



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:59 PM GMT

PDB ID : 4QLA
Title : Crystal structure of juvenile hormone epoxide hydrolase from the silkworm *Bombyx mori*
Authors : Zhou, K.; Jia, N.; Hu, C.; Jiang, Y.L.; Yang, J.P.; Chen, Y.; Li, S.; Zhou, C.Z.
Deposited on : 2014-06-11
Resolution : 2.30 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

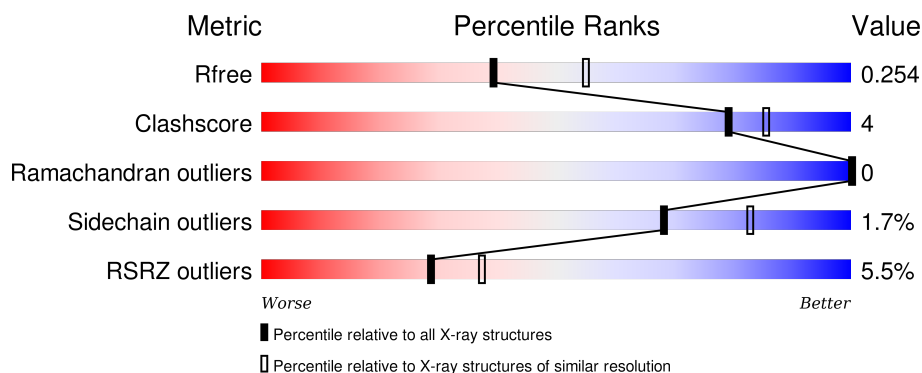
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

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	447	
1	B	447	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6682 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 Juvenile hormone epoxide hydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	424	Total	C	N	O	S	0	0	0
			3419	2245	556	608	10			
1	B	380	Total	C	N	O	S	0	0	0
			3072	2015	506	541	10			

There are 26 discrepancies between the modelled and reference sequences:

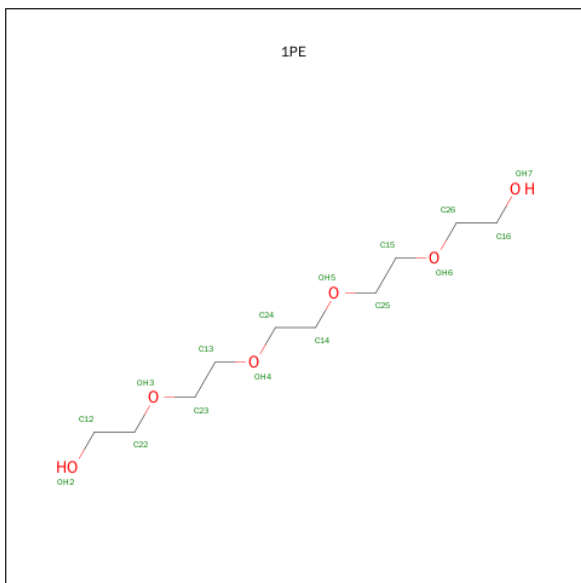
Chain	Residue	Modelled	Actual	Comment	Reference
A	15	MET	-	EXPRESSION TAG	UNP H9JWL6
A	16	HIS	-	EXPRESSION TAG	UNP H9JWL6
A	17	HIS	-	EXPRESSION TAG	UNP H9JWL6
A	18	HIS	-	EXPRESSION TAG	UNP H9JWL6
A	19	HIS	-	EXPRESSION TAG	UNP H9JWL6
A	20	HIS	-	EXPRESSION TAG	UNP H9JWL6
A	21	HIS	-	EXPRESSION TAG	UNP H9JWL6
A	22	GLY	-	EXPRESSION TAG	UNP H9JWL6
A	142	GLY	ASP	ENGINEERED MUTATION	UNP H9JWL6
A	336	LEU	PHE	ENGINEERED MUTATION	UNP H9JWL6
A	402	HIS	TYR	ENGINEERED MUTATION	UNP H9JWL6
A	414	MET	LEU	ENGINEERED MUTATION	UNP H9JWL6
A	433	ALA	SER	ENGINEERED MUTATION	UNP H9JWL6
B	15	MET	-	EXPRESSION TAG	UNP H9JWL6
B	16	HIS	-	EXPRESSION TAG	UNP H9JWL6
B	17	HIS	-	EXPRESSION TAG	UNP H9JWL6
B	18	HIS	-	EXPRESSION TAG	UNP H9JWL6
B	19	HIS	-	EXPRESSION TAG	UNP H9JWL6
B	20	HIS	-	EXPRESSION TAG	UNP H9JWL6
B	21	HIS	-	EXPRESSION TAG	UNP H9JWL6
B	22	GLY	-	EXPRESSION TAG	UNP H9JWL6
B	142	GLY	ASP	ENGINEERED MUTATION	UNP H9JWL6
B	336	LEU	PHE	ENGINEERED MUTATION	UNP H9JWL6
B	402	HIS	TYR	ENGINEERED MUTATION	UNP H9JWL6
B	414	MET	LEU	ENGINEERED MUTATION	UNP H9JWL6

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Chain	Residue	Modelled	Actual	Comment	Reference
B	433	ALA	SER	ENGINEERED MUTATION	UNP H9JWL6

- Molecule 2 is PENTAETHYLENE GLYCOL (three-letter code: 1PE) (formula: $C_{10}H_{22}O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			16	10	6		
2	B	1	Total	C	O	0	0
			16	10	6		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	81	Total	O	0	0
			81	81		
3	B	78	Total	O	0	0
			78	78		

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	161.40Å 51.47Å 124.45Å 90.00° 122.68° 90.00°	Depositor
Resolution (Å)	39.39 – 2.30 39.39 – 2.30	Depositor EDS
% Data completeness (in resolution range)	98.9 (39.39-2.30) 98.9 (39.39-2.30)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.23 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.203 , 0.251 0.204 , 0.254	Depositor DCC
R_{free} test set	1907 reflections (5.25%)	DCC
Wilson B-factor (Å ²)	33.5	Xtriage
Anisotropy	0.709	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 37.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 38204 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6682	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.04% 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

Bond lengths and bond angles in the following residue types are not validated in this section: 1PE

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	0.55	6/3520 (0.2%)	0.55	0/4776
1	B	0.55	0/3162	0.57	0/4283
All	All	0.55	6/6682 (0.1%)	0.56	0/9059

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	228	TRP	CD2-CE2	5.46	1.48	1.41
1	A	38	TRP	CD2-CE2	5.20	1.47	1.41
1	A	154	TRP	CD2-CE2	5.12	1.47	1.41
1	A	348	TRP	CD2-CE2	5.12	1.47	1.41
1	A	361	TRP	CD2-CE2	5.03	1.47	1.41
1	A	102	TRP	CD2-CE2	5.03	1.47	1.41

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	3419	0	3435	21	0
1	B	3072	0	3081	26	0
2	A	16	0	22	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	16	0	22	0	0
3	A	81	0	0	2	0
3	B	78	0	0	2	0
All	All	6682	0	6560	46	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 (46) 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:172:SER:O	1:A:173:LYS:HB2	1.74	0.87
1:B:118:TYR:HB3	1:B:135:ILE:HD11	1.72	0.71
1:A:411:ILE:C	1:A:411:ILE:HD12	2.14	0.68
1:B:152:HIS:CE1	1:B:157:SER:HA	2.30	0.66
1:B:118:TYR:HB3	1:B:135:ILE:CD1	2.29	0.62
1:A:411:ILE:O	1:A:411:ILE:HD12	2.02	0.60
1:B:74:ARG:HD2	1:B:99:LEU:HD13	1.84	0.58
1:A:251:MET:O	1:A:251:MET:HG3	2.02	0.58
1:A:176:ASN:HB2	3:A:660:HOH:O	2.06	0.55
1:A:311:ILE:HD13	1:B:95:LEU:HD11	1.91	0.53
1:B:210:LYS:HE2	1:B:243:GLU:OE2	2.08	0.53
1:A:107:PRO:HB2	1:A:110:GLU:HB2	1.91	0.52
1:A:136:THR:HG22	1:A:180:GLU:OE2	2.09	0.52
1:A:153:GLY:HA3	1:A:227:ASP:HB3	1.94	0.50
1:B:409:PRO:HA	1:B:412:LEU:HD12	1.94	0.49
1:B:401:LYS:HA	1:B:406:TYR:HE1	1.77	0.49
1:B:200:LEU:HD12	1:B:204:GLU:CB	2.43	0.48
1:A:139:VAL:HG13	1:A:143:VAL:HB	1.97	0.47
1:B:154:TRP:CE2	1:B:155:PRO:HB3	2.50	0.47
1:B:346:TYR:CD2	1:B:347:ARG:HG2	2.50	0.47
1:B:163:GLU:CG	3:B:653:HOH:O	2.64	0.46
1:B:252:ALA:HB1	1:B:412:LEU:HD21	1.98	0.45
1:B:120:HIS:CE1	1:B:133:MET:HG2	2.51	0.45
1:B:310:GLY:HA3	1:B:367:VAL:HG22	1.99	0.45
1:A:346:TYR:CD2	1:A:347:ARG:HG2	2.52	0.45
1:A:120:HIS:CE1	1:A:133:MET:HG2	2.52	0.44
1:A:200:LEU:CD1	1:A:204:GLU:HB3	2.47	0.44
1:A:333:ASN:HB3	1:A:336:LEU:HG	2.00	0.44
1:B:154:TRP:CG	1:B:155:PRO:HA	2.53	0.43
1:B:224:GLN:HG2	1:B:248:HIS:HB3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:298:TYR:CG	1:A:299:MET:N	2.87	0.43
1:B:200:LEU:HD12	1:B:204:GLU:HB2	2.01	0.43
1:A:269:PHE:HA	1:A:270:PRO:HD3	1.79	0.42
1:A:310:GLY:HA3	1:A:367:VAL:HG22	2.00	0.42
1:A:200:LEU:HD12	1:A:204:GLU:HB2	2.00	0.42
1:B:200:LEU:HD12	1:B:204:GLU:HB3	2.02	0.42
1:A:327:LYS:HE2	3:A:680:HOH:O	2.19	0.42
1:B:346:TYR:HD2	1:B:347:ARG:HG2	1.84	0.42
1:B:407:GLN:HB3	1:B:411:ILE:HD11	2.03	0.41
1:A:407:GLN:OE1	1:A:411:ILE:HD11	2.20	0.41
1:B:373:TYR:O	1:B:377:PHE:HD1	2.04	0.41
1:A:408:PRO:HG2	1:A:411:ILE:HG23	2.03	0.41
1:B:153:GLY:HA3	1:B:227:ASP:HB3	2.02	0.40
1:B:302:GLN:HG2	1:B:370:MET:HG2	2.03	0.40
1:B:327:LYS:HE2	3:B:673:HOH:O	2.20	0.40
1:B:161:PHE:O	1:B:165:ILE:HD12	2.22	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	422/447 (94%)	410 (97%)	12 (3%)	0	100	100
1	B	374/447 (84%)	359 (96%)	15 (4%)	0	100	100
All	All	796/894 (89%)	769 (97%)	27 (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	A	365/387 (94%)	358 (98%)	7 (2%)	65	81
1	B	326/387 (84%)	321 (98%)	5 (2%)	72	85
All	All	691/774 (89%)	679 (98%)	12 (2%)	68	83

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	44	LEU
1	A	139	VAL
1	A	196	VAL
1	A	210	LYS
1	A	307	ASP
1	A	411	ILE
1	A	415	LYS
1	B	113	LYS
1	B	139	VAL
1	B	165	ILE
1	B	307	ASP
1	B	425	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](#)

2 ligands are modelled in this entry.

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	1PE	A	501	-	15,15,15	0.49	0	14,14,14	0.28	0
2	1PE	B	501	-	15,15,15	0.52	0	14,14,14	0.22	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	1PE	A	501	-	-	0/13/13/13	0/0/0/0
2	1PE	B	501	-	-	0/13/13/13	0/0/0/0

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 [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	424/447 (94%)	0.15	19 (4%) 37 46	23, 38, 71, 94	1 (0%)
1	B	380/447 (85%)	0.15	25 (6%) 22 29	23, 35, 83, 129	1 (0%)
All	All	804/894 (89%)	0.15	44 (5%) 29 37	23, 36, 73, 129	2 (0%)

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	456	ASN	8.4
1	B	406	TYR	7.5
1	A	269	PHE	6.5
1	B	409	PRO	6.2
1	B	382	PHE	5.6
1	B	293	LEU	5.2
1	A	278	LEU	4.8
1	A	276	PRO	4.7
1	B	408	PRO	4.4
1	B	411	ILE	4.1
1	A	277	GLU	4.1
1	A	261	PHE	4.1
1	B	292	LEU	4.1
1	B	410	CYS	4.1
1	B	455	LYS	4.0
1	A	173	LYS	3.7
1	B	427	ASP	3.4
1	A	454	LEU	3.3
1	A	142	GLY	3.3
1	A	456	ASN	3.3
1	B	389	ILE	3.3
1	B	415	LYS	3.3
1	A	275	GLU	2.9
1	A	383	ASP	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	139	VAL	2.8
1	B	141	LYS	2.8
1	B	291	THR	2.7
1	A	270	PRO	2.7
1	A	274	VAL	2.6
1	A	257	PRO	2.5
1	A	455	LYS	2.5
1	B	47	LYS	2.5
1	B	140	PRO	2.4
1	B	251	MET	2.3
1	B	412	LEU	2.2
1	B	401	LYS	2.2
1	B	414	MET	2.2
1	B	252	ALA	2.2
1	B	138	LYS	2.2
1	A	242	LYS	2.1
1	A	141	LYS	2.1
1	B	46	GLN	2.1
1	A	379	SER	2.0
1	A	96	ASP	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	1PE	B	501	16/16	0.78	0.19	0.74	58,63,73,74	0
2	1PE	A	501	16/16	0.68	0.17	-0.24	63,68,73,73	0

6.5 Other polymers [i](#)

There are no such residues in this entry.