



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 03:54 PM GMT

PDB ID : 4DTZ
Title : cytochrome P450 BM3h-8C8 MRI sensor bound to dopamine
Authors : Brustad, E.M.; Lelyveld, V.S.; Snow, C.D.; Crook, N.; Martinez, F.M.; Scholl, T.J.; Jasanoff, A.; Arnold, F.H.
Deposited on : 2012-02-21
Resolution : 1.55 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 (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

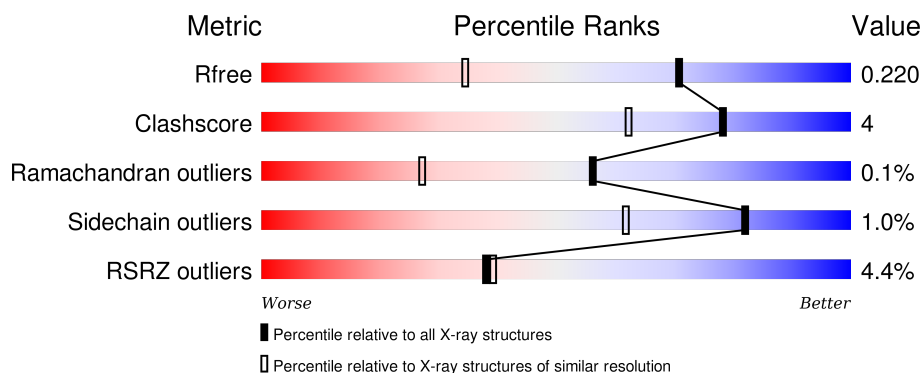
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.55 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	1665 (1.58-1.54)
Clashscore	102246	1014 (1.56-1.56)
Ramachandran outliers	100387	1704 (1.58-1.54)
Sidechain outliers	100360	1702 (1.58-1.54)
RSRZ outliers	91569	1668 (1.58-1.54)

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 ≥ 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	469	<div> <div>4%</div> <div>86%</div> <div>10%</div> <div>.</div> </div>
1	B	469	<div> <div>5%</div> <div>84%</div> <div>12%</div> <div>.</div> </div>

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
3	LDP	A	501	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8549 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 cytochrome P450 BM3 variant 8C8.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	453	Total	C	N	O	S	0	7	0
			3678	2349	629	682	18			
1	B	453	Total	C	N	O	S	0	6	0
			3665	2342	623	682	18			

There are 22 discrepancies between the modelled and reference sequences:

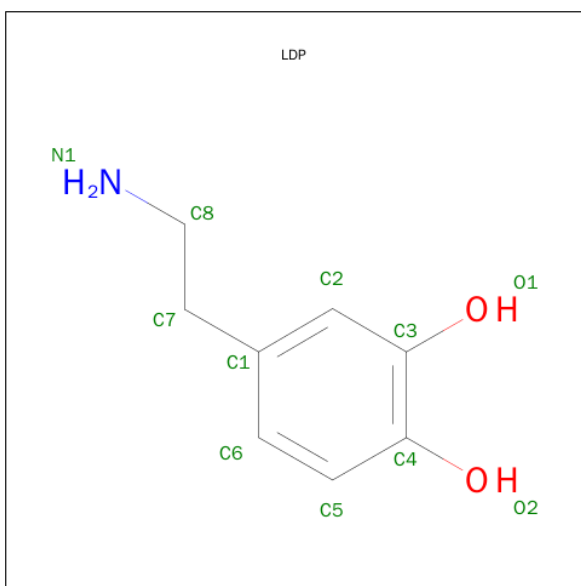
Chain	Residue	Modelled	Actual	Comment	Reference
A	75	PRO	LEU	ENGINEERED MUTATION	UNP P14779
A	189	ARG	GLN	ENGINEERED MUTATION	UNP P14779
A	263	ALA	ILE	ENGINEERED MUTATION	UNP P14779
A	268	ALA	THR	ENGINEERED MUTATION	UNP P14779
A	286	GLU	VAL	ENGINEERED MUTATION	UNP P14779
A	464	HIS	-	EXPRESSION TAG	UNP P14779
A	465	HIS	-	EXPRESSION TAG	UNP P14779
A	466	HIS	-	EXPRESSION TAG	UNP P14779
A	467	HIS	-	EXPRESSION TAG	UNP P14779
A	468	HIS	-	EXPRESSION TAG	UNP P14779
A	469	HIS	-	EXPRESSION TAG	UNP P14779
B	75	PRO	LEU	ENGINEERED MUTATION	UNP P14779
B	189	ARG	GLN	ENGINEERED MUTATION	UNP P14779
B	263	ALA	ILE	ENGINEERED MUTATION	UNP P14779
B	268	ALA	THR	ENGINEERED MUTATION	UNP P14779
B	286	GLU	VAL	ENGINEERED MUTATION	UNP P14779
B	464	HIS	-	EXPRESSION TAG	UNP P14779
B	465	HIS	-	EXPRESSION TAG	UNP P14779
B	466	HIS	-	EXPRESSION TAG	UNP P14779
B	467	HIS	-	EXPRESSION TAG	UNP P14779
B	468	HIS	-	EXPRESSION TAG	UNP P14779
B	469	HIS	-	EXPRESSION TAG	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 43	C 34	Fe 1	N 4	O 4	0	0
2	B	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 3 is L-DOPAMINE (three-letter code: LDP) (formula: $C_8H_{11}NO_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			11	8	1	2		
3	B	1	Total	C	N	O	0	0
			11	8	1	2		

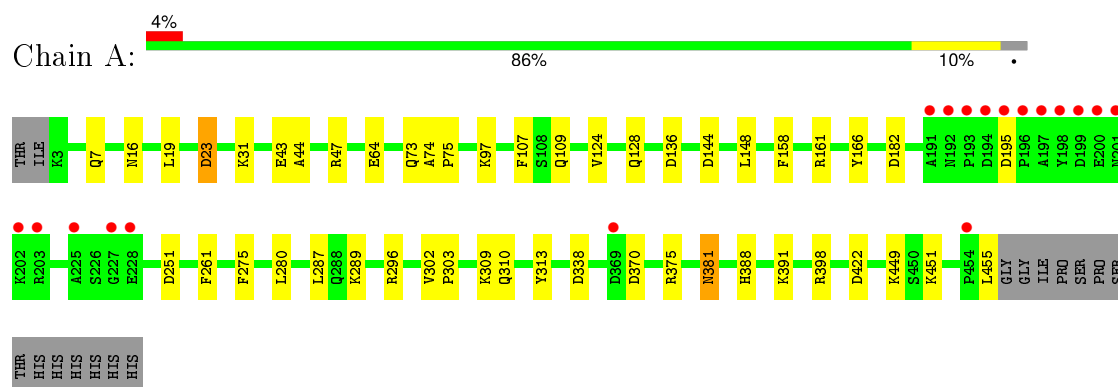
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	573	Total 573	O 573	0	0
4	B	525	Total 525	O 525	0	0

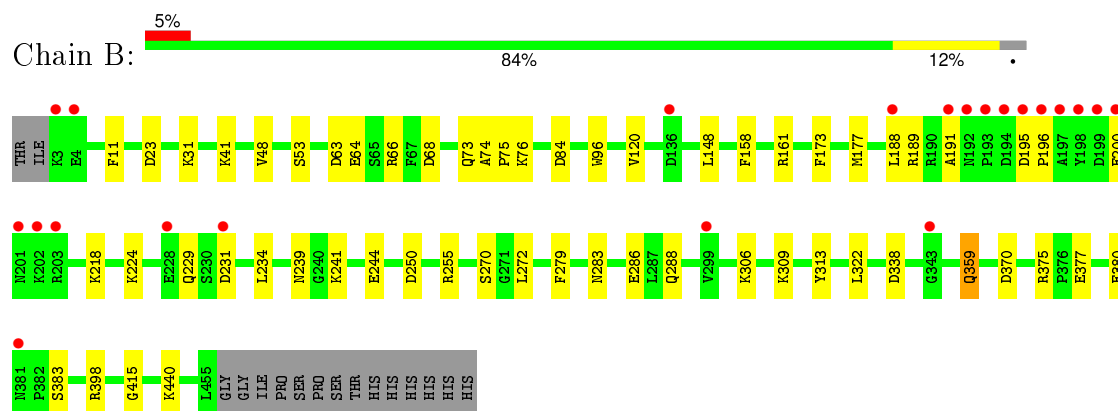
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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: cytochrome P450 BM3 variant 8C8



- Molecule 1: cytochrome P450 BM3 variant 8C8



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	58.70 Å 146.37 Å 63.98 Å 90.00° 97.59° 90.00°	Depositor
Resolution (Å)	37.39 – 1.55 37.39 – 1.55	Depositor EDS
% Data completeness (in resolution range)	95.6 (37.39-1.55) 95.6 (37.39-1.55)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.14 (at 1.55 Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.177 , 0.222 0.175 , 0.220	Depositor DCC
R_{free} test set	7801 reflections (5.56%)	DCC
Wilson B-factor (Å ²)	17.0	Xtriage
Anisotropy	0.052	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 49.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Outliers	0 of 147676 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	8549	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.33% 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, 0.375 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, LDP

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	1.26	3/3785 (0.1%)	1.11	10/5116 (0.2%)
1	B	1.29	12/3769 (0.3%)	1.17	17/5095 (0.3%)
All	All	1.27	15/7554 (0.2%)	1.14	27/10211 (0.3%)

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	288	GLN	CG-CD	9.67	1.73	1.51
1	A	107	PHE	CE1-CZ	7.28	1.51	1.37
1	B	173	PHE	CE2-CZ	6.52	1.49	1.37
1	B	279	PHE	CD1-CE1	6.42	1.52	1.39
1	B	377	GLU	CG-CD	6.21	1.61	1.51
1	B	359	GLN	CB-CG	-6.18	1.35	1.52
1	B	64	GLU	CD-OE2	-5.86	1.19	1.25
1	A	64	GLU	CD-OE1	5.73	1.31	1.25
1	B	288	GLN	CD-OE1	5.72	1.36	1.24
1	B	313	TYR	CD1-CE1	5.54	1.47	1.39
1	B	415	GLY	C-O	5.32	1.32	1.23
1	B	309	LYS	CD-CE	5.22	1.64	1.51
1	B	120	VAL	CB-CG2	5.13	1.63	1.52
1	A	166	TYR	CE2-CZ	5.12	1.45	1.38
1	B	11	PHE	CE2-CZ	5.10	1.47	1.37

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	370	ASP	CB-CG-OD1	9.17	126.55	118.30
1	A	161	ARG	NE-CZ-NH1	8.82	124.71	120.30
1	B	84	ASP	CB-CG-OD2	-7.63	111.43	118.30
1	A	19	LEU	CB-CG-CD1	-7.56	98.16	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	66	ARG	NE-CZ-NH2	-7.30	116.65	120.30
1	B	398	ARG	NE-CZ-NH2	-7.30	116.65	120.30
1	A	338	ASP	CB-CG-OD1	7.03	124.62	118.30
1	B	84	ASP	CB-CG-OD1	6.79	124.42	118.30
1	A	275	PHE	CB-CG-CD2	-6.58	116.19	120.80
1	B	66	ARG	NE-CZ-NH1	6.35	123.47	120.30
1	B	338	ASP	CB-CG-OD1	6.27	123.94	118.30
1	A	251	ASP	CB-CG-OD1	6.18	123.86	118.30
1	A	251	ASP	CB-CG-OD2	-5.96	112.93	118.30
1	B	375	ARG	NE-CZ-NH2	5.96	123.28	120.30
1	A	144	ASP	CB-CG-OD2	-5.88	113.00	118.30
1	A	398	ARG	NE-CZ-NH2	-5.87	117.36	120.30
1	B	234	LEU	CB-CG-CD2	-5.79	101.16	111.00
1	B	161	ARG	NE-CZ-NH2	-5.75	117.42	120.30
1	B	63	ASP	CB-CG-OD2	-5.74	113.13	118.30
1	B	177	MET	CG-SD-CE	-5.61	91.23	100.20
1	B	63	ASP	CB-CG-OD1	5.56	123.30	118.30
1	B	68	ASP	CB-CG-OD2	-5.56	113.30	118.30
1	B	96	TRP	CA-CB-CG	-5.55	103.16	113.70
1	A	261	PHE	CB-CG-CD2	-5.21	117.16	120.80
1	B	255	ARG	NE-CZ-NH2	-5.18	117.71	120.30
1	B	250	ASP	CB-CG-OD2	-5.09	113.72	118.30
1	A	23	ASP	CB-CG-OD2	-5.05	113.76	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	3678	0	3655	28	1
1	B	3665	0	3633	26	0
2	A	43	0	30	0	0
2	B	43	0	30	0	0
3	A	11	0	9	0	0
3	B	11	0	9	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	573	0	0	9	1
4	B	525	0	0	12	2
All	All	8549	0	7366	53	2

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 (53) 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:B:229:GLN:HE22	1:B:239:ASN:HD21	1.08	0.98
1:A:74:ALA:HB3	1:A:75:PRO:HD3	1.54	0.90
1:A:296[B]:ARG:NH2	4:A:866:HOH:O	2.06	0.87
1:A:16:ASN:HD22	1:A:43:GLU:H	1.21	0.86
1:B:244:GLU:HG3	4:B:1119:HOH:O	1.76	0.85
1:B:23:ASP:OD2	4:B:702:HOH:O	1.93	0.85
1:B:74:ALA:HB3	1:B:75:PRO:HD3	1.57	0.84
1:A:7:GLN:HE21	1:A:16:ASN:HD21	1.25	0.84
1:B:380[A]:GLU:OE2	4:B:1116:HOH:O	1.99	0.80
1:A:136:ASP:OD2	1:B:218:LYS:NZ	2.18	0.74
1:A:370:ASP:OD2	1:A:375[A]:ARG:NH2	2.21	0.74
1:B:31:LYS:HE2	4:B:712:HOH:O	1.96	0.66
1:B:74:ALA:HB3	1:B:75:PRO:CD	2.29	0.62
1:A:310:GLN:HG2	4:A:1106:HOH:O	2.00	0.62
1:B:41:LYS:HE2	1:B:48:VAL:HG21	1.82	0.61
1:A:289:LYS:NZ	1:A:289:LYS:HB2	2.16	0.60
1:A:73:GLN:NE2	4:A:1034:HOH:O	2.19	0.59
1:A:388:HIS:HA	1:A:391:LYS:HD3	1.84	0.59
1:A:97:LYS:HB2	4:A:743:HOH:O	2.02	0.59
1:B:73:GLN:HG3	1:B:188:LEU:HD22	1.85	0.58
1:B:76:LYS:HD2	4:B:1082:HOH:O	2.04	0.58
1:A:381:ASN:ND2	4:A:885:HOH:O	2.37	0.57
1:A:7:GLN:HE21	1:A:16:ASN:ND2	2.00	0.56
1:A:280:LEU:HB3	1:A:287:LEU:HD13	1.88	0.55
1:B:229:GLN:HE22	1:B:239:ASN:ND2	1.90	0.54
1:A:289:LYS:HB2	1:A:289:LYS:HZ1	1.70	0.54
1:A:289:LYS:NZ	1:A:289:LYS:CB	2.71	0.53
1:A:74:ALA:HB3	1:A:75:PRO:CD	2.35	0.53
1:B:73:GLN:CG	1:B:188:LEU:HD22	2.40	0.51
1:B:224:LYS:NZ	4:B:870:HOH:O	2.41	0.51
1:B:229:GLN:NE2	1:B:239:ASN:HD21	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:47:ARG:HD2	4:A:988:HOH:O	2.09	0.51
1:B:306:LYS:HG2	4:B:697:HOH:O	2.11	0.50
1:A:128:GLN:NE2	4:A:822:HOH:O	2.45	0.49
1:B:200:GLU:HA	1:B:200:GLU:OE1	2.13	0.49
1:A:289:LYS:HD2	1:A:313:TYR:CE1	2.49	0.48
1:A:109:GLN:HE22	1:A:309:LYS:HZ2	1.61	0.48
1:A:449:LYS:HE3	4:A:1109:HOH:O	2.15	0.47
1:B:440:LYS:HE2	4:B:1057:HOH:O	2.13	0.47
1:A:124:VAL:HG13	1:A:455:LEU:HD13	1.95	0.47
1:A:422:ASP:OD1	1:A:451:LYS:NZ	2.46	0.46
1:B:195:ASP:OD2	1:B:196:PRO:HD2	2.14	0.46
1:A:16:ASN:ND2	1:A:43:GLU:H	2.01	0.46
1:B:189:ARG:HD2	4:B:1054:HOH:O	2.16	0.46
1:A:44:ALA:HB3	1:A:47:ARG:HG3	2.00	0.44
1:B:270[B]:SER:OG	4:B:891:HOH:O	2.21	0.43
1:B:53:SER:HB3	1:B:359:GLN:HB3	2.01	0.42
1:A:31:LYS:HE2	4:A:680:HOH:O	2.19	0.41
1:B:283:ASN:HB3	1:B:286:GLU:OE1	2.20	0.41
1:B:272:LEU:HD13	1:B:322:LEU:HG	2.01	0.41
1:A:302:VAL:HA	1:A:303:PRO:HD3	1.96	0.41
1:B:241:LYS:HE3	4:B:901:HOH:O	2.21	0.41
1:B:231:ASP:CG	4:B:1092:HOH:O	2.59	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:23:ASP:OD2	4:B:1077:HOH:O[2_646]	2.05	0.15
4:A:1065:HOH:O	4:B:1044:HOH:O[2_545]	2.10	0.10

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	458/469 (98%)	446 (97%)	12 (3%)	0	100	100
1	B	457/469 (97%)	444 (97%)	12 (3%)	1 (0%)	52	25
All	All	915/938 (98%)	890 (97%)	24 (3%)	1 (0%)	56	27

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	191	ALA

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	399/409 (98%)	394 (99%)	5 (1%)	76	51
1	B	397/409 (97%)	394 (99%)	3 (1%)	86	71
All	All	796/818 (97%)	788 (99%)	8 (1%)	82	62

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	148	LEU
1	A	158	PHE
1	A	182	ASP
1	A	195	ASP
1	A	381	ASN
1	B	148	LEU
1	B	158	PHE
1	B	383	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	16	ASN
1	A	109	GLN

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Mol	Chain	Res	Type
1	A	381	ASN
1	B	229	GLN

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 ⓘ

4 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	500	1,3	30,50,50	2.67	8 (26%)	24,82,82	2.51	13 (54%)
3	LDP	A	501	2	10,11,11	1.60	2 (20%)	13,14,14	1.90	2 (15%)
2	HEM	B	500	1,3	30,50,50	2.23	9 (30%)	24,82,82	2.79	12 (50%)
3	LDP	B	501	2	10,11,11	1.89	3 (30%)	13,14,14	2.01	5 (38%)

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	500	1,3	-	0/10/54/54	0/0/8/8
3	LDP	A	501	2	-	0/3/3/3	0/1/1/1
2	HEM	B	500	1,3	-	0/10/54/54	0/0/8/8
3	LDP	B	501	2	-	0/3/3/3	0/1/1/1

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	500	HEM	C3B-C4B	-10.65	1.42	1.51
2	B	500	HEM	C3B-C4B	-6.81	1.45	1.51
2	A	500	HEM	C3D-C4D	-5.99	1.43	1.51
2	A	500	HEM	C2D-C3D	-4.18	1.42	1.54
2	B	500	HEM	C3D-C4D	-3.72	1.46	1.51
2	B	500	HEM	C2C-C1C	-3.70	1.45	1.52
2	B	500	HEM	C2D-C1D	-3.23	1.41	1.51
2	A	500	HEM	C2B-C1B	-2.46	1.43	1.51
2	A	500	HEM	C2D-C1D	-2.40	1.44	1.51
2	B	500	HEM	C2B-C1B	-2.18	1.44	1.51
2	A	500	HEM	CHC-C4B	-2.09	1.33	1.38
3	B	501	LDP	O2-C4	2.10	1.40	1.36
2	A	500	HEM	CMC-C2C	2.12	1.58	1.53
3	B	501	LDP	C2-C1	2.30	1.43	1.39
2	B	500	HEM	C3C-CAC	2.32	1.55	1.51
2	B	500	HEM	CHC-C1C	2.50	1.42	1.36
2	B	500	HEM	CAA-C2A	2.65	1.56	1.52
3	A	501	LDP	C3-C4	2.72	1.44	1.40
2	A	500	HEM	CMA-C3A	2.75	1.57	1.51
3	A	501	LDP	C5-C6	3.29	1.44	1.38
2	B	500	HEM	C4C-NC	3.70	1.40	1.36
3	B	501	LDP	C3-C4	4.45	1.47	1.40

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	501	LDP	C8-C7-C1	-4.82	102.93	112.82
3	B	501	LDP	C6-C5-C4	-4.61	115.77	120.49
2	B	500	HEM	CBA-CAA-C2A	-4.48	104.50	112.53
2	B	500	HEM	CAA-CBA-CGA	-4.44	104.62	112.75
2	B	500	HEM	C4B-CHC-C1C	-4.17	118.85	125.82
2	A	500	HEM	CMA-C3A-C4A	-3.00	123.40	128.36
2	A	500	HEM	CBA-CAA-C2A	-2.98	107.18	112.53
2	B	500	HEM	CMA-C3A-C4A	-2.88	123.59	128.36
2	B	500	HEM	CBD-CAD-C3D	-2.87	105.20	113.55

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	500	HEM	C1D-CHD-C4C	-2.85	121.06	125.82
2	A	500	HEM	CBD-CAD-C3D	-2.62	105.94	113.55
3	B	501	LDP	C3-C2-C1	-2.18	118.19	120.83
3	B	501	LDP	C2-C3-C4	2.01	121.64	119.82
2	A	500	HEM	CMA-C3A-C2A	2.08	129.58	125.24
2	A	500	HEM	CMD-C2D-C3D	2.17	123.93	114.35
2	B	500	HEM	C2D-C3D-C4D	2.38	105.53	101.50
3	B	501	LDP	C5-C6-C1	2.39	124.30	121.04
2	A	500	HEM	C3C-CAC-CBC	2.40	128.14	124.46
2	B	500	HEM	CHD-C1D-ND	2.49	130.52	124.52
2	A	500	HEM	C2D-C3D-C4D	2.49	105.73	101.50
2	A	500	HEM	CMB-C2B-C3B	2.72	123.31	116.53
2	A	500	HEM	C3B-C4B-CHC	3.27	127.77	123.16
2	B	500	HEM	CMC-C2C-C3C	3.48	125.23	116.53
2	A	500	HEM	C3B-CAB-CBB	3.49	129.80	124.46
2	B	500	HEM	CAD-C3D-C4D	3.60	125.18	112.47
3	B	501	LDP	C8-C7-C1	3.64	120.29	112.82
3	A	501	LDP	C2-C3-C4	3.67	123.14	119.82
2	A	500	HEM	CAD-C3D-C4D	3.74	125.67	112.47
2	B	500	HEM	CMB-C2B-C3B	4.56	127.91	116.53
2	A	500	HEM	CAD-C3D-C2D	5.20	128.18	113.22
2	A	500	HEM	CMC-C2C-C3C	5.36	129.91	116.53
2	B	500	HEM	CAD-C3D-C2D	5.59	129.29	113.22

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	453/469 (96%)	0.02	18 (3%) 42 44	9, 17, 33, 55	0
1	B	453/469 (96%)	0.04	22 (4%) 33 34	9, 17, 37, 70	0
All	All	906/938 (96%)	0.03	40 (4%) 38 39	9, 17, 35, 70	0

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	191	ALA	8.1
1	B	196	PRO	7.7
1	B	197	ALA	7.3
1	A	196	PRO	6.8
1	B	198	TYR	6.7
1	A	191	ALA	6.0
1	A	199	ASP	5.6
1	A	198	TYR	5.3
1	A	192	ASN	5.0
1	A	200	GLU	4.1
1	B	195	ASP	4.1
1	B	192	ASN	3.9
1	A	197	ALA	3.9
1	A	194	ASP	3.7
1	B	3	LYS	3.6
1	B	194	ASP	3.5
1	B	193	PRO	3.4
1	A	203	ARG	3.2
1	A	195	ASP	3.2
1	B	203	ARG	3.1
1	B	201	ASN	2.9
1	A	202	LYS	2.9
1	B	200	GLU	2.9
1	A	193	PRO	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	188	LEU	2.8
1	B	202	LYS	2.8
1	A	201	ASN	2.7
1	B	136	ASP	2.6
1	B	343	GLY	2.4
1	A	225	ALA	2.4
1	B	4	GLU	2.4
1	A	369	ASP	2.3
1	B	299	VAL	2.2
1	A	227	GLY	2.2
1	A	454	PRO	2.2
1	B	228	GLU	2.1
1	A	228	GLU	2.1
1	B	199	ASP	2.1
1	B	231	ASP	2.1
1	B	381	ASN	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	LDP	A	501	11/11	0.97	0.14	2.76	10,12,14,15	0
3	LDP	B	501	11/11	0.95	0.12	1.98	11,13,16,22	0
2	HEM	A	500	43/43	0.98	0.12	0.87	6,10,12,16	0
2	HEM	B	500	43/43	0.98	0.10	0.69	6,9,13,14	0

6.5 Other polymers

There are no such residues in this entry.