



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:35 AM GMT

PDB ID : 3BDZ
Title : The Role of Asn 242 in P450cin
Authors : Mehareenna, Y.T.; Poulos, T.L.
Deposited on : 2007-11-15
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 (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

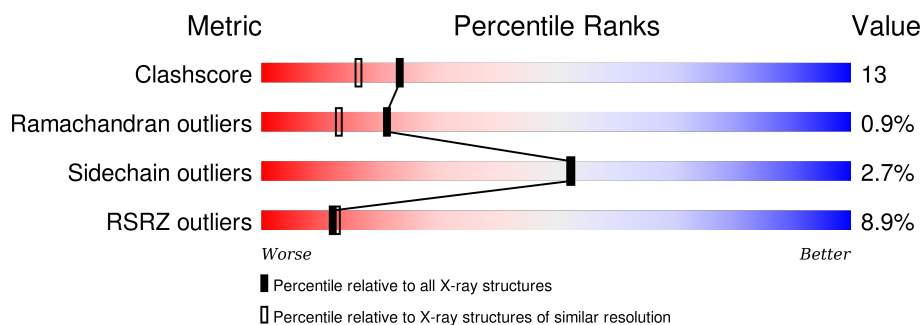
1 Overall quality at a glance

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))
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 ≥ 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	397	<div> <div>5%</div> <div>80%</div> <div>18%</div> <div>.</div> </div>
1	B	397	<div> <div>13%</div> <div>80%</div> <div>18%</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	MLI	A	500[A]	-	-	-	X
3	MLI	A	500[B]	-	-	X	X

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6933 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.

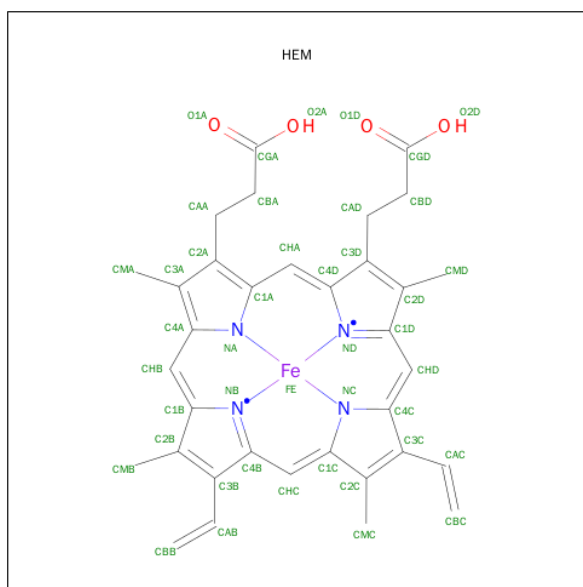
- Molecule 1 is a protein called P450cin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	397	Total	C	N	O	S	0	0	0
			3146	2006	548	581	11			
1	B	397	Total	C	N	O	S	0	0	0
			3146	2006	548	581	11			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	242	ALA	ASN	ENGINEERED	UNP Q8VQF6
B	242	ALA	ASN	ENGINEERED	UNP Q8VQF6

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



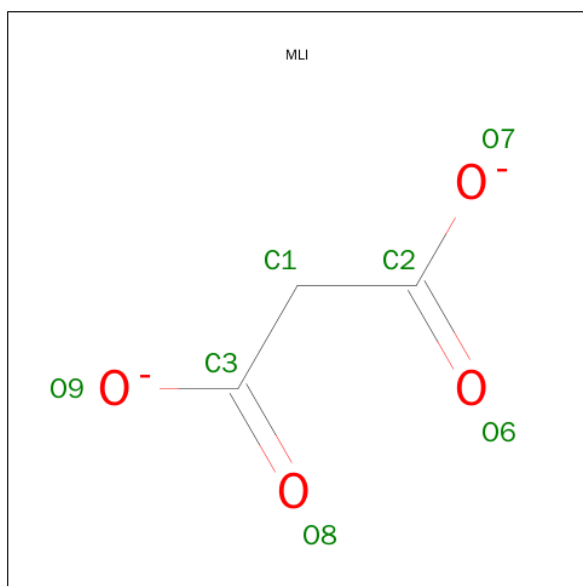
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

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

- Molecule 3 is MALONATE ION (three-letter code: MLI) (formula: $C_3H_2O_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	1
			14	6	8		
3	B	1	Total	C	O	0	1
			14	6	8		

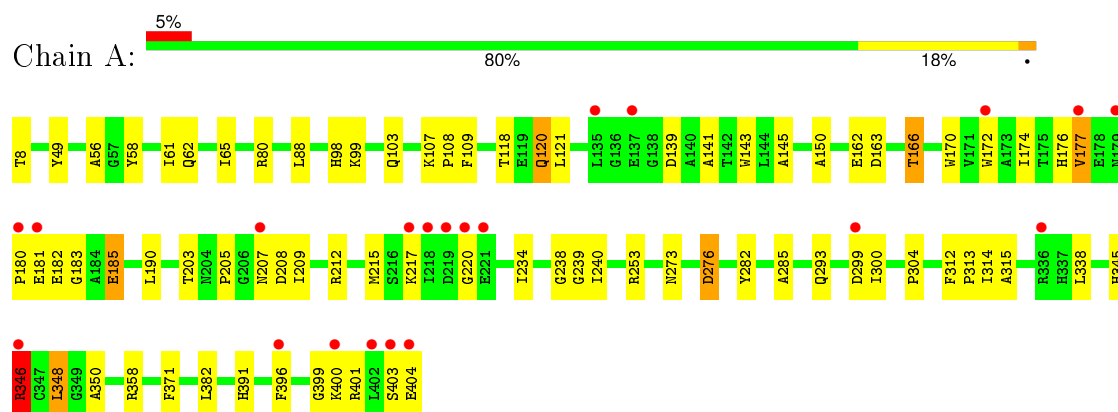
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	289	Total	O	0	0
			289	289		
4	B	238	Total	O	0	0
			238	238		

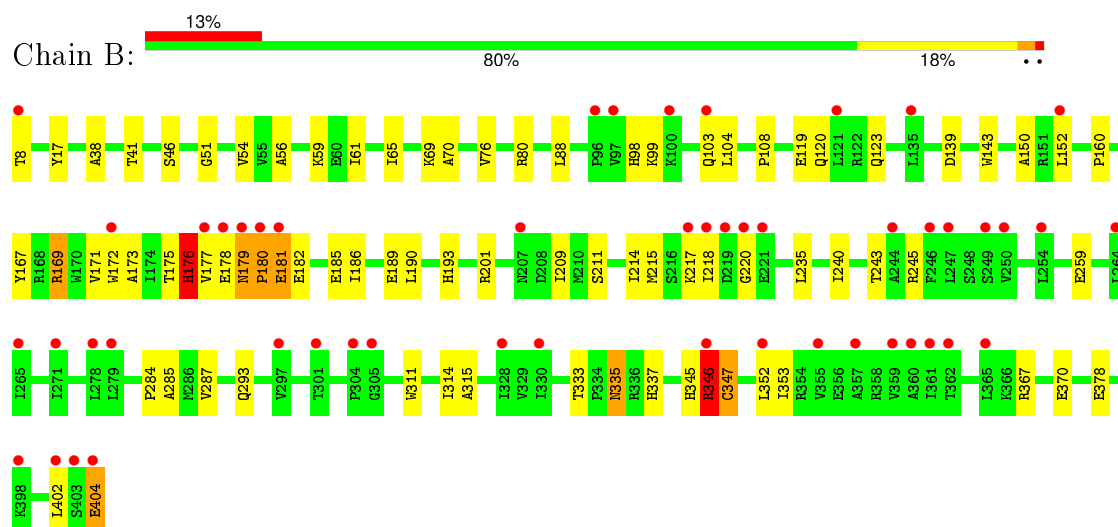
3 Residue-property plots [i](#)

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: P450cin



• Molecule 1: P450cin



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	63.33Å 68.23Å 103.56Å 90.00° 95.53° 90.00°	Depositor
Resolution (Å)	46.30 – 2.00 46.30 – 2.00	Depositor EDS
% Data completeness (in resolution range)	94.5 (46.30-2.00) 94.5 (46.30-2.00)	Depositor EDS
R_{merge}	0.20	Depositor
R_{sym}	0.20	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.95 (at 2.00Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.181 , 0.227 0.244 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	25.2	Xtriage
Anisotropy	0.393	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 56.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	2 of 56408 reflections (0.004%)	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6933	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.03% 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, MLI

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	0.68	0/3226	0.75	2/4386 (0.0%)
1	B	0.69	0/3226	0.77	3/4386 (0.1%)
All	All	0.69	0/6452	0.76	5/8772 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	276	ASP	CB-CG-OD1	-5.92	112.97	118.30
1	B	285	ALA	N-CA-C	5.76	126.56	111.00
1	A	285	ALA	N-CA-C	5.55	125.99	111.00
1	B	169	ARG	NE-CZ-NH1	-5.36	117.62	120.30
1	B	176	HIS	N-CA-C	5.00	124.51	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	49	TYR	Sidechain

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	3146	0	3098	83	0
1	B	3146	0	3098	82	0
2	A	43	0	30	7	0
2	B	43	0	30	8	0
3	A	14	0	4	3	0
3	B	14	0	4	0	0
4	A	289	0	0	13	0
4	B	238	0	0	12	0
All	All	6933	0	6264	169	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 13.

All (169) 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:172:TRP:CZ3	1:A:177:VAL:CG2	2.10	1.33
1:A:172:TRP:CZ3	1:A:177:VAL:HG22	1.71	1.24
1:A:172:TRP:CE3	4:A:778:HOH:O	1.96	1.14
1:A:172:TRP:HE3	4:A:778:HOH:O	1.27	1.13
1:A:172:TRP:CH2	1:A:177:VAL:HG22	1.90	1.05
1:B:172:TRP:CZ3	1:B:177:VAL:HG22	1.95	1.00
1:A:172:TRP:CE3	1:A:177:VAL:HG23	1.97	0.99
1:A:172:TRP:CZ3	1:A:177:VAL:HG23	2.03	0.94
1:A:172:TRP:CE3	1:A:177:VAL:CG2	2.54	0.87
1:B:172:TRP:CZ3	1:B:177:VAL:CG2	2.58	0.84
1:B:172:TRP:CH2	1:B:177:VAL:HG22	2.19	0.76
1:A:98:HIS:HE1	2:A:450:HEM:O1D	1.69	0.75
1:B:123:GLN:HG3	4:B:724:HOH:O	1.88	0.73
1:B:243:THR:HG22	1:B:353:ILE:HD13	1.70	0.73
1:A:181:GLU:O	1:A:185:GLU:HB3	1.88	0.73
1:B:345:HIS:O	1:B:346:ARG:O	2.09	0.70
3:A:500[B]:MLI:O6	4:A:685:HOH:O	2.10	0.69
1:B:347:CYS:HB2	2:B:450:HEM:NA	2.08	0.68
1:B:345:HIS:O	1:B:346:ARG:C	2.31	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:403:SER:O	1:A:404:GLU:HB3	1.95	0.66
1:A:293:GLN:HA	1:A:304:PRO:HG3	1.78	0.66
1:A:98:HIS:CE1	2:A:450:HEM:O1D	2.49	0.66
2:B:450:HEM:HMB1	2:B:450:HEM:HBB2	1.78	0.66
1:B:172:TRP:CE3	1:B:177:VAL:HG22	2.30	0.66
1:B:346:ARG:NH2	4:B:665:HOH:O	2.28	0.65
1:B:98:HIS:HE1	2:B:450:HEM:O1D	1.79	0.65
1:A:401:ARG:NH2	1:A:404:GLU:HG2	2.11	0.65
1:B:80:ARG:HG2	1:B:80:ARG:HH11	1.61	0.65
1:A:172:TRP:CH2	1:A:177:VAL:CG2	2.65	0.64
1:A:172:TRP:CZ3	1:A:177:VAL:HG21	2.27	0.64
1:A:120:GLN:HB3	4:A:771:HOH:O	1.96	0.64
1:B:177:VAL:HG11	1:B:182:GLU:CD	2.18	0.63
1:B:370:GLU:HB3	4:B:663:HOH:O	1.99	0.62
1:B:217:LYS:HE3	1:B:220:GLY:O	2.00	0.61
1:A:109:PHE:CD2	1:A:348:LEU:HD13	2.35	0.61
1:A:107:LYS:HB2	1:A:108:PRO:HD3	1.82	0.61
2:A:450:HEM:HBB2	2:A:450:HEM:HMB1	1.84	0.60
1:A:239:GLY:N	2:A:450:HEM:HAC	2.16	0.60
1:B:104:LEU:HD11	1:B:218:ILE:HD11	1.82	0.60
1:B:99:LYS:HE3	1:B:103:GLN:NE2	2.15	0.60
1:A:177:VAL:O	1:A:177:VAL:HG12	2.00	0.60
1:B:345:HIS:HD2	2:B:450:HEM:O2D	1.84	0.59
1:A:346:ARG:NH1	1:A:350:ALA:CB	2.66	0.59
1:B:347:CYS:HB2	2:B:450:HEM:C1A	2.38	0.59
1:A:177:VAL:HG11	1:A:183:GLY:N	2.17	0.59
1:B:178:GLU:HG3	4:B:728:HOH:O	2.03	0.58
1:B:108:PRO:HB2	1:B:209:ILE:CD1	2.34	0.57
1:B:108:PRO:CB	1:B:209:ILE:HD12	2.33	0.57
1:A:190:LEU:C	1:A:190:LEU:HD23	2.25	0.57
1:B:345:HIS:HE1	4:B:521:HOH:O	1.88	0.57
1:A:107:LYS:HG2	4:A:657:HOH:O	2.05	0.56
1:A:345:HIS:O	1:A:346:ARG:C	2.42	0.56
1:B:152:LEU:HD21	1:B:352:LEU:HD11	1.86	0.56
1:B:173:ALA:HA	1:B:177:VAL:HG23	1.88	0.55
1:A:391:HIS:HE1	4:A:662:HOH:O	1.89	0.55
1:A:162:GLU:H	1:A:162:GLU:CD	2.10	0.55
1:A:346:ARG:NH1	1:A:350:ALA:HB1	2.21	0.55
1:B:178:GLU:O	1:B:180:PRO:N	2.39	0.55
1:A:358:ARG:HD3	4:A:621:HOH:O	2.07	0.55
1:B:17:TYR:OH	1:B:176:HIS:HD2	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:172:TRP:CZ3	1:B:177:VAL:HG21	2.41	0.54
1:B:119:GLU:HG3	4:B:721:HOH:O	2.07	0.54
1:B:345:HIS:C	1:B:346:ARG:O	2.45	0.54
1:A:80:ARG:CZ	4:A:569:HOH:O	2.55	0.54
1:B:108:PRO:HB2	1:B:209:ILE:HD12	1.88	0.54
1:B:169:ARG:NH1	1:B:189:GLU:OE1	2.34	0.53
1:A:166:THR:HG21	4:A:774:HOH:O	2.07	0.52
1:A:345:HIS:O	1:A:346:ARG:O	2.27	0.52
1:B:175:THR:O	1:B:176:HIS:HB2	2.08	0.52
1:B:335:ASN:ND2	1:B:337:HIS:HD2	2.08	0.52
1:B:378:GLU:HA	1:B:378:GLU:OE1	2.08	0.52
1:A:203:THR:O	1:A:205:PRO:HD3	2.10	0.51
1:B:172:TRP:CE3	1:B:177:VAL:CG2	2.92	0.51
1:A:98:HIS:C	1:A:98:HIS:CD2	2.83	0.51
1:B:175:THR:O	1:B:176:HIS:CB	2.59	0.51
1:B:190:LEU:HD23	1:B:190:LEU:C	2.31	0.51
1:A:109:PHE:CG	1:A:348:LEU:HD13	2.47	0.50
1:A:234:ILE:HD11	3:A:500[B]:MLI:O8	2.12	0.50
1:A:253:ARG:HD3	1:A:282:TYR:OH	2.11	0.50
1:A:401:ARG:NH2	1:A:404:GLU:CG	2.74	0.50
1:A:346:ARG:HH11	1:A:350:ALA:CB	2.25	0.49
1:B:98:HIS:CE1	2:B:450:HEM:O1D	2.62	0.49
1:B:172:TRP:CH2	1:B:177:VAL:HG13	2.47	0.49
1:A:150:ALA:HA	1:A:240:ILE:HG12	1.94	0.49
1:A:399:GLY:C	1:A:400:LYS:HD2	2.32	0.49
1:A:62:GLN:HA	1:A:62:GLN:OE1	2.11	0.49
1:B:69:LYS:HG2	4:B:624:HOH:O	2.12	0.49
1:A:109:PHE:CD2	1:A:348:LEU:CD1	2.96	0.49
1:B:139:ASP:O	1:B:143:TRP:HB3	2.13	0.49
1:B:209:ILE:HD13	4:B:667:HOH:O	2.13	0.49
1:B:217:LYS:HE3	1:B:220:GLY:C	2.34	0.48
1:B:160:PRO:HD2	1:B:167:TYR:OH	2.14	0.48
1:A:346:ARG:HH12	1:A:350:ALA:HB1	1.78	0.48
1:A:166:THR:HG22	4:A:719:HOH:O	2.13	0.48
1:A:56:ALA:HB1	1:A:314:ILE:HG21	1.94	0.48
1:B:186:ILE:HD12	1:B:186:ILE:N	2.29	0.48
1:B:120:GLN:HG2	4:B:722:HOH:O	2.14	0.47
1:A:176:HIS:O	1:A:177:VAL:C	2.52	0.47
1:A:293:GLN:CA	1:A:304:PRO:HG3	2.43	0.47
1:A:205:PRO:HG3	1:A:215:MET:CE	2.44	0.47
1:B:150:ALA:HB1	4:B:708:HOH:O	2.13	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:203:THR:C	1:A:205:PRO:HD3	2.34	0.47
1:B:181:GLU:O	1:B:185:GLU:HB2	2.13	0.47
1:A:345:HIS:O	2:A:450:HEM:HAA2	2.14	0.47
1:B:175:THR:O	1:B:176:HIS:CD2	2.68	0.47
1:B:179:ASN:O	1:B:182:GLU:HB3	2.15	0.47
1:B:335:ASN:C	1:B:335:ASN:HD22	2.18	0.46
1:A:61:ILE:HD12	1:A:315:ALA:HB2	1.97	0.46
1:B:88:LEU:HA	4:B:584:HOH:O	2.16	0.46
1:A:170:TRP:O	1:A:174:ILE:HG13	2.15	0.46
1:A:177:VAL:CG1	1:A:182:GLU:HB3	2.46	0.46
1:B:180:PRO:O	1:B:182:GLU:N	2.49	0.46
1:A:176:HIS:NE2	1:A:382:LEU:HD11	2.31	0.46
1:B:311:TRP:HB3	1:B:314:ILE:HG12	1.98	0.45
1:A:88:LEU:HA	4:A:595:HOH:O	2.16	0.45
1:B:38:ALA:HA	1:B:41:THR:O	2.16	0.45
1:B:70:ALA:HA	1:B:293:GLN:HG2	1.97	0.45
1:B:404:GLU:O	1:B:404:GLU:OE2	2.34	0.45
1:B:61:ILE:HD12	1:B:315:ALA:HB2	1.99	0.45
1:B:108:PRO:CB	1:B:209:ILE:CD1	2.95	0.45
1:A:99:LYS:O	1:A:103:GLN:HG3	2.17	0.45
1:A:172:TRP:CD2	1:A:177:VAL:HG23	2.50	0.44
1:B:171:VAL:HG21	1:B:240:ILE:HD11	1.98	0.44
1:B:367:ARG:CZ	1:B:402:LEU:HG	2.47	0.44
1:A:172:TRP:CZ3	4:A:778:HOH:O	2.40	0.44
1:B:245:ARG:HA	1:B:245:ARG:NE	2.32	0.44
1:B:189:GLU:O	1:B:193:HIS:HD2	2.00	0.44
1:A:62:GLN:NE2	1:A:338:LEU:HD21	2.32	0.44
1:B:54:VAL:HG13	1:B:311:TRP:CD1	2.53	0.44
1:B:99:LYS:O	1:B:103:GLN:HG3	2.18	0.44
1:B:259:GLU:OE2	4:B:683:HOH:O	2.20	0.44
1:A:207:ASN:HA	1:A:212:ARG:HH21	1.82	0.44
1:B:182:GLU:O	1:B:186:ILE:HD13	2.18	0.43
1:A:98:HIS:ND1	1:A:345:HIS:CE1	2.86	0.43
1:B:347:CYS:HB2	2:B:450:HEM:C4A	2.53	0.43
2:A:450:HEM:HBB2	2:A:450:HEM:CMB	2.48	0.43
1:A:139:ASP:O	1:A:143:TRP:HB3	2.19	0.43
1:A:177:VAL:HG11	1:A:182:GLU:HB3	2.00	0.43
1:A:400:LYS:N	1:A:400:LYS:HD2	2.34	0.43
1:B:56:ALA:HB1	1:B:314:ILE:HG21	2.00	0.43
1:B:211:SER:O	1:B:215:MET:HG3	2.19	0.42
1:B:17:TYR:OH	1:B:176:HIS:CD2	2.71	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:141:ALA:HA	1:A:145:ALA:HB3	2.02	0.42
1:A:181:GLU:CA	1:A:181:GLU:OE2	2.68	0.42
1:A:176:HIS:O	1:A:177:VAL:O	2.38	0.42
1:B:46:SER:O	1:B:51:GLY:HA2	2.19	0.42
1:B:98:HIS:CE1	1:B:345:HIS:CD2	3.07	0.42
1:B:76:VAL:HG23	1:B:287:VAL:CG2	2.49	0.42
1:A:118:THR:O	1:A:121:LEU:HB2	2.19	0.42
1:A:181:GLU:OE2	1:A:181:GLU:HA	2.19	0.42
1:B:284:PRO:HD2	2:B:450:HEM:HMB3	2.02	0.41
1:B:169:ARG:HH12	1:B:189:GLU:CD	2.22	0.41
1:B:201:ARG:HB2	1:B:214:ILE:CD1	2.49	0.41
1:B:180:PRO:C	1:B:182:GLU:N	2.73	0.41
1:A:163:ASP:HB3	1:A:166:THR:HG23	2.02	0.41
1:A:108:PRO:HB3	1:A:209:ILE:HG12	2.01	0.41
1:A:238:GLY:HA2	3:A:500[A]:MLI:O6	2.21	0.41
1:A:299:ASP:C	1:A:300:ILE:HD12	2.41	0.41
1:A:312:PHE:HB2	1:A:313:PRO:HD3	2.02	0.41
1:A:371:PHE:HB3	1:A:396:PHE:HB3	2.03	0.41
1:A:239:GLY:CA	2:A:450:HEM:HAC	2.51	0.41
1:A:217:LYS:HD3	1:A:220:GLY:O	2.20	0.41
1:A:273:ASN:HB2	4:A:582:HOH:O	2.21	0.41
1:B:80:ARG:CG	1:B:80:ARG:HH11	2.30	0.41
1:B:217:LYS:HE2	1:B:217:LYS:HB3	1.94	0.41
1:A:58:TYR:HA	1:A:315:ALA:HB1	2.03	0.40
1:A:208:ASP:O	1:A:212:ARG:HG2	2.22	0.40
1:A:345:HIS:C	1:A:346:ARG:O	2.58	0.40
1:B:108:PRO:HB2	1:B:209:ILE:HD11	2.02	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	395/397 (100%)	380 (96%)	12 (3%)	3 (1%)	24	15
1	B	395/397 (100%)	379 (96%)	12 (3%)	4 (1%)	19	11
All	All	790/794 (100%)	759 (96%)	24 (3%)	7 (1%)	21	13

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	177	VAL
1	A	346	ARG
1	B	176	HIS
1	B	181	GLU
1	B	346	ARG
1	B	180	PRO
1	A	180	PRO

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	332/332 (100%)	324 (98%)	8 (2%)	57	58
1	B	332/332 (100%)	322 (97%)	10 (3%)	48	47
All	All	664/664 (100%)	646 (97%)	18 (3%)	52	52

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	THR
1	A	65	ILE
1	A	120	GLN
1	A	166	THR
1	A	185	GLU
1	A	276	ASP
1	A	346	ARG
1	A	348	LEU
1	B	8	THR

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Mol	Chain	Res	Type
1	B	59	LYS
1	B	65	ILE
1	B	179	ASN
1	B	235	LEU
1	B	333	THR
1	B	335	ASN
1	B	346	ARG
1	B	347	CYS
1	B	404	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	98	HIS
1	A	204	ASN
1	A	391	HIS
1	B	66	GLN
1	B	98	HIS
1	B	103	GLN
1	B	176	HIS
1	B	179	ASN
1	B	193	HIS
1	B	335	ASN
1	B	345	HIS

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

6 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	450	1,3	30,50,50	2.55	8 (26%)	24,82,82	2.49	8 (33%)
3	MLI	A	500[A]	2	0,6,6	0.00	-	0,7,7	0.00	-
3	MLI	A	500[B]	-	0,6,6	0.00	-	0,7,7	0.00	-
2	HEM	B	450	1,3	30,50,50	2.89	9 (30%)	24,82,82	2.39	8 (33%)
3	MLI	B	500[A]	2	0,6,6	0.00	-	0,7,7	0.00	-
3	MLI	B	500[B]	-	0,6,6	0.00	-	0,7,7	0.00	-

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	450	1,3	-	0/10/54/54	0/0/8/8
3	MLI	A	500[A]	2	-	0/0/4/4	0/0/0/0
3	MLI	A	500[B]	-	-	0/0/4/4	0/0/0/0
2	HEM	B	450	1,3	-	0/10/54/54	0/0/8/8
3	MLI	B	500[A]	2	-	0/0/4/4	0/0/0/0
3	MLI	B	500[B]	-	-	0/0/4/4	0/0/0/0

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	450	HEM	C3B-C4B	-7.57	1.45	1.51
2	A	450	HEM	C2D-C3D	-7.18	1.33	1.54
2	A	450	HEM	C3C-CAC	-7.14	1.37	1.51
2	B	450	HEM	C3C-CAC	-6.80	1.38	1.51
2	B	450	HEM	C3B-CAB	-5.80	1.40	1.51
2	A	450	HEM	C3B-C4B	-5.41	1.47	1.51
2	B	450	HEM	C2D-C3D	-5.33	1.38	1.54
2	B	450	HEM	C2C-C1C	-5.25	1.42	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	450	HEM	C3B-CAB	-4.77	1.42	1.51
2	A	450	HEM	C3D-C4D	-3.49	1.47	1.51
2	A	450	HEM	C2D-C1D	-2.24	1.44	1.51
2	A	450	HEM	CAA-C2A	2.02	1.55	1.52
2	B	450	HEM	CHD-C4C	2.15	1.41	1.36
2	B	450	HEM	CAA-C2A	2.31	1.56	1.52
2	B	450	HEM	CMB-C2B	2.63	1.59	1.53
2	A	450	HEM	CHD-C4C	2.81	1.43	1.36
2	B	450	HEM	C1C-NC	3.71	1.40	1.36

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	450	HEM	C3C-CAC-CBC	-4.23	117.96	124.46
2	B	450	HEM	CBA-CAA-C2A	-3.57	106.14	112.53
2	A	450	HEM	CBD-CAD-C3D	-2.55	106.14	113.55
2	B	450	HEM	CMA-C3A-C4A	-2.08	124.92	128.36
2	A	450	HEM	C2D-C3D-C4D	2.12	105.09	101.50
2	B	450	HEM	CAD-C3D-C4D	2.73	122.11	112.47
2	A	450	HEM	CMD-C2D-C3D	3.28	128.88	114.35
2	B	450	HEM	CMC-C2C-C3C	3.56	125.41	116.53
2	B	450	HEM	CMB-C2B-C3B	3.99	126.48	116.53
2	A	450	HEM	CAD-C3D-C4D	4.29	127.59	112.47
2	B	450	HEM	CMD-C2D-C3D	4.32	133.45	114.35
2	A	450	HEM	CMB-C2B-C3B	4.35	127.39	116.53
2	B	450	HEM	C2D-C3D-C4D	4.44	109.02	101.50
2	A	450	HEM	CAD-C3D-C2D	4.78	126.96	113.22
2	B	450	HEM	CAD-C3D-C2D	5.43	128.82	113.22
2	A	450	HEM	CMC-C2C-C3C	5.98	131.45	116.53

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 18 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	450	HEM	7	0
3	A	500[A]	MLI	1	0
3	A	500[B]	MLI	2	0
2	B	450	HEM	8	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 ⓘ

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	397/397 (100%)	0.56	21 (5%)	30 32	42, 52, 74, 89	0
1	B	397/397 (100%)	0.95	50 (12%)	5 5	43, 54, 73, 89	0
All	All	794/794 (100%)	0.75	71 (8%)	12 13	42, 53, 74, 89	0

All (71) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	404	GLU	8.2
1	A	220	GLY	7.4
1	B	218	ILE	6.6
1	B	177	VAL	6.5
1	A	172	TRP	6.3
1	B	404	GLU	6.1
1	A	219	ASP	5.9
1	B	172	TRP	5.9
1	B	180	PRO	5.8
1	B	219	ASP	5.8
1	B	220	GLY	5.7
1	A	181	GLU	5.3
1	B	8	THR	5.2
1	A	177	VAL	5.1
1	A	218	ILE	4.8
1	B	221	GLU	4.5
1	A	180	PRO	4.1
1	A	179	ASN	4.1
1	B	181	GLU	4.0
1	B	246	PHE	3.7
1	B	179	ASN	3.7
1	B	250	VAL	3.5
1	B	265	ILE	3.4
1	B	178	GLU	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	217	LYS	3.3
1	A	403	SER	3.3
1	B	217	LYS	3.2
1	B	361	ILE	3.2
1	B	96	PRO	3.1
1	A	135	LEU	3.1
1	B	100	LYS	3.1
1	B	304	PRO	3.0
1	B	271	ILE	3.0
1	B	360	ALA	3.0
1	B	254	LEU	3.0
1	A	299	ASP	2.9
1	B	403	SER	2.9
1	B	402	LEU	2.8
1	A	221	GLU	2.7
1	A	346	ARG	2.7
1	B	355	VAL	2.6
1	A	402	LEU	2.5
1	B	247	LEU	2.5
1	B	121	LEU	2.5
1	B	398	LYS	2.5
1	A	400	LYS	2.4
1	B	346	ARG	2.4
1	B	244	ALA	2.4
1	B	362	THR	2.4
1	A	207	ASN	2.3
1	B	135	LEU	2.3
1	B	328	ILE	2.3
1	B	365	LEU	2.3
1	B	357	ALA	2.3
1	B	297	VAL	2.3
1	B	103	GLN	2.3
1	B	352	LEU	2.3
1	B	249	SER	2.2
1	B	279	LEU	2.2
1	B	359	VAL	2.2
1	B	97	VAL	2.2
1	B	207	ASN	2.2
1	A	336	ARG	2.1
1	B	152	LEU	2.1
1	B	264	LEU	2.1
1	B	330	ILE	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	137	GLU	2.1
1	B	301	THR	2.1
1	B	278	LEU	2.1
1	B	305	GLY	2.1
1	A	396	PHE	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 [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	MLI	A	500[A]	7/7	0.86	0.24	9.91	21,27,30,31	7
3	MLI	A	500[B]	7/7	0.86	0.24	6.08	9,13,20,23	7
3	MLI	B	500[B]	7/7	0.87	0.19	1.76	24,27,28,28	7
3	MLI	B	500[A]	7/7	0.87	0.19	1.21	26,29,30,31	7
2	HEM	A	450	43/43	0.91	0.14	0.00	42,45,47,55	0
2	HEM	B	450	43/43	0.92	0.15	-0.38	45,47,52,55	0

6.5 Other polymers [i](#)

There are no such residues in this entry.