



# Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 08:57 am BST

PDB ID : 2AB0  
Title : Crystal Structure of E. coli protein YajL (ThiJ)  
Authors : Wilson, M.A.; Ringe, D.; Petsko, G.A.  
Deposited on : 2005-07-14  
Resolution : 1.10 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

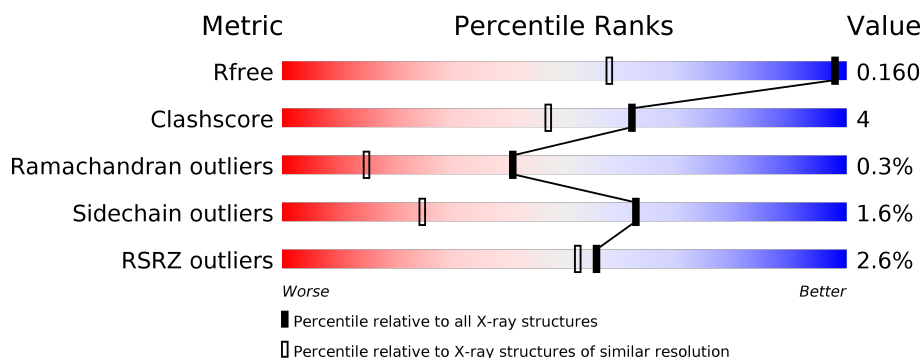
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.10 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	1619 (1.14-1.06)
Clashscore	141614	1671 (1.14-1.06)
Ramachandran outliers	138981	1615 (1.14-1.06)
Sidechain outliers	138945	1613 (1.14-1.06)
RSRZ outliers	127900	1588 (1.14-1.06)

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	205	<div> <div style="width: 100%; height: 10px; background-color: red; position: relative;"> <span style="position: absolute; top: -10px; left: 0;">%</span> </div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> <span>81%</span> <span>13%</span> <span>5%</span> </div> </div>
1	B	205	<div> <div style="width: 100%; height: 10px; background-color: red; position: relative;"> <span style="position: absolute; top: -10px; left: 0;">3%</span> </div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> <span>84%</span> <span>9%</span> <span>• 5%</span> </div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3538 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 YajL.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	195	Total	C	N	O	S	0	15	0
			1572	1002	264	298	8			
1	B	195	Total	C	N	O	S	0	9	0
			1526	972	260	287	7			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	197	GLN	-	EXPRESSION TAG	UNP Q46948
A	198	LEU	-	EXPRESSION TAG	UNP Q46948
A	199	GLU	-	EXPRESSION TAG	UNP Q46948
A	200	HIS	-	EXPRESSION TAG	UNP Q46948
A	201	HIS	-	EXPRESSION TAG	UNP Q46948
A	202	HIS	-	EXPRESSION TAG	UNP Q46948
A	203	HIS	-	EXPRESSION TAG	UNP Q46948
A	204	HIS	-	EXPRESSION TAG	UNP Q46948
A	205	HIS	-	EXPRESSION TAG	UNP Q46948
B	197	GLN	-	EXPRESSION TAG	UNP Q46948
B	198	LEU	-	EXPRESSION TAG	UNP Q46948
B	199	GLU	-	EXPRESSION TAG	UNP Q46948
B	200	HIS	-	EXPRESSION TAG	UNP Q46948
B	201	HIS	-	EXPRESSION TAG	UNP Q46948
B	202	HIS	-	EXPRESSION TAG	UNP Q46948
B	203	HIS	-	EXPRESSION TAG	UNP Q46948
B	204	HIS	-	EXPRESSION TAG	UNP Q46948
B	205	HIS	-	EXPRESSION TAG	UNP Q46948

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	188	Total	O	0	0
			188	188		

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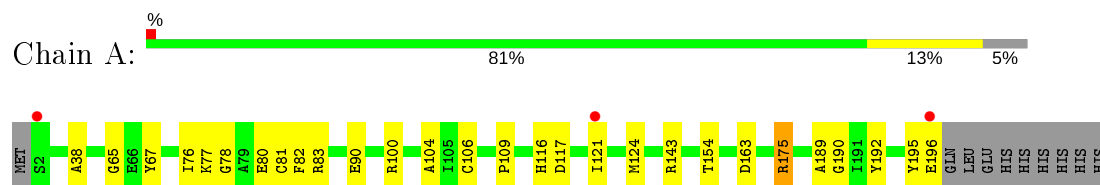
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	252	Total	O	0	0
			252	252		

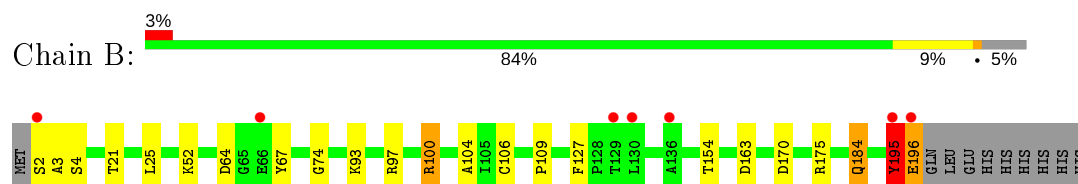
### 3 Residue-property plots [i](#)

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: YajL



#### • Molecule 1: YajL



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	42.72Å 78.47Å 99.95Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	100.00 – 1.10 42.15 – 1.10	Depositor EDS
% Data completeness (in resolution range)	96.6 (100.00-1.10) 91.0 (42.15-1.10)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.66 (at 1.10Å)	Xtriage
Refinement program	SHELXL-97	Depositor
R, $R_{free}$	0.136 , 0.171 0.157 , 0.160	Depositor DCC
$R_{free}$ test set	6609 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	10.1	Xtriage
Anisotropy	0.421	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 47.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	3538	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	16.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 17.77% 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

### 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	0.54	0/1597	1.11	15/2173 (0.7%)
1	B	0.52	0/1550	1.04	6/2104 (0.3%)
All	All	0.53	0/3147	1.07	21/4277 (0.5%)

There are no bond length outliers.

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	121	ILE	C-N-CA	-7.58	106.39	122.30
1	A	189	ALA	C-N-CA	-7.44	106.68	122.30
1	A	121	ILE	O-C-N	-7.23	110.92	123.20
1	A	124	MET	CA-CB-CG	6.96	125.13	113.30
1	A	175[A]	ARG	NE-CZ-NH1	6.40	123.50	120.30
1	A	175[B]	ARG	NE-CZ-NH1	6.40	123.50	120.30
1	B	100	ARG	NE-CZ-NH2	-6.20	117.20	120.30
1	A	82	PHE	CB-CG-CD1	6.07	125.05	120.80
1	A	175[A]	ARG	NE-CZ-NH2	-6.05	117.28	120.30
1	A	175[B]	ARG	NE-CZ-NH2	-6.05	117.28	120.30
1	A	189	ALA	O-C-N	-6.03	112.94	123.20
1	B	184	GLN	CB-CG-CD	5.69	126.41	111.60
1	A	192	TYR	CB-CG-CD1	5.67	124.40	121.00
1	B	195	TYR	CB-CG-CD2	5.63	124.38	121.00
1	A	117	ASP	CB-CG-OD2	-5.51	113.34	118.30
1	A	143	ARG	NE-CZ-NH2	-5.44	117.58	120.30
1	B	64	ASP	CB-CG-OD1	5.31	123.08	118.30
1	B	195	TYR	C-N-CA	5.12	134.49	121.70
1	A	190	GLY	N-CA-C	5.09	125.83	113.10
1	B	195	TYR	CB-CG-CD1	-5.09	117.95	121.00
1	A	65	GLY	O-C-N	-5.06	114.61	122.70

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	1572	0	1607	13	0
1	B	1526	0	1573	15	0
2	A	188	0	0	5	0
2	B	252	0	0	2	0
All	All	3538	0	3180	28	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 (28) 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:77[A]:LYS:HD3	2:A:296:HOH:O	1.85	0.77
1:A:195:TYR:CE2	1:A:196:GLU:HG2	2.35	0.62
1:A:90:GLU:HG2	2:A:271:HOH:O	2.00	0.60
1:A:38:ALA:HB3	1:A:81[B]:CYS:SG	2.41	0.60
1:B:195:TYR:O	1:B:196:GLU:HB2	2.05	0.56
1:A:76[B]:ILE:HD11	2:A:302:HOH:O	2.06	0.54
1:B:175[B]:ARG:NH1	1:B:195:TYR:O	2.49	0.46
1:A:106[A]:CYS:O	1:A:109:PRO:HD2	2.16	0.45
1:A:175[B]:ARG:NH1	2:A:384:HOH:O	2.49	0.45
1:B:93:LYS:O	1:B:97:ARG:HG2	2.17	0.45
1:B:2:SER:OG	1:B:3:ALA:N	2.49	0.45
1:A:76[B]:ILE:HD11	1:A:80:GLU:OE1	2.17	0.44
1:B:52:LYS:NZ	2:B:388:HOH:O	2.49	0.44
1:A:104:ALA:O	1:A:154:THR:HA	2.18	0.44
1:A:77[B]:LYS:NZ	2:A:292:HOH:O	2.50	0.44
1:A:67[A]:TYR:O	1:A:100:ARG:HD2	2.18	0.43
1:B:21:THR:O	1:B:25:LEU:HG	2.18	0.43
1:B:175[A]:ARG:HG3	2:B:442:HOH:O	2.17	0.43
1:B:104:ALA:O	1:B:154:THR:HA	2.19	0.42
1:A:83:ARG:O	1:A:116:HIS:HE1	2.03	0.42
1:B:106[B]:CYS:SG	1:B:127:PHE:HB2	2.59	0.42
1:B:4[A]:SER:OG	1:B:67:TYR:HA	2.19	0.42
1:B:170:ASP:OD1	1:B:175[A]:ARG:HG3	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:74:GLY:HA3	1:B:106[A]:CYS:HB3	2.03	0.41
1:B:106[B]:CYS:SG	1:B:127:PHE:CD1	3.14	0.41
1:B:106[B]:CYS:O	1:B:109:PRO:HD2	2.21	0.40
1:B:67:TYR:O	1:B:100:ARG:HD2	2.20	0.40
1:A:78:GLY:O	1:A:81[B]:CYS:HB3	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	208/205 (102%)	202 (97%)	6 (3%)	0	100	100
1	B	202/205 (98%)	196 (97%)	5 (2%)	1 (0%)	29	6
All	All	410/410 (100%)	398 (97%)	11 (3%)	1 (0%)	41	17

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	195	TYR

### 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.

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	170/165 (103%)	169 (99%)	1 (1%)	86	61
1	B	164/165 (99%)	160 (98%)	4 (2%)	49	11
All	All	334/330 (101%)	329 (98%)	5 (2%)	62	27

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	163	ASP
1	B	163	ASP
1	B	184	GLN
1	B	195	TYR
1	B	196	GLU

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

Mol	Chain	Res	Type
1	A	116	HIS
1	B	123	ASN
1	B	184	GLN

### 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

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, 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	195/205 (95%)	-0.24	3 (1%) 73 69	7, 12, 27, 52	0
1	B	195/205 (95%)	0.11	7 (3%) 42 39	7, 11, 30, 81	0
All	All	390/410 (95%)	-0.07	10 (2%) 56 52	7, 11, 28, 81	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	195	TYR	9.3
1	B	196	GLU	5.7
1	B	129	THR	4.5
1	B	130	LEU	4.1
1	B	66	GLU	2.8
1	B	136	ALA	2.8
1	A	196	GLU	2.6
1	B	2	SER	2.6
1	A	2	SER	2.5
1	A	121	ILE	2.4

### 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

There are no such residues in this entry.