



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 3ZGD
Title : crystal structure of a KEAP1 mutant
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Deposited on : 2012-12-17
Resolution : 1.98 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 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

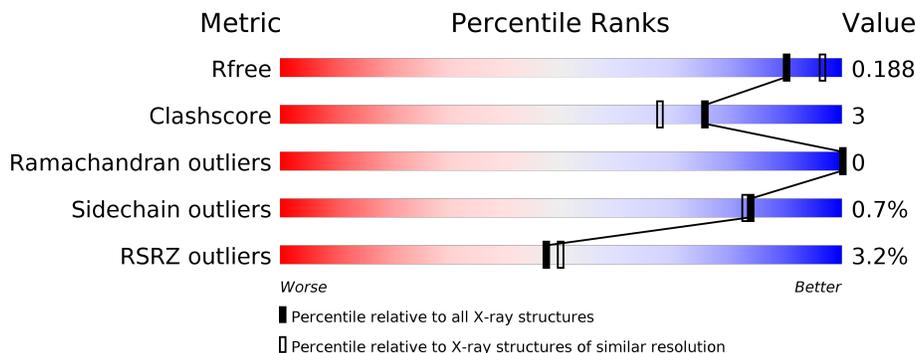
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.98 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	11647 (2.00-1.96)
Clashscore	141614	1014 (1.98-1.98)
Ramachandran outliers	138981	1006 (1.98-1.98)
Sidechain outliers	138945	1006 (1.98-1.98)
RSRZ outliers	127900	11410 (2.00-1.96)

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	310	
1	B	310	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	ACT	B	1611	-	-	X	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 5099 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 KELCH-LIKE ECH-ASSOCIATED PROTEIN 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	285	2197	1365	401	416	15	0	1	0
1	B	285	2194	1363	399	417	15	0	1	0

There are 46 discrepancies between the modelled and reference sequences:

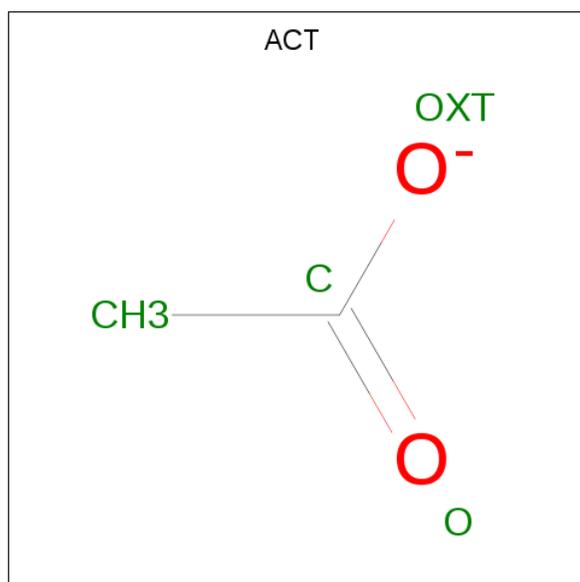
Chain	Residue	Modelled	Actual	Comment	Reference
A	300	MET	-	expression tag	UNP Q14145
A	301	GLY	-	expression tag	UNP Q14145
A	302	SER	-	expression tag	UNP Q14145
A	303	SER	-	expression tag	UNP Q14145
A	304	HIS	-	expression tag	UNP Q14145
A	305	HIS	-	expression tag	UNP Q14145
A	306	HIS	-	expression tag	UNP Q14145
A	307	HIS	-	expression tag	UNP Q14145
A	308	HIS	-	expression tag	UNP Q14145
A	309	HIS	-	expression tag	UNP Q14145
A	310	SER	-	expression tag	UNP Q14145
A	311	SER	-	expression tag	UNP Q14145
A	312	GLY	-	expression tag	UNP Q14145
A	313	LEU	-	expression tag	UNP Q14145
A	314	VAL	-	expression tag	UNP Q14145
A	315	PRO	-	expression tag	UNP Q14145
A	316	ARG	-	expression tag	UNP Q14145
A	317	GLY	-	expression tag	UNP Q14145
A	318	SER	-	expression tag	UNP Q14145
A	319	HIS	-	expression tag	UNP Q14145
A	320	MET	-	expression tag	UNP Q14145
A	540	ALA	GLU	engineered mutation	UNP Q14145
A	542	ALA	GLU	engineered mutation	UNP Q14145
B	300	MET	-	expression tag	UNP Q14145
B	301	GLY	-	expression tag	UNP Q14145

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Chain	Residue	Modelled	Actual	Comment	Reference
B	302	SER	-	expression tag	UNP Q14145
B	303	SER	-	expression tag	UNP Q14145
B	304	HIS	-	expression tag	UNP Q14145
B	305	HIS	-	expression tag	UNP Q14145
B	306	HIS	-	expression tag	UNP Q14145
B	307	HIS	-	expression tag	UNP Q14145
B	308	HIS	-	expression tag	UNP Q14145
B	309	HIS	-	expression tag	UNP Q14145
B	310	SER	-	expression tag	UNP Q14145
B	311	SER	-	expression tag	UNP Q14145
B	312	GLY	-	expression tag	UNP Q14145
B	313	LEU	-	expression tag	UNP Q14145
B	314	VAL	-	expression tag	UNP Q14145
B	315	PRO	-	expression tag	UNP Q14145
B	316	ARG	-	expression tag	UNP Q14145
B	317	GLY	-	expression tag	UNP Q14145
B	318	SER	-	expression tag	UNP Q14145
B	319	HIS	-	expression tag	UNP Q14145
B	320	MET	-	expression tag	UNP Q14145
B	540	ALA	GLU	engineered mutation	UNP Q14145
B	542	ALA	GLU	engineered mutation	UNP Q14145

- Molecule 2 is ACETATE ION (three-letter code: ACT) (formula: C₂H₃O₂).



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

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0
2	B	1	Total C O 4 2 2	0	0
2	B	1	Total C O 4 2 2	0	0
2	B	1	Total C O 4 2 2	0	0
2	B	1	Total C O 4 2 2	0	0
2	B	1	Total C O 4 2 2	0	0
2	B	1	Total C O 4 2 2	0	0

- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Na 1 1	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	348	Total O 348 348	0	0
4	B	319	Total O 319 319	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	75.72Å 75.77Å 202.04Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	37.86 – 1.98 37.86 – 1.98	Depositor EDS
% Data completeness (in resolution range)	99.9 (37.86-1.98) 99.9 (37.86-1.98)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.91 (at 1.98Å)	Xtrriage
Refinement program	BUSTER 2.11.2	Depositor
R, R_{free}	0.164 , 0.181 0.168 , 0.188	Depositor DCC
R_{free} test set	4014 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å ²)	39.6	Xtrriage
Anisotropy	0.135	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 54.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	0.025 for k,h,-l	Xtrriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	5099	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.52% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² 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

5.1 Standard geometry

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

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	0.42	0/2250	0.63	0/3062
1	B	0.42	0/2247	0.63	0/3059
All	All	0.42	0/4497	0.63	0/6121

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

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	2197	0	2093	13	0
1	B	2194	0	2086	14	0
2	A	16	0	12	0	0
2	B	24	0	18	5	0
3	A	1	0	0	0	0
4	A	348	0	0	0	0
4	B	319	0	0	1	0
All	All	5099	0	4209	27	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 3.

All (27) 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:493:GLU:H	2:B:1611:ACT:H1	1.37	0.89
1:A:466:ALA:HB1	1:A:514:VAL:HG23	1.74	0.69
1:B:466:ALA:HB1	1:B:514:VAL:HG23	1.76	0.67
1:A:330:THR:HG21	1:A:597:MET:HE3	1.87	0.56
1:B:330:THR:HG21	1:B:597:MET:HE3	1.86	0.56
2:B:1611:ACT:H2	4:B:2182:HOH:O	2.10	0.50
1:B:492:PRO:HD2	2:B:1611:ACT:H3	1.94	0.49
1:B:426:TYR:CZ	1:B:442:ARG:HD3	2.47	0.49
1:A:362:ARG:NH2	1:A:394:ASP:OD2	2.47	0.47
1:B:362:ARG:HG3	1:B:365:LEU:HD13	1.96	0.47
1:B:515:LEU:HD22	1:B:566:ILE:HG13	1.99	0.45
1:A:472:LEU:HB3	1:A:490:TYR:HB3	1.98	0.45
1:B:369:VAL:HG21	1:B:609:THR:HB	1.98	0.44
1:A:466:ALA:HB1	1:A:514:VAL:CG2	2.47	0.44
1:B:472:LEU:HB3	1:B:490:TYR:HB3	1.99	0.44
1:B:493:GLU:N	2:B:1611:ACT:H1	2.19	0.44
1:A:515:LEU:HD22	1:A:566:ILE:HG13	1.99	0.43
1:B:493:GLU:H	2:B:1611:ACT:CH3	2.20	0.43
1:A:369:VAL:HG21	1:A:609:THR:HB	2.00	0.43
1:A:426:TYR:CZ	1:A:442[B]:ARG:HD3	2.54	0.42
1:A:362:ARG:HD2	1:A:378:GLY:O	2.19	0.42
1:A:336:ARG:CZ	1:B:336:ARG:HD3	2.49	0.42
1:A:426:TYR:CZ	1:A:442[A]:ARG:HD3	2.55	0.41
1:A:365:LEU:H	1:A:365:LEU:HD23	1.86	0.41
1:B:365:LEU:HD23	1:B:365:LEU:H	1.85	0.40
1:A:416:ILE:HD11	1:A:427:ALA:HB1	2.03	0.40
1:B:416:ILE:HD11	1:B:427:ALA:HB1	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	284/310 (92%)	277 (98%)	7 (2%)	0	100	100
1	B	284/310 (92%)	277 (98%)	7 (2%)	0	100	100
All	All	568/620 (92%)	554 (98%)	14 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	230/250 (92%)	228 (99%)	2 (1%)	78	77
1	B	230/250 (92%)	229 (100%)	1 (0%)	91	90
All	All	460/500 (92%)	457 (99%)	3 (1%)	84	83

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	528	GLN
1	A	557	LEU
1	B	557	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 11 ligands modelled in this entry, 1 is monoatomic - leaving 10 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
2	ACT	A	1611	-	1,3,3	4.29	1 (100%)	0,3,3	0.00	-
2	ACT	B	1615	-	1,3,3	4.49	1 (100%)	0,3,3	0.00	-
2	ACT	A	1612	-	1,3,3	4.52	1 (100%)	0,3,3	0.00	-
2	ACT	A	1610	-	1,3,3	5.01	1 (100%)	0,3,3	0.00	-
2	ACT	B	1612	-	1,3,3	5.53	1 (100%)	0,3,3	0.00	-
2	ACT	B	1611	-	1,3,3	1.26	0	0,3,3	0.00	-
2	ACT	A	1613	-	1,3,3	5.62	1 (100%)	0,3,3	0.00	-
2	ACT	B	1613	-	1,3,3	3.72	1 (100%)	0,3,3	0.00	-
2	ACT	B	1610	-	1,3,3	4.14	1 (100%)	0,3,3	0.00	-
2	ACT	B	1614	-	1,3,3	5.11	1 (100%)	0,3,3	0.00	-

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1613	ACT	CH3-C	5.62	1.55	1.48
2	B	1612	ACT	CH3-C	5.53	1.55	1.48
2	B	1614	ACT	CH3-C	5.11	1.55	1.48
2	A	1610	ACT	CH3-C	5.01	1.55	1.48
2	A	1612	ACT	CH3-C	4.52	1.54	1.48
2	B	1615	ACT	CH3-C	4.49	1.54	1.48
2	A	1611	ACT	CH3-C	4.29	1.54	1.48
2	B	1610	ACT	CH3-C	4.14	1.54	1.48
2	B	1613	ACT	CH3-C	3.72	1.53	1.48

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1611	ACT	5	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	285/310 (91%)	0.06	6 (2%) 63 65	31, 40, 58, 79	1 (0%)
1	B	285/310 (91%)	0.18	12 (4%) 36 38	32, 40, 57, 82	1 (0%)
All	All	570/620 (91%)	0.12	18 (3%) 47 50	31, 40, 58, 82	2 (0%)

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	465	VAL	4.6
1	B	349	ASP	3.4
1	B	512	VAL	3.3
1	B	418	VAL	3.2
1	A	325	GLY	3.0
1	A	349	ASP	2.9
1	B	463	VAL	2.9
1	B	416	ILE	2.8
1	A	557	LEU	2.7
1	A	348	SER	2.6
1	B	466	ALA	2.4
1	A	587	ASP	2.3
1	B	428	VAL	2.1
1	B	468	LEU	2.1
1	B	464	GLY	2.1
1	A	559	ILE	2.1
1	B	514	VAL	2.1
1	B	552	HIS	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 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
2	ACT	B	1610	4/4	0.85	0.23	54,56,57,59	0
2	ACT	B	1611	4/4	0.90	0.16	33,40,41,47	0
2	ACT	B	1615	4/4	0.94	0.15	44,50,54,56	0
2	ACT	B	1614	4/4	0.94	0.16	45,49,50,54	0
2	ACT	A	1613	4/4	0.96	0.23	50,51,52,53	0
2	ACT	A	1612	4/4	0.97	0.11	44,45,45,55	0
2	ACT	B	1613	4/4	0.97	0.11	44,45,47,50	0
2	ACT	B	1612	4/4	0.98	0.14	34,34,35,39	0
3	NA	A	1614	1/1	0.98	0.04	49,49,49,49	0
2	ACT	A	1611	4/4	0.98	0.13	47,48,50,58	0
2	ACT	A	1610	4/4	0.98	0.09	37,38,39,40	0

6.5 Other polymers [i](#)

There are no such residues in this entry.