



Full wwPDB X-ray Structure Validation Report ⓘ

May 25, 2020 – 10:04 pm BST

PDB ID : 4IFL
Title : Crystal Structures of apo Keap1, Keap1-peptide, and Keap1-compound complexes
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Deposited on : 2012-12-14
Resolution : 1.80 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
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

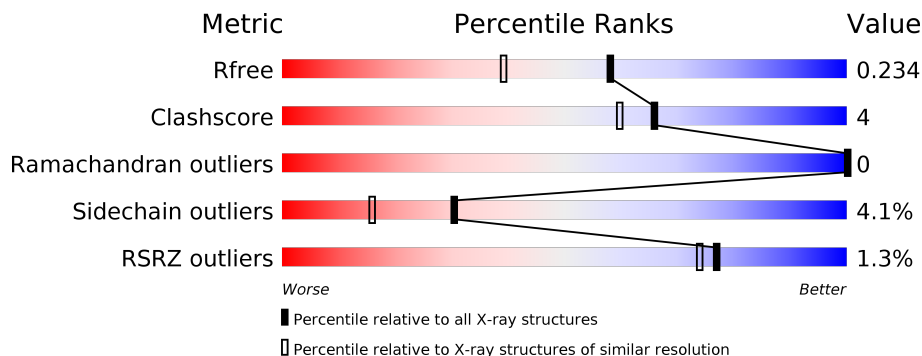
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

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	X	291	
2	P	16	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2623 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	X	285	2190	1361	396	418	15	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
X	319	SER	-	EXPRESSION TAG	UNP Q14145
X	320	SER	-	EXPRESSION TAG	UNP Q14145

- Molecule 2 is a protein called Nrf2 peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	P	16	132	86	18	28	0	0	0

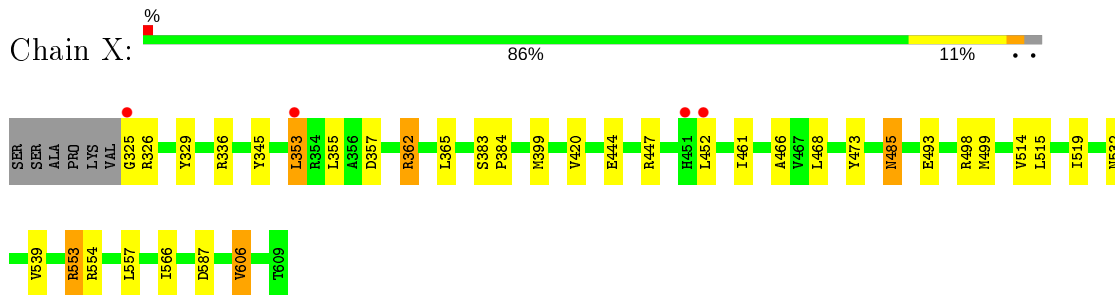
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	X	283	Total 283	O 283	0	0
3	P	18	Total 18	O 18	0	0

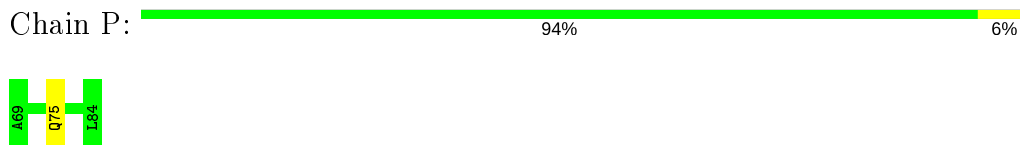
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: kelch-like ECH-associated protein 1



- Molecule 2: Nrf2 peptide



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	77.50Å 92.03Å 46.16Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.66 – 1.80 39.66 – 1.80	Depositor EDS
% Data completeness (in resolution range)	88.4 (39.66-1.80) 88.4 (39.66-1.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.53 (at 1.81Å)	Xtrriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.190 , 0.235 0.188 , 0.234	Depositor DCC
R_{free} test set	1394 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	29.9	Xtrriage
Anisotropy	0.043	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 42.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.37$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	2623	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.41% 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](#)

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	X	1.35	6/2243 (0.3%)	1.07	6/3053 (0.2%)
2	P	1.47	0/134	1.04	0/178
All	All	1.36	6/2377 (0.3%)	1.07	6/3231 (0.2%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	X	473	TYR	CD1-CE1	6.44	1.49	1.39
1	X	325	GLY	N-CA	5.80	1.54	1.46
1	X	345	TYR	CD1-CE1	-5.33	1.31	1.39
1	X	606	VAL	CB-CG1	-5.22	1.41	1.52
1	X	420	VAL	CB-CG1	5.05	1.63	1.52
1	X	473	TYR	CE1-CZ	5.01	1.45	1.38

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	499	MET	CG-SD-CE	-7.43	88.31	100.20
1	X	336	ARG	CB-CA-C	-6.49	97.41	110.40
1	X	554	ARG	NE-CZ-NH1	-6.06	117.27	120.30
1	X	357	ASP	CB-CG-OD1	5.67	123.40	118.30
1	X	498	ARG	NE-CZ-NH2	-5.16	117.72	120.30
1	X	587	ASP	CB-CG-OD1	5.12	122.91	118.30

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	X	2190	0	2078	17	0
2	P	132	0	117	1	0
3	P	18	0	0	1	0
3	X	283	0	0	2	0
All	All	2623	0	2195	18	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 4.

All (18) 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:X:466:ALA:HB1	1:X:514:VAL:HG23	1.69	0.75
1:X:383:SER:HB2	1:X:384:PRO:HD2	1.74	0.69
1:X:485:ASN:HD22	1:X:485:ASN:H	1.47	0.63
1:X:553:ARG:HD2	3:X:972:HOH:O	1.99	0.62
2:P:75:GLN:HG3	3:P:117:HOH:O	2.09	0.52
1:X:515:LEU:HD22	1:X:566:ILE:HG13	1.93	0.49
1:X:326:ARG:HG2	1:X:326:ARG:HH21	1.79	0.48
1:X:468:LEU:CD2	1:X:539:VAL:HG23	2.45	0.46
1:X:452:LEU:HD22	3:X:932:HOH:O	2.15	0.46
1:X:444:GLU:OE1	1:X:447:ARG:HD3	2.16	0.45
1:X:362:ARG:HG3	1:X:365:LEU:HD13	1.99	0.45
1:X:468:LEU:HD21	1:X:539:VAL:HG23	2.01	0.43
1:X:353:LEU:HD13	1:X:355:LEU:HD21	2.01	0.42
1:X:485:ASN:HD22	1:X:485:ASN:N	2.11	0.42
1:X:329:TYR:O	1:X:606:VAL:HA	2.19	0.41
1:X:383:SER:HB2	1:X:384:PRO:CD	2.46	0.40
1:X:329:TYR:O	1:X:606:VAL:HG23	2.20	0.40
1:X:365:LEU:HD23	1:X:365:LEU:H	1.86	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	X	283/291 (97%)	275 (97%)	8 (3%)	0	100	100
2	P	14/16 (88%)	14 (100%)	0	0	100	100
All	All	297/307 (97%)	289 (97%)	8 (3%)	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	X	230/236 (98%)	220 (96%)	10 (4%)	29	14
2	P	13/13 (100%)	13 (100%)	0	100	100
All	All	243/249 (98%)	233 (96%)	10 (4%)	30	16

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	X	353	LEU
1	X	362	ARG
1	X	399	MET
1	X	461	ILE
1	X	485	ASN
1	X	493	GLU
1	X	519	ILE
1	X	532	ASN
1	X	553	ARG
1	X	557	LEU

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

Mol	Chain	Res	Type
1	X	414	ASN
1	X	485	ASN
1	X	532	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 [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	X	285/291 (97%)	-0.07	4 (1%) 75 72	19, 25, 36, 52	0
2	P	16/16 (100%)	-0.17	0 100 100	22, 28, 37, 38	0
All	All	301/307 (98%)	-0.07	4 (1%) 77 74	19, 26, 36, 52	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	X	452	LEU	3.2
1	X	353	LEU	2.5
1	X	451	HIS	2.5
1	X	325	GLY	2.3

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.