



# wwPDB X-ray Structure Validation Summary Report ⓘ

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PDB ID : 6DEV  
Title : Human caspase-6 E35K  
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Deposited on : 2018-05-13  
Resolution : 2.35 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](mailto: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.

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

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	rb-20031633
Percentile statistics	:	20171227.v01 (using entries in the PDB archive December 27th 2017)
Refmac	:	5.8.0158
CCP4	:	7.0 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20031633

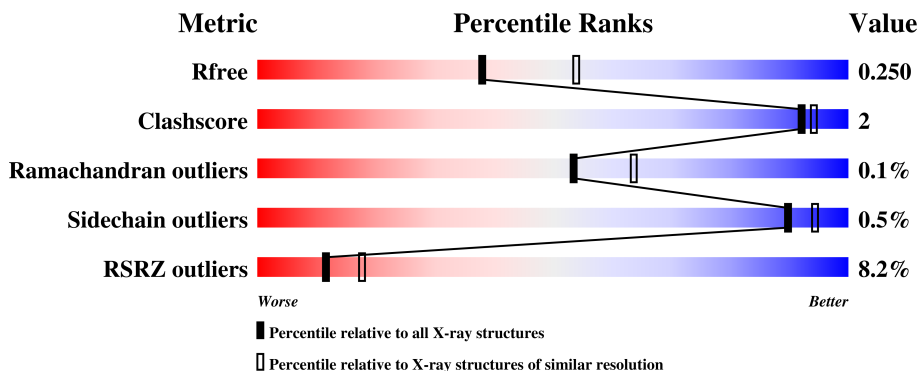
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.35 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))
$R_{free}$	111664	1763 (2.36-2.32)
Clashscore	122126	1858 (2.36-2.32)
Ramachandran outliers	120053	1834 (2.36-2.32)
Sidechain outliers	120020	1835 (2.36-2.32)
RSRZ outliers	108989	1737 (2.36-2.32)

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  $\geq 3$ , 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.

Mol	Chain	Length	Quality of chain
1	A	302	<div> <div>6%</div> <div>72%</div> <div>25%</div> </div>
1	B	302	<div> <div>4%</div> <div>71%</div> <div>25%</div> </div>
1	C	302	<div> <div>7%</div> <div>71%</div> <div>25%</div> </div>
1	D	302	<div> <div>7%</div> <div>71%</div> <div>26%</div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 14598 atoms, of which 7173 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 Caspase-6.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	225	Total	C	H	N	O	S	0	0	0
			3613	1161	1800	316	322	14			
1	B	225	Total	C	H	N	O	S	0	0	0
			3610	1161	1797	315	323	14			
1	C	225	Total	C	H	N	O	S	0	0	0
			3617	1164	1800	316	322	15			
1	D	223	Total	C	H	N	O	S	0	0	0
			3573	1151	1776	311	321	14			

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	initiating methionine	UNP P55212
A	1	ALA	-	expression tag	UNP P55212
A	35	LYS	GLU	engineered mutation	UNP P55212
A	294	LEU	-	expression tag	UNP P55212
A	295	GLU	-	expression tag	UNP P55212
A	296	HIS	-	expression tag	UNP P55212
A	297	HIS	-	expression tag	UNP P55212
A	298	HIS	-	expression tag	UNP P55212
A	299	HIS	-	expression tag	UNP P55212
A	300	HIS	-	expression tag	UNP P55212
A	301	HIS	-	expression tag	UNP P55212
B	0	MET	-	initiating methionine	UNP P55212
B	1	ALA	-	expression tag	UNP P55212
B	35	LYS	GLU	engineered mutation	UNP P55212
B	294	LEU	-	expression tag	UNP P55212
B	295	GLU	-	expression tag	UNP P55212
B	296	HIS	-	expression tag	UNP P55212
B	297	HIS	-	expression tag	UNP P55212
B	298	HIS	-	expression tag	UNP P55212
B	299	HIS	-	expression tag	UNP P55212
B	300	HIS	-	expression tag	UNP P55212

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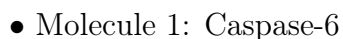
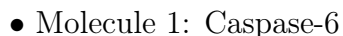
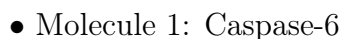
Chain	Residue	Modelled	Actual	Comment	Reference
B	301	HIS	-	expression tag	UNP P55212
C	0	MET	-	initiating methionine	UNP P55212
C	1	ALA	-	expression tag	UNP P55212
C	35	LYS	GLU	engineered mutation	UNP P55212
C	294	LEU	-	expression tag	UNP P55212
C	295	GLU	-	expression tag	UNP P55212
C	296	HIS	-	expression tag	UNP P55212
C	297	HIS	-	expression tag	UNP P55212
C	298	HIS	-	expression tag	UNP P55212
C	299	HIS	-	expression tag	UNP P55212
C	300	HIS	-	expression tag	UNP P55212
C	301	HIS	-	expression tag	UNP P55212
D	0	MET	-	initiating methionine	UNP P55212
D	1	ALA	-	expression tag	UNP P55212
D	35	LYS	GLU	engineered mutation	UNP P55212
D	294	LEU	-	expression tag	UNP P55212
D	295	GLU	-	expression tag	UNP P55212
D	296	HIS	-	expression tag	UNP P55212
D	297	HIS	-	expression tag	UNP P55212
D	298	HIS	-	expression tag	UNP P55212
D	299	HIS	-	expression tag	UNP P55212
D	300	HIS	-	expression tag	UNP P55212
D	301	HIS	-	expression tag	UNP P55212

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	68	Total O 68 68	0	0
2	B	60	Total O 60 60	0	0
2	C	29	Total O 29 29	0	0
2	D	28	Total O 28 28	0	0



- Molecule 1: Caspase-6





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	64.39Å 91.33Å 86.68Å 90.00° 91.80° 90.00°	Depositor
Resolution (Å)	36.47 – 2.35 45.66 – 2.35	Depositor EDS
% Data completeness (in resolution range)	84.0 (36.47-2.35) 78.9 (45.66-2.35)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.44 (at 2.34Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155)	Depositor
R, $R_{free}$	0.205 , 0.248 0.206 , 0.250	Depositor DCC
$R_{free}$ test set	2022 reflections (5.72%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	28.9	Xtriage
Anisotropy	0.111	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.39 , 31.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.075 for h,-k,-l	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	14598	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.87% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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](#)

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	0.25	0/1855	0.44	0/2492
1	B	0.25	0/1855	0.42	0/2493
1	C	0.24	0/1860	0.41	0/2498
1	D	0.25	0/1839	0.41	0/2471
All	All	0.25	0/7409	0.42	0/9954

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 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	1813	1800	1799	7	0
1	B	1813	1797	1796	9	0
1	C	1817	1800	1799	6	0
1	D	1797	1776	1775	5	0
2	A	68	0	0	3	0
2	B	60	0	0	2	0
2	C	29	0	0	1	0
2	D	28	0	0	1	0
All	All	7425	7173	7169	24	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 2.

The worst 5 of 24 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:144:LYS:NZ	1:B:202:ALA:O	2.07	0.88
1:A:51:ASN:ND2	2:A:402:HOH:O	2.22	0.71
1:B:43:ARG:NH2	2:B:402:HOH:O	2.28	0.67
1:A:51:ASN:ND2	1:A:53:GLU:OE2	2.28	0.66
1:A:247:GLU:OE1	2:A:401:HOH:O	2.12	0.66

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	219/302 (72%)	212 (97%)	7 (3%)	0	100	100
1	B	219/302 (72%)	211 (96%)	8 (4%)	0	100	100
1	C	221/302 (73%)	215 (97%)	5 (2%)	1 (0%)	31	34
1	D	217/302 (72%)	212 (98%)	5 (2%)	0	100	100
All	All	876/1208 (72%)	850 (97%)	25 (3%)	1 (0%)	53	64

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	276	PRO

### 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	197/263 (75%)	196 (100%)	1 (0%)	90	94
1	B	197/263 (75%)	197 (100%)	0	100	100
1	C	197/263 (75%)	195 (99%)	2 (1%)	78	86
1	D	195/263 (74%)	194 (100%)	1 (0%)	90	94
All	All	786/1052 (75%)	782 (100%)	4 (0%)	90	94

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	57	TRP
1	C	217	TYR
1	C	266	ASP
1	D	57	TRP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	161	GLN
1	D	58	HIS

### 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 carbohydrates 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 [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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	225/302 (74%)	0.34	18 (8%) 12 19	13, 34, 75, 85	0
1	B	225/302 (74%)	0.37	13 (5%) 23 33	14, 37, 73, 85	0
1	C	225/302 (74%)	0.52	22 (9%) 7 12	27, 46, 82, 100	0
1	D	223/302 (73%)	0.58	21 (9%) 8 14	24, 46, 82, 97	0
All	All	898/1208 (74%)	0.45	74 (8%) 11 18	13, 41, 80, 100	0

The worst 5 of 74 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	98	LEU	7.7
1	A	57	TRP	7.4
1	A	214	GLU	7.0
1	C	61	LEU	6.5
1	B	292	SER	6.2

### 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](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.