



# wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 10, 2020 – 08:29 AM BST

PDB ID : 6IAK  
Title : The crystal structure of the chicken CREB3 bZIP  
Authors : Sabaratnam, K.; Renner, M.  
Deposited on : 2018-11-26  
Resolution : 3.95 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 : 2.13.1  
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) : 2.13.1

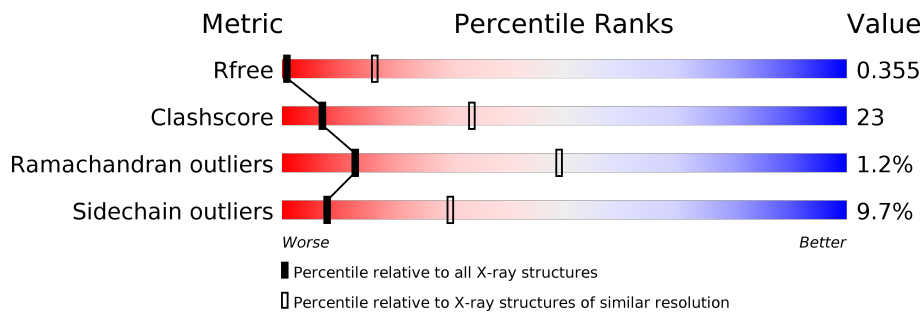
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.95 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	1025 (4.22-3.70)
Clashscore	141614	1085 (4.22-3.70)
Ramachandran outliers	138981	1047 (4.22-3.70)
Sidechain outliers	138945	1039 (4.22-3.70)

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 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  $\leq 5\%$

Mol	Chain	Length	Quality of chain
1	A	441	7% 7% • 85%
1	B	441	7% 6% • 86%
1	C	441	7% 7% • 86%
1	D	441	7% 6% 86%
1	E	441	9% 6% 85%
1	F	441	9% 6% • 84%
1	G	441	6% • 92%

## 2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 3563 atoms, of which 0 are hydrogens and 0 are deuteriums.

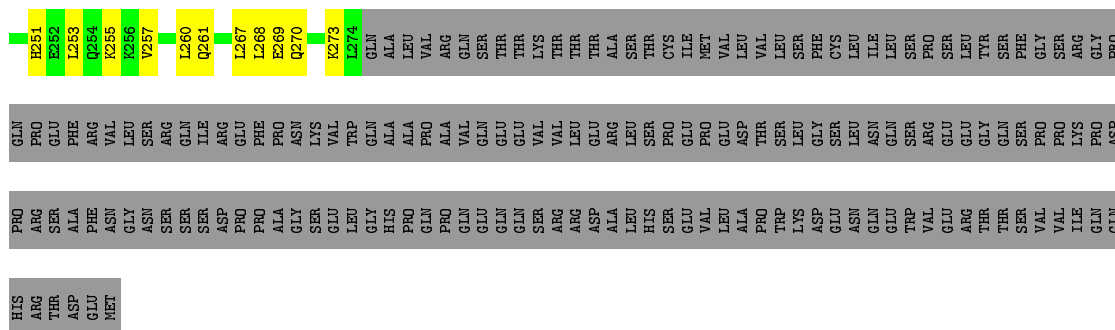
In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Uncharacterized protein.

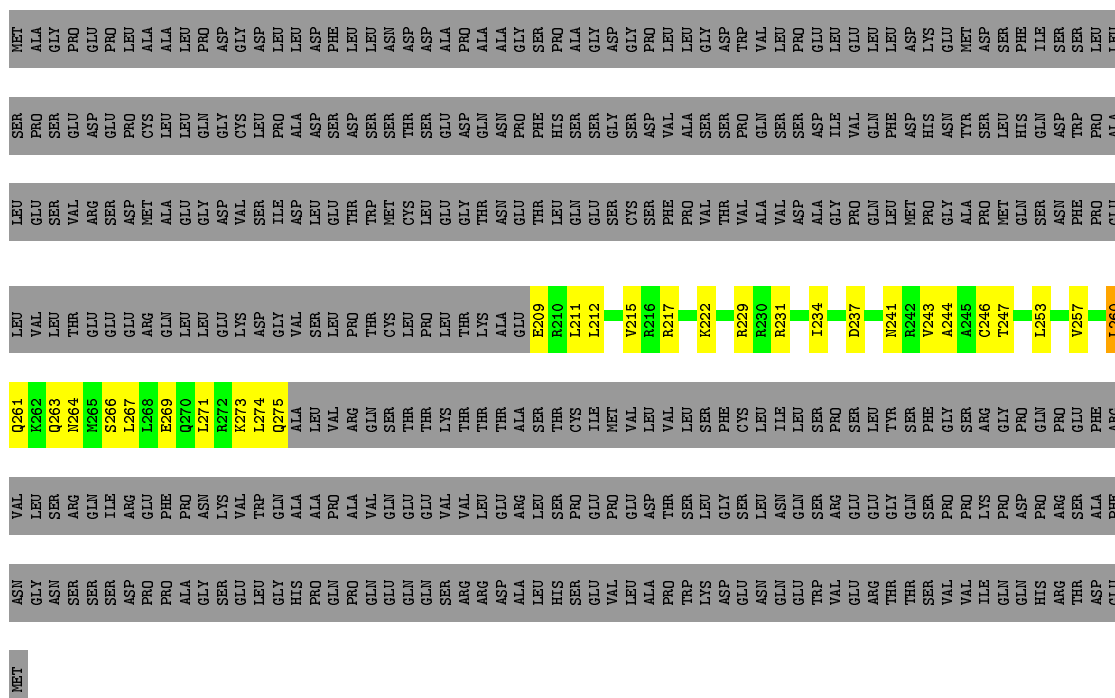
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	68	Total	C	N	O	S	0	0	0
			567	349	120	96	2			
1	B	62	Total	C	N	O	S	0	0	0
			511	313	106	90	2			
1	C	61	Total	C	N	O	S	0	0	0
			509	310	108	89	2			
1	D	62	Total	C	N	O	S	0	0	0
			518	317	110	89	2			
1	E	67	Total	C	N	O	S	0	0	0
			563	345	119	97	2			
1	F	71	Total	C	N	O	S	0	0	0
			593	363	124	104	2			
1	G	37	Total	C	N	O	S	0	0	0
			302	180	65	56	1			



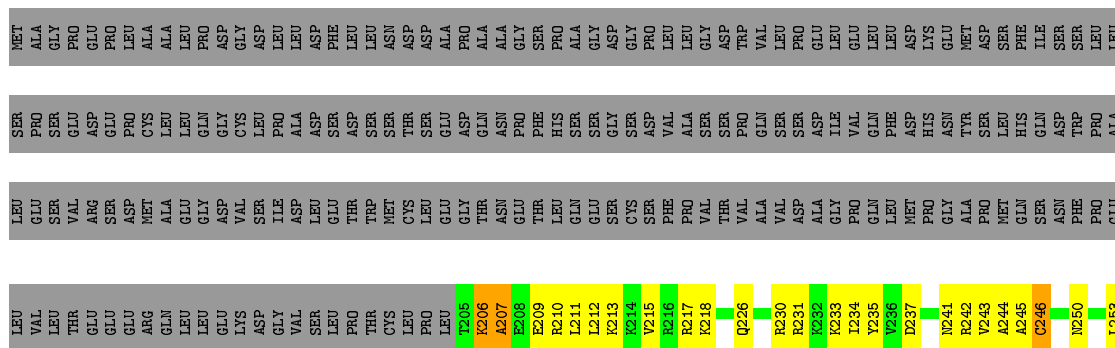




• Molecule 1: Uncharacterized protein



• Molecule 1: Uncharacterized protein





## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	137.90Å 167.06Å 115.46Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	51.64 – 3.95 51.64 – 3.95	Depositor EDS
% Data completeness (in resolution range)	99.8 (51.64-3.95) 99.9 (51.64-3.95)	Depositor EDS
$R_{merge}$	0.30	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.92 (at 4.00Å)	Xtrriage
Refinement program	BUSTER 2.10.3	Depositor
R, $R_{free}$	0.261 , 0.300 0.312 , 0.355	Depositor DCC
$R_{free}$ test set	605 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	126.6	Xtrriage
Anisotropy	0.469	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 168.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.89	EDS
Total number of atoms	3563	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	158.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.96% 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	0.74	0/568	1.10	3/751 (0.4%)
1	B	0.76	0/512	1.07	2/679 (0.3%)
1	C	0.78	0/510	1.07	2/675 (0.3%)
1	D	0.81	0/519	0.99	0/685
1	E	0.75	0/564	1.05	3/745 (0.4%)
1	F	0.77	0/594	1.03	1/785 (0.1%)
1	G	0.70	0/303	1.07	1/402 (0.2%)
All	All	0.76	0/3570	1.06	12/4722 (0.3%)

There are no bond length outliers.

The worst 5 of 12 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	B	276	ALA	N-CA-CB	-6.40	101.14	110.10
1	A	252	GLU	N-CA-CB	6.15	121.67	110.60
1	F	207	ALA	C-N-CA	5.93	136.52	121.70
1	E	212	LEU	N-CA-C	-5.62	95.81	111.00
1	E	266	SER	N-CA-CB	-5.54	102.19	110.50

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	567	0	626	45	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	511	0	552	34	0
1	C	509	0	550	40	0
1	D	518	0	568	38	0
1	E	563	0	616	18	0
1	F	593	0	647	38	0
1	G	302	0	309	3	0
All	All	3563	0	3868	169	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 23.

The worst 5 of 169 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:F:211:LEU:CD1	1:F:212:LEU:HD12	1.12	1.56
1:F:211:LEU:CD1	1:F:212:LEU:CD1	1.74	1.55
1:F:211:LEU:HD11	1:F:212:LEU:CD1	1.34	1.39
1:C:253:LEU:HG	1:D:253:LEU:HD22	1.28	1.15
1:F:211:LEU:HD11	1:F:212:LEU:HD13	1.32	1.12

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	66/441 (15%)	59 (89%)	7 (11%)	0	100	100
1	B	60/441 (14%)	55 (92%)	3 (5%)	2 (3%)	4	30
1	C	59/441 (13%)	57 (97%)	1 (2%)	1 (2%)	9	42
1	D	60/441 (14%)	56 (93%)	3 (5%)	1 (2%)	9	42
1	E	65/441 (15%)	55 (85%)	10 (15%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	F	69/441 (16%)	62 (90%)	6 (9%)	1 (1%)	11	45
1	G	35/441 (8%)	31 (89%)	4 (11%)	0	100	100
All	All	414/3087 (13%)	375 (91%)	34 (8%)	5 (1%)	13	49

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	206	LYS
1	B	275	GLN
1	C	216	ARG
1	B	277	LEU
1	D	215	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	62/392 (16%)	56 (90%)	6 (10%)	8	30
1	B	56/392 (14%)	50 (89%)	6 (11%)	6	27
1	C	56/392 (14%)	51 (91%)	5 (9%)	9	35
1	D	57/392 (14%)	53 (93%)	4 (7%)	15	43
1	E	62/392 (16%)	56 (90%)	6 (10%)	8	30
1	F	65/392 (17%)	61 (94%)	4 (6%)	18	46
1	G	32/392 (8%)	25 (78%)	7 (22%)	1	6
All	All	390/2744 (14%)	352 (90%)	38 (10%)	8	30

5 of 38 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	216	ARG
1	E	209	GLU
1	G	250	ASN

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Mol	Chain	Res	Type
1	D	251	HIS
1	E	237	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 25 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	249	GLN
1	D	226	GLN
1	F	275	GLN
1	C	250	ASN
1	D	264	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 monosaccharides 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

Unable to reproduce the depositors R factor - this section is therefore empty.

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

Unable to reproduce the depositors R factor - this section is therefore empty.

### 6.3 Carbohydrates

Unable to reproduce the depositors R factor - this section is therefore empty.

### 6.4 Ligands

Unable to reproduce the depositors R factor - this section is therefore empty.

### 6.5 Other polymers

Unable to reproduce the depositors R factor - this section is therefore empty.