



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:50 PM GMT

PDB ID : 4APY  
Title : Ethylene glycol-bound form of P450 CYP125A3 from Mycobacterium smegmatis  
Authors : Frank, D.J.; Garcia Fernandez, E.; Kells, P.M.; Garcia Lopez, J.L.; Podust, L.M.; Ortiz de Montellano, P.R.  
Deposited on : 2012-04-09  
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](mailto: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.

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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  
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

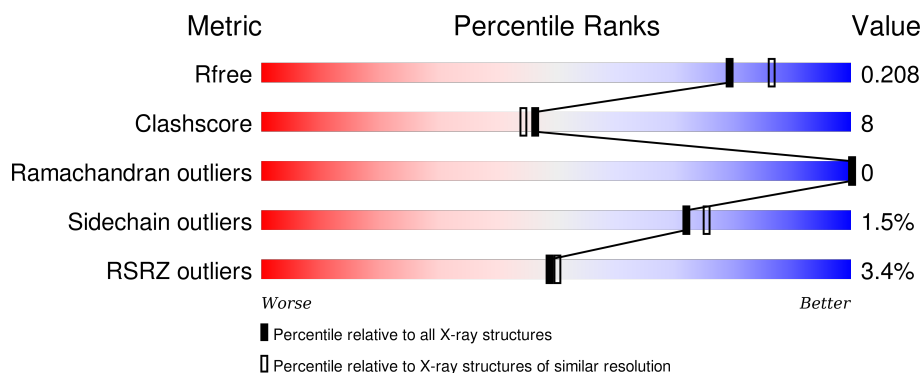
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*


The reported resolution of this entry is 2.00 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	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (2.00-2.00)

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. 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	433	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	EDO	A	1416	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	EDO	A	1419	-	-	-	X
2	EDO	A	1421	-	-	-	X
2	EDO	A	1422	-	-	-	X
2	EDO	A	1424	-	-	-	X
2	EDO	A	1425	-	-	X	-
2	EDO	A	1427	-	-	-	X
2	EDO	A	1428	-	-	-	X

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 3686 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.

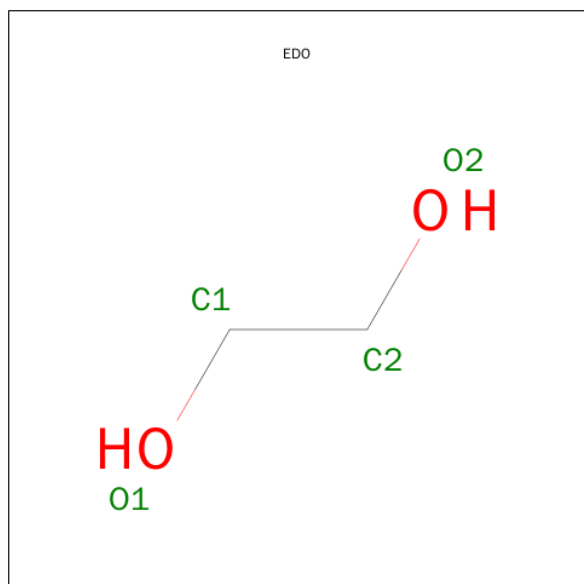
- Molecule 1 is a protein called P 450 HEME-THIOLATE PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	414	3259	2049	565	631	14	0	6	0

There are 6 discrepancies between the modelled and reference sequences:

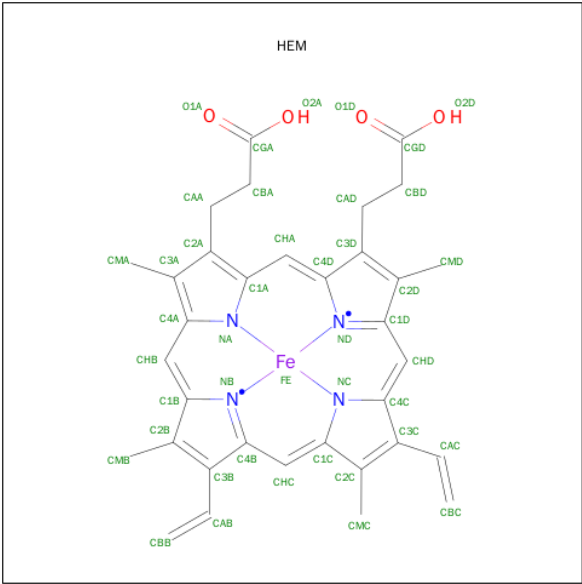
Chain	Residue	Modelled	Actual	Comment	Reference
A	428	HIS	-	EXPRESSION TAG	UNP A0R4Y3
A	429	HIS	-	EXPRESSION TAG	UNP A0R4Y3
A	430	HIS	-	EXPRESSION TAG	UNP A0R4Y3
A	431	HIS	-	EXPRESSION TAG	UNP A0R4Y3
A	432	HIS	-	EXPRESSION TAG	UNP A0R4Y3
A	433	HIS	-	EXPRESSION TAG	UNP A0R4Y3

- Molecule 2 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C<sub>2</sub>H<sub>6</sub>O<sub>2</sub>).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0

- Molecule 3 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: C<sub>34</sub>H<sub>32</sub>FeN<sub>4</sub>O<sub>4</sub>).



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

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Cl	0	0
			1	1		

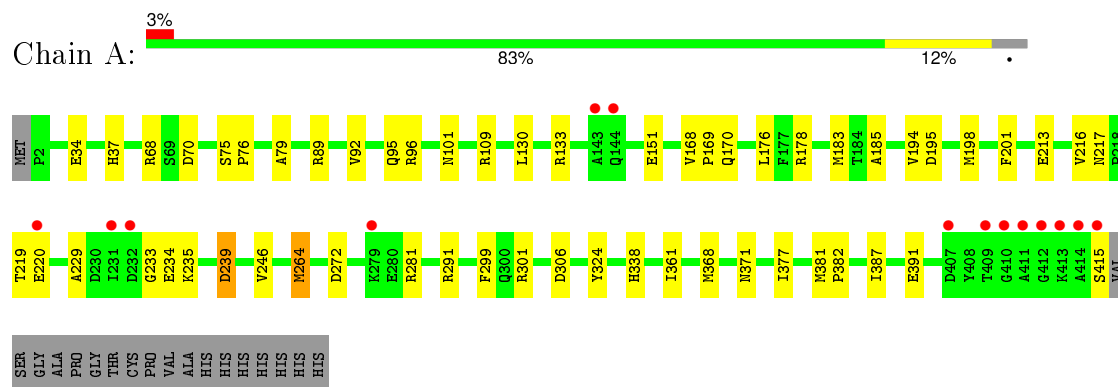
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	339	Total	O	0	0
			339	339		

### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: P450 HEME-THIOLATE PROTEIN



## 4 Data and refinement statistics

Property	Value	Source
Space group	I 41	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	108.19Å 108.19Å 120.23Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	76.50 – 2.00 54.09 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.2 (76.50-2.00) 99.3 (54.09-2.00)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.89 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, $R_{free}$	0.164 , 0.206 0.166 , 0.208	Depositor DCC
$R_{free}$ test set	2341 reflections (5.33%)	DCC
Wilson B-factor (Å <sup>2</sup> )	38.5	Xtriage
Anisotropy	0.088	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 44.4	EDS
Estimated twinning fraction	0.035 for -h,k,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 46295 reflections	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	3686	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.33% 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, 0.375 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: HEM, EDO, CL

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	1.14	3/3332 (0.1%)	1.01	10/4523 (0.2%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	324	TYR	CD1-CE1	6.66	1.49	1.39
1	A	201	PHE	CB-CG	6.04	1.61	1.51
1	A	151	GLU	CB-CG	5.44	1.62	1.52

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	68	ARG	NE-CZ-NH1	7.29	123.94	120.30
1	A	239	ASP	CB-CG-OD1	-7.08	111.93	118.30
1	A	89	ARG	NE-CZ-NH2	-6.94	116.83	120.30
1	A	178	ARG	NE-CZ-NH1	6.11	123.36	120.30
1	A	291	ARG	NE-CZ-NH2	-6.10	117.25	120.30
1	A	306	ASP	CB-CG-OD1	6.01	123.71	118.30
1	A	264	MET	CG-SD-CE	-5.54	91.34	100.20
1	A	109	ARG	CG-CD-NE	-5.42	100.42	111.80
1	A	70	ASP	CB-CG-OD1	5.33	123.10	118.30
1	A	301	ARG	NE-CZ-NH2	-5.00	117.80	120.30

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	3259	0	3138	49	0
2	A	44	0	65	9	0
3	A	43	0	30	2	0
4	A	1	0	0	0	0
5	A	339	0	0	6	0
All	All	3686	0	3233	50	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 8.

All (50) 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:219:THR:HG22	1:A:220:GLU:H	1.23	1.00
1:A:219:THR:HG22	1:A:220:GLU:N	1.84	0.90
1:A:195:ASP:HB3	2:A:1425:EDO:H12	1.55	0.87
3:A:1418:HEM:O1A	5:A:2277:HOH:O	1.91	0.87
1:A:264:MET:HE3	1:A:377:ILE:HG21	1.59	0.84
1:A:219:THR:CG2	1:A:220:GLU:H	1.92	0.83
1:A:185:ALA:HB2	1:A:194:VAL:CG2	2.10	0.82
1:A:185:ALA:HB2	1:A:194:VAL:HG21	1.64	0.79
1:A:264:MET:CE	1:A:377:ILE:HG21	2.13	0.79
1:A:281:ARG:O	2:A:1417:EDO:H22	1.85	0.76
1:A:198:MET:H	2:A:1425:EDO:H11	1.53	0.73
1:A:133:ARG:HH22	1:A:170:GLN:HE22	1.35	0.72
1:A:391:GLU:OE1	5:A:2327:HOH:O	2.07	0.72
1:A:34:GLU:OE1	1:A:37:HIS:HD2	1.72	0.71
1:A:195:ASP:HB3	2:A:1425:EDO:C1	2.19	0.70
1:A:216:VAL:HG23	1:A:217:ASN:ND2	2.10	0.67
1:A:264:MET:CE	1:A:377:ILE:HD13	2.25	0.67
1:A:133:ARG:NH2	1:A:170:GLN:HE22	1.95	0.65
1:A:272:ASP:HB2	5:A:2254:HOH:O	1.95	0.65
1:A:37:HIS:HE1	5:A:2013:HOH:O	1.80	0.64
1:A:216:VAL:HG23	1:A:217:ASN:HD22	1.66	0.61
1:A:195:ASP:CB	2:A:1425:EDO:H12	2.30	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:185:ALA:CB	1:A:194:VAL:CG2	2.80	0.58
1:A:183:MET:HE3	1:A:246:VAL:HG13	1.85	0.57
1:A:219:THR:CG2	1:A:220:GLU:N	2.54	0.56
1:A:361:ILE:HD12	3:A:1418:HEM:HBD2	1.88	0.55
1:A:387:ILE:HG22	1:A:387:ILE:O	2.07	0.54
1:A:234:GLU:OE1	1:A:234:GLU:HA	2.08	0.54
1:A:198:MET:N	2:A:1425:EDO:H11	2.21	0.53
1:A:183:MET:CE	1:A:246:VAL:HG13	2.38	0.53
1:A:79:ALA:H	1:A:101:ASN:ND2	2.06	0.53
1:A:133:ARG:HH22	1:A:170:GLN:NE2	2.04	0.51
1:A:264:MET:HE2	1:A:377:ILE:HD13	1.91	0.51
1:A:168:VAL:HG21	1:A:176:LEU:HD22	1.94	0.50
1:A:92:VAL:O	1:A:95:GLN:HG2	2.13	0.48
1:A:229:ALA:HB1	1:A:233:GLY:HA2	1.96	0.47
1:A:387:ILE:CG2	1:A:387:ILE:O	2.64	0.46
1:A:185:ALA:CB	1:A:194:VAL:HG22	2.46	0.45
1:A:133:ARG:HB3	5:A:2176:HOH:O	2.17	0.44
1:A:338:HIS:ND1	2:A:1420:EDO:H12	2.33	0.44
1:A:76:PRO:HA	2:A:1422:EDO:H11	1.98	0.44
1:A:213:GLU:O	1:A:216:VAL:HG22	2.18	0.44
1:A:381:MET:N	1:A:382:PRO:CD	2.81	0.44
1:A:75:SER:N	1:A:76:PRO:CD	2.81	0.43
1:A:96:ARG:HA	1:A:101:ASN:HB2	2.00	0.42
1:A:76:PRO:O	2:A:1422:EDO:C1	2.67	0.42
1:A:130:LEU:HD12	1:A:368:MET:CE	2.50	0.42
1:A:130:LEU:HD12	1:A:368:MET:HE3	2.02	0.42
1:A:371:ASN:HB3	5:A:2322:HOH:O	2.21	0.41
1:A:34:GLU:OE1	1:A:37:HIS:CD2	2.62	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	418/433 (96%)	404 (97%)	14 (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	341/357 (96%)	336 (98%)	5 (2%)	72	75

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

Mol	Chain	Res	Type
1	A	169	PRO
1	A	235	LYS
1	A	239	ASP
1	A	299	PHE
1	A	415	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	37	HIS
1	A	101	ASN
1	A	170	GLN
1	A	217	ASN
1	A	364	ASN
1	A	371	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 carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 13 ligands modelled in this entry, 1 is monoatomic - leaving 12 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$
2	EDO	A	1416	-	3,3,3	0.83	0	2,2,2	0.18	0
2	EDO	A	1417	-	3,3,3	0.94	0	2,2,2	0.62	0
3	HEM	A	1418	1,2	30,50,50	2.49	9 (30%)	24,82,82	2.50	11 (45%)
2	EDO	A	1419	-	3,3,3	1.14	0	2,2,2	0.59	0
2	EDO	A	1420	-	3,3,3	0.37	0	2,2,2	2.06	1 (50%)
2	EDO	A	1421	-	3,3,3	0.44	0	2,2,2	0.67	0
2	EDO	A	1422	-	3,3,3	0.40	0	2,2,2	1.77	1 (50%)
2	EDO	A	1423	3	3,3,3	0.91	0	2,2,2	0.84	0
2	EDO	A	1424	-	3,3,3	0.63	0	2,2,2	0.66	0
2	EDO	A	1425	-	3,3,3	0.55	0	2,2,2	0.15	0
2	EDO	A	1427	-	3,3,3	0.51	0	2,2,2	0.48	0
2	EDO	A	1428	-	3,3,3	0.57	0	2,2,2	0.17	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	EDO	A	1416	-	-	0/1/1/1	0/0/0/0
2	EDO	A	1417	-	-	0/1/1/1	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	HEM	A	1418	1,2	-	0/10/54/54	0/0/8/8
2	EDO	A	1419	-	-	0/1/1/1	0/0/0/0
2	EDO	A	1420	-	-	0/1/1/1	0/0/0/0
2	EDO	A	1421	-	-	0/1/1/1	0/0/0/0
2	EDO	A	1422	-	-	0/1/1/1	0/0/0/0
2	EDO	A	1423	3	-	0/1/1/1	0/0/0/0
2	EDO	A	1424	-	-	0/1/1/1	0/0/0/0
2	EDO	A	1425	-	-	0/1/1/1	0/0/0/0
2	EDO	A	1427	-	-	0/1/1/1	0/0/0/0
2	EDO	A	1428	-	-	0/1/1/1	0/0/0/0

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1418	HEM	C3B-C4B	-8.07	1.44	1.51
3	A	1418	HEM	C3D-C4D	-5.49	1.44	1.51
3	A	1418	HEM	C2D-C1D	-3.46	1.40	1.51
3	A	1418	HEM	C2C-C1C	-3.46	1.46	1.52
3	A	1418	HEM	C2B-C1B	-2.46	1.43	1.51
3	A	1418	HEM	C1C-NC	2.16	1.38	1.36
3	A	1418	HEM	FE-ND	2.54	2.10	1.97
3	A	1418	HEM	C3B-CAB	3.37	1.57	1.51
3	A	1418	HEM	CAA-C2A	3.44	1.57	1.52

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1418	HEM	CMA-C3A-C4A	-5.11	119.92	128.36
3	A	1418	HEM	CBA-CAA-C2A	-3.14	106.90	112.53
2	A	1420	EDO	O1-C1-C2	-2.82	92.33	112.54
3	A	1418	HEM	CBD-CAD-C3D	-2.46	106.41	113.55
2	A	1422	EDO	O2-C2-C1	-2.44	95.03	112.54
3	A	1418	HEM	CMA-C3A-C2A	2.27	129.98	125.24
3	A	1418	HEM	C3C-CAC-CBC	2.31	128.00	124.46
3	A	1418	HEM	C2D-C3D-C4D	2.41	105.58	101.50
3	A	1418	HEM	CMD-C2D-C3D	2.67	126.14	114.35
3	A	1418	HEM	CMC-C2C-C3C	3.34	124.88	116.53
3	A	1418	HEM	CAD-C3D-C2D	4.32	125.63	113.22
3	A	1418	HEM	CAD-C3D-C4D	4.62	128.77	112.47
3	A	1418	HEM	CMB-C2B-C3B	4.85	128.63	116.53

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1417	EDO	1	0
3	A	1418	HEM	2	0
2	A	1420	EDO	1	0
2	A	1422	EDO	2	0
2	A	1425	EDO	5	0

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

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	414/433 (95%)	-0.31	14 (3%)	49 50	26, 36, 59, 88	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	414	ALA	6.9
1	A	411	ALA	6.2
1	A	413	LYS	4.7
1	A	412	GLY	4.5
1	A	415	SER	4.4
1	A	231	ILE	4.0
1	A	410	GLY	3.1
1	A	220	GLU	2.4
1	A	409	THR	2.3
1	A	144	GLN	2.2
1	A	143	ALA	2.2
1	A	407	ASP	2.1
1	A	279	LYS	2.0
1	A	232	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

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
2	EDO	A	1428	4/4	0.89	0.31	9.11	57,62,64,68	0
2	EDO	A	1424	4/4	0.78	0.18	8.10	70,73,74,75	0
2	EDO	A	1419	4/4	0.78	0.17	6.92	51,58,58,62	0
2	EDO	A	1421	4/4	0.93	0.20	4.52	48,52,53,53	0
2	EDO	A	1422	4/4	0.80	0.19	4.10	54,57,59,63	0
2	EDO	A	1416	4/4	0.92	0.14	3.36	58,58,60,60	0
2	EDO	A	1427	4/4	0.96	0.13	2.03	43,46,52,53	0
2	EDO	A	1417	4/4	0.85	0.13	1.17	44,47,49,51	0
3	HEM	A	1418	43/43	0.99	0.12	0.82	23,27,31,39	0
2	EDO	A	1423	4/4	0.96	0.10	-0.83	39,40,41,46	0
4	CL	A	1426	1/1	0.96	0.23	-	67,67,67,67	1
2	EDO	A	1425	4/4	0.86	0.18	-	29,41,42,49	4
2	EDO	A	1420	4/4	0.89	0.15	-	49,56,56,62	0

## 6.5 Other polymers

There are no such residues in this entry.