



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:23 PM GMT

PDB ID : 4I6E
Title : A vertebrate cryptochrome
Authors : Xing, W.; Busino, L.; Hinds, T.R.; Marionni, S.T.; Saifee, N.H.; Bush, M.F.;
Pagano, M.; Zheng, N.
Deposited on : 2012-11-29
Resolution : 2.70 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
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : trunk26865

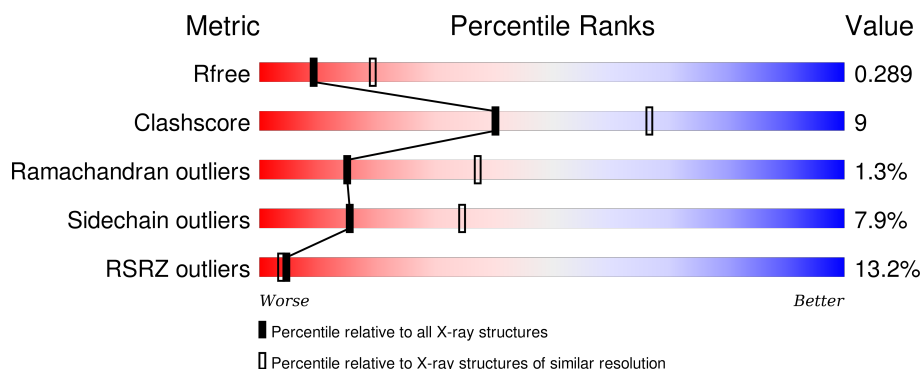
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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}	91344	2103 (2.70-2.70)
Clashscore	102246	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)
RSRZ outliers	91569	2107 (2.70-2.70)

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 ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	512	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3822 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.

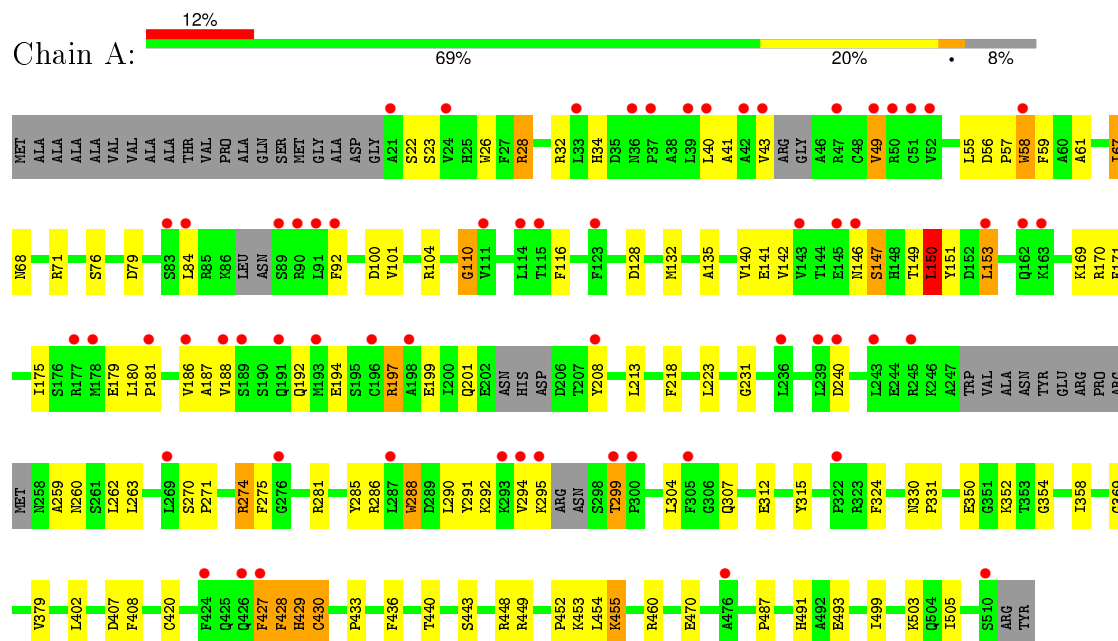
- Molecule 1 is a protein called Cryptochrome-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	471	Total	C	N	O	S	0	0	0
			3821	2460	669	673	19			

- Molecule 2 is water.

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

- Molecule 1: Cryptochrome-2



4 Data and refinement statistics

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, α , β , γ	68.94Å 68.94Å 127.53Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.74 – 2.70 48.74 – 2.40	Depositor EDS
% Data completeness (in resolution range)	100.0 (48.74-2.70) 99.9 (48.74-2.40)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.10 (at 2.39Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.1_1168)	Depositor
R, R_{free}	0.244 , 0.287 0.245 , 0.289	Depositor DCC
R_{free} test set	1377 reflections (8.41%)	DCC
Wilson B-factor (Å ²)	59.4	Xtriage
Anisotropy	0.119	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 62.5	EDS
Estimated twinning fraction	0.044 for h,-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 23256 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	3822	wwPDB-VP
Average B, all atoms (Å ²)	111.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.73% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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	0.44	0/3928	0.64	1/5323 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	150	LEU	CA-CB-CG	5.11	127.04	115.30

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	3821	0	3757	66	0
2	A	1	0	0	0	0
All	All	3822	0	3757	66	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 9.

All (66) 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:428:PHE:HA	1:A:429:HIS:HB2	1.67	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:354:GLY:HA3	1:A:487:PRO:HA	1.70	0.72
1:A:150:LEU:N	1:A:312:GLU:OE2	2.23	0.70
1:A:194:GLU:HA	1:A:197:ARG:HE	1.61	0.66
1:A:199:GLU:OE1	1:A:201:GLN:NE2	2.29	0.66
1:A:443:SER:O	1:A:448:ARG:NH2	2.28	0.66
1:A:132:MET:HG2	1:A:142:VAL:HG11	1.79	0.63
1:A:194:GLU:HB3	1:A:197:ARG:HH21	1.64	0.63
1:A:259:ALA:HA	1:A:262:LEU:HD23	1.81	0.62
1:A:436:PHE:O	1:A:440:THR:HG23	1.99	0.61
1:A:56:ASP:HB3	1:A:59:PHE:HB3	1.82	0.61
1:A:23:SER:HB2	1:A:49:VAL:HG23	1.85	0.59
1:A:149:THR:HG21	1:A:153:LEU:HD21	1.85	0.59
1:A:171:PHE:O	1:A:175:ILE:HG12	2.04	0.57
1:A:354:GLY:HA3	1:A:487:PRO:CA	2.34	0.56
1:A:40:LEU:HD23	1:A:188:VAL:HG12	1.88	0.55
1:A:34:HIS:HA	1:A:187:ALA:HB2	1.89	0.55
1:A:194:GLU:HA	1:A:197:ARG:NE	2.21	0.55
1:A:58:TRP:HA	1:A:61:ALA:HB3	1.90	0.53
1:A:429:HIS:O	1:A:430:CYS:HB2	2.08	0.53
1:A:448:ARG:HG2	1:A:455:LYS:HA	1.92	0.52
1:A:433:PRO:HD3	1:A:491:HIS:CE1	2.45	0.51
1:A:67:ILE:HG23	1:A:218:PHE:CD2	2.46	0.51
1:A:26:TRP:CD2	1:A:28:ARG:HD2	2.46	0.51
1:A:420:CYS:SG	1:A:427:PHE:HA	2.50	0.51
1:A:428:PHE:HA	1:A:429:HIS:CB	2.31	0.50
1:A:146:ASN:O	1:A:281:ARG:NH2	2.43	0.50
1:A:240:ASP:OD2	1:A:286:ARG:NH2	2.45	0.50
1:A:291:TYR:CE2	1:A:295:LYS:HD2	2.48	0.49
1:A:315:TYR:OH	1:A:407:ASP:OD1	2.09	0.47
1:A:26:TRP:HB3	1:A:116:PHE:HB3	1.97	0.47
1:A:135:ALA:HB1	1:A:140:VAL:HB	1.96	0.47
1:A:32:ARG:NH2	1:A:147:SER:HB2	2.29	0.47
1:A:56:ASP:HA	1:A:57:PRO:HD3	1.79	0.47
1:A:101:VAL:HG22	1:A:104:ARG:HH21	1.80	0.47
1:A:76:SER:HA	1:A:231:GLY:HA2	1.98	0.46
1:A:201:GLN:OE1	1:A:201:GLN:N	2.47	0.46
1:A:350:GLU:HB3	1:A:352:LYS:HE3	1.97	0.46
1:A:151:TYR:OH	1:A:175:ILE:HG23	2.17	0.45
1:A:175:ILE:HD13	1:A:175:ILE:HA	1.81	0.45
1:A:292:LYS:HD3	1:A:292:LYS:O	2.16	0.45
1:A:67:ILE:HD12	1:A:68:ASN:H	1.82	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:22:SER:OG	1:A:22:SER:O	2.35	0.45
1:A:151:TYR:CZ	1:A:181:PRO:HG3	2.52	0.45
1:A:324:PHE:HB3	1:A:408:PHE:CZ	2.52	0.44
1:A:179:GLU:H	1:A:179:GLU:CD	2.20	0.44
1:A:288:TRP:HZ2	1:A:299:THR:HG22	1.83	0.44
1:A:128:ASP:O	1:A:132:MET:HG3	2.18	0.43
1:A:181:PRO:O	1:A:285:TYR:OH	2.23	0.43
1:A:452:PRO:C	1:A:454:LEU:H	2.21	0.43
1:A:67:ILE:H	1:A:67:ILE:HG13	1.48	0.43
1:A:71:ARG:HD3	1:A:213:LEU:HD11	1.99	0.43
1:A:499:ILE:HA	1:A:499:ILE:HD13	1.83	0.42
1:A:330:ASN:HA	1:A:331:PRO:HD3	1.94	0.42
1:A:213:LEU:O	1:A:218:PHE:HB2	2.19	0.42
1:A:84:LEU:HA	1:A:84:LEU:HD23	1.91	0.42
1:A:274:ARG:HD3	1:A:275:PHE:CD1	2.55	0.42
1:A:92:PHE:HA	1:A:208:TYR:HB3	2.01	0.42
1:A:270:SER:OG	1:A:271:PRO:HD3	2.19	0.41
1:A:358:ILE:HG23	1:A:379:VAL:HG12	2.02	0.41
1:A:369:GLY:HA2	1:A:402:LEU:O	2.20	0.41
1:A:43:VAL:HG12	1:A:192:GLN:NE2	2.35	0.41
1:A:22:SER:OG	1:A:110:GLY:O	2.30	0.41
1:A:43:VAL:HG12	1:A:192:GLN:HE21	1.86	0.41
1:A:290:LEU:O	1:A:294:VAL:HG23	2.20	0.41
1:A:150:LEU:HB2	1:A:312:GLU:OE2	2.21	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	459/512 (90%)	410 (89%)	43 (9%)	6 (1%)	15 37

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	429	HIS
1	A	430	CYS
1	A	41	ALA
1	A	100	ASP
1	A	197	ARG
1	A	110	GLY

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	404/434 (93%)	372 (92%)	32 (8%)	15	34

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	ARG
1	A	49	VAL
1	A	55	LEU
1	A	58	TRP
1	A	67	ILE
1	A	79	ASP
1	A	141	GLU
1	A	147	SER
1	A	150	LEU
1	A	153	LEU
1	A	169	LYS
1	A	170	ARG
1	A	180	LEU
1	A	186	VAL
1	A	223	LEU
1	A	260	ASN
1	A	263	LEU
1	A	274	ARG
1	A	288	TRP

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Mol	Chain	Res	Type
1	A	299	THR
1	A	304	LEU
1	A	307	GLN
1	A	427	PHE
1	A	428	PHE
1	A	449	ARG
1	A	453	LYS
1	A	455	LYS
1	A	460	ARG
1	A	470	GLU
1	A	493	GLU
1	A	503	LYS
1	A	505	ILE

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](#)

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 ⓘ

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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	471/512 (91%)	0.79	62 (13%) 4 4	43, 110, 168, 201	0

All (62) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	24	VAL	8.0
1	A	33	LEU	7.6
1	A	245	ARG	7.1
1	A	193	MET	6.3
1	A	143	VAL	5.9
1	A	111	VAL	5.8
1	A	424	PHE	5.7
1	A	39	LEU	5.5
1	A	188	VAL	5.2
1	A	92	PHE	5.1
1	A	177	ARG	4.8
1	A	89	SER	4.6
1	A	191	GLN	4.6
1	A	186	VAL	4.5
1	A	36	ASN	4.5
1	A	269	LEU	4.4
1	A	51	CYS	4.3
1	A	91	LEU	4.1
1	A	510	SER	4.1
1	A	47	ARG	4.0
1	A	153	LEU	3.9
1	A	243	LEU	3.8
1	A	83	SER	3.8
1	A	21	ALA	3.7
1	A	300	PRO	3.6
1	A	240	ASP	3.6
1	A	49	VAL	3.5

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Mol	Chain	Res	Type	RSRZ
1	A	293	LYS	3.3
1	A	236	LEU	3.3
1	A	84	LEU	3.1
1	A	198	ALA	3.1
1	A	189	SER	3.0
1	A	305	PHE	3.0
1	A	114	LEU	3.0
1	A	40	LEU	3.0
1	A	162	GLN	2.9
1	A	178	MET	2.9
1	A	115	THR	2.8
1	A	50	ARG	2.7
1	A	299	THR	2.7
1	A	163	LYS	2.7
1	A	427	PHE	2.7
1	A	37	PRO	2.7
1	A	208	TYR	2.6
1	A	146	ASN	2.6
1	A	476	ALA	2.6
1	A	145	GLU	2.6
1	A	42	ALA	2.5
1	A	43	VAL	2.5
1	A	52	VAL	2.4
1	A	426	GLN	2.4
1	A	295	LYS	2.4
1	A	276	GLY	2.4
1	A	322	PRO	2.3
1	A	181	PRO	2.2
1	A	58	TRP	2.2
1	A	123	PHE	2.1
1	A	239	LEU	2.1
1	A	294	VAL	2.1
1	A	90	ARG	2.1
1	A	287	LEU	2.0
1	A	196	CYS	2.0

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

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.