



# Full wwPDB X-ray Structure Validation Report ⓘ

May 23, 2020 – 02:30 pm BST

PDB ID : 5NLB  
Title : Crystal structure of human CUL3 N-terminal domain bound to KEAP1 BTB and 3-box  
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Deposited on : 2017-04-04  
Resolution : 3.45 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](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	:	2.11
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

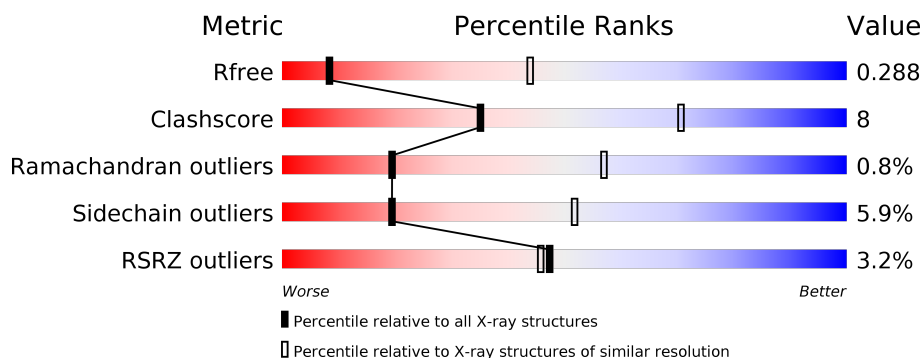
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.45 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	1291 (3.52-3.40)
Clashscore	141614	1372 (3.52-3.40)
Ramachandran outliers	138981	1337 (3.52-3.40)
Sidechain outliers	138945	1338 (3.52-3.40)
RSRZ outliers	127900	1205 (3.52-3.40)

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  $\leq 5\%$ . 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	154	<div> <div>3%</div> <div> <div></div> <div>70%</div> <div>29%</div> <div>.</div> </div> </div>
2	B	356	<div> <div>3%</div> <div> <div></div> <div>76%</div> <div>21%</div> <div>..</div> </div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3882 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
1	A	154	Total	C	N	O	S	0	0	0
			1039	664	172	195	8			

- Molecule 2 is a protein called Cullin-3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	348	Total	C	N	O	S	0	0	0
			2843	1786	500	534	23			

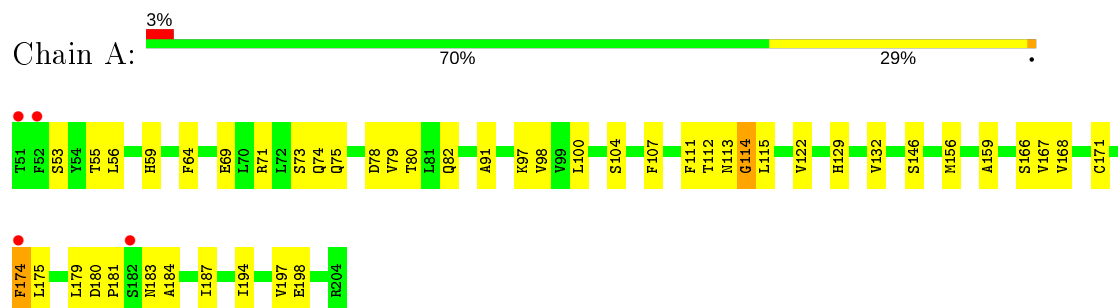
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	342	ARG	ILE	engineered mutation	UNP Q13618
B	346	ASP	LEU	engineered mutation	UNP Q13618

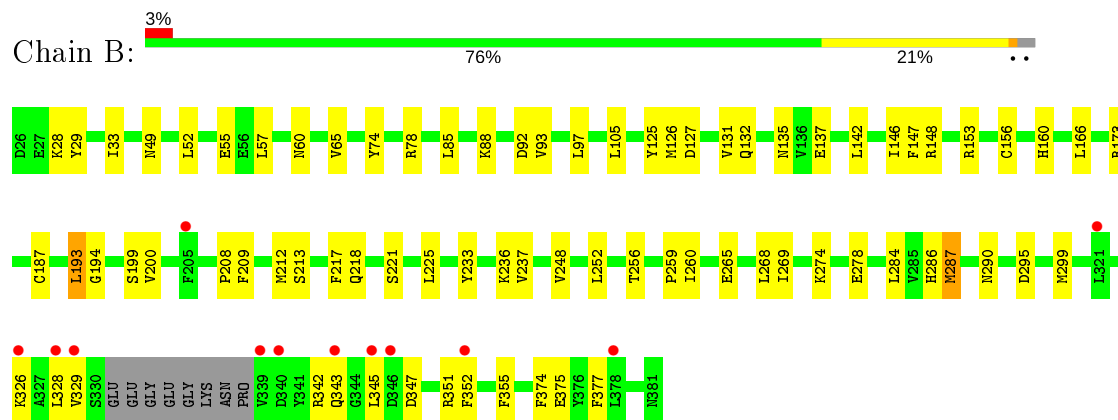
### 3 Residue-property plots [i](#)

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: Cullin-3



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	41.00Å 233.38Å 164.71Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.38 – 3.45 40.38 – 3.45	Depositor EDS
% Data completeness (in resolution range)	99.6 (40.38-3.45) 99.7 (40.38-3.45)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.78 (at 3.48Å)	Xtriage
Refinement program	PHENIX (1.11.1 _2575: ???)	Depositor
R, $R_{free}$	0.238 , 0.288 0.238 , 0.288	Depositor DCC
$R_{free}$ test set	521 reflections (4.78%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	140.2	Xtriage
Anisotropy	0.462	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 125.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.45$ , $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	3882	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	167.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.11% 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	0.52	0/1053	0.69	0/1424
2	B	0.56	0/2883	0.80	3/3870 (0.1%)
All	All	0.55	0/3936	0.77	3/5294 (0.1%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	225	LEU	CA-CB-CG	6.06	129.23	115.30
2	B	193	LEU	CB-CG-CD2	-5.08	102.37	111.00
2	B	187	CYS	CA-CB-SG	5.04	123.06	114.00

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	1039	0	872	20	0
2	B	2843	0	2830	44	1
All	All	3882	0	3702	62	1

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 8.

All (62) 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:156:MET:HE1	1:A:175:LEU:HD11	1.64	0.78
1:A:184:ALA:HA	1:A:187:ILE:HD12	1.65	0.77
2:B:28:LYS:HG3	2:B:29:TYR:HD1	1.58	0.69
2:B:148:ARG:HB2	2:B:193:LEU:HD22	1.75	0.68
1:A:115:LEU:HB3	2:B:52:LEU:O	1.95	0.66
2:B:342:ARG:HA	2:B:345:LEU:HD12	1.76	0.66
2:B:74:TYR:OH	2:B:137:GLU:O	2.14	0.65
1:A:171:CYS:HA	1:A:174:PHE:HB3	1.77	0.65
1:A:71:ARG:O	1:A:74:GLN:N	2.30	0.64
1:A:56:LEU:HD22	1:A:59:HIS:HB2	1.78	0.64
1:A:73:SER:HB2	1:A:75:GLN:HG2	1.79	0.63
2:B:209:PHE:CE2	2:B:260:ILE:HD13	2.34	0.63
2:B:221:SER:HB3	2:B:268:LEU:HD23	1.80	0.63
2:B:153:ARG:HH12	2:B:200:VAL:HA	1.67	0.59
2:B:213:SER:HB3	2:B:217:PHE:CE2	2.39	0.58
1:A:82:GLN:HA	1:A:91:ALA:O	2.04	0.58
2:B:265:GLU:HA	2:B:269:ILE:HD12	1.85	0.57
2:B:49:ASN:HA	2:B:52:LEU:HD12	1.87	0.56
1:A:78:ASP:OD1	1:A:97:LYS:N	2.24	0.56
2:B:347:ASP:O	2:B:351:ARG:HB2	2.05	0.56
2:B:166:LEU:HD11	2:B:208:PRO:HB2	1.88	0.55
2:B:65:VAL:HB	2:B:125:TYR:HD2	1.71	0.55
2:B:88:LYS:NZ	2:B:92:ASP:OD2	2.39	0.55
1:A:183:ASN:O	1:A:187:ILE:HG13	2.08	0.54
1:A:100:LEU:HB3	1:A:107:PHE:CD2	2.42	0.53
1:A:159:ALA:HB3	1:A:168:VAL:HG23	1.90	0.53
1:A:171:CYS:HA	1:A:174:PHE:CB	2.38	0.53
2:B:65:VAL:HG21	2:B:126:MET:HB2	1.91	0.52
2:B:85:LEU:HD12	2:B:146:ILE:HG22	1.92	0.52
2:B:173:ARG:NH2	2:B:213:SER:OG	2.38	0.51
1:A:79:VAL:HG21	1:A:111:PHE:HE1	1.75	0.51
2:B:166:LEU:HD22	2:B:212:MET:CE	2.41	0.49
2:B:233:TYR:O	2:B:237:VAL:HG23	2.12	0.49
2:B:256:THR:O	2:B:259:PRO:HD2	2.12	0.49
2:B:218:GLN:NE2	2:B:221:SER:OG	2.46	0.49
2:B:287:MET:CE	2:B:295:ASP:HB3	2.44	0.48
2:B:28:LYS:HG3	2:B:29:TYR:CD1	2.44	0.48
2:B:284:LEU:HD12	2:B:299:MET:SD	2.53	0.47
2:B:374:PHE:HA	2:B:377:PHE:CZ	2.49	0.47
2:B:33:ILE:HG23	2:B:60:ASN:HB3	1.96	0.46
2:B:286:HIS:CE1	2:B:290:ASN:ND2	2.84	0.46
2:B:256:THR:C	2:B:259:PRO:HD2	2.35	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:97:LEU:O	2:B:160:HIS:NE2	2.49	0.46
2:B:248:VAL:HG22	2:B:252:LEU:HD12	1.98	0.45
1:A:100:LEU:HB3	1:A:107:PHE:HD2	1.83	0.44
1:A:180:ASP:HB3	1:A:183:ASN:H	1.82	0.44
1:A:129:HIS:O	1:A:132:VAL:HG22	2.18	0.44
2:B:131:VAL:HG23	2:B:132:GLN:HE21	1.81	0.44
2:B:352:PHE:HA	2:B:355:PHE:HD2	1.83	0.44
2:B:274:LYS:HE3	2:B:278:GLU:OE2	2.19	0.43
2:B:93:VAL:HG11	2:B:105:LEU:HD13	2.00	0.43
2:B:52:LEU:HD23	2:B:52:LEU:HA	1.78	0.43
2:B:78:ARG:HD3	2:B:142:LEU:HD11	2.00	0.43
2:B:166:LEU:HD22	2:B:212:MET:HE2	2.00	0.42
1:A:115:LEU:HG	2:B:52:LEU:HB2	2.01	0.42
2:B:65:VAL:HB	2:B:125:TYR:CD2	2.53	0.42
1:A:112:THR:O	1:A:114:GLY:N	2.53	0.42
2:B:85:LEU:HD13	2:B:147:PHE:HA	2.02	0.41
2:B:326:LYS:HA	2:B:329:VAL:HG22	2.03	0.41
2:B:57:LEU:HD23	2:B:57:LEU:HA	1.88	0.41
1:A:174:PHE:CD1	1:A:174:PHE:C	2.94	0.41
2:B:194:GLY:HA2	2:B:200:VAL:HG23	2.01	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:236:LYS:NZ	2:B:375:GLU:OE1[1_655]	2.19	0.01

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

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	152/154 (99%)	124 (82%)	24 (16%)	4 (3%)	5 32

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	344/356 (97%)	333 (97%)	11 (3%)	0	100	100
All	All	496/510 (97%)	457 (92%)	35 (7%)	4 (1%)	19	57

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	114	GLY
1	A	113	ASN
1	A	194	ILE
1	A	197	VAL

### 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	77/131 (59%)	62 (80%)	15 (20%)	1	6
2	B	315/324 (97%)	307 (98%)	8 (2%)	47	75
All	All	392/455 (86%)	369 (94%)	23 (6%)	19	52

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	53	SER
1	A	55	THR
1	A	64	PHE
1	A	69	GLU
1	A	80	THR
1	A	98	VAL
1	A	104	SER
1	A	122	VAL
1	A	146	SER
1	A	166	SER
1	A	167	VAL
1	A	174	PHE

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Mol	Chain	Res	Type
1	A	179	LEU
1	A	181	PRO
1	A	198	GLU
2	B	55	GLU
2	B	127	ASP
2	B	135	ASN
2	B	156	CYS
2	B	199	SER
2	B	287	MET
2	B	328	LEU
2	B	343	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	132	GLN
2	B	218	GLN
2	B	286	HIS
2	B	290	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, 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	154/154 (100%)	-0.11	4 (2%) 56 53	130, 163, 238, 259	0
2	B	348/356 (97%)	0.18	12 (3%) 45 43	123, 163, 219, 245	0
All	All	502/510 (98%)	0.09	16 (3%) 47 46	123, 163, 227, 259	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	328	LEU	10.4
2	B	329	VAL	4.2
2	B	352	PHE	3.3
2	B	343	GLN	3.1
2	B	321	LEU	2.9
2	B	340	ASP	2.9
2	B	345	LEU	2.8
2	B	326	LYS	2.7
1	A	52	PHE	2.7
1	A	182	SER	2.4
1	A	51	THR	2.4
2	B	346	ASP	2.3
2	B	205	PHE	2.3
2	B	339	VAL	2.2
2	B	378	LEU	2.1
1	A	174	PHE	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](#)

There are no ligands in this entry.

### 6.5 Other polymers [i](#)

There are no such residues in this entry.