



wwPDB X-ray Structure Validation Summary Report ⓘ

May 21, 2020 – 09:25 pm BST

PDB ID : 6DC6
Title : Crystal structure of human ubiquitin activating enzyme E1 (Uba1) in complex with ubiquitin
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Deposited on : 2018-05-04
Resolution : 3.14 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

<https://www.wwpdb.org/validation/2017/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 : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

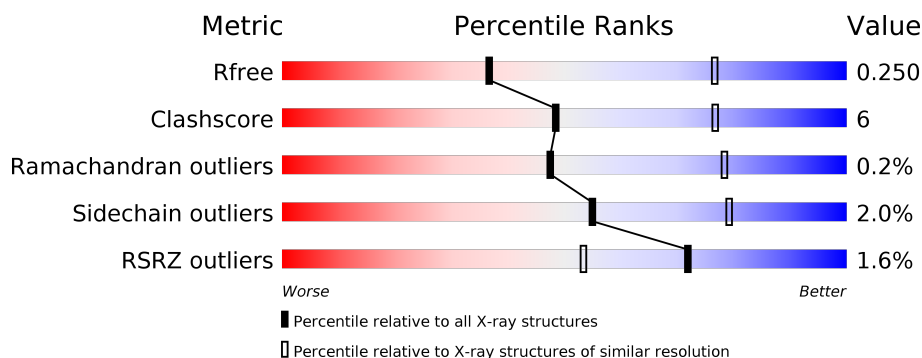
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.14 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}	130704	1626 (3.18-3.10)
Clashscore	141614	1735 (3.18-3.10)
Ramachandran outliers	138981	1677 (3.18-3.10)
Sidechain outliers	138945	1677 (3.18-3.10)
RSRZ outliers	127900	1588 (3.18-3.10)

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 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	1015	<div> <div>%</div> <div> <div></div> <div>81%</div> <div>16%</div> <div>•</div> </div> </div>
1	C	1015	<div> <div>2%</div> <div> <div></div> <div>80%</div> <div>18%</div> <div>••</div> </div> </div>
2	B	80	<div> <div>%</div> <div> <div></div> <div>83%</div> <div>13%</div> <div>5%</div> </div> </div>
2	D	80	<div> <div>5%</div> <div> <div></div> <div>88%</div> <div>8%</div> <div>5%</div> </div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 16888 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-like modifier-activating enzyme 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	992	Total	C	N	O	S	0	0	0
			7819	4965	1340	1475	39			
1	C	996	Total	C	N	O	S	0	0	0
			7843	4978	1345	1481	39			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	44	GLY	-	expression tag	UNP P22314
A	45	ALA	-	expression tag	UNP P22314
A	46	MET	-	expression tag	UNP P22314
A	47	GLY	-	expression tag	UNP P22314
A	48	SER	-	expression tag	UNP P22314
A	632	ALA	CYS	engineered mutation	UNP P22314
C	44	GLY	-	expression tag	UNP P22314
C	45	ALA	-	expression tag	UNP P22314
C	46	MET	-	expression tag	UNP P22314
C	47	GLY	-	expression tag	UNP P22314
C	48	SER	-	expression tag	UNP P22314
C	632	ALA	CYS	engineered mutation	UNP P22314

- Molecule 2 is a protein called Ubiquitin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	76	Total	C	N	O	S	0	0	0
			602	378	105	118	1			
2	D	76	Total	C	N	O	S	0	0	0
			602	378	105	118	1			

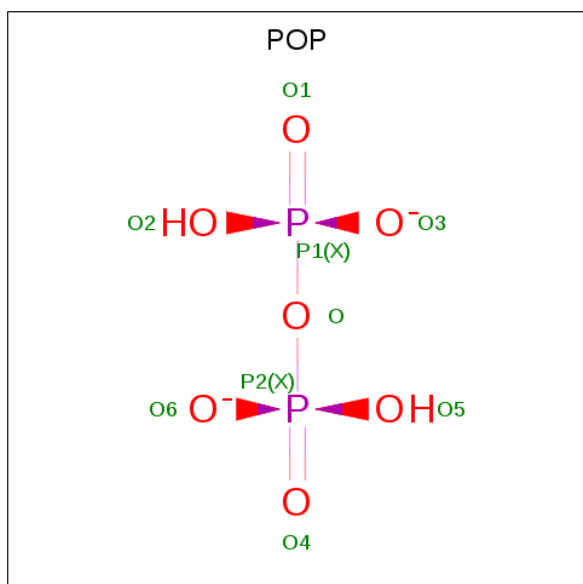
There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-3	GLY	-	expression tag	UNP P62979
B	-2	ALA	-	expression tag	UNP P62979
B	-1	MET	-	expression tag	UNP P62979
B	0	ALA	-	expression tag	UNP P62979
D	-3	GLY	-	expression tag	UNP P62979
D	-2	ALA	-	expression tag	UNP P62979
D	-1	MET	-	expression tag	UNP P62979
D	0	ALA	-	expression tag	UNP P62979

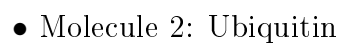
- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	2	Total Mg 2 2	0	0
3	C	2	Total Mg 2 2	0	0

- Molecule 4 is PYROPHOSPHATE 2- (three-letter code: POP) (formula: $\text{H}_2\text{O}_7\text{P}_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total O P 9 7 2	0	0
4	C	1	Total O P 9 7 2	0	0



- Molecule 2: Ubiquitin



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	105.00Å 70.44Å 188.85Å 90.00° 95.19° 90.00°	Depositor
Resolution (Å)	41.70 – 3.14 41.70 – 3.14	Depositor EDS
% Data completeness (in resolution range)	94.5 (41.70-3.14) 92.2 (41.70-3.14)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.78 (at 3.12Å)	Xtriage
Refinement program	PHENIX 1.10.1 _2155	Depositor
R, R_{free}	0.212 , 0.250 0.212 , 0.250	Depositor DCC
R_{free} test set	2000 reflections (4.35%)	wwPDB-VP
Wilson B-factor (Å ²)	65.0	Xtriage
Anisotropy	0.553	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 54.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	16888	wwPDB-VP
Average B, all atoms (Å ²)	90.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.28% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MG, POP

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	0.26	0/7994	0.44	1/10845 (0.0%)
1	C	0.26	0/8018	0.43	1/10878 (0.0%)
2	B	0.25	0/608	0.45	0/816
2	D	0.24	0/608	0.44	0/816
All	All	0.26	0/17228	0.44	2/23355 (0.0%)

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	C	0	1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	492	LEU	CA-CB-CG	7.18	131.81	115.30
1	C	679	LEU	CA-CB-CG	5.23	127.34	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	238	ALA	Peptide

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	7819	0	7696	96	0
1	C	7843	0	7717	102	0
2	B	602	0	629	9	0
2	D	602	0	629	6	0
3	A	2	0	0	0	0
3	C	2	0	0	0	0
4	A	9	0	0	1	0
4	C	9	0	0	2	0
All	All	16888	0	16671	202	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.

The worst 5 of 202 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:683:GLN:H	1:A:684:PRO:HD2	1.54	0.72
1:A:83:LEU:HD22	1:A:191:THR:HG23	1.72	0.72
1:A:80:LEU:HD11	1:A:101:LEU:HD22	1.72	0.72
1:A:57:ARG:NH1	4:A:1103:POP:O1	2.23	0.71
1:C:322:LYS:HB3	1:C:325:ARG:HG3	1.72	0.71

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	988/1015 (97%)	926 (94%)	60 (6%)	2 (0%)	47	78
1	C	992/1015 (98%)	933 (94%)	56 (6%)	3 (0%)	41	72
2	B	74/80 (92%)	72 (97%)	2 (3%)	0	100	100
2	D	74/80 (92%)	71 (96%)	3 (4%)	0	100	100
All	All	2128/2190 (97%)	2002 (94%)	121 (6%)	5 (0%)	47	78

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	1045	GLY
1	C	150	PRO
1	A	683	GLN
1	C	229	PRO
1	A	229	PRO

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	851/868 (98%)	839 (99%)	12 (1%)	67	85
1	C	853/868 (98%)	829 (97%)	24 (3%)	43	71
2	B	68/69 (99%)	68 (100%)	0	100	100
2	D	68/69 (99%)	68 (100%)	0	100	100
All	All	1840/1874 (98%)	1804 (98%)	36 (2%)	55	79

5 of 36 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	223	MET
1	C	288	ARG
1	C	966	ASN
1	C	239	ARG
1	C	345	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	514	ASN
1	A	914	GLN
1	C	374	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 for Mogul analysis.

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	$\# Z > 2$
4	POP	C	1103	3	6,8,8	0.74	0	13,13,13	1.28	1 (7%)
4	POP	A	1103	3	6,8,8	0.73	0	13,13,13	1.34	1 (7%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	POP	C	1103	3	-	0/6/6/6	-
4	POP	A	1103	3	-	0/6/6/6	-

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1103	POP	P2-O-P1	-3.64	120.34	132.83
4	C	1103	POP	P2-O-P1	-3.55	120.64	132.83

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	1103	POP	2	0
4	A	1103	POP	1	0

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	992/1015 (97%)	-0.18	6 (0%) 89 80	37, 74, 119, 166	0
1	C	996/1015 (98%)	0.03	23 (2%) 60 40	50, 95, 156, 189	0
2	B	76/80 (95%)	-0.07	1 (1%) 77 61	44, 87, 125, 135	0
2	D	76/80 (95%)	0.15	4 (5%) 26 12	56, 104, 143, 151	0
All	All	2140/2190 (97%)	-0.06	34 (1%) 72 53	37, 86, 144, 189	0

The worst 5 of 34 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	278	CYS	6.4
1	C	279	ASP	4.2
1	C	371	PRO	3.5
1	C	968	GLU	3.3
2	D	61	ILE	3.3

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 carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	MG	A	1101	1/1	0.93	0.19	36,36,36,36	0
3	MG	C	1101	1/1	0.93	0.21	33,33,33,33	0
3	MG	A	1102	1/1	0.96	0.27	23,23,23,23	0
3	MG	C	1102	1/1	0.97	0.21	30,30,30,30	0
4	POP	A	1103	9/9	0.97	0.20	3,57,79,86	0
4	POP	C	1103	9/9	0.97	0.22	26,45,90,99	0

6.5 Other polymers [i](#)

There are no such residues in this entry.