



wwPDB X-ray Structure Validation Summary Report ⓘ

Oct 4, 2023 – 10:38 PM EDT

PDB ID : 6VK6
Title : Crystal Structure of Methylosinus trichosporium OB3b Soluble Methane Monooxygenase Hydroxylase
Authors : Jones, J.C.; Banerjee, R.; Shi, K.; Aihara, H.; Lipscomb, J.D.
Deposited on : 2020-01-18
Resolution : 1.52 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 : **FAILED**
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtrriage (Phenix) : 1.13
EDS : **FAILED**
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.35.1

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 1.52 Å.

There are no overall percentile quality scores available for this entry.

MolProbity and EDS failed to run properly - the sequence quality summary graphics cannot be shown.

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 18856 atoms, of which 8741 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 Methane monooxygenase component A alpha chain.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	515	8291	2719	4045	734	780	13	0	17	0

- Molecule 2 is a protein called Methane monooxygenase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
2	B	387	6221	2039	3026	560	591	5	0	10	0

- Molecule 3 is a protein called Methane monooxygenase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
3	C	168	2782	880	1412	235	254	1	0	2	0

- Molecule 4 is FE (III) ION (three-letter code: FE) (formula: Fe) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	Fe	0	0
			2	2		

- Molecule 5 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	
			Total	C	H			O
5	A	1	10	2	6	2	0	0
5	A	1	10	2	6	2	0	0
5	A	1	10	2	6	2	0	0
5	A	1	10	2	6	2	0	0
5	A	1	20	4	12	4	0	1
5	A	1	10	2	6	2	0	0
5	A	1	10	2	6	2	0	0
5	A	1	10	2	6	2	0	0
5	A	1	10	2	6	2	0	0
5	A	1	10	2	6	2	0	0
5	A	1	10	2	6	2	0	0
5	A	1	10	2	6	2	0	0
5	A	1	10	2	6	2	0	0
5	A	1	10	2	6	2	0	0

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

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	C	1	Total	C	H	O	0	0
			10	2	6	2		
5	C	1	Total	C	H	O	0	0
			10	2	6	2		
5	C	1	Total	C	H	O	0	0
			10	2	6	2		
5	C	1	Total	C	H	O	0	0
			10	2	6	2		
5	C	1	Total	C	H	O	0	0
			10	2	6	2		
5	C	1	Total	C	H	O	0	0
			10	2	6	2		

- Molecule 6 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	2	Total	Cl	0	0
			2	2		
6	C	2	Total	Cl	0	0
			2	2		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	495	Total	O	0	2
			497	497		
7	B	423	Total	O	0	0
			423	423		
7	C	206	Total	O	0	0
			206	206		

MolProbity and EDS failed to run properly - this section is therefore empty.

3 Data and refinement statistics

EDS failed to run properly - this section is therefore incomplete.

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	63.07Å 292.63Å 141.79Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	65.01 – 1.52	Depositor
% Data completeness (in resolution range)	93.5 (65.01-1.52)	Depositor
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.47 (at 1.52Å)	Xtrriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.146 , 0.164	Depositor
Wilson B-factor (Å ²)	18.4	Xtrriage
Anisotropy	0.419	Xtrriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
Total number of atoms	18856	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.42% 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.

4 Model quality [i](#)

4.1 Standard geometry [i](#)

MolProbity failed to run properly - this section is therefore empty.

4.2 Too-close contacts [i](#)

MolProbity failed to run properly - this section is therefore empty.

4.3 Torsion angles [i](#)

4.3.1 Protein backbone [i](#)

MolProbity failed to run properly - this section is therefore empty.

4.3.2 Protein sidechains [i](#)

MolProbity failed to run properly - this section is therefore empty.

4.3.3 RNA [i](#)

MolProbity failed to run properly - this section is therefore empty.

4.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

4.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

4.6 Ligand geometry [i](#)

Of 49 ligands modelled in this entry, 6 are monoatomic - leaving 43 for Mogul analysis.

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
5	EDO	A	810	-	3,3,3	0.51	0	2,2,2	0.12	0
5	EDO	A	804	-	3,3,3	0.47	0	2,2,2	0.55	0
5	EDO	A	817	-	3,3,3	0.49	0	2,2,2	0.23	0
5	EDO	A	821	-	3,3,3	0.47	0	2,2,2	0.42	0
5	EDO	A	822	-	3,3,3	0.50	0	2,2,2	0.26	0
5	EDO	A	818	-	3,3,3	0.47	0	2,2,2	0.36	0
5	EDO	B	410	-	3,3,3	0.48	0	2,2,2	0.26	0
5	EDO	A	807[B]	-	3,3,3	0.48	0	2,2,2	0.27	0
5	EDO	A	807[A]	-	3,3,3	0.47	0	2,2,2	0.30	0
5	EDO	A	805	-	3,3,3	0.48	0	2,2,2	0.24	0
5	EDO	C	203	-	3,3,3	0.46	0	2,2,2	0.36	0
5	EDO	B	414	-	3,3,3	0.50	0	2,2,2	0.20	0
5	EDO	C	205	-	3,3,3	0.45	0	2,2,2	0.37	0
5	EDO	A	816	-	3,3,3	0.47	0	2,2,2	0.43	0
5	EDO	A	823	-	3,3,3	0.48	0	2,2,2	0.22	0
5	EDO	B	407	-	3,3,3	0.46	0	2,2,2	0.35	0
5	EDO	A	803	-	3,3,3	0.47	0	2,2,2	0.51	0
5	EDO	A	814	-	3,3,3	0.46	0	2,2,2	0.30	0
5	EDO	C	208	-	3,3,3	0.48	0	2,2,2	0.33	0
5	EDO	A	825	-	3,3,3	0.45	0	2,2,2	0.50	0
5	EDO	C	206	-	3,3,3	0.47	0	2,2,2	0.41	0
5	EDO	A	819	-	3,3,3	0.49	0	2,2,2	0.10	0
5	EDO	B	404	-	3,3,3	0.46	0	2,2,2	0.40	0
5	EDO	B	413	-	3,3,3	0.47	0	2,2,2	0.39	0
5	EDO	A	820	-	3,3,3	0.49	0	2,2,2	0.24	0
5	EDO	A	809	-	3,3,3	0.47	0	2,2,2	0.32	0
5	EDO	A	824	-	3,3,3	0.49	0	2,2,2	0.15	0
5	EDO	B	412	-	3,3,3	0.47	0	2,2,2	0.27	0
5	EDO	C	209	-	3,3,3	0.47	0	2,2,2	0.28	0
5	EDO	B	411	-	3,3,3	0.49	0	2,2,2	0.34	0
5	EDO	A	808	-	3,3,3	0.48	0	2,2,2	0.22	0
5	EDO	A	815	-	3,3,3	0.47	0	2,2,2	0.42	0
5	EDO	C	204	-	3,3,3	0.47	0	2,2,2	0.47	0
5	EDO	A	812	4	3,3,3	0.49	0	2,2,2	0.32	0
5	EDO	A	811	-	3,3,3	0.47	0	2,2,2	0.40	0
5	EDO	A	813	-	3,3,3	0.47	0	2,2,2	0.33	0
5	EDO	C	207	-	3,3,3	0.47	0	2,2,2	0.28	0
5	EDO	B	406	-	3,3,3	0.48	0	2,2,2	0.24	0
5	EDO	A	806	-	3,3,3	0.47	0	2,2,2	0.35	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	EDO	B	403	-	3,3,3	0.49	0	2,2,2	0.36	0
5	EDO	B	408	-	3,3,3	0.47	0	2,2,2	0.17	0
5	EDO	B	409	-	3,3,3	0.48	0	2,2,2	0.48	0
5	EDO	B	405	-	3,3,3	0.47	0	2,2,2	0.30	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
5	EDO	A	810	-	-	0/1/1/1	-
5	EDO	A	804	-	-	0/1/1/1	-
5	EDO	A	817	-	-	0/1/1/1	-
5	EDO	A	821	-	-	1/1/1/1	-
5	EDO	A	822	-	-	1/1/1/1	-
5	EDO	A	818	-	-	0/1/1/1	-
5	EDO	B	410	-	-	1/1/1/1	-
5	EDO	A	807[B]	-	-	1/1/1/1	-
5	EDO	A	807[A]	-	-	0/1/1/1	-
5	EDO	A	805	-	-	0/1/1/1	-
5	EDO	C	203	-	-	0/1/1/1	-
5	EDO	B	414	-	-	0/1/1/1	-
5	EDO	C	205	-	-	1/1/1/1	-
5	EDO	A	816	-	-	1/1/1/1	-
5	EDO	A	823	-	-	0/1/1/1	-
5	EDO	B	407	-	-	0/1/1/1	-
5	EDO	A	803	-	-	0/1/1/1	-
5	EDO	A	814	-	-	0/1/1/1	-
5	EDO	C	208	-	-	1/1/1/1	-
5	EDO	A	825	-	-	0/1/1/1	-
5	EDO	C	206	-	-	0/1/1/1	-
5	EDO	A	819	-	-	0/1/1/1	-
5	EDO	B	404	-	-	0/1/1/1	-
5	EDO	B	413	-	-	0/1/1/1	-
5	EDO	A	820	-	-	1/1/1/1	-
5	EDO	A	809	-	-	0/1/1/1	-
5	EDO	A	824	-	-	1/1/1/1	-
5	EDO	B	412	-	-	0/1/1/1	-
5	EDO	C	209	-	-	0/1/1/1	-
5	EDO	B	411	-	-	1/1/1/1	-
5	EDO	A	808	-	-	1/1/1/1	-
5	EDO	A	815	-	-	0/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	EDO	C	204	-	-	1/1/1/1	-
5	EDO	A	812	4	-	0/1/1/1	-
5	EDO	A	811	-	-	0/1/1/1	-
5	EDO	A	813	-	-	0/1/1/1	-
5	EDO	C	207	-	-	0/1/1/1	-
5	EDO	B	406	-	-	0/1/1/1	-
5	EDO	A	806	-	-	0/1/1/1	-
5	EDO	B	403	-	-	0/1/1/1	-
5	EDO	B	408	-	-	0/1/1/1	-
5	EDO	B	409	-	-	0/1/1/1	-
5	EDO	B	405	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

5 of 13 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	807[B]	EDO	O1-C1-C2-O2
5	B	410	EDO	O1-C1-C2-O2
5	A	822	EDO	O1-C1-C2-O2
5	B	405	EDO	O1-C1-C2-O2
5	C	205	EDO	O1-C1-C2-O2

There are no ring outliers.

No monomer is involved in short contacts.

4.7 Other polymers [i](#)

There are no such residues in this entry.

4.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

5 Fit of model and data

5.1 Protein, DNA and RNA chains

EDS failed to run properly - this section is therefore empty.

5.2 Non-standard residues in protein, DNA, RNA chains

EDS failed to run properly - this section is therefore empty.

5.3 Carbohydrates

EDS failed to run properly - this section is therefore empty.

5.4 Ligands

EDS failed to run properly - this section is therefore empty.

5.5 Other polymers

EDS failed to run properly - this section is therefore empty.