this section will therefore be empty.

6.5 Other polymers

EDS was not executed - this section will therefore be empty.