



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 18, 2017 – 01:09 AM EST

PDB ID : 5E7Y
Title : Crystal structure of P450 BM3 heme domain M7 variant
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Deposited on : 2015-10-13
Resolution : 2.00 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 : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20028442
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) : rb-20028442

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	GOL	A	502	-	-	-	X
3	GOL	A	503	-	-	-	X
3	GOL	A	504	-	-	-	X
3	GOL	A	505	-	-	-	X
3	GOL	A	506	-	-	-	X
3	GOL	B	502	-	-	-	X
3	GOL	B	503	-	-	-	X
3	GOL	B	504	-	-	-	X

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 7952 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.

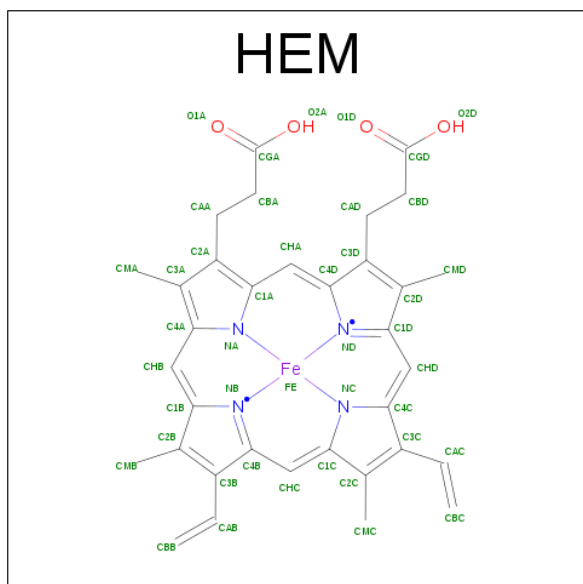
- Molecule 1 is a protein called Bifunctional P-450/NADPH-P450 reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	451	Total	C	N	O	S	0	13	0
			3701	2364	626	695	16			
1	B	452	Total	C	N	O	S	0	11	0
			3708	2366	631	695	16			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	87	ALA	PHE	engineered mutation	UNP P14779
A	281	GLY	VAL	engineered mutation	UNP P14779
A	354	SER	MET	engineered mutation	UNP P14779
B	87	ALA	PHE	engineered mutation	UNP P14779
B	281	GLY	VAL	engineered mutation	UNP P14779
B	354	SER	MET	engineered mutation	UNP P14779

- Molecule 2 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	B	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	217	Total 217	O 217	0	0
4	B	192	Total 192	O 192	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	59.51Å 153.89Å 61.59Å 90.00° 94.97° 90.00°	Depositor
Resolution (Å)	20.00 – 2.00 19.87 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.7 (20.00-2.00) 99.8 (19.87-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.64 (at 2.01Å)	Xtriage
Refinement program	REFMAC 5.8.0103	Depositor
R, R_{free}	0.188 , 0.220 0.193 , 0.222	Depositor DCC
R_{free} test set	3704 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	30.6	Xtriage
Anisotropy	0.138	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 46.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.032 for l,-k,h	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7952	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.42% 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.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, HEM

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	0/3821	0.74	1/5167 (0.0%)
1	B	0.54	0/3816	0.73	0/5157
All	All	0.54	0/7637	0.74	1/10324 (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	242	ASP	CB-CG-OD1	5.26	123.04	118.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	3701	0	3687	11	0
1	B	3708	0	3672	16	0
2	A	43	0	30	1	0
2	B	43	0	30	1	0
3	A	30	0	40	0	0
3	B	18	0	24	0	0
4	A	217	0	0	2	0
4	B	192	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	7952	0	7483	29	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 2.

All (29) 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:27:GLN:HA	1:A:30:MET:HE2	1.53	0.91
1:A:27:GLN:HA	1:A:30:MET:CE	2.05	0.86
1:B:158[B]:PHE:CE1	1:B:258:ILE:HG23	2.22	0.74
1:B:45:PRO:C	1:B:46:GLY:N	2.49	0.65
1:B:46:GLY:O	1:B:47:ARG:N	2.33	0.61
1:A:128[B]:GLN:NE2	4:A:601:HOH:O	2.25	0.61
1:B:158[B]:PHE:CZ	1:B:258:ILE:HG23	2.38	0.59
1:B:157:GLY:C	1:B:158[B]:PHE:CD1	2.77	0.59
1:B:46:GLY:C	1:B:47:ARG:N	2.55	0.58
1:A:27:GLN:HA	1:A:30:MET:HE3	1.86	0.57
1:B:128[B]:GLN:OE1	1:B:128[B]:GLN:HA	2.04	0.56
1:A:27:GLN:CA	1:A:30:MET:HE2	2.32	0.54
1:B:309:LYS:HA	1:B:309:LYS:HE3	1.92	0.50
1:A:266:HIS:NE2	1:A:267:GLU:HG3	2.28	0.49
1:B:267:GLU:OE1	4:B:601:HOH:O	2.20	0.48
2:B:501:HEM:HBC2	2:B:501:HEM:HMC2	1.96	0.48
1:B:388:HIS:HA	1:B:391:LYS:HD3	1.96	0.48
2:A:501:HEM:HMC2	2:A:501:HEM:HBC2	1.97	0.46
1:B:158[B]:PHE:CD1	1:B:258:ILE:HG12	2.51	0.45
1:A:388:HIS:HA	1:A:391:LYS:HD3	1.97	0.45
1:A:120:VAL:HG11	1:A:302[B]:VAL:HG13	1.99	0.45
1:A:118:MET:HG2	4:A:794:HOH:O	2.17	0.43
1:B:45:PRO:C	1:B:46:GLY:CA	2.87	0.43
1:A:158:PHE:CE2	1:A:219:ILE:HD13	2.55	0.42
1:B:154:GLY:HA2	1:B:158[B]:PHE:CE2	2.55	0.41
1:A:272:LEU:HD13	1:A:322:LEU:HG	2.02	0.41
1:B:266:HIS:NE2	1:B:267:GLU:HG3	2.35	0.41
1:B:272:LEU:HD13	1:B:322:LEU:HG	2.02	0.41
1:B:112:MET:SD	1:B:405:PHE:HA	2.62	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	462/471 (98%)	451 (98%)	11 (2%)	0	100	100
1	B	458/471 (97%)	445 (97%)	13 (3%)	0	100	100
All	All	920/942 (98%)	896 (97%)	24 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	406/410 (99%)	394 (97%)	12 (3%)	48	47
1	B	404/410 (98%)	393 (97%)	11 (3%)	52	52
All	All	810/820 (99%)	787 (97%)	23 (3%)	65	50

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	21	ASN
1	A	110[A]	GLN
1	A	110[B]	GLN
1	A	136[A]	ASP
1	A	136[B]	ASP
1	A	148	LEU
1	A	158	PHE
1	A	167[A]	ARG

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Mol	Chain	Res	Type
1	A	167[B]	ARG
1	A	277[A]	LEU
1	A	277[B]	LEU
1	A	440	LYS
1	B	20	LEU
1	B	136[A]	ASP
1	B	136[B]	ASP
1	B	148	LEU
1	B	158[A]	PHE
1	B	158[B]	PHE
1	B	167	ARG
1	B	169[A]	GLN
1	B	169[B]	GLN
1	B	285[A]	HIS
1	B	285[B]	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	319	ASN

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](#)

10 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	HEM	A	501	1,4	24,50,50	0.79	1 (4%)	16,82,82	1.15	2 (12%)
3	GOL	A	502	-	5,5,5	0.22	0	5,5,5	0.16	0
3	GOL	A	503	-	5,5,5	0.35	0	5,5,5	0.32	0
3	GOL	A	504	-	5,5,5	0.30	0	5,5,5	0.37	0
3	GOL	A	505	-	5,5,5	0.35	0	5,5,5	0.38	0
3	GOL	A	506	-	5,5,5	0.33	0	5,5,5	0.36	0
2	HEM	B	501	1,4	24,50,50	0.69	1 (4%)	16,82,82	1.06	1 (6%)
3	GOL	B	502	-	5,5,5	0.31	0	5,5,5	0.29	0
3	GOL	B	503	-	5,5,5	0.31	0	5,5,5	0.57	0
3	GOL	B	504	-	5,5,5	0.37	0	5,5,5	0.44	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	HEM	A	501	1,4	-	0/6/54/54	0/0/8/8
3	GOL	A	502	-	-	0/4/4/4	0/0/0/0
3	GOL	A	503	-	-	0/4/4/4	0/0/0/0
3	GOL	A	504	-	-	0/4/4/4	0/0/0/0
3	GOL	A	505	-	-	0/4/4/4	0/0/0/0
3	GOL	A	506	-	-	0/4/4/4	0/0/0/0
2	HEM	B	501	1,4	-	0/6/54/54	0/0/8/8
3	GOL	B	502	-	-	0/4/4/4	0/0/0/0
3	GOL	B	503	-	-	0/4/4/4	0/0/0/0
3	GOL	B	504	-	-	0/4/4/4	0/0/0/0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	HEM	C1B-NB	-2.19	1.33	1.36
2	B	501	HEM	C1B-NB	-2.07	1.34	1.36

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	HEM	CBD-CAD-C3D	-2.48	108.12	112.47
2	A	501	HEM	CMA-C3A-C4A	-2.35	124.31	128.31
2	B	501	HEM	CBD-CAD-C3D	-2.12	108.75	112.47

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	HEM	1	0
2	B	501	HEM	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	B	2

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	B	46:GLY	C	47:ARG	N	2.55
1	B	45:PRO	C	46:GLY	N	2.49

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	451/471 (95%)	-0.11	12 (2%) 58 58	20, 36, 66, 91	0
1	B	452/471 (95%)	0.01	22 (4%) 33 35	24, 40, 67, 92	0
All	All	903/942 (95%)	-0.05	34 (3%) 44 45	20, 38, 66, 92	0

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	46	GLY	5.2
1	A	13	GLU	4.4
1	B	191	ALA	4.1
1	B	192	ASN	3.9
1	B	197	ALA	3.2
1	B	196	PRO	3.1
1	A	383	SER	3.0
1	A	21	ASN	3.0
1	B	194	ASP	3.0
1	A	428	ASN	3.0
1	B	94	LYS	3.0
1	B	231	ASP	2.9
1	A	196	PRO	2.9
1	B	136[A]	ASP	2.8
1	B	227	GLY	2.7
1	B	244	GLU	2.7
1	B	228	GLU	2.6
1	B	195	ASP	2.5
1	B	343	GLY	2.5
1	A	47	ARG	2.4
1	A	197	ALA	2.4
1	B	428	ASN	2.3
1	B	169[A]	GLN	2.3
1	B	259	ILE	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	247	GLU	2.3
1	A	136[A]	ASP	2.3
1	A	380	GLU	2.3
1	A	191	ALA	2.3
1	A	23[A]	ASP	2.3
1	B	199	ASP	2.2
1	B	189	GLN	2.2
1	A	430	GLU	2.2
1	B	193	PRO	2.1
1	B	426	HIS	2.1

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
3	GOL	B	504	6/6	0.82	0.28	12.79	40,58,59,60	0
3	GOL	A	505	6/6	0.86	0.24	12.38	36,49,53,54	0
3	GOL	A	503	6/6	0.87	0.33	8.92	52,64,66,75	0
3	GOL	A	506	6/6	0.80	0.22	4.86	42,61,68,72	0
3	GOL	A	504	6/6	0.83	0.25	4.43	41,63,73,77	0
3	GOL	B	503	6/6	0.73	0.27	3.65	52,63,70,70	0
3	GOL	A	502	6/6	0.87	0.14	3.34	50,53,59,66	0
3	GOL	B	502	6/6	0.82	0.19	2.50	60,61,67,69	0
2	HEM	A	501	43/43	0.98	0.08	-0.71	17,20,24,25	0
2	HEM	B	501	43/43	0.98	0.08	-0.88	22,25,32,34	0

6.5 Other polymers

There are no such residues in this entry.