



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 17, 2017 – 10:28 AM EDT

PDB ID : 4ZGC
Title : Crystal Structure Analysis of Kelch protein (with disulfide bond) from Plasmodium falciparum
Authors : Jiang, D.Q.; Tempel, W.; Loppnau, P.; Graslund, S.; He, H.; Ravichandran, M.; Seitova, A.; Arrowsmith, C.H.; Edwards, A.M.; Bountra, C.; El Bakkouri, M.; Senisterra, G.; Osman, K.T.; Lovato, D.V.; Hui, R.; Hutchinson, A.; Lin, Y.H.; Structural Genomics Consortium (SGC)
Deposited on : unknown
Resolution : 2.50 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

<http://wwpdb.org/validation/2016/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)	:	1.9-1692
EDS	:	rb-20029824
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20029824

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	UNX	A	801	-	-	-	X
2	UNX	A	803	-	-	-	X
2	UNX	A	820	-	-	-	X

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5898 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 protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	372	Total	C	N	O	S	0	4	0
			2877	1854	473	534	16			
1	B	373	Total	C	N	O	S	0	4	0
			2861	1837	480	528	16			

- Molecule 2 is UNKNOWN ATOM OR ION (three-letter code: UNX) (formula: X).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	12	Total	X	0	0
			12	12		
2	A	24	Total	X	0	0
			24	24		

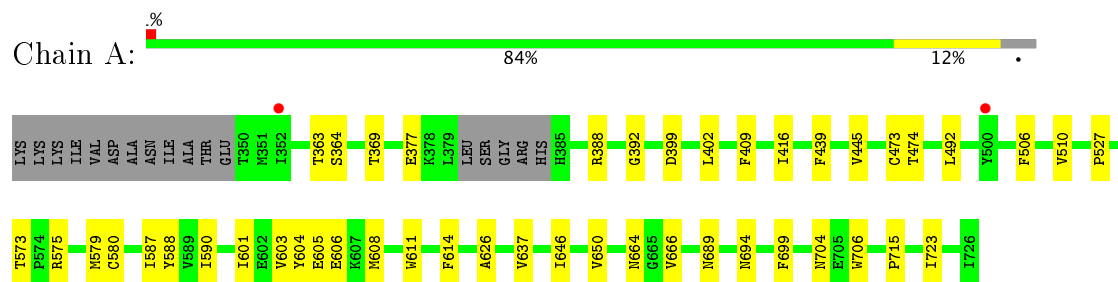
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	62	Total	O	0	0
			62	62		
3	B	62	Total	O	0	0
			62	62		

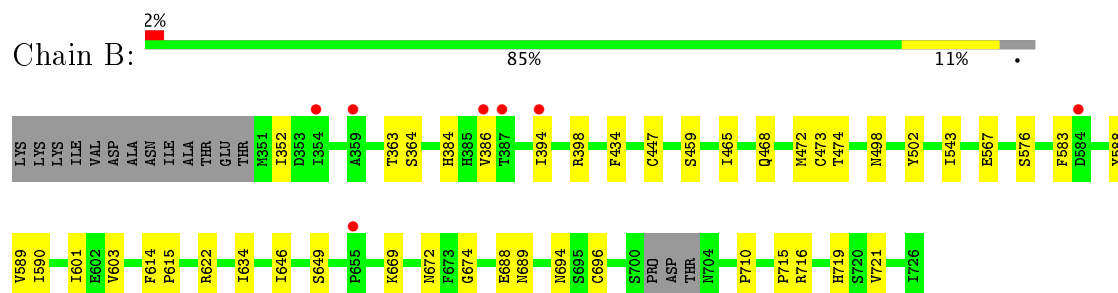
3 Residue-property plots [i](#)

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 protein



- Molecule 1: Kelch protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	62.94Å 117.63Å 74.51Å 90.00° 113.77° 90.00°	Depositor
Resolution (Å)	36.49 – 2.50 36.49 – 2.50	Depositor EDS
% Data completeness (in resolution range)	98.2 (36.49-2.50) 98.2 (36.49-2.50)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.33 (at 2.51Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, R_{free}	0.192 , 0.253 0.204 , 0.264	Depositor DCC
R_{free} test set	1668 reflections (5.22%)	DCC
Wilson B-factor (Å ²)	35.6	Xtriage
Anisotropy	0.423	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 44.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.033 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5898	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.98% 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](#)

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

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	0.46	0/2958	0.59	0/4020
1	B	0.44	0/2939	0.58	0/3995
All	All	0.45	0/5897	0.58	0/8015

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	2877	0	2684	26	0
1	B	2861	0	2649	27	0
2	A	24	0	0	0	0
2	B	12	0	0	0	0
3	A	62	0	0	1	0
3	B	62	0	0	0	0
All	All	5898	0	5333	53	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 5.

All (53) 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:384:HIS:CD2	1:B:384:HIS:CB	2.59	0.85
1:B:567:GLU:CG	1:B:567:GLU:OE1	2.36	0.73
1:B:465:ILE:O	1:B:468:GLN:NE2	2.20	0.72
1:B:384:HIS:ND1	1:B:384:HIS:CB	2.52	0.71
1:B:567:GLU:CG	1:B:567:GLU:OE2	2.39	0.69
1:A:369:THR:HG22	1:A:377:GLU:HA	1.80	0.63
1:A:506:PHE:HB3	1:A:527:PRO:HG3	1.82	0.62
1:A:637[A]:VAL:HG12	1:A:650:VAL:HG22	1.82	0.61
1:B:672:ASN:ND2	1:B:688:GLU:OE2	2.31	0.60
1:B:646:ILE:HD13	1:B:689:ASN:HA	1.83	0.58
1:B:567:GLU:OE1	1:B:567:GLU:OE2	2.23	0.56
1:A:445:VAL:HG22	1:A:723:ILE:HD12	1.88	0.55
1:B:386:VAL:HG21	1:B:394:ILE:HG12	1.88	0.54
1:B:590:ILE:HG23	1:B:601:ILE:HG12	1.90	0.53
1:B:398:ARG:HD2	1:B:434:PHE:CE2	2.45	0.52
1:A:646:ILE:HD13	1:A:689:ASN:HA	1.92	0.51
1:B:622:ARG:NH2	1:B:649:SER:OG	2.41	0.48
1:A:587:ILE:HB	1:A:604:TYR:HB3	1.96	0.48
1:B:543:ILE:HG21	1:B:589:VAL:HG21	1.95	0.48
1:A:580:CYS:HA	1:A:588:TYR:O	2.14	0.47
1:A:664:ASN:HB2	3:A:936:HOH:O	2.15	0.47
1:A:604:TYR:HB2	1:A:611:TRP:CZ3	2.50	0.47
1:B:498:ASN:HB2	1:B:502:TYR:O	2.14	0.46
1:B:614:PHE:HA	1:B:615:PRO:HD3	1.79	0.46
1:A:416:ILE:HD12	1:A:439:PHE:HB3	1.98	0.46
1:B:447[A]:CYS:SG	1:B:721:VAL:HG22	2.56	0.46
1:B:459:SER:HB2	1:B:472:MET:HE1	1.99	0.45
1:A:445:VAL:HG22	1:A:723:ILE:CD1	2.47	0.45
1:A:694:ASN:HB3	1:A:715:PRO:HG3	1.99	0.45
1:A:492:LEU:O	1:A:510:VAL:HA	2.17	0.44
1:B:473:CYS:SG	1:B:474:THR:N	2.87	0.44
1:A:573:THR:O	1:A:575:ARG:HG3	2.18	0.44
1:A:603:VAL:HG23	1:A:614:PHE:HB2	1.99	0.44
1:A:388:ARG:NH1	1:A:392:GLY:O	2.43	0.44
1:B:696[B]:CYS:HB3	1:B:710:PRO:O	2.17	0.44
1:A:473:CYS:SG	1:A:474:THR:N	2.90	0.43
1:A:363:THR:OG1	1:A:364:SER:N	2.51	0.43
1:A:699:PHE:HB2	1:A:706:TRP:CZ3	2.54	0.43
1:A:399:ASP:OD2	1:A:402:LEU:HG	2.19	0.43
1:A:579:MET:HG2	1:A:626:ALA:HA	2.01	0.43
1:B:694:ASN:HB3	1:B:715:PRO:HG3	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:409:PHE:CZ	1:A:439:PHE:HA	2.54	0.43
1:B:674:GLY:H	1:B:719:HIS:CE1	2.37	0.43
1:B:588:TYR:CD2	1:B:603:VAL:HG22	2.54	0.42
1:A:699:PHE:CE1	1:A:704:ASN:HA	2.54	0.42
1:B:363:THR:OG1	1:B:364:SER:N	2.53	0.42
1:A:604:TYR:CE2	1:A:606:GLU:HA	2.55	0.42
1:B:352:ILE:N	1:B:363:THR:O	2.51	0.41
1:B:447[B]:CYS:SG	1:B:716:ARG:HD2	2.60	0.41
1:A:605:GLU:HG3	1:A:608:MET:HE2	2.03	0.41
1:B:583:PHE:CE1	1:B:634:ILE:HG12	2.55	0.41
1:A:590:ILE:HG23	1:A:601:ILE:HG12	2.01	0.41
1:B:669:LYS:HB2	1:B:689:ASN:ND2	2.36	0.41

There are no symmetry-related clashes.

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	372/389 (96%)	358 (96%)	14 (4%)	0	100	100
1	B	373/389 (96%)	355 (95%)	18 (5%)	0	100	100
All	All	745/778 (96%)	713 (96%)	32 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	289/340 (85%)	288 (100%)	1 (0%)	94	98
1	B	282/340 (83%)	281 (100%)	1 (0%)	93	98
All	All	571/680 (84%)	569 (100%)	2 (0%)	93	98

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	666	VAL
1	B	576	SER

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

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 ⓘ

Of 36 ligands modelled in this entry, 36 are unknown - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	372/389 (95%)	-0.22	2 (0%) 90 91	19, 32, 56, 70	3 (0%)
1	B	373/389 (95%)	-0.06	7 (1%) 67 69	20, 37, 63, 81	3 (0%)
All	All	745/778 (95%)	-0.14	9 (1%) 79 80	19, 35, 60, 81	6 (0%)

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	386	VAL	3.6
1	B	354	ILE	2.8
1	B	359	ALA	2.4
1	A	500	TYR	2.4
1	B	387	THR	2.3
1	B	394	ILE	2.3
1	A	352	ILE	2.2
1	B	584	ASP	2.1
1	B	655	PRO	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. LLDF column lists the quality of electron

density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
2	UNX	A	820	1/1	0.91	0.64	17.71	16,16,16,16	0
2	UNX	A	803	1/1	0.81	0.61	16.61	29,29,29,29	0
2	UNX	A	801	1/1	0.83	0.33	6.74	21,21,21,21	0
2	UNX	A	822	1/1	0.84	0.18	-	29,29,29,29	0
2	UNX	A	811	1/1	0.85	0.68	-	23,23,23,23	0
2	UNX	B	802	1/1	0.93	0.12	-	24,24,24,24	0
2	UNX	A	813	1/1	0.91	0.58	-	32,32,32,32	0
2	UNX	B	804	1/1	0.91	0.18	-	26,26,26,26	0
2	UNX	A	804	1/1	0.58	0.22	-	25,25,25,25	0
2	UNX	B	806	1/1	0.89	0.33	-	34,34,34,34	0
2	UNX	A	810	1/1	0.97	0.07	-	23,23,23,23	0
2	UNX	A	808	1/1	0.80	0.22	-	23,23,23,23	0
2	UNX	A	814	1/1	0.87	0.49	-	30,30,30,30	0
2	UNX	A	805	1/1	0.89	0.44	-	33,33,33,33	0
2	UNX	A	807	1/1	0.91	0.11	-	30,30,30,30	0
2	UNX	A	815	1/1	0.33	0.91	-	48,48,48,48	0
2	UNX	B	808	1/1	0.83	0.28	-	26,26,26,26	0
2	UNX	B	803	1/1	0.91	0.35	-	29,29,29,29	0
2	UNX	A	812	1/1	0.77	0.17	-	33,33,33,33	0
2	UNX	A	823	1/1	0.81	0.31	-	28,28,28,28	0
2	UNX	A	809	1/1	0.75	0.26	-	37,37,37,37	0
2	UNX	B	812	1/1	0.77	0.17	-	32,32,32,32	0
2	UNX	B	805	1/1	0.91	0.36	-	28,28,28,28	0
2	UNX	A	802	1/1	0.95	0.32	-	23,23,23,23	0
2	UNX	A	818	1/1	0.93	0.37	-	28,28,28,28	0
2	UNX	B	811	1/1	0.88	0.30	-	27,27,27,27	0
2	UNX	B	801	1/1	0.87	0.28	-	23,23,23,23	0
2	UNX	A	806	1/1	0.96	0.14	-	34,34,34,34	0
2	UNX	B	807	1/1	0.92	0.53	-	37,37,37,37	0
2	UNX	B	809	1/1	0.96	0.48	-	29,29,29,29	0
2	UNX	A	817	1/1	0.91	0.46	-	35,35,35,35	0
2	UNX	B	810	1/1	0.71	0.47	-	31,31,31,31	0
2	UNX	A	821	1/1	0.86	0.20	-	23,23,23,23	0
2	UNX	A	816	1/1	0.69	0.32	-	36,36,36,36	0
2	UNX	A	824	1/1	0.86	0.20	-	31,31,31,31	0
2	UNX	A	819	1/1	0.85	0.17	-	27,27,27,27	0

6.5 Other polymers [i](#)

There are no such residues in this entry.