



# Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 11:14 PM GMT

PDB ID : 1J51  
Title : CRYSTAL STRUCTURE OF CYTOCHROME P450CAM MUTANT (F87W/Y96F/V247L/C334A)WITH 1,3,5-TRICHLOROBENZENE  
Authors : Chen, X.; Christopher, A.; Jones, J.; Guo, Q.; Xu, F.; Cao, R.; Wong, L-L.; Rao, Z.  
Deposited on : 2002-01-05  
Resolution : 2.20 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

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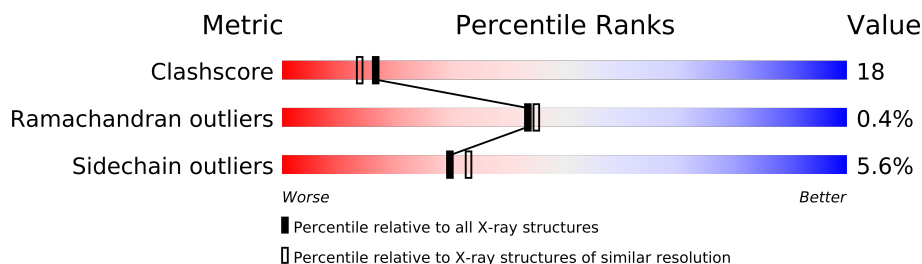
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 21963  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.20 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	79885	3751 (2.20-2.20)
Ramachandran outliers	78287	3681 (2.20-2.20)
Sidechain outliers	78261	3682 (2.20-2.20)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	414	
1	B	414	
1	C	414	
1	D	414	

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 13468 atoms, of which 0 are hydrogen and 0 are deuterium.

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 P450CAM.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	405	Total	C	N	O	S	0	0	0
			3209	2036	561	595	17			
1	B	405	Total	C	N	O	S	0	0	0
			3209	2036	561	595	17			
1	C	405	Total	C	N	O	S	0	0	0
			3209	2036	561	595	17			
1	D	405	Total	C	N	O	S	0	0	0
			3209	2036	561	595	17			

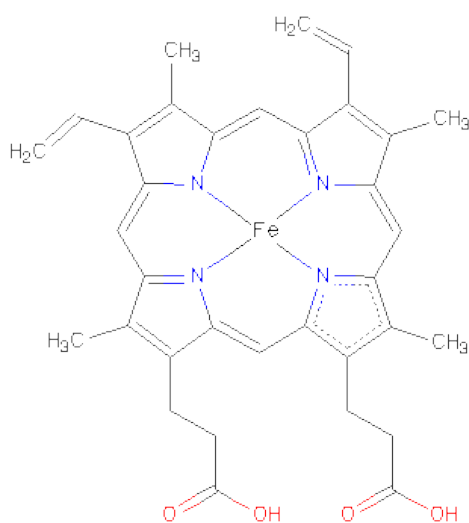
There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	87	TRP	PHE	ENGINEERED	UNP P00183
A	96	PHE	TYR	ENGINEERED	UNP P00183
A	247	LEU	VAL	ENGINEERED	UNP P00183
A	334	ALA	CYS	ENGINEERED	UNP P00183
B	87	TRP	PHE	ENGINEERED	UNP P00183
B	96	PHE	TYR	ENGINEERED	UNP P00183
B	247	LEU	VAL	ENGINEERED	UNP P00183
B	334	ALA	CYS	ENGINEERED	UNP P00183
C	87	TRP	PHE	ENGINEERED	UNP P00183
C	96	PHE	TYR	ENGINEERED	UNP P00183
C	247	LEU	VAL	ENGINEERED	UNP P00183
C	334	ALA	CYS	ENGINEERED	UNP P00183
D	87	TRP	PHE	ENGINEERED	UNP P00183
D	96	PHE	TYR	ENGINEERED	UNP P00183
D	247	LEU	VAL	ENGINEERED	UNP P00183
D	334	ALA	CYS	ENGINEERED	UNP P00183

- Molecule 2 is POTASSIUM ION (three-letter code: K) (formula: K).

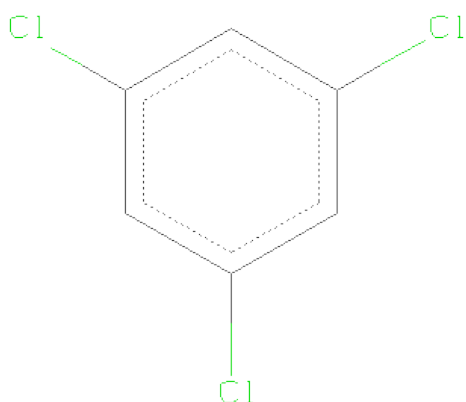
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total K 1 1	0	0
2	A	1	Total K 1 1	0	0
2	D	1	Total K 1 1	0	0
2	C	1	Total K 1 1	0	0

- Molecule 3 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula:  $C_{34}H_{32}FeN_4O_4$ ).



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

- Molecule 4 is 1,3,5-TRICHLORO-BENZENE (three-letter code: TCZ) (formula:  $C_6H_3Cl_3$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	Cl	0	0
			9	6	3		
4	B	1	Total	C	Cl	0	0
			9	6	3		
4	C	1	Total	C	Cl	0	0
			9	6	3		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	136	Total	O	0	0
			136	136		
5	B	80	Total	O	0	0
			80	80		
5	C	121	Total	O	0	0
			121	121		
5	D	92	Total	O	0	0
			92	92		

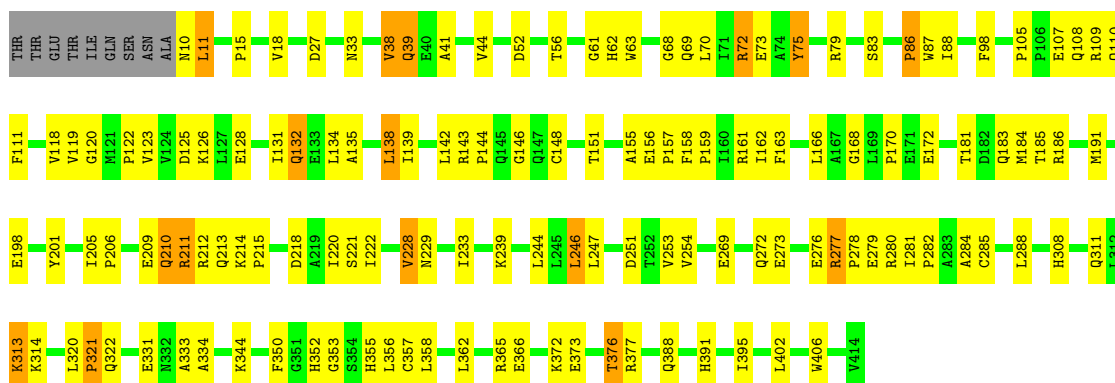
### 3 Residue-property plots

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.

Note EDS was not executed.

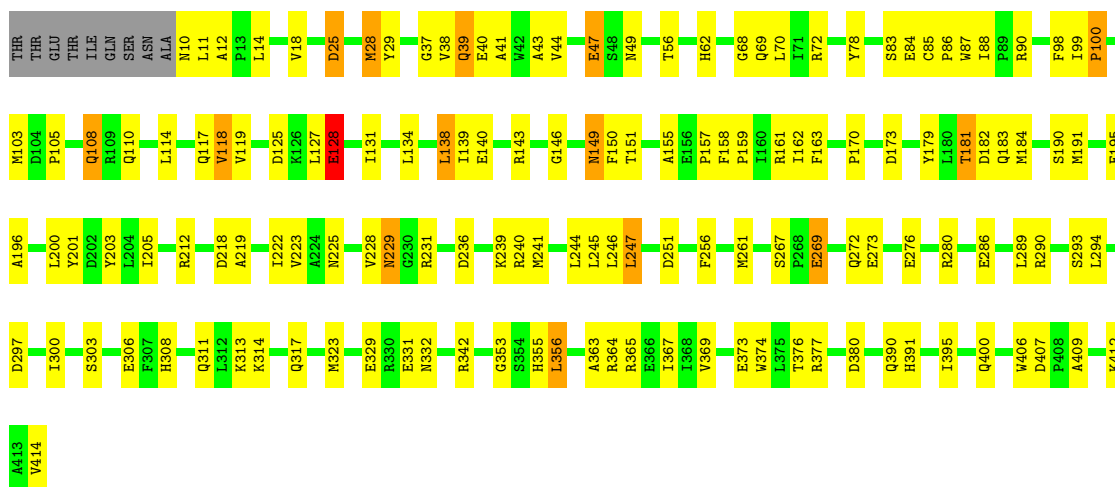
#### • Molecule 1: CYTOCHROME P450CAM

Chain A: 



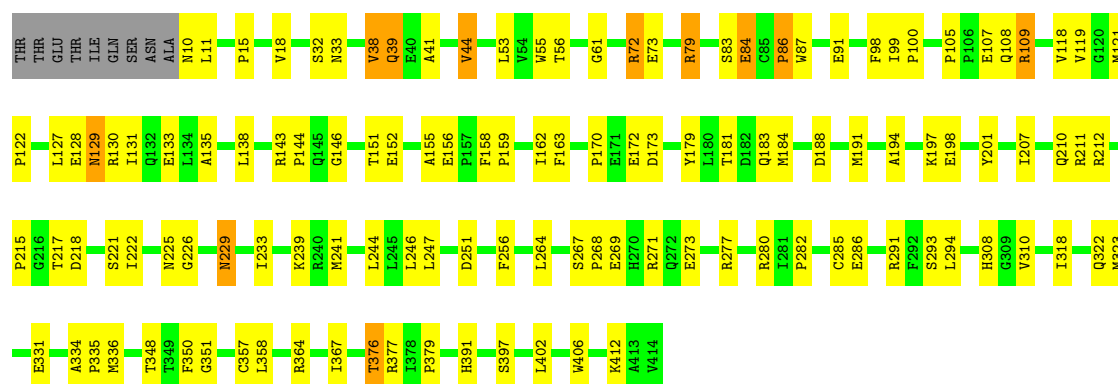
#### • Molecule 1: CYTOCHROME P450CAM

Chain B: 



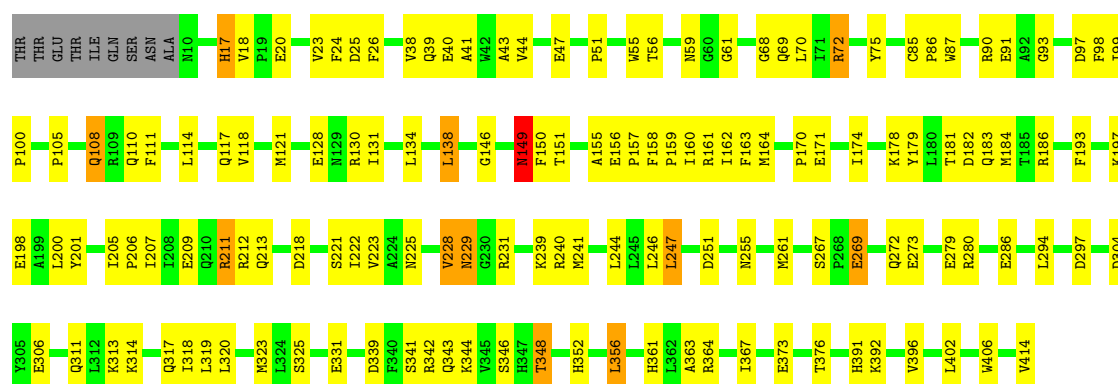
#### • Molecule 1: CYTOCHROME P450CAM

Chain C: 



# Molecule 1: CYTOCHROME P450CAM

Chain D:



## 4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	66.87Å 62.44Å 95.52Å 89.98° 90.30° 90.09°	Depositor
Resolution (Å)	500.00 – 2.20	Depositor
% Data completeness (in resolution range)	99.8 (500.00-2.20)	Depositor
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.186 , 0.258	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	13468	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	21.0	wwPDB-VP

## 5 Model quality

### 5.1 Standard geometry

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

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	0.63	0/3289	0.83	4/4470 (0.1%)
1	B	0.63	0/3289	0.80	1/4470 (0.0%)
1	C	0.65	0/3289	0.82	2/4470 (0.0%)
1	D	0.62	0/3289	0.80	3/4470 (0.1%)
All	All	0.63	0/13156	0.81	10/17880 (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
1	C	0	1
All	All	0	2

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	72	ARG	NE-CZ-NH1	-7.51	116.54	120.30
1	A	72	ARG	NE-CZ-NH1	-6.66	116.97	120.30
1	B	247	LEU	N-CA-C	-5.75	95.48	111.00
1	A	109	ARG	NE-CZ-NH2	-5.70	117.45	120.30
1	D	149	ASN	N-CA-C	-5.68	95.66	111.00
1	A	109	ARG	NE-CZ-NH1	5.59	123.10	120.30
1	D	320	LEU	N-CA-C	-5.54	96.05	111.00
1	A	105	PRO	N-CA-C	-5.42	98.01	112.10
1	C	105	PRO	N-CA-C	-5.20	98.59	112.10
1	D	247	LEU	N-CA-C	-5.17	97.05	111.00

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	75	TYR	Sidechain
1	C	179	TYR	Sidechain

## 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3209	0	3159	115	0
1	B	3209	0	3159	136	0
1	C	3209	0	3159	92	0
1	D	3209	0	3159	126	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	43	0	30	2	0
3	B	43	0	30	2	0
3	C	43	0	30	1	0
3	D	43	0	30	2	0
4	A	9	0	3	0	0
4	B	9	0	3	1	0
4	C	9	0	3	0	0
5	A	136	0	0	4	0
5	B	80	0	0	7	0
5	C	121	0	0	4	0
5	D	92	0	0	8	0
All	All	13468	0	12765	465	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 18.

All (465) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:210:GLN:HE21	1:A:210:GLN:N	1.33	1.26
1:C:131:ILE:HG12	1:C:162:ILE:HD13	1.28	1.15
1:D:207:ILE:HG23	1:D:211:ARG:NH1	1.65	1.10
1:B:163:PHE:HE2	1:B:246:LEU:HD23	1.25	1.02
1:B:191:MET:HE2	1:B:196:ALA:HA	1.42	1.01
1:D:110:GLN:HB3	1:D:228:VAL:HG23	1.47	0.95
1:B:181:THR:HG22	5:B:2452:HOH:O	1.68	0.94
1:A:210:GLN:N	1:A:210:GLN:NE2	2.18	0.91
1:D:200:LEU:HD11	1:D:246:LEU:HD12	1.53	0.90
1:D:72:ARG:HG3	1:D:72:ARG:HH11	1.34	0.90
1:D:131:ILE:HG12	1:D:162:ILE:HD13	1.55	0.88
1:B:131:ILE:HG12	1:B:162:ILE:HD13	1.55	0.88
1:B:163:PHE:CE2	1:B:246:LEU:HD23	2.09	0.88
1:D:117:GLN:HB2	5:D:4469:HOH:O	1.73	0.87
1:C:87:TRP:HZ3	1:C:184:MET:O	1.60	0.84
1:D:376:THR:HG22	1:D:414:VAL:HG21	1.59	0.84
1:C:163:PHE:CE2	1:C:246:LEU:HD23	2.12	0.84
1:C:131:ILE:HG12	1:C:162:ILE:CD1	2.08	0.84
1:B:99:ILE:HD11	1:B:240:ARG:HH21	1.42	0.84
1:D:207:ILE:HG23	1:D:211:ARG:HH11	1.44	0.83
1:D:163:PHE:HE2	1:D:246:LEU:HD23	1.43	0.82
1:A:210:GLN:HE21	1:A:210:GLN:H	1.24	0.82
1:A:72:ARG:HH12	1:A:331:GLU:CD	1.83	0.82
1:A:72:ARG:NH1	1:A:331:GLU:OE2	2.12	0.82
1:D:151:THR:O	1:D:156:GLU:HG3	1.81	0.80
1:B:414:VAL:HG12	1:B:414:VAL:O	1.81	0.80
1:C:163:PHE:HE2	1:C:246:LEU:HD23	1.47	0.79
1:A:277:ARG:HA	1:A:277:ARG:HE	1.47	0.78
1:C:87:TRP:CZ3	1:C:184:MET:O	2.37	0.78
1:C:72:ARG:HH12	1:C:331:GLU:CD	1.88	0.77
1:D:207:ILE:CG2	1:D:211:ARG:NH1	2.48	0.77
1:C:129:ASN:H	1:C:129:ASN:HD22	1.31	0.77
1:C:129:ASN:N	1:C:129:ASN:HD22	1.81	0.77
1:C:376:THR:HG22	1:C:377:ARG:HG3	1.68	0.76
1:C:376:THR:O	1:C:412:LYS:HE3	1.86	0.75
1:A:158:PHE:HB3	1:A:159:PRO:HD3	1.67	0.75
1:D:97:ASP:O	1:D:240:ARG:HD2	1.87	0.74
1:B:118:VAL:HG22	5:B:2471:HOH:O	1.86	0.74
1:B:182:ASP:OD1	5:B:2452:HOH:O	2.06	0.74
1:A:87:TRP:HZ3	1:A:184:MET:O	1.71	0.72
1:A:212:ARG:NH2	1:A:233:ILE:O	2.22	0.72
1:C:212:ARG:NH2	1:C:233:ILE:O	2.23	0.72
1:B:56:THR:HG21	1:B:62:HIS:CE1	2.25	0.72

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:131:ILE:HG12	1:A:162:ILE:HD13	1.71	0.72
1:D:131:ILE:HA	1:D:162:ILE:HD11	1.72	0.71
1:D:72:ARG:HD3	1:D:352:HIS:CE1	2.26	0.71
1:C:212:ARG:HA	1:C:225:ASN:HD21	1.54	0.71
1:A:181:THR:HG23	1:A:247:LEU:HD12	1.73	0.71
1:B:87:TRP:CZ3	1:B:184:MET:O	2.43	0.71
1:D:273:GLU:CG	1:D:280:ARG:HH12	2.03	0.70
1:D:228:VAL:O	1:D:231:ARG:HG3	1.90	0.70
1:C:33:ASN:HB3	1:C:41:ALA:HA	1.71	0.70
1:A:172:GLU:HB3	1:D:272:GLN:NE2	2.06	0.70
1:B:163:PHE:HE2	1:B:246:LEU:CD2	2.01	0.70
1:A:277:ARG:HG3	1:A:280:ARG:NH2	2.07	0.70
1:B:110:GLN:CD	1:B:229:ASN:HB2	2.11	0.70
1:A:206:PRO:O	1:A:210:GLN:NE2	2.25	0.69
1:B:110:GLN:NE2	1:B:229:ASN:HB2	2.07	0.69
1:B:62:HIS:CD2	1:B:88:ILE:HD13	2.28	0.69
1:B:376:THR:HG22	1:B:414:VAL:HG21	1.75	0.68
1:A:211:ARG:HG2	1:A:221:SER:OG	1.94	0.68
1:B:87:TRP:HZ3	1:B:184:MET:O	1.76	0.68
1:B:228:VAL:O	1:B:231:ARG:HG3	1.93	0.68
1:C:181:THR:HG23	1:C:247:LEU:HD12	1.76	0.68
1:B:131:ILE:HA	1:B:162:ILE:HD11	1.75	0.67
1:D:44:VAL:O	1:D:47:GLU:HG2	1.94	0.67
1:D:72:ARG:HG3	1:D:72:ARG:NH1	2.07	0.67
1:B:110:GLN:HG3	1:B:229:ASN:HB2	1.76	0.67
1:A:210:GLN:HE21	1:A:210:GLN:CA	2.06	0.67
1:A:278:PRO:O	1:A:281:ILE:HG13	1.95	0.67
1:C:41:ALA:O	1:C:44:VAL:HG13	1.94	0.67
1:A:209:GLU:HB3	1:A:210:GLN:HE22	1.60	0.67
1:B:313:LYS:HG3	1:B:314:LYS:N	2.09	0.66
1:D:414:VAL:HG12	1:D:414:VAL:O	1.93	0.66
1:C:211:ARG:HG3	1:C:221:SER:OG	1.96	0.65
1:B:72:ARG:HH12	1:B:331:GLU:CD	1.98	0.65
1:C:291:ARG:CZ	1:C:336:MET:HE3	2.26	0.65
1:C:72:ARG:NH1	1:C:331:GLU:OE2	2.28	0.65
1:D:261:MET:HE2	1:D:261:MET:HA	1.79	0.65
1:B:128:GLU:OE1	1:B:128:GLU:HA	1.96	0.65
1:A:311:GLN:NE2	5:A:1475:HOH:O	2.27	0.65
1:B:99:ILE:HD11	1:B:240:ARG:NH2	2.12	0.65
1:B:212:ARG:HA	1:B:225:ASN:HD21	1.62	0.64
1:C:358:LEU:HD12	3:C:417:HEM:HMD1	1.79	0.64
1:D:376:THR:CG2	1:D:414:VAL:HG21	2.27	0.64

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:87:TRP:CZ3	1:A:184:MET:O	2.50	0.64
1:B:303:SER:N	5:B:2509:HOH:O	2.30	0.64
1:B:294:LEU:HD23	1:B:294:LEU:H	1.62	0.64
1:A:143:ARG:HB3	1:A:144:PRO:CD	2.27	0.64
1:A:166:LEU:HD11	5:A:1581:HOH:O	1.96	0.64
1:B:390:GLN:HG3	1:B:400:GLN:HB2	1.80	0.64
1:D:150:PHE:CE2	1:D:155:ALA:HB2	2.33	0.64
1:D:98:PHE:HB3	1:D:244:LEU:HB2	1.80	0.63
1:B:244:LEU:O	1:B:247:LEU:O	2.17	0.63
1:A:277:ARG:HB2	1:A:280:ARG:HH21	1.64	0.63
1:B:390:GLN:HE21	1:B:400:GLN:HG3	1.64	0.63
1:B:151:THR:HA	1:B:155:ALA:HB3	1.81	0.63
1:B:110:GLN:CG	1:B:229:ASN:HB2	2.29	0.63
1:B:223:VAL:HG22	1:B:241:MET:HE2	1.81	0.63
1:A:181:THR:CG2	1:A:251:ASP:HB2	2.30	0.62
1:D:17:HIS:CD2	1:D:313:LYS:HB2	2.35	0.62
1:A:353:GLY:O	1:A:356:LEU:HD23	1.99	0.62
1:D:207:ILE:HG23	1:D:211:ARG:HH12	1.62	0.62
1:D:269:GLU:H	1:D:269:GLU:CD	2.03	0.62
1:C:107:GLU:CD	1:C:107:GLU:H	2.03	0.62
1:D:348:THR:HG22	1:D:352:HIS:HB2	1.82	0.61
1:B:98:PHE:HB3	1:B:244:LEU:HB2	1.82	0.61
1:A:282:PRO:O	1:A:285:CYS:HB3	1.99	0.61
1:D:223:VAL:HG22	1:D:241:MET:HE2	1.82	0.61
1:A:172:GLU:CB	1:D:272:GLN:HE22	2.13	0.61
1:A:15:PRO:HB2	1:A:18:VAL:HG23	1.81	0.61
1:D:69:GLN:NE2	1:D:70:LEU:HD12	2.16	0.61
1:D:273:GLU:HG2	1:D:280:ARG:NH1	2.16	0.61
1:C:256:PHE:CE1	1:C:367:ILE:HD13	2.35	0.61
1:D:87:TRP:HZ3	1:D:184:MET:O	1.84	0.61
1:A:281:ILE:HD12	1:A:372:LYS:HD2	1.83	0.60
1:D:158:PHE:HB3	1:D:159:PRO:HD3	1.83	0.60
1:A:98:PHE:HB3	1:A:244:LEU:HB2	1.81	0.60
1:B:38:VAL:HG12	1:B:391:HIS:ND1	2.16	0.60
1:A:155:ALA:O	1:A:159:PRO:HD2	2.02	0.60
1:C:109:ARG:HD3	5:C:3504:HOH:O	2.01	0.60
1:C:79:ARG:HG3	1:C:79:ARG:HH11	1.66	0.60
1:B:158:PHE:HB3	1:B:159:PRO:HD3	1.83	0.60
1:C:163:PHE:HE2	1:C:246:LEU:CD2	2.14	0.59
1:B:68:GLY:HA3	1:B:331:GLU:OE2	2.03	0.59
1:A:205:ILE:HB	1:A:206:PRO:HD3	1.84	0.59
1:B:363:ALA:O	1:B:367:ILE:HG13	2.03	0.59

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:273:GLU:HG3	1:A:280:ARG:NH2	2.17	0.58
1:B:72:ARG:HH11	1:B:72:ARG:HG3	1.68	0.58
1:A:181:THR:HG23	1:A:247:LEU:CD1	2.34	0.58
1:D:229:ASN:O	1:D:229:ASN:ND2	2.36	0.58
1:B:41:ALA:O	1:B:44:VAL:HG22	2.04	0.58
1:C:218:ASP:O	1:C:222:ILE:HG12	2.03	0.58
1:B:191:MET:HB2	1:B:195:GLU:HB2	1.86	0.58
1:C:18:VAL:HG11	1:C:55:TRP:CG	2.39	0.58
1:A:128:GLU:HA	1:A:128:GLU:OE2	2.03	0.58
1:D:87:TRP:CZ3	1:D:184:MET:O	2.57	0.57
1:A:151:THR:O	1:A:156:GLU:HG3	2.04	0.57
1:D:279:GLU:HG3	5:D:4488:HOH:O	2.03	0.57
1:D:267:SER:HB2	1:D:269:GLU:OE1	2.04	0.57
1:C:38:VAL:HG22	1:C:391:HIS:ND1	2.18	0.57
1:D:56:THR:O	1:D:61:GLY:HA2	2.03	0.57
1:D:17:HIS:HE2	1:D:313:LYS:HB3	1.69	0.57
1:C:268:PRO:O	1:C:271:ARG:HB2	2.05	0.57
1:D:186:ARG:HD2	1:D:392:LYS:HB3	1.86	0.57
1:D:41:ALA:O	1:D:44:VAL:HG22	2.05	0.57
1:B:14:LEU:HD11	1:B:18:VAL:CG1	2.35	0.57
1:A:33:ASN:HB3	1:A:41:ALA:HA	1.86	0.56
1:D:273:GLU:CG	1:D:280:ARG:NH1	2.67	0.56
1:C:294:LEU:HD23	1:C:294:LEU:H	1.69	0.56
1:D:130:ARG:HG2	5:D:4445:HOH:O	2.06	0.56
1:B:110:GLN:NE2	1:B:229:ASN:CB	2.67	0.56
1:B:294:LEU:N	1:B:294:LEU:HD23	2.20	0.56
1:A:142:LEU:HD22	1:A:148:CYS:HB3	1.87	0.56
1:C:282:PRO:O	1:C:285:CYS:HB3	2.05	0.56
1:B:353:GLY:O	1:B:356:LEU:HD22	2.05	0.56
1:A:122:PRO:O	1:A:126:LYS:HE3	2.06	0.56
1:B:99:ILE:CD1	1:B:240:ARG:HH21	2.16	0.56
1:B:99:ILE:HG23	1:B:100:PRO:HA	1.87	0.55
1:B:228:VAL:HG22	1:B:228:VAL:O	2.07	0.55
1:D:373:GLU:HA	1:D:373:GLU:OE2	2.06	0.55
1:D:110:GLN:HG3	1:D:229:ASN:HB2	1.89	0.55
1:D:68:GLY:HA3	1:D:331:GLU:OE2	2.05	0.55
1:C:98:PHE:HB3	1:C:244:LEU:HB2	1.87	0.55
1:B:269:GLU:CD	1:B:269:GLU:H	2.07	0.55
1:B:191:MET:CE	1:B:196:ALA:HA	2.26	0.55
1:A:206:PRO:O	1:A:210:GLN:CD	2.45	0.55
1:B:269:GLU:OE1	1:B:269:GLU:N	2.40	0.55
1:B:236:ASP:O	1:B:240:ARG:HG3	2.06	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:181:THR:HG21	1:A:251:ASP:HB2	1.87	0.55
1:A:172:GLU:CB	1:D:272:GLN:NE2	2.68	0.55
1:B:56:THR:HG21	1:B:62:HIS:NE2	2.21	0.55
1:C:181:THR:HG23	1:C:247:LEU:CD1	2.37	0.55
1:A:41:ALA:O	1:A:44:VAL:HG22	2.06	0.55
1:D:306:GLU:HG2	1:D:311:GLN:HG2	1.88	0.54
1:D:306:GLU:HG3	1:D:311:GLN:NE2	2.22	0.54
1:A:172:GLU:OE2	1:D:272:GLN:NE2	2.40	0.54
1:A:62:HIS:CD2	1:A:88:ILE:HD13	2.43	0.54
1:A:134:LEU:O	1:A:138:LEU:HB2	2.08	0.54
1:D:131:ILE:HG12	1:D:162:ILE:CD1	2.33	0.54
1:C:84:GLU:HG3	5:C:3503:HOH:O	2.08	0.54
1:C:15:PRO:HB2	1:C:18:VAL:HG23	1.88	0.54
1:A:39:GLN:NE2	1:A:39:GLN:H	2.05	0.54
1:B:306:GLU:HB2	1:B:311:GLN:NE2	2.23	0.54
1:C:379:PRO:HD3	1:C:412:LYS:HE2	1.90	0.54
1:D:110:GLN:HB3	1:D:228:VAL:CG2	2.32	0.53
1:D:231:ARG:HD2	1:D:231:ARG:O	2.07	0.53
1:B:244:LEU:HD11	4:B:2450:TCZ:H6	1.91	0.53
1:D:341:SER:O	1:D:342:ARG:C	2.47	0.53
1:B:170:PRO:HG2	1:B:173:ASP:OD1	2.08	0.53
1:B:37:GLY:HA3	1:B:40:GLU:OE2	2.09	0.53
1:B:414:VAL:CG1	1:B:414:VAL:O	2.54	0.53
1:D:40:GLU:O	1:D:43:ALA:HB3	2.07	0.53
1:C:56:THR:O	1:C:61:GLY:HA2	2.09	0.53
1:C:280:ARG:HA	5:C:3525:HOH:O	2.09	0.53
1:D:72:ARG:HH12	1:D:331:GLU:CD	2.12	0.53
1:A:18:VAL:HG13	1:A:63:TRP:CH2	2.44	0.53
1:D:90:ARG:HB2	1:D:317:GLN:OE1	2.08	0.53
1:A:118:VAL:HG23	1:A:119:VAL:HG13	1.91	0.53
1:C:83:SER:O	1:C:86:PRO:HD3	2.09	0.53
1:D:294:LEU:HD23	1:D:294:LEU:H	1.74	0.52
1:D:75:TYR:HB3	1:D:352:HIS:O	2.09	0.52
1:A:143:ARG:HB3	1:A:144:PRO:HD3	1.92	0.52
1:B:143:ARG:HH11	1:B:143:ARG:HG2	1.73	0.52
1:C:121:MET:HB3	1:C:122:PRO:HD3	1.90	0.52
1:D:38:VAL:HG12	1:D:391:HIS:ND1	2.23	0.52
1:D:17:HIS:NE2	1:D:313:LYS:HB2	2.25	0.52
1:D:197:LYS:HE3	1:D:198:GLU:OE2	2.10	0.52
1:B:118:VAL:CG2	1:B:219:ALA:HA	2.39	0.52
1:B:223:VAL:CG2	1:B:241:MET:HE2	2.40	0.52
1:D:17:HIS:NE2	1:D:313:LYS:CB	2.73	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:111:PHE:CD2	1:A:228:VAL:HG11	2.45	0.52
1:B:377:ARG:HH11	1:B:377:ARG:HG3	1.75	0.52
1:C:39:GLN:H	1:C:39:GLN:NE2	2.08	0.52
1:C:143:ARG:HB3	1:C:144:PRO:HD3	1.90	0.52
1:A:110:GLN:NE2	1:A:229:ASN:OD1	2.42	0.52
1:A:218:ASP:O	1:A:222:ILE:HG12	2.10	0.51
1:B:251:ASP:HB2	5:B:2452:HOH:O	2.09	0.51
1:A:172:GLU:HB2	1:D:272:GLN:HE22	1.73	0.51
1:C:229:ASN:O	1:C:229:ASN:OD1	2.28	0.51
1:B:56:THR:HG22	1:B:62:HIS:O	2.09	0.51
1:C:229:ASN:O	1:C:229:ASN:CG	2.49	0.51
1:B:376:THR:CG2	1:B:414:VAL:HG21	2.40	0.51
1:B:161:ARG:NH1	5:B:2494:HOH:O	2.43	0.51
1:A:69:GLN:O	1:A:73:GLU:HG3	2.11	0.51
1:D:105:PRO:HG3	1:D:108:GLN:NE2	2.24	0.51
1:C:146:GLY:HA2	1:C:406:TRP:NE1	2.25	0.51
1:A:183:GLN:HG3	1:A:191:MET:HG2	1.92	0.51
1:D:228:VAL:O	1:D:229:ASN:HB3	2.11	0.51
1:B:200:LEU:O	1:B:203:TYR:HB3	2.11	0.50
1:A:284:ALA:O	1:A:288:LEU:HG	2.11	0.50
1:D:251:ASP:OD2	1:D:255:ASN:ND2	2.40	0.50
1:D:201:TYR:CE2	1:D:239:LYS:HE3	2.46	0.50
1:B:146:GLY:HA2	1:B:406:TRP:CD1	2.46	0.50
1:D:151:THR:HA	1:D:155:ALA:HB3	1.93	0.50
1:B:365:ARG:HD3	1:B:369:VAL:HG23	1.94	0.50
1:C:129:ASN:H	1:C:129:ASN:ND2	2.05	0.50
1:A:388:GLN:NE2	5:A:1578:HOH:O	2.44	0.50
1:C:318:ILE:O	1:C:318:ILE:HG23	2.12	0.50
1:A:209:GLU:HB3	1:A:210:GLN:NE2	2.24	0.50
1:B:72:ARG:NH1	1:B:72:ARG:HG3	2.27	0.50
1:D:205:ILE:HB	1:D:206:PRO:HD3	1.93	0.50
1:B:157:PRO:O	1:B:161:ARG:HG3	2.12	0.50
1:B:179:TYR:CE1	1:B:183:GLN:NE2	2.80	0.49
1:B:191:MET:HE2	1:B:196:ALA:CA	2.30	0.49
1:D:184:MET:HG2	1:D:193:PHE:CE1	2.48	0.49
1:B:119:VAL:HG11	1:B:245:LEU:HD22	1.94	0.49
1:C:291:ARG:CZ	1:C:336:MET:CE	2.89	0.49
1:D:356:LEU:HG	5:D:4455:HOH:O	2.11	0.49
1:A:186:ARG:HH22	1:A:251:ASP:CG	2.15	0.49
1:D:160:ILE:O	1:D:164:MET:HG2	2.13	0.49
1:D:297:ASP:OD2	3:D:417:HEM:O2A	2.30	0.49
1:A:157:PRO:O	1:A:161:ARG:HB2	2.12	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:218:ASP:OD1	1:A:220:ILE:HB	2.12	0.49
1:B:179:TYR:OH	1:B:190:SER:CB	2.61	0.49
1:B:150:PHE:CZ	1:B:261:MET:HG3	2.47	0.49
1:B:293:SER:O	1:B:323:MET:HG3	2.12	0.49
1:D:318:ILE:HG23	1:D:318:ILE:O	2.13	0.49
1:B:72:ARG:NH1	1:B:331:GLU:OE2	2.36	0.49
1:D:105:PRO:HG3	1:D:108:GLN:HE22	1.78	0.49
1:D:201:TYR:CD2	1:D:239:LYS:HE3	2.48	0.49
1:A:377:ARG:HH11	1:A:377:ARG:HG3	1.78	0.49
1:B:39:GLN:NE2	1:B:39:GLN:H	2.11	0.49
1:D:178:LYS:HE3	1:D:182:ASP:OD1	2.13	0.49
1:D:17:HIS:HE2	1:D:313:LYS:CB	2.25	0.49
1:D:99:ILE:HD11	1:D:240:ARG:CZ	2.43	0.48
1:A:253:VAL:O	1:A:254:VAL:C	2.51	0.48
1:B:228:VAL:CG2	1:B:228:VAL:O	2.61	0.48
1:B:231:ARG:O	1:B:231:ARG:HD2	2.12	0.48
1:C:109:ARG:HH11	1:C:109:ARG:HG3	1.77	0.48
1:B:286:GLU:OE2	1:B:364:ARG:NE	2.35	0.48
1:A:107:GLU:H	1:A:107:GLU:CD	2.16	0.48
1:C:181:THR:CG2	1:C:251:ASP:HB2	2.42	0.48
1:B:273:GLU:OE1	1:B:280:ARG:NH2	2.45	0.48
1:D:163:PHE:CE2	1:D:246:LEU:HD23	2.34	0.48
1:A:272:GLN:NE2	1:A:276:GLU:HG3	2.29	0.48
1:B:114:LEU:O	1:B:117:GLN:HB3	2.12	0.48
1:A:210:GLN:H	1:A:210:GLN:NE2	2.00	0.48
1:D:211:ARG:HG3	1:D:221:SER:OG	2.14	0.48
1:B:218:ASP:O	1:B:222:ILE:HG12	2.14	0.48
1:A:277:ARG:CA	1:A:277:ARG:HE	2.24	0.48
1:C:294:LEU:HD23	1:C:294:LEU:N	2.28	0.48
1:A:142:LEU:HD22	1:A:148:CYS:CB	2.42	0.48
1:A:355:HIS:HD1	3:A:417:HEM:CGD	2.27	0.48
1:B:134:LEU:HG	1:B:138:LEU:HD22	1.96	0.48
1:C:273:GLU:HG3	1:C:280:ARG:HH21	1.79	0.48
1:D:149:ASN:ND2	1:D:402:LEU:H	2.12	0.48
1:A:132:GLN:HB3	1:A:132:GLN:HE21	1.39	0.48
1:D:91:GLU:OE1	5:D:4479:HOH:O	2.20	0.47
1:B:78:TYR:CE1	1:B:105:PRO:HB2	2.49	0.47
1:A:52:ASP:HB3	1:A:70:LEU:HD22	1.97	0.47
1:D:294:LEU:HD23	1:D:294:LEU:N	2.28	0.47
1:B:11:LEU:HG	1:B:25:ASP:OD1	2.13	0.47
1:C:334:ALA:N	1:C:335:PRO:CD	2.77	0.47
1:C:156:GLU:OE1	5:C:3477:HOH:O	2.20	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:118:VAL:HG23	1:C:119:VAL:HG13	1.95	0.47
1:A:218:ASP:OD2	1:A:221:SER:OG	2.23	0.47
1:D:157:PRO:O	1:D:161:ARG:HG3	2.15	0.47
1:B:373:GLU:OE2	1:B:373:GLU:HA	2.15	0.47
1:C:239:LYS:HE2	1:C:239:LYS:HB3	1.60	0.47
1:D:72:ARG:NH1	1:D:72:ARG:CG	2.73	0.47
1:D:244:LEU:O	1:D:247:LEU:O	2.33	0.47
1:B:272:GLN:O	1:B:276:GLU:HG3	2.14	0.47
1:D:171:GLU:O	1:D:174:ILE:HD12	2.14	0.47
1:B:103:MET:O	1:B:355:HIS:CE1	2.68	0.47
1:A:132:GLN:OE1	1:A:373:GLU:OE1	2.33	0.47
1:C:158:PHE:HB3	1:C:159:PRO:HD3	1.97	0.47
1:B:256:PHE:CE2	1:B:289:LEU:HA	2.49	0.47
1:A:125:ASP:O	1:A:128:GLU:HB2	2.15	0.47
1:D:325:SER:HB2	1:D:348:THR:OG1	2.14	0.47
1:A:73:GLU:CD	1:A:308:HIS:HE2	2.17	0.47
1:D:161:ARG:NH1	5:D:4440:HOH:O	2.47	0.47
1:B:90:ARG:HB2	1:B:317:GLN:OE1	2.14	0.46
1:C:129:ASN:ND2	1:C:129:ASN:N	2.53	0.46
1:A:87:TRP:CZ2	1:A:395:ILE:HG21	2.50	0.46
1:D:306:GLU:CG	1:D:311:GLN:HG2	2.45	0.46
1:C:155:ALA:O	1:C:159:PRO:CD	2.63	0.46
1:A:213:GLN:C	1:A:215:PRO:HD3	2.35	0.46
1:D:304:ASP:N	1:D:314:LYS:HB2	2.30	0.46
1:A:122:PRO:O	1:A:126:LYS:HG2	2.16	0.46
1:B:179:TYR:OH	1:B:190:SER:HB3	2.16	0.46
1:D:323:MET:O	1:D:323:MET:HG2	2.15	0.46
1:C:127:LEU:O	1:C:131:ILE:HG13	2.16	0.46
1:A:281:ILE:CD1	1:A:372:LYS:HD2	2.46	0.46
1:B:39:GLN:HE21	1:B:39:GLN:H	1.62	0.46
1:B:70:LEU:HD11	1:B:308:HIS:CE1	2.51	0.46
1:A:228:VAL:O	1:A:229:ASN:OD1	2.33	0.46
1:A:56:THR:O	1:A:61:GLY:HA2	2.16	0.46
1:C:379:PRO:CD	1:C:412:LYS:HE2	2.46	0.46
1:D:273:GLU:CD	1:D:280:ARG:HH12	2.18	0.46
1:B:40:GLU:O	1:B:43:ALA:HB3	2.16	0.46
1:D:297:ASP:O	1:D:319:LEU:HD12	2.16	0.46
1:B:47:GLU:HG2	1:B:49:ASN:HD21	1.80	0.46
1:A:277:ARG:CG	1:A:280:ARG:NH2	2.79	0.46
1:B:139:ILE:HD13	1:B:374:TRP:HA	1.97	0.46
1:D:24:PHE:O	1:D:26:PHE:N	2.47	0.46
1:C:264:LEU:HA	1:C:264:LEU:HD23	1.78	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:114:LEU:O	1:D:117:GLN:HB3	2.16	0.46
1:C:91:GLU:CD	1:C:91:GLU:H	2.19	0.46
1:B:201:TYR:HB3	1:B:239:LYS:HD2	1.97	0.45
1:B:229:ASN:C	1:B:229:ASN:HD22	2.20	0.45
1:C:109:ARG:HH11	1:C:109:ARG:CG	2.29	0.45
1:D:121:MET:HE1	1:D:361:HIS:HB3	1.98	0.45
1:C:181:THR:HG21	1:C:251:ASP:HB2	1.99	0.45
1:D:212:ARG:HA	1:D:225:ASN:HD21	1.81	0.45
1:C:207:ILE:O	1:C:211:ARG:HB2	2.16	0.45
1:C:143:ARG:HB3	1:C:144:PRO:CD	2.46	0.45
1:A:214:LYS:N	1:A:215:PRO:HD3	2.32	0.45
1:D:149:ASN:HD21	1:D:402:LEU:H	1.64	0.45
1:A:279:GLU:N	1:A:279:GLU:OE2	2.38	0.45
1:C:293:SER:O	1:C:323:MET:HG3	2.17	0.45
1:A:27:ASP:OD1	1:A:27:ASP:C	2.55	0.45
1:D:59:ASN:ND2	5:D:4508:HOH:O	2.32	0.45
1:C:226:GLY:O	1:C:233:ILE:HG22	2.17	0.44
1:D:286:GLU:OE2	1:D:364:ARG:NE	2.28	0.44
1:C:286:GLU:OE2	1:C:364:ARG:NE	2.31	0.44
1:A:135:ALA:O	1:A:139:ILE:HG13	2.17	0.44
1:D:110:GLN:CG	1:D:229:ASN:HB2	2.47	0.44
1:B:201:TYR:CG	1:B:239:LYS:HD2	2.53	0.44
1:B:85:CYS:HB3	1:B:300:ILE:HB	2.00	0.44
1:C:183:GLN:HG3	1:C:191:MET:HG2	1.98	0.44
1:B:28:MET:HG3	1:B:29:TYR:CD1	2.53	0.44
1:A:172:GLU:HB3	1:D:272:GLN:HE22	1.75	0.44
1:B:294:LEU:N	1:B:294:LEU:CD2	2.81	0.44
1:A:358:LEU:HD12	3:A:417:HEM:HMD3	1.99	0.44
1:B:105:PRO:HG3	1:B:108:GLN:HE22	1.83	0.44
1:C:99:ILE:HA	1:C:100:PRO:HA	1.72	0.44
1:C:322:GLN:HB3	1:C:348:THR:O	2.17	0.44
1:D:18:VAL:HG21	1:D:55:TRP:CD2	2.53	0.44
1:A:344:LYS:HA	1:A:344:LYS:HD2	1.68	0.43
1:B:131:ILE:HA	1:B:162:ILE:CD1	2.43	0.43
1:A:168:GLY:O	1:A:211:ARG:NH2	2.44	0.43
1:D:108:GLN:HB3	1:D:108:GLN:HE21	1.59	0.43
1:B:84:GLU:HG2	1:B:84:GLU:O	2.16	0.43
1:B:118:VAL:HG13	1:B:119:VAL:HG13	2.00	0.43
1:D:339:ASP:OD2	1:D:341:SER:HB2	2.18	0.43
1:A:155:ALA:O	1:A:159:PRO:CD	2.66	0.43
1:C:197:LYS:HE2	1:C:197:LYS:HB3	1.83	0.43
1:A:83:SER:O	1:A:86:PRO:HD3	2.18	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:68:GLY:HA3	1:A:331:GLU:OE2	2.18	0.43
1:A:75:TYR:CZ	1:A:320:LEU:HB2	2.53	0.43
1:C:170:PRO:HB2	1:C:172:GLU:HG2	2.01	0.43
1:C:53:LEU:HD23	1:C:310:VAL:HG21	2.00	0.43
1:B:191:MET:HE3	1:B:195:GLU:C	2.39	0.43
1:B:11:LEU:O	1:B:12:ALA:C	2.57	0.43
1:A:146:GLY:HA2	1:A:406:TRP:NE1	2.34	0.43
1:A:277:ARG:CB	1:A:280:ARG:HH21	2.31	0.43
1:B:355:HIS:O	1:B:356:LEU:C	2.57	0.43
1:B:108:GLN:HB3	1:B:108:GLN:HE21	1.51	0.43
1:C:350:PHE:HB3	1:C:357:CYS:HB3	1.99	0.43
1:A:313:LYS:O	1:A:314:LYS:C	2.57	0.43
1:B:83:SER:O	1:B:86:PRO:HD3	2.18	0.43
1:B:56:THR:CG2	1:B:62:HIS:CE1	2.99	0.42
1:C:135:ALA:HB2	1:C:158:PHE:CE1	2.54	0.42
1:B:140:GLU:HA	1:B:140:GLU:OE2	2.18	0.42
1:D:164:MET:HE3	1:D:170:PRO:O	2.18	0.42
1:B:251:ASP:CG	1:B:251:ASP:O	2.56	0.42
1:C:172:GLU:HG3	1:C:173:ASP:N	2.35	0.42
1:C:130:ARG:NH2	1:C:133:GLU:OE1	2.52	0.42
1:D:179:TYR:O	1:D:183:GLN:HG2	2.18	0.42
1:C:322:GLN:HG2	1:C:351:GLY:HA2	2.02	0.42
1:B:390:GLN:HE21	1:B:400:GLN:CG	2.31	0.42
1:B:377:ARG:HG3	1:B:377:ARG:NH1	2.33	0.42
1:A:376:THR:HG22	1:A:377:ARG:HG3	2.01	0.42
1:C:151:THR:O	1:C:156:GLU:HG3	2.20	0.42
1:D:313:LYS:O	1:D:314:LYS:C	2.58	0.42
1:A:350:PHE:HB3	1:A:357:CYS:HB3	2.01	0.42
1:A:163:PHE:CE2	1:A:246:LEU:HD22	2.55	0.42
3:D:417:HEM:HBB2	3:D:417:HEM:HMB1	2.01	0.42
1:D:100:PRO:HB3	1:D:111:PHE:HB2	2.02	0.42
1:B:267:SER:HB2	1:B:269:GLU:OE1	2.20	0.42
1:A:362:LEU:O	1:A:366:GLU:HG3	2.20	0.42
1:D:344:LYS:HA	1:D:344:LYS:HD2	1.75	0.42
3:B:417:HEM:HMB1	3:B:417:HEM:HBB2	2.00	0.42
1:C:294:LEU:CD2	1:C:294:LEU:N	2.83	0.42
1:D:317:GLN:HG2	5:D:4494:HOH:O	2.20	0.42
1:A:38:VAL:HG22	1:A:391:HIS:HB3	2.01	0.42
1:B:303:SER:HA	1:B:314:LYS:HD2	2.02	0.42
1:B:212:ARG:HA	1:B:225:ASN:ND2	2.31	0.42
1:C:215:PRO:HB3	1:C:225:ASN:OD1	2.19	0.41
1:C:99:ILE:O	1:C:241:MET:HA	2.20	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:33:ASN:ND2	5:A:1586:HOH:O	2.49	0.41
1:D:55:TRP:C	1:D:55:TRP:CD1	2.93	0.41
1:D:251:ASP:O	1:D:396:VAL:HG11	2.20	0.41
1:B:297:ASP:OD2	3:B:417:HEM:O2A	2.38	0.41
1:B:290:ARG:HD2	1:B:332:ASN:ND2	2.35	0.41
1:D:163:PHE:HE2	1:D:246:LEU:CD2	2.24	0.41
1:A:170:PRO:HB2	1:A:172:GLU:HG2	2.01	0.41
1:B:149:ASN:C	1:B:149:ASN:HD22	2.23	0.41
1:A:288:LEU:HD23	1:A:288:LEU:HA	1.83	0.41
1:A:146:GLY:HA2	1:A:406:TRP:CD1	2.55	0.41
1:D:85:CYS:SG	1:D:93:GLY:HA3	2.60	0.41
1:B:149:ASN:C	1:B:149:ASN:ND2	2.74	0.41
1:A:205:ILE:CB	1:A:206:PRO:HD3	2.50	0.41
1:B:118:VAL:CG2	5:B:2471:HOH:O	2.56	0.41
1:B:239:LYS:HB3	1:B:239:LYS:HE2	1.78	0.41
1:A:321:PRO:O	1:A:322:GLN:C	2.58	0.41
1:C:73:GLU:CD	1:C:308:HIS:HE2	2.22	0.41
1:A:10:ASN:CG	1:A:11:LEU:H	2.24	0.41
1:B:267:SER:C	1:B:269:GLU:OE1	2.59	0.41
1:C:194:ALA:O	1:C:198:GLU:HG2	2.21	0.41
1:D:146:GLY:HA2	1:D:406:TRP:CD1	2.56	0.41
1:D:17:HIS:NE2	1:D:313:LYS:HB3	2.33	0.41
1:C:201:TYR:HB3	1:C:239:LYS:HD2	2.02	0.41
1:C:10:ASN:N	1:C:10:ASN:OD1	2.54	0.41
1:D:363:ALA:O	1:D:367:ILE:HG13	2.21	0.41
1:B:407:ASP:OD1	1:B:409:ALA:HB3	2.21	0.41
1:A:333:ALA:O	1:A:334:ALA:C	2.59	0.41
1:B:99:ILE:CG2	1:B:100:PRO:HA	2.51	0.41
1:A:185:THR:HG21	1:A:251:ASP:OD1	2.21	0.41
1:B:125:ASP:O	1:B:128:GLU:HB2	2.21	0.41
1:A:75:TYR:HB3	1:A:352:HIS:O	2.21	0.41
1:D:218:ASP:O	1:D:222:ILE:HG12	2.21	0.41
1:C:267:SER:HB3	1:C:269:GLU:OE1	2.21	0.41
1:B:127:LEU:O	1:B:128:GLU:C	2.60	0.40
1:D:134:LEU:HG	1:D:138:LEU:HD22	2.03	0.40
1:D:110:GLN:CB	1:D:228:VAL:HG23	2.33	0.40
1:B:99:ILE:HD11	1:B:240:ARG:HE	1.85	0.40
1:D:23:VAL:HA	1:D:55:TRP:O	2.21	0.40
1:B:303:SER:HA	1:B:314:LYS:CD	2.51	0.40
1:C:294:LEU:HD12	1:C:397:SER:O	2.21	0.40
1:A:120:GLY:O	1:A:123:VAL:N	2.54	0.40
1:A:201:TYR:HB3	1:A:239:LYS:HD2	2.02	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:188:ASP:C	1:C:188:ASP:OD2	2.60	0.40
1:B:87:TRP:CZ2	1:B:395:ILE:HG21	2.57	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	403/414 (97%)	378 (94%)	24 (6%)	1 (0%)	56	62
1	B	403/414 (97%)	379 (94%)	21 (5%)	3 (1%)	30	28
1	C	403/414 (97%)	383 (95%)	20 (5%)	0	100	100
1	D	403/414 (97%)	370 (92%)	30 (7%)	3 (1%)	30	28
All	All	1612/1656 (97%)	1510 (94%)	95 (6%)	7 (0%)	43	45

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	25	ASP
1	B	25	ASP
1	D	51	PRO
1	B	28	MET
1	D	128	GLU
1	A	321	PRO
1	B	128	GLU

### 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	349/357 (98%)	330 (95%)	19 (5%)	31	35
1	B	349/357 (98%)	330 (95%)	19 (5%)	31	35
1	C	349/357 (98%)	329 (94%)	20 (6%)	29	32
1	D	349/357 (98%)	329 (94%)	20 (6%)	29	32
All	All	1396/1428 (98%)	1318 (94%)	78 (6%)	30	33

All (78) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	LEU
1	A	38	VAL
1	A	39	GLN
1	A	79	ARG
1	A	86	PRO
1	A	108	GLN
1	A	132	GLN
1	A	138	LEU
1	A	198	GLU
1	A	210	GLN
1	A	211	ARG
1	A	228	VAL
1	A	246	LEU
1	A	269	GLU
1	A	277	ARG
1	A	313	LYS
1	A	365	ARG
1	A	376	THR
1	A	402	LEU
1	B	10	ASN
1	B	39	GLN
1	B	47	GLU
1	B	69	GLN
1	B	100	PRO
1	B	108	GLN
1	B	118	VAL
1	B	128	GLU
1	B	138	LEU
1	B	149	ASN
1	B	181	THR
1	B	205	ILE
1	B	229	ASN
1	B	269	GLU

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Mol	Chain	Res	Type
1	B	329	GLU
1	B	342	ARG
1	B	356	LEU
1	B	380	ASP
1	B	412	LYS
1	C	11	LEU
1	C	32	SER
1	C	38	VAL
1	C	39	GLN
1	C	44	VAL
1	C	79	ARG
1	C	84	GLU
1	C	86	PRO
1	C	108	GLN
1	C	109	ARG
1	C	128	GLU
1	C	129	ASN
1	C	138	LEU
1	C	152	GLU
1	C	210	GLN
1	C	217	THR
1	C	229	ASN
1	C	277	ARG
1	C	376	THR
1	C	402	LEU
1	D	17	HIS
1	D	20	GLU
1	D	39	GLN
1	D	72	ARG
1	D	86	PRO
1	D	108	GLN
1	D	118	VAL
1	D	138	LEU
1	D	149	ASN
1	D	181	THR
1	D	209	GLU
1	D	211	ARG
1	D	213	GLN
1	D	228	VAL
1	D	229	ASN
1	D	269	GLU
1	D	343	GLN

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Mol	Chain	Res	Type
1	D	346	SER
1	D	348	THR
1	D	356	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (46) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	21	HIS
1	A	39	GLN
1	A	46	GLN
1	A	108	GLN
1	A	129	ASN
1	A	132	GLN
1	A	145	GLN
1	A	210	GLN
1	A	213	GLN
1	A	272	GLN
1	A	388	GLN
1	B	39	GLN
1	B	46	GLN
1	B	49	ASN
1	B	69	GLN
1	B	108	GLN
1	B	117	GLN
1	B	132	GLN
1	B	149	ASN
1	B	210	GLN
1	B	225	ASN
1	B	229	ASN
1	B	308	HIS
1	B	311	GLN
1	B	390	GLN
1	B	400	GLN
1	C	10	ASN
1	C	30	ASN
1	C	39	GLN
1	C	46	GLN
1	C	108	GLN
1	C	110	GLN
1	C	129	ASN
1	C	132	GLN
1	C	145	GLN

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Mol	Chain	Res	Type
1	C	213	GLN
1	C	225	ASN
1	C	272	GLN
1	C	311	GLN
1	D	39	GLN
1	D	46	GLN
1	D	108	GLN
1	D	117	GLN
1	D	132	GLN
1	D	149	ASN
1	D	225	ASN

### 5.3.3 RNA ⓘ

There are no RNA chains 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 11 ligands modelled in this entry, 4 are monoatomic - leaving 7 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	$\# Z  > 2$
4	TCZ	A	1450	-	9,9,9	2.52	7 (77%)	12,12,12	0.95	1 (8%)
3	HEM	A	417	1	49,50,50	3.24	24 (48%)	46,82,82	2.59	18 (39%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
4	TCZ	B	2450	-	9,9,9	2.61	6 (66%)	12,12,12	1.09	0
3	HEM	B	417	1	49,50,50	5.29	24 (48%)	46,82,82	2.39	17 (36%)
4	TCZ	C	3450	-	9,9,9	2.58	6 (66%)	12,12,12	1.00	0
3	HEM	C	417	1	49,50,50	3.98	22 (44%)	46,82,82	2.48	19 (41%)
3	HEM	D	417	1,5	49,50,50	4.73	22 (44%)	46,82,82	2.37	16 (34%)

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
4	TCZ	A	1450	-	-	0/0/0/0	0/1/1/1
3	HEM	A	417	1	-	0/14/114/114	0/0/8/8
4	TCZ	B	2450	-	-	0/0/0/0	0/1/1/1
3	HEM	B	417	1	-	0/14/114/114	0/0/8/8
4	TCZ	C	3450	-	-	0/0/0/0	0/1/1/1
3	HEM	C	417	1	-	0/14/114/114	0/0/8/8
3	HEM	D	417	1,5	-	0/14/114/114	0/0/8/8

All (111) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	417	HEM	C2D-C1D	20.17	1.49	1.44
3	D	417	HEM	C2D-C1D	19.88	1.49	1.44
3	B	417	HEM	C3D-C4D	18.49	1.49	1.44
3	C	417	HEM	C2D-C1D	15.21	1.48	1.44
3	B	417	HEM	C2B-C1B	15.10	1.48	1.44
3	D	417	HEM	C2B-C1B	13.69	1.48	1.44
3	A	417	HEM	C2D-C1D	12.48	1.47	1.44
3	D	417	HEM	C3D-C4D	12.32	1.47	1.44
3	C	417	HEM	C2B-C1B	12.20	1.47	1.44
3	B	417	HEM	C3B-C2B	-6.90	1.31	1.43
3	D	417	HEM	C3B-C2B	-6.73	1.32	1.43
3	B	417	HEM	C4A-C3A	6.61	1.48	1.40
3	C	417	HEM	C3B-C2B	-6.60	1.32	1.43
3	B	417	HEM	C3C-C2C	-6.54	1.32	1.43
3	C	417	HEM	C4A-C3A	6.54	1.48	1.40
3	D	417	HEM	C4A-C3A	6.36	1.48	1.40
3	C	417	HEM	C3D-C4D	6.24	1.46	1.44
3	A	417	HEM	C4A-C3A	6.14	1.47	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	417	HEM	C3C-CAC	6.12	1.59	1.40
3	D	417	HEM	C3C-C2C	-6.06	1.33	1.43
3	B	417	HEM	C3C-CAC	5.99	1.59	1.40
3	A	417	HEM	C3C-C2C	-5.85	1.33	1.43
3	A	417	HEM	C3D-C2D	-5.62	1.33	1.43
3	D	417	HEM	C3C-CAC	5.61	1.58	1.40
3	C	417	HEM	C3C-C2C	-5.46	1.34	1.43
3	C	417	HEM	C3D-C2D	-5.43	1.34	1.43
3	D	417	HEM	C3B-CAB	5.42	1.57	1.40
3	A	417	HEM	C3B-C2B	-5.42	1.34	1.43
3	A	417	HEM	C3C-CAC	5.21	1.56	1.40
3	B	417	HEM	C3B-CAB	5.10	1.56	1.40
3	C	417	HEM	C3B-CAB	5.05	1.56	1.40
3	D	417	HEM	C3D-C2D	-4.83	1.35	1.43
3	A	417	HEM	C3B-CAB	4.76	1.55	1.40
3	D	417	HEM	CBD-CGD	4.63	1.62	1.50
3	B	417	HEM	C3B-C4B	4.45	1.49	1.44
3	D	417	HEM	C3B-C4B	4.38	1.49	1.44
3	B	417	HEM	CAA-C2A	4.33	1.59	1.52
3	B	417	HEM	C3D-C2D	-4.03	1.36	1.43
3	C	417	HEM	C3B-C4B	4.00	1.49	1.44
3	C	417	HEM	CAA-C2A	3.96	1.59	1.52
3	A	417	HEM	C2B-C1B	3.88	1.45	1.44
3	B	417	HEM	CBD-CGD	3.84	1.60	1.50
3	A	417	HEM	C3B-C4B	3.84	1.49	1.44
3	C	417	HEM	CBD-CGD	3.80	1.60	1.50
3	D	417	HEM	CHD-C4C	3.77	1.43	1.36
3	B	417	HEM	CHA-C4D	3.66	1.41	1.35
3	A	417	HEM	CBD-CGD	3.53	1.59	1.50
3	C	417	HEM	CBA-CGA	3.49	1.59	1.50
4	C	3450	TCZ	C2-C3	3.41	1.44	1.38
4	A	1450	TCZ	C4-C3	3.40	1.44	1.38
4	B	2450	TCZ	C4-C3	3.39	1.44	1.38
4	B	2450	TCZ	C4-C5	3.37	1.44	1.38
3	B	417	HEM	CBA-CGA	3.35	1.59	1.50
3	C	417	HEM	CMC-C2C	3.31	1.57	1.47
3	B	417	HEM	CHD-C4C	3.28	1.42	1.36
3	B	417	HEM	CMD-C2D	3.28	1.57	1.47
4	B	2450	TCZ	C2-C1	3.27	1.44	1.38
3	A	417	HEM	CMC-C2C	3.26	1.57	1.47
4	C	3450	TCZ	C4-C5	3.19	1.43	1.38
3	A	417	HEM	CMD-C2D	3.18	1.57	1.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	417	HEM	FE-ND	3.17	2.09	1.97
4	A	1450	TCZ	C4-C5	3.16	1.43	1.38
3	B	417	HEM	FE-NA	3.12	2.05	1.92
4	C	3450	TCZ	C2-C1	3.11	1.43	1.38
3	B	417	HEM	FE-ND	3.09	2.09	1.97
3	D	417	HEM	CBA-CGA	3.08	1.58	1.50
3	A	417	HEM	CBA-CGA	3.07	1.58	1.50
3	A	417	HEM	CAA-C2A	3.07	1.57	1.52
3	D	417	HEM	CAA-C2A	3.05	1.57	1.52
3	C	417	HEM	CMB-C2B	3.03	1.56	1.47
3	A	417	HEM	CMB-C2B	2.97	1.56	1.47
3	D	417	HEM	FE-ND	2.94	2.08	1.97
3	A	417	HEM	FE-NA	2.94	2.05	1.92
4	A	1450	TCZ	C1-CL1	-2.89	1.67	1.74
3	B	417	HEM	CMC-C2C	2.85	1.56	1.47
4	B	2450	TCZ	C1-CL1	-2.83	1.67	1.74
3	D	417	HEM	CMD-C2D	2.79	1.56	1.47
3	B	417	HEM	C4A-NA	2.70	1.41	1.36
3	D	417	HEM	CMB-C2B	2.69	1.55	1.47
3	D	417	HEM	CMC-C2C	2.69	1.55	1.47
4	C	3450	TCZ	C6-C1	2.69	1.43	1.38
4	A	1450	TCZ	C2-C3	2.67	1.43	1.38
4	C	3450	TCZ	C4-C3	2.66	1.43	1.38
4	C	3450	TCZ	C1-CL1	-2.66	1.68	1.74
3	C	417	HEM	FE-ND	2.66	2.07	1.97
3	A	417	HEM	CHA-C4D	2.64	1.39	1.35
3	C	417	HEM	CMD-C2D	2.63	1.55	1.47
4	B	2450	TCZ	C2-C3	2.62	1.42	1.38
4	A	1450	TCZ	C6-C1	2.59	1.42	1.38
4	B	2450	TCZ	C6-C1	2.58	1.42	1.38
3	C	417	HEM	O2A-CGA	-2.57	1.21	1.30
3	B	417	HEM	FE-NB	2.55	2.07	1.97
3	D	417	HEM	FE-NA	2.54	2.03	1.92
3	A	417	HEM	CHB-C1B	2.51	1.39	1.35
3	B	417	HEM	CMB-C2B	2.44	1.55	1.47
3	D	417	HEM	CHB-C1B	2.42	1.39	1.35
3	C	417	HEM	FE-NA	2.40	2.02	1.92
3	D	417	HEM	C4A-NA	2.39	1.41	1.36
3	A	417	HEM	C3D-C4D	2.33	1.45	1.44
3	A	417	HEM	CHD-C4C	2.33	1.40	1.36
3	C	417	HEM	CHA-C4D	2.30	1.39	1.35
3	C	417	HEM	FE-NC	2.25	2.06	1.97

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1450	TCZ	C2-C1	2.24	1.42	1.38
3	B	417	HEM	O2D-CGD	-2.11	1.22	1.30
3	A	417	HEM	C1C-NC	2.08	1.41	1.38
4	A	1450	TCZ	C3-CL3	-2.07	1.69	1.74
3	B	417	HEM	CAD-C3D	2.05	1.58	1.51
3	A	417	HEM	C4A-NA	2.05	1.40	1.36
3	D	417	HEM	FE-NB	2.05	2.05	1.97
3	A	417	HEM	CHC-C1C	2.03	1.39	1.36
3	C	417	HEM	CHD-C4C	2.02	1.39	1.36

All (71) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	417	HEM	C3B-C4B-NB	-9.28	107.36	114.00
3	C	417	HEM	C3B-C4B-NB	-7.83	108.40	114.00
3	D	417	HEM	C3B-C4B-NB	-7.63	108.54	114.00
3	B	417	HEM	C3B-C4B-NB	-7.25	108.81	114.00
3	B	417	HEM	C4D-ND-C1D	5.29	110.58	105.16
3	A	417	HEM	C4D-ND-C1D	4.80	110.08	105.16
3	A	417	HEM	O2D-CGD-O1D	4.79	135.47	123.30
3	C	417	HEM	C4D-ND-C1D	4.78	110.05	105.16
3	D	417	HEM	C2D-C1D-ND	-4.70	107.38	112.93
3	D	417	HEM	C4D-ND-C1D	4.70	109.97	105.16
3	C	417	HEM	C2D-C1D-ND	-4.64	107.45	112.93
3	B	417	HEM	C2D-C1D-ND	-4.61	107.48	112.93
3	A	417	HEM	C2D-C1D-ND	-4.52	107.59	112.93
3	B	417	HEM	O2D-CGD-O1D	4.50	134.75	123.30
3	C	417	HEM	O2D-CGD-O1D	4.37	134.42	123.30
3	C	417	HEM	CHD-C1D-ND	4.35	128.20	124.58
3	D	417	HEM	O2D-CGD-O1D	4.28	134.19	123.30
3	B	417	HEM	C4A-CHB-C1B	-4.01	122.19	127.47
3	A	417	HEM	O2A-CGA-O1A	4.00	133.47	123.30
3	A	417	HEM	CHD-C1D-ND	3.96	127.88	124.58
3	A	417	HEM	C4A-CHB-C1B	-3.90	122.33	127.47
3	D	417	HEM	O2A-CGA-O1A	3.83	133.05	123.30
3	D	417	HEM	CHD-C1D-ND	3.83	127.77	124.58
3	B	417	HEM	CMA-C3A-C4A	-3.81	122.76	128.62
3	B	417	HEM	CHD-C1D-ND	3.78	127.73	124.58
3	D	417	HEM	C4A-CHB-C1B	-3.78	122.50	127.47
3	B	417	HEM	O2A-CGA-O1A	3.75	132.85	123.30
3	C	417	HEM	CMA-C3A-C4A	-3.71	122.92	128.62
3	C	417	HEM	O2A-CGA-O1A	3.67	132.64	123.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	417	HEM	C4A-CHB-C1B	-3.52	122.84	127.47
3	A	417	HEM	O1D-CGD-CBD	-3.46	111.13	123.03
3	C	417	HEM	CBD-CAD-C3D	3.45	121.91	114.37
3	D	417	HEM	CMA-C3A-C4A	-3.43	123.34	128.62
3	D	417	HEM	O1D-CGD-CBD	-3.38	111.41	123.03
3	B	417	HEM	O1D-CGD-CBD	-3.21	111.99	123.03
3	A	417	HEM	CMA-C3A-C4A	-3.20	123.70	128.62
3	C	417	HEM	O1D-CGD-CBD	-3.13	112.25	123.03
3	D	417	HEM	CBD-CAD-C3D	3.10	121.13	114.37
3	C	417	HEM	CHA-C4D-ND	3.09	128.55	124.31
3	A	417	HEM	CHA-C4D-ND	3.01	128.44	124.31
3	B	417	HEM	CHB-C4A-NA	2.98	129.56	124.58
3	C	417	HEM	CBA-CAA-C2A	2.96	117.91	112.69
3	C	417	HEM	C1A-CHA-C4D	-2.79	123.80	127.47
3	C	417	HEM	CMA-C3A-C2A	2.78	130.19	124.94
3	C	417	HEM	CHB-C4A-NA	2.78	129.22	124.58
3	D	417	HEM	CHB-C4A-NA	2.77	129.21	124.58
3	B	417	HEM	CMA-C3A-C2A	2.76	130.15	124.94
3	D	417	HEM	CMA-C3A-C2A	2.75	130.12	124.94
3	A	417	HEM	CBD-CAD-C3D	2.72	120.32	114.37
3	B	417	HEM	CBD-CAD-C3D	2.70	120.27	114.37
3	A	417	HEM	CHC-C4B-NB	2.58	126.73	124.58
3	A	417	HEM	CMB-C2B-C3B	2.57	132.21	126.16
3	D	417	HEM	O2A-CGA-CBA	-2.54	105.22	114.22
3	A	417	HEM	CHB-C4A-NA	2.53	128.80	124.58
3	C	417	HEM	O2A-CGA-CBA	-2.49	105.43	114.22
3	A	417	HEM	O2A-CGA-CBA	-2.47	105.47	114.22
3	B	417	HEM	O2A-CGA-CBA	-2.42	105.67	114.22
3	B	417	HEM	C3A-C4A-CHB	-2.40	121.44	126.00
3	D	417	HEM	CHA-C4D-ND	2.40	127.61	124.31
3	A	417	HEM	CBA-CAA-C2A	2.39	116.91	112.69
3	B	417	HEM	CBA-CAA-C2A	2.38	116.89	112.69
3	C	417	HEM	CHC-C4B-NB	2.32	126.51	124.58
3	C	417	HEM	CMB-C2B-C3B	2.29	131.55	126.16
3	A	417	HEM	CMA-C3A-C2A	2.26	129.20	124.94
4	A	1450	TCZ	C5-C6-C1	2.24	119.93	117.47
3	A	417	HEM	C1A-CHA-C4D	-2.16	124.63	127.47
3	D	417	HEM	C1A-CHA-C4D	-2.15	124.64	127.47
3	D	417	HEM	CMB-C2B-C3B	2.09	131.08	126.16
3	B	417	HEM	CHA-C4D-ND	2.07	127.15	124.31
3	C	417	HEM	C3A-C4A-CHB	-2.07	122.08	126.00
3	B	417	HEM	CMB-C2B-C3B	2.05	131.00	126.16

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 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 ⓘ

EDS was not executed - this section will therefore be empty.

### 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

### 6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

### 6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

### 6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.