



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 19, 2024 – 05:27 PM EDT

PDB ID : 1WU2
Title : Crystal Structure of molybdopterin biosynthesis moeA protein from *Pyrococcus horikoshii* OT3
Authors : Lokanath, N.K.; Kunishima, N.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)
Deposited on : 2004-11-30
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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

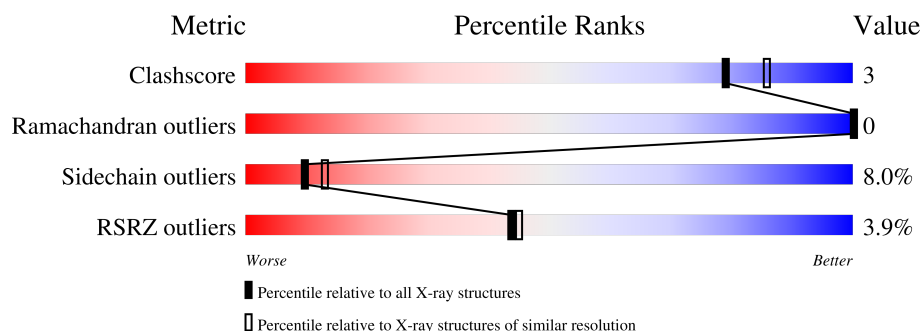
1 Overall quality at a glance

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))
Clashscore	180529	6698 (2.30-2.30)
Ramachandran outliers	177936	6640 (2.30-2.30)
Sidechain outliers	177891	6640 (2.30-2.30)
RSRZ outliers	164620	5963 (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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 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	396	<div> <div>4%</div> <div> <div></div> <div>83%</div> <div>12%</div> <div>• •</div> </div> </div>
1	B	396	<div> <div>4%</div> <div> <div></div> <div>84%</div> <div>10%</div> <div>• •</div> </div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6401 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 molybdopterin biosynthesis moeA protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	383	Total	C	N	O	S	Se	0	0	0
			3009	1943	494	565	1	6			
1	B	379	Total	C	N	O	S	Se	0	0	0
			2969	1917	487	558	1	6			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	MET	modified residue	GB 14591417
A	2	MSE	MET	modified residue	GB 14591417
A	118	MSE	MET	modified residue	GB 14591417
A	121	MSE	MET	modified residue	GB 14591417
A	168	MSE	MET	modified residue	GB 14591417
A	215	MSE	MET	modified residue	GB 14591417
A	296	MSE	MET	modified residue	GB 14591417
A	318	MSE	MET	modified residue	GB 14591417
B	1	MSE	MET	modified residue	GB 14591417
B	2	MSE	MET	modified residue	GB 14591417
B	118	MSE	MET	modified residue	GB 14591417
B	121	MSE	MET	modified residue	GB 14591417
B	168	MSE	MET	modified residue	GB 14591417
B	215	MSE	MET	modified residue	GB 14591417
B	296	MSE	MET	modified residue	GB 14591417
B	318	MSE	MET	modified residue	GB 14591417

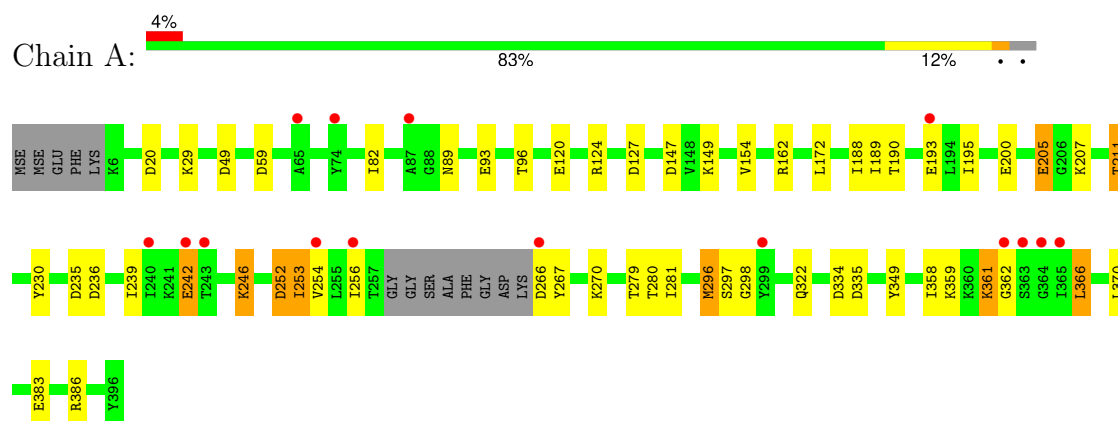
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	226	Total	O	0	0
			226	226		
2	B	197	Total	O	0	0
			197	197		

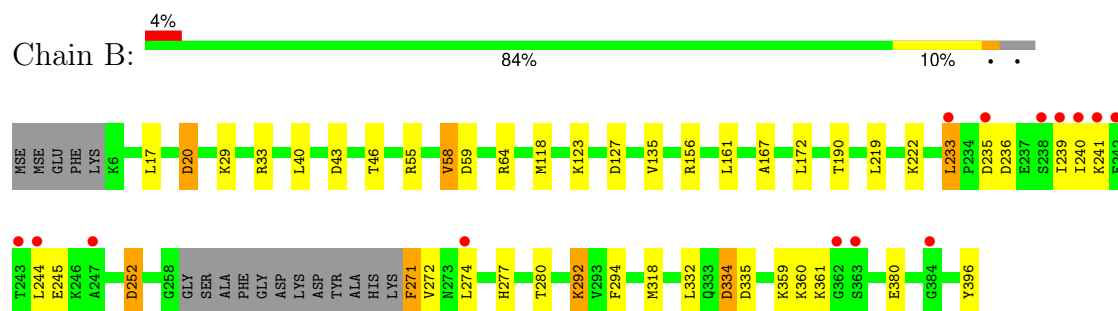
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: molybdopterin biosynthesis moeA protein



- Molecule 1: molybdopterin biosynthesis moeA protein



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	173.15Å 65.91Å 76.68Å 90.00° 102.17° 90.00°	Depositor
Resolution (Å)	20.00 – 2.30 20.00 – 2.30	Depositor EDS
% Data completeness (in resolution range)	98.1 (20.00-2.30) 97.9 (20.00-2.30)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.50 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.1.27	Depositor
R, R_{free}	0.233 , 0.282 0.237 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	33.7	Xtriage
Anisotropy	0.507	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 29.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	6401	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.66% 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

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	0.30	0/3056	0.66	11/4117 (0.3%)
1	B	0.29	0/3014	0.62	7/4060 (0.2%)
All	All	0.29	0/6070	0.64	18/8177 (0.2%)

There are no bond length outliers.

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	127	ASP	CB-CG-OD2	6.16	123.84	118.30
1	B	127	ASP	CB-CG-OD2	6.09	123.78	118.30
1	B	334	ASP	CB-CG-OD2	5.79	123.51	118.30
1	A	266	ASP	CB-CG-OD2	5.77	123.50	118.30
1	A	59	ASP	CB-CG-OD2	5.76	123.48	118.30
1	A	335	ASP	CB-CG-OD2	5.76	123.48	118.30
1	A	252	ASP	CB-CG-OD2	5.50	123.25	118.30
1	B	59	ASP	CB-CG-OD2	5.50	123.25	118.30
1	B	20	ASP	CB-CG-OD2	5.49	123.24	118.30
1	A	147	ASP	CB-CG-OD2	5.46	123.21	118.30
1	A	20	ASP	CB-CG-OD2	5.41	123.17	118.30
1	B	252	ASP	CB-CG-OD2	5.23	123.01	118.30
1	A	49	ASP	CB-CG-OD2	5.21	122.99	118.30
1	A	235	ASP	CB-CG-OD2	5.19	122.97	118.30
1	B	335	ASP	CB-CG-OD2	5.13	122.92	118.30
1	A	334	ASP	CB-CG-OD2	5.06	122.85	118.30
1	B	43	ASP	CB-CG-OD2	5.02	122.82	118.30
1	A	236	ASP	CB-CG-OD2	5.01	122.81	118.30

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	3009	0	3105	19	0
1	B	2969	0	3070	14	0
2	A	226	0	0	0	0
2	B	197	0	0	1	0
All	All	6401	0	6175	32	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 3.

All (32) 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:361:LYS:CB	1:A:362:GLY:HA3	2.18	0.74
1:A:361:LYS:HB3	1:A:362:GLY:HA3	1.73	0.70
1:A:239:ILE:O	1:A:242:GLU:HG2	1.97	0.64
1:A:256:ILE:HG23	1:A:298:GLY:HA2	1.79	0.64
1:A:230:TYR:CE2	1:A:246:LYS:HG3	2.38	0.58
1:B:190:THR:HG22	1:B:233:LEU:HD21	1.90	0.52
1:B:46:THR:HG23	1:B:172:LEU:HG	1.92	0.51
1:A:361:LYS:CG	1:A:362:GLY:HA3	2.41	0.50
1:A:211:THR:HG23	1:B:167:ALA:HB1	1.94	0.49
1:A:242:GLU:C	1:A:246:LYS:HZ2	2.18	0.47
1:B:222:LYS:NZ	1:B:396:TYR:OXT	2.48	0.47
1:B:274:LEU:HD21	1:B:277:HIS:HB2	1.95	0.47
1:A:242:GLU:O	1:A:246:LYS:HD3	2.16	0.46
1:A:349:TYR:HB2	1:A:358:ILE:HD11	1.98	0.46
1:A:361:LYS:HB3	1:A:362:GLY:CA	2.43	0.45
1:A:361:LYS:HG2	1:A:362:GLY:HA3	1.97	0.45
1:B:20:ASP:O	1:B:292:LYS:NZ	2.50	0.45
1:A:205:GLU:HB3	1:A:207:LYS:HZ2	1.83	0.44
1:A:188:ILE:HD12	1:A:253:ILE:HD11	2.01	0.43
1:A:366:LEU:HD13	1:A:370:LEU:HD13	2.00	0.43
1:B:64:ARG:NH1	2:B:438:HOH:O	2.52	0.43
1:B:235:ASP:HA	1:B:239:ILE:HD11	2.01	0.43
1:A:254:VAL:HG13	1:A:296:MSE:SE	2.69	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:58:VAL:HG21	1:B:118:MSE:HG2	2.02	0.42
1:B:280:THR:HG21	1:B:380:GLU:OE2	2.18	0.42
1:A:154:VAL:HG21	1:A:172:LEU:HD11	2.02	0.42
1:B:40:LEU:HD13	1:B:161:LEU:HD21	2.02	0.42
1:B:17:LEU:HD11	1:B:294:PHE:CE2	2.54	0.42
1:A:279:THR:HG23	1:A:281:ILE:H	1.86	0.41
1:B:252:ASP:CB	1:B:318:MSE:HE1	2.51	0.41
1:A:242:GLU:OE2	1:A:246:LYS:NZ	2.40	0.41
1:B:240:ILE:HG23	1:B:271:PHE:N	2.36	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	379/396 (96%)	364 (96%)	15 (4%)	0	100	100
1	B	375/396 (95%)	362 (96%)	13 (4%)	0	100	100
All	All	754/792 (95%)	726 (96%)	28 (4%)	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	325/326 (100%)	294 (90%)	31 (10%)	7	8
1	B	321/326 (98%)	300 (94%)	21 (6%)	14	20
All	All	646/652 (99%)	594 (92%)	52 (8%)	10	13

All (52) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	29	LYS
1	A	82	ILE
1	A	89	ASN
1	A	93	GLU
1	A	96	THR
1	A	120	GLU
1	A	124	ARG
1	A	149	LYS
1	A	162	ARG
1	A	189	ILE
1	A	190	THR
1	A	193	GLU
1	A	195	ILE
1	A	200	GLU
1	A	205	GLU
1	A	211	THR
1	A	242	GLU
1	A	246	LYS
1	A	252	ASP
1	A	253	ILE
1	A	267	TYR
1	A	270	LYS
1	A	280	THR
1	A	296	MSE
1	A	297	SER
1	A	322	GLN
1	A	359	LYS
1	A	361	LYS
1	A	366	LEU
1	A	383	GLU
1	A	386	ARG
1	B	29	LYS
1	B	33	ARG
1	B	55	ARG
1	B	58	VAL

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Mol	Chain	Res	Type
1	B	123	LYS
1	B	135	VAL
1	B	156	ARG
1	B	219	LEU
1	B	233	LEU
1	B	236	ASP
1	B	241	LYS
1	B	244	LEU
1	B	245	GLU
1	B	271	PHE
1	B	272	VAL
1	B	292	LYS
1	B	332	LEU
1	B	334	ASP
1	B	359	LYS
1	B	360	LYS
1	B	361	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	217	GLN
1	A	322	GLN
1	A	352	ASN
1	B	273	ASN
1	B	352	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry

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

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	377/396 (95%)	0.37	15 (3%) 43 44	23, 35, 57, 63	0
1	B	373/396 (94%)	0.41	14 (3%) 44 46	24, 37, 55, 77	0
All	All	750/792 (94%)	0.39	29 (3%) 44 45	23, 36, 57, 77	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	363	SER	4.7
1	B	240	ILE	4.6
1	A	363	SER	4.2
1	A	362	GLY	4.0
1	B	241	LYS	4.0
1	B	244	LEU	3.8
1	A	243	THR	3.3
1	B	362	GLY	3.3
1	A	65	ALA	3.1
1	B	239	ILE	3.0
1	A	299	TYR	2.8
1	B	233	LEU	2.8
1	A	364	GLY	2.6
1	B	247	ALA	2.5
1	B	274	LEU	2.5
1	A	365	ILE	2.5
1	B	238	SER	2.5
1	B	243	THR	2.3
1	B	242	GLU	2.3
1	A	240	ILE	2.3
1	A	266	ASP	2.2
1	A	242	GLU	2.2
1	A	87	ALA	2.2
1	A	74	TYR	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	256	ILE	2.1
1	A	254	VAL	2.0
1	A	193	GLU	2.0
1	B	235	ASP	2.0
1	B	384	GLY	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 monosaccharides 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.