



# Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 09:43 pm BST

PDB ID : 4EGP  
Title : The X-ray crystal structure of CYP199A4 in complex with 2-naphthoic acid  
Authors : Zhou, W.; Bell, S.G.; Yang, W.; Zhou, R.M.; Tan, A.B.H.; Wong, L.-L.  
Deposited on : 2012-03-31  
Resolution : 3.00 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

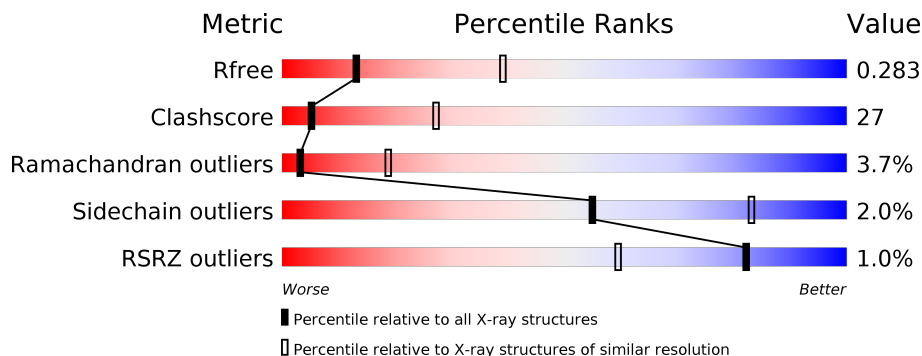
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.00 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	2092 (3.00-3.00)
Clashscore	141614	2416 (3.00-3.00)
Ramachandran outliers	138981	2333 (3.00-3.00)
Sidechain outliers	138945	2336 (3.00-3.00)
RSRZ outliers	127900	1990 (3.00-3.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. 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	410	
1	B	410	

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	FIV	A	502	-	-	X	-

## 2 Entry composition [i](#)

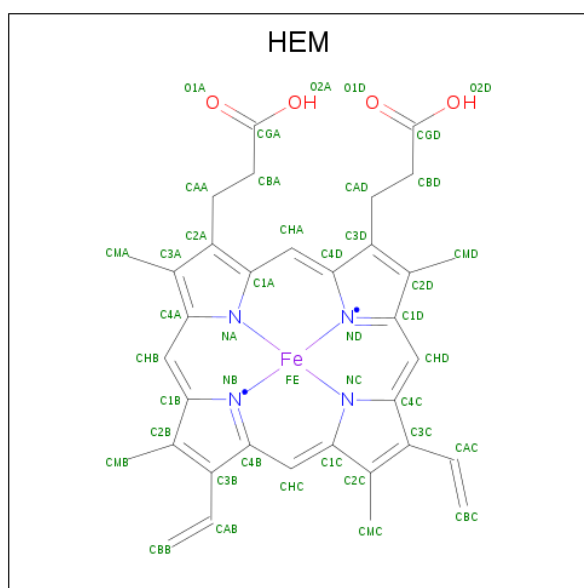
There are 7 unique types of molecules in this entry. The entry contains 6211 atoms, of which 0 are hydrogens and 0 are deuteriums.

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.

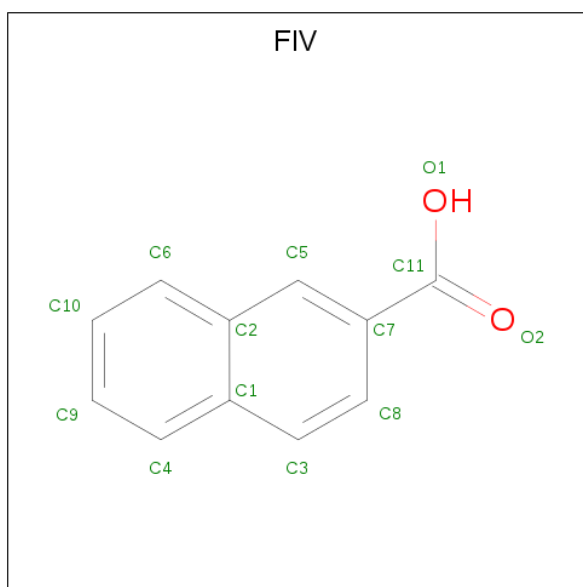
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	393	Total 3024	C 1914	N 534	O 564	S 12	0	1	0
1	B	393	Total 3021	C 1912	N 534	O 564	S 11	0	0	0

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



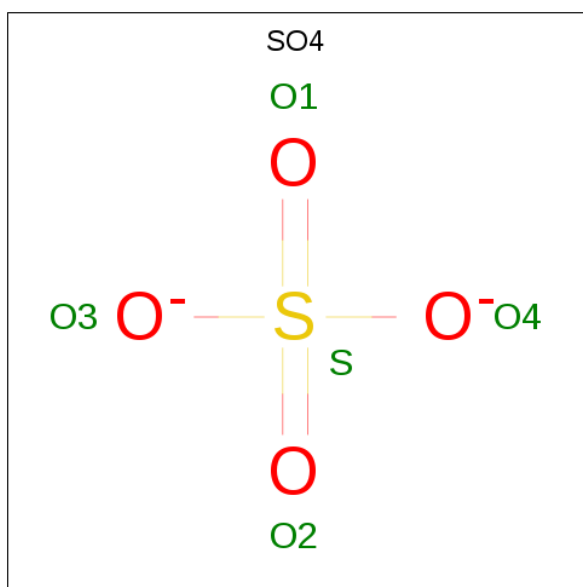
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	Fe	N	O		
2	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	B	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 3 is naphthalene-2-carboxylic acid (three-letter code: FIV) (formula:  $C_{11}H_8O_2$ ).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	C O	0	0
			13	11 2		
3	A	1	Total	C O	0	0
			13	11 2		
3	B	1	Total	C O	0	0
			13	11 2		

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



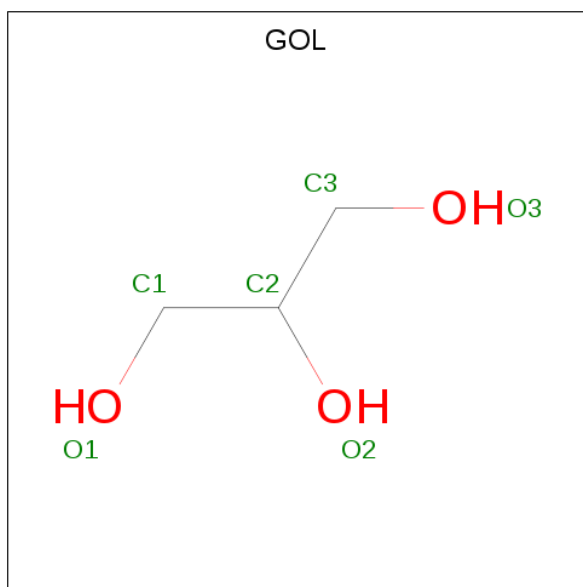
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	O S	0	0
			5	4 1		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total O S 5 4 1	0	0
4	B	1	Total O S 5 4 1	0	0

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



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

- Molecule 6 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

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

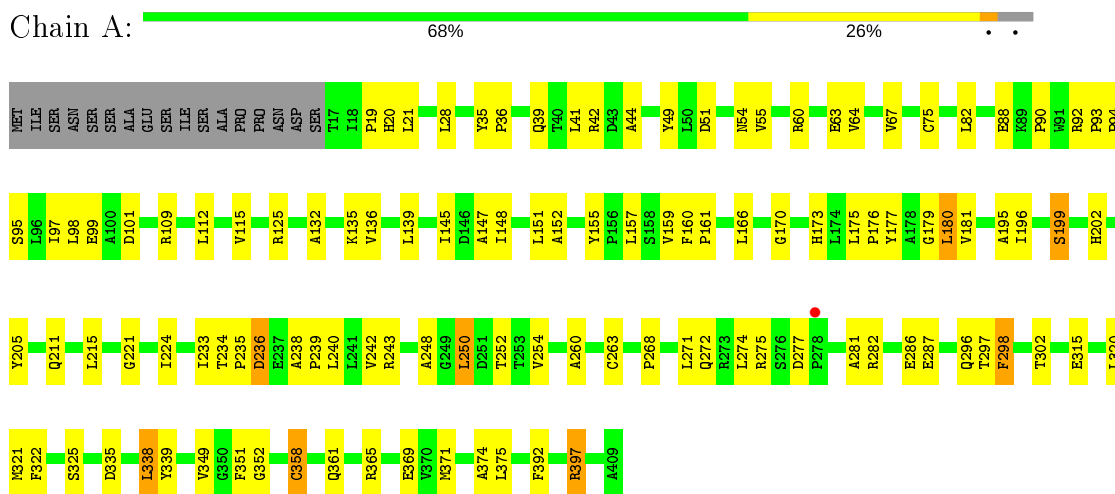
- Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	11	Total O 11 11	0	0
7	B	8	Total O 8 8	0	0

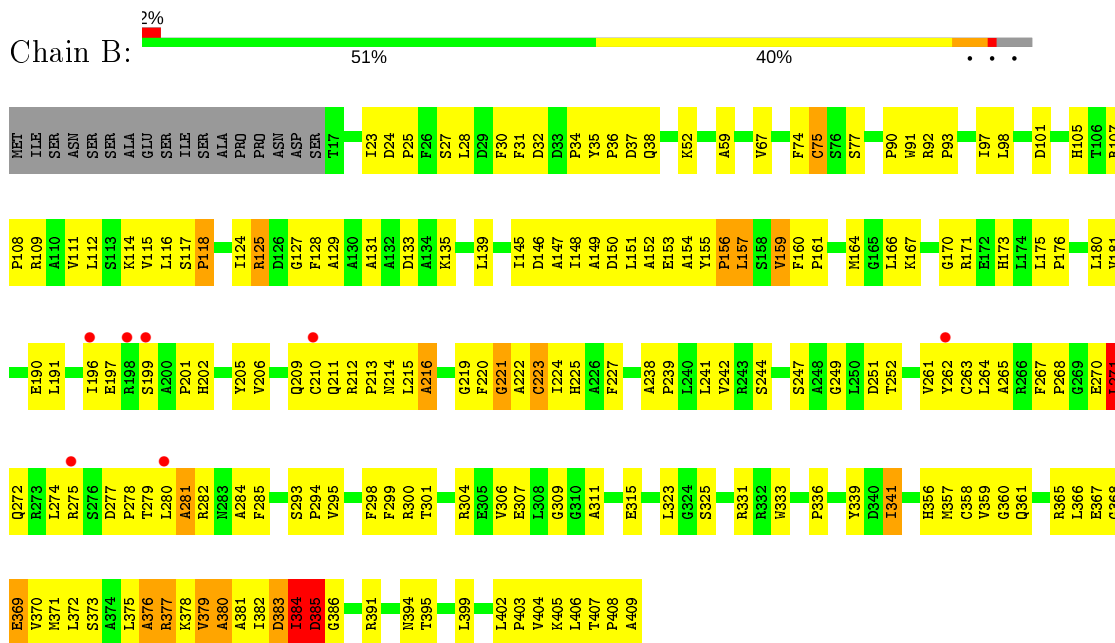
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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



- Molecule 1: Cytochrome P450



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	143.54Å 172.21Å 106.26Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.94 – 3.00 48.94 – 3.00	Depositor EDS
% Data completeness (in resolution range)	99.0 (48.94-3.00) 99.0 (48.94-3.00)	Depositor EDS
$R_{merge}$	0.14	Depositor
$R_{sym}$	0.09	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.37 (at 3.01Å)	Xtrriage
Refinement program	REFMAC 5.5.0109	Depositor
R, $R_{free}$	0.230 , 0.287 0.234 , 0.283	Depositor DCC
$R_{free}$ test set	1346 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	59.8	Xtrriage
Anisotropy	0.686	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 36.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.45$ , $\langle L^2 \rangle = 0.28$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	6211	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	62.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.11% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, FIV, SO4, HEM, CL

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z  > 5$	RMSZ	# $ Z  > 5$
1	A	0.68	0/3099	0.82	3/4217 (0.1%)
1	B	0.69	1/3093 (0.0%)	0.84	5/4209 (0.1%)
All	All	0.69	1/6192 (0.0%)	0.83	8/8426 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	223	CYS	CB-SG	-5.26	1.73	1.81

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	385	ASP	N-CA-CB	-10.96	90.87	110.60
1	A	281	ALA	CB-CA-C	-7.86	98.31	110.10
1	B	384	ILE	CB-CA-C	-5.77	100.05	111.60
1	A	180	LEU	CA-CB-CG	5.70	128.41	115.30
1	B	293	SER	N-CA-C	5.42	125.64	111.00
1	B	271	LEU	CA-CB-CG	5.25	127.37	115.30
1	B	251	ASP	CB-CA-C	-5.17	100.06	110.40
1	A	338	LEU	CA-CB-CG	5.07	126.96	115.30

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within

the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3024	0	2998	99	0
1	B	3021	0	2993	227	0
2	A	43	0	30	3	0
2	B	43	0	30	8	0
3	A	26	0	14	7	0
3	B	13	0	7	4	0
4	A	10	0	0	0	0
4	B	5	0	0	0	0
5	A	6	0	8	0	0
6	A	1	0	0	0	0
7	A	11	0	0	1	0
7	B	8	0	0	1	0
All	All	6211	0	6080	327	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 27.

All (327) 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:381:ALA:CB	1:B:407:THR:HB	1.34	1.54
1:B:376:ALA:O	1:B:377:ARG:HG3	1.36	1.25
1:B:381:ALA:CB	1:B:407:THR:CB	2.20	1.17
1:B:381:ALA:HB3	1:B:407:THR:HB	1.30	1.13
1:B:384:ILE:HG13	1:B:385:ASP:N	1.58	1.10
1:B:379:VAL:HG22	1:B:380:ALA:H	1.05	1.09
1:B:381:ALA:HB3	1:B:407:THR:CB	1.85	1.05
1:B:384:ILE:CG1	1:B:385:ASP:H	1.66	1.04
1:B:379:VAL:HG22	1:B:380:ALA:N	1.67	1.04
1:B:244:SER:HB2	3:B:502:FIV:H2	1.36	1.03
1:B:381:ALA:HB1	1:B:407:THR:CB	1.85	1.03
1:B:384:ILE:O	1:B:385:ASP:HB2	1.23	1.03
1:B:384:ILE:HD12	1:B:403:PRO:O	1.56	1.02
1:B:384:ILE:O	1:B:385:ASP:CB	2.11	0.99
1:B:281:ALA:O	1:B:372:LEU:HD22	1.67	0.95
1:B:384:ILE:HG13	1:B:385:ASP:H	0.78	0.93
1:B:384:ILE:N	1:B:404:VAL:HG13	1.84	0.92
1:B:381:ALA:HB1	1:B:407:THR:HB	0.92	0.92
1:B:244:SER:HB2	3:B:502:FIV:C8	2.00	0.92
1:A:92:ARG:HH22	3:A:502:FIV:H2	1.34	0.92

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:379:VAL:HA	1:B:409:ALA:HB2	1.52	0.91
1:B:382:ILE:O	1:B:383:ASP:HB2	1.70	0.90
1:B:379:VAL:CG2	1:B:380:ALA:H	1.85	0.89
1:B:376:ALA:O	1:B:377:ARG:CG	2.19	0.89
1:B:281:ALA:O	1:B:372:LEU:CD2	2.21	0.88
1:B:383:ASP:O	1:B:384:ILE:HB	1.73	0.87
1:B:382:ILE:HG23	1:B:404:VAL:HG11	1.55	0.87
1:B:147:ALA:HB3	1:B:402:LEU:O	1.75	0.85
1:B:282:ARG:HH22	1:B:365:ARG:HH11	1.24	0.84
1:B:24:ASP:OD2	1:B:52:LYS:HE3	1.77	0.83
1:B:381:ALA:HB3	1:B:407:THR:CA	2.08	0.83
1:B:167:LYS:HE3	1:B:209:GLN:HE22	1.44	0.81
1:A:75:CYS:HB2	1:A:101:ASP:OD2	1.81	0.80
1:A:39:GLN:HA	1:A:42:ARG:HH21	1.46	0.80
1:B:384:ILE:HA	1:B:404:VAL:HA	1.64	0.80
1:B:270:GLU:HB3	1:B:341:ILE:HG12	1.62	0.80
1:A:236:ASP:O	1:A:240:LEU:HD13	1.81	0.80
1:B:383:ASP:O	1:B:384:ILE:CB	2.30	0.79
1:B:282:ARG:NH2	1:B:365:ARG:HH11	1.79	0.79
1:B:166:LEU:HD22	1:B:209:GLN:HE21	1.46	0.78
1:B:277:ASP:O	1:B:280:LEU:HD12	1.83	0.78
1:A:20:HIS:HD2	1:A:49:TYR:CD2	2.02	0.77
1:B:384:ILE:H	1:B:404:VAL:HG13	1.49	0.77
1:A:20:HIS:CD2	1:A:49:TYR:CD2	2.73	0.76
1:B:151:LEU:HD21	1:B:371:MET:CE	2.16	0.76
1:B:265:ALA:HB1	1:B:383:ASP:HA	1.66	0.75
1:A:240:LEU:HD12	1:A:240:LEU:N	2.02	0.75
1:B:265:ALA:HB2	1:B:382:ILE:HG22	1.68	0.74
1:B:127:GLY:O	1:B:131:ALA:HB3	1.87	0.73
1:A:135:LYS:O	1:A:139:LEU:HG	1.89	0.73
1:B:274:LEU:HD11	1:B:281:ALA:H	1.54	0.72
1:B:155:TYR:N	1:B:156:PRO:HD2	2.05	0.72
1:B:210:CYS:O	1:B:215:LEU:HD11	1.89	0.72
1:B:161:PRO:HA	1:B:164:MET:HB2	1.73	0.71
1:A:397:ARG:HH11	1:A:397:ARG:CG	2.03	0.71
1:B:117:SER:HB2	1:B:118:PRO:HD2	1.73	0.71
1:B:378:LYS:HG2	1:B:378:LYS:O	1.90	0.70
1:B:196:ILE:HG22	1:B:197:GLU:HG3	1.74	0.69
1:B:278:PRO:C	1:B:280:LEU:H	1.95	0.69
1:B:252:THR:HB	2:B:501:HEM:HBB2	1.74	0.69
1:B:145:ILE:HD13	1:B:406:LEU:HD11	1.73	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:265:ALA:HB2	1:B:382:ILE:CG2	2.23	0.69
1:B:384:ILE:CD1	1:B:403:PRO:O	2.37	0.69
1:A:151:LEU:HD11	1:A:371:MET:SD	2.34	0.68
1:B:384:ILE:CA	1:B:404:VAL:HA	2.23	0.68
1:B:220:PHE:O	1:B:221:GLY:C	2.29	0.68
1:A:20:HIS:CD2	1:A:51:ASP:OD1	2.46	0.68
1:B:156:PRO:HG2	1:B:157:LEU:H	1.59	0.68
1:B:151:LEU:HD21	1:B:371:MET:HE3	1.77	0.67
1:B:267:PHE:N	1:B:268:PRO:HD3	2.10	0.67
1:B:376:ALA:C	1:B:377:ARG:HG3	2.14	0.66
1:A:19:PRO:HG3	1:A:44:ALA:O	1.94	0.66
1:B:285:PHE:HB2	1:B:372:LEU:HD11	1.77	0.66
1:B:280:LEU:O	1:B:282:ARG:N	2.29	0.66
1:B:157:LEU:HD11	1:B:175:LEU:HD11	1.77	0.66
1:B:382:ILE:O	1:B:383:ASP:CB	2.39	0.66
1:B:381:ALA:HB3	1:B:407:THR:C	2.17	0.65
1:B:220:PHE:O	1:B:224:ILE:N	2.20	0.65
1:B:114:LYS:O	1:B:117:SER:HB3	1.95	0.65
1:B:373:SER:HB3	1:B:377:ARG:HH21	1.62	0.65
1:B:145:ILE:CD1	1:B:406:LEU:HD11	2.27	0.65
1:B:133:ASP:HB3	1:B:378:LYS:NZ	2.12	0.65
1:B:220:PHE:O	1:B:222:ALA:N	2.30	0.64
1:B:129:ALA:HB2	1:B:370:VAL:HG11	1.78	0.64
1:B:282:ARG:NH2	1:B:365:ARG:NH1	2.45	0.64
1:B:220:PHE:O	1:B:223:CYS:N	2.31	0.64
1:B:125:ARG:HB3	7:B:606:HOH:O	1.99	0.63
1:B:215:LEU:O	1:B:216:ALA:C	2.36	0.63
1:A:20:HIS:HD2	1:A:51:ASP:OD1	1.80	0.63
1:B:373:SER:HB3	1:B:377:ARG:NH2	2.14	0.63
1:B:111:VAL:O	1:B:115:VAL:HG23	1.98	0.62
1:B:146:ASP:O	1:B:148:ILE:N	2.30	0.62
1:A:21:LEU:HD13	1:A:41:LEU:HD23	1.82	0.62
1:B:282:ARG:HH22	1:B:365:ARG:NH1	1.96	0.61
1:B:35:TYR:HD2	1:B:38:GLN:OE1	1.82	0.61
1:B:151:LEU:HD21	1:B:371:MET:HE2	1.82	0.61
1:A:240:LEU:CD1	1:A:240:LEU:N	2.62	0.61
1:B:212:ARG:N	1:B:213:PRO:HD2	2.14	0.61
1:B:381:ALA:CB	1:B:407:THR:CG2	2.79	0.61
1:A:320:LEU:HD12	1:A:321:MET:H	1.66	0.61
1:A:215:LEU:HD22	1:A:221:GLY:HA3	1.82	0.60
1:B:379:VAL:O	1:B:380:ALA:HB2	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:238:ALA:HB3	1:A:239:PRO:HD3	1.84	0.60
1:A:28:LEU:HD22	1:A:392:PHE:CD1	2.37	0.60
1:B:105:HIS:HE1	2:B:501:HEM:O1D	1.85	0.60
1:A:139:LEU:HD13	1:A:145:ILE:HD13	1.83	0.60
1:A:157:LEU:O	1:A:161:PRO:HD2	2.01	0.60
1:A:97:ILE:HD13	1:A:112:LEU:HD12	1.84	0.59
1:A:180:LEU:HD21	1:A:199:SER:HB2	1.84	0.59
1:B:155:TYR:N	1:B:156:PRO:CD	2.65	0.59
1:B:384:ILE:HG23	1:B:385:ASP:N	2.18	0.59
1:A:181:VAL:HG11	3:A:502:FIV:H3	1.83	0.59
1:B:268:PRO:HA	1:B:271:LEU:HD23	1.85	0.59
1:B:368:GLY:O	1:B:372:LEU:HD13	2.03	0.59
1:A:39:GLN:CA	1:A:42:ARG:HH21	2.13	0.58
1:B:379:VAL:HG23	1:B:407:THR:O	2.03	0.58
1:A:240:LEU:CD1	1:A:240:LEU:H	2.15	0.58
1:B:154:ALA:N	1:B:156:PRO:HD2	2.19	0.58
1:B:166:LEU:HD22	1:B:209:GLN:NE2	2.19	0.58
1:B:381:ALA:O	1:B:407:THR:N	2.34	0.57
1:B:281:ALA:O	1:B:284:ALA:HB3	2.04	0.57
1:B:379:VAL:CG2	1:B:407:THR:O	2.53	0.57
1:A:173:HIS:C	1:A:176:PRO:HD2	2.24	0.57
1:B:77:SER:HB3	1:B:101:ASP:CG	2.25	0.57
1:B:34:PRO:O	1:B:38:GLN:HG3	2.05	0.57
1:A:196:ILE:O	1:A:199:SER:HB3	2.05	0.56
1:A:272:GLN:HE22	1:A:275:ARG:HH11	1.53	0.56
1:A:335:ASP:HB3	1:A:338:LEU:HG	1.86	0.56
1:B:97:ILE:HG23	1:B:98:LEU:H	1.71	0.56
1:A:268:PRO:O	1:A:271:LEU:HB3	2.04	0.56
1:B:295:VAL:HG11	1:B:298:PHE:CE1	2.40	0.56
1:B:384:ILE:HA	1:B:404:VAL:HG22	1.87	0.56
1:A:250:LEU:HD12	1:A:254:VAL:HG23	1.88	0.56
1:B:369:GLU:O	1:B:369:GLU:HG2	2.05	0.55
1:A:211:GLN:NE2	7:A:606:HOH:O	2.37	0.55
1:A:351:PHE:HB3	1:A:358:CYS:HB3	1.89	0.55
1:B:278:PRO:HB3	1:B:376:ALA:HB1	1.88	0.55
1:B:148:ILE:HG12	1:B:402:LEU:HB3	1.88	0.55
1:A:95:SER:HB3	1:A:98:LEU:H	1.72	0.55
1:B:219:GLY:O	1:B:220:PHE:HD1	1.89	0.55
1:B:124:ILE:HG13	1:B:125:ARG:N	2.22	0.54
1:B:378:LYS:CG	1:B:378:LYS:O	2.55	0.54
1:B:170:GLY:HA3	1:B:205:TYR:CZ	2.42	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:177:TYR:O	1:A:180:LEU:HB3	2.08	0.54
1:A:260:ALA:O	1:A:263:CYS:HB2	2.07	0.54
1:B:304:ARG:O	1:B:315:GLU:HA	2.08	0.54
1:B:91:TRP:CE3	1:B:92:ARG:HB3	2.42	0.54
1:A:139:LEU:HB3	1:A:145:ILE:HG21	1.89	0.54
1:B:128:PHE:CE1	1:B:159:VAL:HG13	2.43	0.54
1:B:274:LEU:HD12	1:B:280:LEU:HB2	1.90	0.53
1:B:265:ALA:CB	1:B:382:ILE:HG22	2.38	0.53
1:B:107:ARG:HB3	1:B:108:PRO:CD	2.38	0.53
1:B:153:GLU:O	1:B:157:LEU:HD22	2.09	0.53
1:B:238:ALA:N	1:B:239:PRO:HD2	2.24	0.53
1:A:63:GLU:O	1:A:67:VAL:HG23	2.08	0.53
1:A:125:ARG:HH22	1:A:369:GLU:CD	2.11	0.53
1:B:285:PHE:HB2	1:B:372:LEU:CD1	2.38	0.53
1:A:97:ILE:HD12	1:A:109:ARG:HA	1.90	0.53
1:A:49:TYR:HE2	1:A:51:ASP:HA	1.74	0.53
1:A:95:SER:HB2	1:A:99:GLU:HG3	1.91	0.53
1:B:379:VAL:CA	1:B:409:ALA:HB2	2.32	0.53
1:A:397:ARG:HH11	1:A:397:ARG:HG3	1.74	0.52
1:B:225:HIS:C	1:B:227:PHE:H	2.13	0.52
1:B:391:ARG:H	1:B:399:LEU:HA	1.73	0.52
1:B:356:HIS:HD1	2:B:501:HEM:CGD	2.23	0.52
1:A:320:LEU:HD12	1:A:321:MET:N	2.25	0.52
1:B:366:LEU:C	1:B:368:GLY:H	2.14	0.51
1:B:97:ILE:HG23	1:B:98:LEU:N	2.25	0.51
1:A:49:TYR:CE2	1:A:51:ASP:HA	2.45	0.51
1:A:397:ARG:HH11	1:A:397:ARG:HG2	1.74	0.51
1:B:149:ALA:HA	1:B:153:GLU:HB2	1.92	0.51
1:B:160:PHE:HB3	1:B:161:PRO:HD3	1.93	0.51
1:B:275:ARG:NH1	1:B:380:ALA:CB	2.74	0.51
1:B:152:ALA:O	1:B:156:PRO:HG3	2.10	0.51
1:B:307:GLU:HA	1:B:311:ALA:O	2.11	0.51
1:A:95:SER:HG	3:A:502:FIV:C11	2.24	0.51
1:B:153:GLU:O	1:B:157:LEU:HD13	2.10	0.51
1:B:153:GLU:C	1:B:157:LEU:HD13	2.32	0.50
1:B:148:ILE:HG12	1:B:402:LEU:CB	2.41	0.50
1:A:173:HIS:O	1:A:176:PRO:HD2	2.11	0.50
1:B:263:CYS:O	1:B:267:PHE:HD2	1.93	0.50
1:B:35:TYR:CD2	1:B:38:GLN:OE1	2.63	0.50
1:B:146:ASP:O	1:B:146:ASP:OD2	2.30	0.50
1:B:384:ILE:CG1	1:B:385:ASP:N	2.32	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:173:HIS:ND1	1:B:201:PRO:O	2.45	0.49
1:B:281:ALA:O	1:B:372:LEU:HD23	2.08	0.49
1:B:147:ALA:CB	1:B:402:LEU:O	2.52	0.49
1:B:357:MET:HG2	1:B:361:GLN:HE22	1.76	0.49
1:B:75:CYS:HB2	1:B:101:ASP:OD2	2.13	0.49
1:B:223:CYS:O	1:B:227:PHE:HD1	1.96	0.49
1:B:221:GLY:HA2	1:B:242:VAL:HG22	1.93	0.49
1:B:274:LEU:HD11	1:B:281:ALA:N	2.26	0.49
1:A:97:ILE:CD1	1:A:109:ARG:HA	2.43	0.49
1:B:278:PRO:C	1:B:280:LEU:N	2.63	0.49
1:A:148:ILE:HA	1:A:152:ALA:HB3	1.95	0.49
1:A:358:CYS:HA	2:A:501:HEM:CHA	2.43	0.49
1:B:215:LEU:O	1:B:216:ALA:O	2.30	0.49
1:A:180:LEU:CD2	1:A:199:SER:HB2	2.43	0.48
1:A:282:ARG:HD2	1:A:369:GLU:OE1	2.13	0.48
1:A:179:GLY:HA3	3:A:503:FIV:H4	1.96	0.48
1:A:115:VAL:HG21	1:A:224:ILE:HG13	1.96	0.48
1:B:109:ARG:NH1	1:B:357:MET:O	2.44	0.48
1:B:264:LEU:HD21	1:B:339:TYR:HE2	1.78	0.48
1:B:300:ARG:HH22	2:B:501:HEM:CGA	2.26	0.48
1:A:176:PRO:HB2	1:A:202:HIS:CE1	2.49	0.47
1:B:274:LEU:CD1	1:B:281:ALA:H	2.23	0.47
1:B:285:PHE:HB2	1:B:372:LEU:HD21	1.96	0.47
1:A:272:GLN:HE22	1:A:275:ARG:NH1	2.11	0.47
1:A:95:SER:OG	3:A:502:FIV:C11	2.61	0.47
1:B:97:ILE:HD13	1:B:112:LEU:HD12	1.96	0.47
1:A:349:VAL:O	1:A:352:GLY:N	2.43	0.47
1:A:175:LEU:N	1:A:176:PRO:CD	2.78	0.47
1:B:267:PHE:N	1:B:268:PRO:CD	2.77	0.47
1:A:252:THR:HB	2:A:501:HEM:C3B	2.49	0.47
1:A:180:LEU:CD1	1:A:195:ALA:O	2.63	0.47
1:B:383:ASP:O	1:B:384:ILE:CG2	2.63	0.47
1:B:261:VAL:C	1:B:263:CYS:H	2.17	0.46
1:B:219:GLY:C	1:B:220:PHE:CD1	2.89	0.46
1:B:220:PHE:C	1:B:222:ALA:N	2.65	0.46
1:B:264:LEU:CD2	1:B:339:TYR:HE2	2.27	0.46
1:B:384:ILE:O	1:B:405:LYS:N	2.49	0.46
1:B:107:ARG:O	1:B:111:VAL:HG23	2.16	0.46
1:A:180:LEU:HD13	1:A:195:ALA:O	2.16	0.46
1:B:375:LEU:C	1:B:377:ARG:H	2.19	0.46
1:B:407:THR:HG23	1:B:408:PRO:HD2	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:263:CYS:HB3	1:A:339:TYR:CD2	2.51	0.46
1:B:59:ALA:HB1	1:B:325:SER:OG	2.16	0.46
1:A:374:ALA:O	1:A:375:LEU:C	2.51	0.46
1:B:373:SER:C	1:B:375:LEU:H	2.19	0.46
1:B:384:ILE:C	1:B:404:VAL:HA	2.35	0.46
1:B:219:GLY:C	1:B:220:PHE:HD1	2.20	0.45
1:B:77:SER:HB3	1:B:101:ASP:OD2	2.16	0.45
1:A:242:VAL:O	1:A:243:ARG:C	2.55	0.45
1:B:180:LEU:HD22	1:B:199:SER:HB2	1.98	0.45
1:B:333:TRP:HB2	1:B:336:PRO:HB3	1.98	0.45
1:A:302:THR:OG1	1:A:315:GLU:HA	2.17	0.45
1:A:42:ARG:HG2	1:A:60:ARG:HG2	1.98	0.45
1:B:156:PRO:CG	1:B:157:LEU:H	2.28	0.45
1:B:275:ARG:HH11	1:B:380:ALA:CB	2.29	0.45
1:B:383:ASP:H	1:B:404:VAL:CG1	2.30	0.45
1:B:190:GLU:HB3	1:B:191:LEU:HD12	1.97	0.45
1:B:358:CYS:SG	2:B:501:HEM:C1C	3.09	0.45
1:B:382:ILE:HG22	1:B:383:ASP:N	2.32	0.45
1:B:384:ILE:CG2	1:B:385:ASP:N	2.74	0.45
1:B:175:LEU:HB2	1:B:176:PRO:HD3	1.99	0.44
1:B:249:GLY:HA2	2:B:501:HEM:C2C	2.53	0.44
1:B:382:ILE:CG2	1:B:404:VAL:HG11	2.37	0.44
1:A:248:ALA:HB2	3:A:502:FIV:C6	2.46	0.44
1:A:234:THR:HB	1:A:235:PRO:HD2	1.99	0.44
1:A:55:VAL:HG21	1:A:82:LEU:CD1	2.48	0.44
1:B:285:PHE:HB2	1:B:372:LEU:CD2	2.48	0.44
1:B:154:ALA:H	1:B:156:PRO:HD2	1.82	0.44
1:B:241:LEU:HA	1:B:241:LEU:HD23	1.85	0.44
1:A:132:ALA:O	1:A:136:VAL:HG23	2.18	0.44
1:B:139:LEU:HD21	1:B:150:ASP:O	2.18	0.44
1:B:224:ILE:HG22	1:B:224:ILE:O	2.17	0.44
1:B:139:LEU:HD22	1:B:150:ASP:HB3	2.00	0.43
1:B:135:LYS:HD2	1:B:135:LYS:HA	1.83	0.43
1:A:160:PHE:HB3	1:A:161:PRO:HD3	2.00	0.43
1:B:139:LEU:CD2	1:B:150:ASP:HB3	2.48	0.43
1:B:384:ILE:HG13	1:B:386:GLY:H	1.82	0.43
1:A:238:ALA:O	1:A:242:VAL:HG23	2.18	0.43
1:B:92:ARG:HG3	1:B:93:PRO:O	2.19	0.43
1:B:24:ASP:HA	1:B:25:PRO:HD3	1.82	0.43
1:A:297:THR:HB	1:A:320:LEU:HD11	2.01	0.43
1:A:93:PRO:HA	1:A:94:PRO:HD3	1.75	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:297:THR:HA	1:A:321:MET:O	2.18	0.43
1:A:181:VAL:CG1	3:A:502:FIV:H3	2.48	0.43
1:B:202:HIS:O	1:B:206:VAL:HG23	2.19	0.43
1:B:295:VAL:HG11	1:B:298:PHE:HE1	1.84	0.43
1:B:67:VAL:HG13	1:B:74:PHE:CG	2.54	0.43
1:A:55:VAL:HG21	1:A:82:LEU:HD11	2.00	0.42
1:B:115:VAL:HG11	1:B:224:ILE:CG1	2.49	0.42
1:B:214:ASN:C	1:B:215:LEU:HG	2.39	0.42
1:B:27:SER:O	1:B:31:PHE:CD1	2.72	0.42
1:A:287:GLU:HG3	1:A:339:TYR:CD1	2.53	0.42
1:A:397:ARG:NH1	1:A:397:ARG:CG	2.71	0.42
1:B:368:GLY:O	1:B:370:VAL:N	2.52	0.42
1:B:30:PHE:HE1	1:B:37:ASP:HB2	1.84	0.42
1:A:88:GLU:O	1:A:90:PRO:HD3	2.19	0.42
1:B:117:SER:CB	1:B:118:PRO:HD2	2.42	0.42
1:A:147:ALA:O	1:A:151:LEU:HB3	2.20	0.42
1:B:133:ASP:HB3	1:B:378:LYS:HZ2	1.81	0.42
1:B:224:ILE:CG2	1:B:224:ILE:O	2.67	0.42
1:B:379:VAL:HG22	1:B:381:ALA:H	1.83	0.42
1:A:92:ARG:HD2	1:A:243:ARG:NH2	2.34	0.42
1:B:383:ASP:H	1:B:404:VAL:HG13	1.84	0.42
1:A:322:PHE:O	1:A:325:SER:HB3	2.20	0.42
1:A:35:TYR:N	1:A:36:PRO:HD2	2.34	0.42
1:B:211:GLN:NE2	1:B:211:GLN:HA	2.35	0.42
1:B:278:PRO:O	1:B:280:LEU:N	2.53	0.42
1:A:298:PHE:O	1:A:321:MET:HB2	2.20	0.42
1:A:97:ILE:HG12	2:A:501:HEM:CGD	2.50	0.42
1:B:379:VAL:CG2	1:B:380:ALA:N	2.42	0.42
1:A:236:ASP:O	1:A:240:LEU:CD1	2.62	0.41
1:A:296:GLN:CG	1:A:397:ARG:HB2	2.50	0.41
1:B:124:ILE:O	1:B:125:ARG:C	2.58	0.41
1:B:299:PHE:CD1	1:B:299:PHE:N	2.89	0.41
1:B:111:VAL:HG13	1:B:227:PHE:CD2	2.55	0.41
1:A:274:LEU:O	1:A:277:ASP:C	2.58	0.41
1:B:368:GLY:C	1:B:370:VAL:H	2.23	0.41
1:B:128:PHE:CD1	1:B:159:VAL:HG13	2.55	0.41
1:B:366:LEU:O	1:B:370:VAL:HG22	2.21	0.41
1:B:145:ILE:HD12	1:B:147:ALA:HA	2.02	0.41
1:B:268:PRO:O	1:B:272:GLN:HG2	2.20	0.41
1:B:360:GLY:HA3	2:B:501:HEM:C3C	2.56	0.41
1:A:49:TYR:OH	1:A:54:ASN:ND2	2.38	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:28:LEU:O	1:B:32:ASP:HB2	2.20	0.41
1:A:112:LEU:HA	1:A:112:LEU:HD23	1.90	0.41
1:A:170:GLY:HA3	1:A:205:TYR:CE1	2.56	0.41
1:B:366:LEU:C	1:B:368:GLY:N	2.74	0.41
1:B:244:SER:CB	3:B:502:FIV:H2	2.27	0.41
1:B:181:VAL:HG21	1:B:247:SER:HB3	2.02	0.41
2:B:501:HEM:C3D	3:B:502:FIV:H8	2.56	0.41
1:B:91:TRP:CZ3	1:B:92:ARG:HB3	2.56	0.41
1:A:286:GLU:OE2	1:A:365:ARG:NE	2.54	0.41
1:B:191:LEU:HD12	1:B:191:LEU:N	2.36	0.41
1:B:151:LEU:CD2	1:B:371:MET:CE	2.95	0.41
1:B:275:ARG:NH1	1:B:380:ALA:HB3	2.36	0.41
1:A:233:ILE:HD12	1:A:238:ALA:HA	2.02	0.41
1:A:20:HIS:O	1:A:21:LEU:HG	2.22	0.40
1:B:223:CYS:O	1:B:227:PHE:CD1	2.75	0.40
1:B:23:ILE:HG22	1:B:24:ASP:N	2.36	0.40
1:B:383:ASP:O	1:B:384:ILE:HG22	2.21	0.40
1:B:116:LEU:HD12	1:B:359:VAL:HG22	2.03	0.40
1:B:404:VAL:HG12	1:B:405:LYS:N	2.36	0.40
1:A:155:TYR:O	1:A:159:VAL:HG22	2.22	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	392/410 (96%)	347 (88%)	42 (11%)	3 (1%)	19 57
1	B	391/410 (95%)	299 (76%)	66 (17%)	26 (7%)	1 6
All	All	783/820 (96%)	646 (82%)	108 (14%)	29 (4%)	3 19

All (29) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	281	ALA
1	B	377	ARG
1	B	380	ALA
1	B	383	ASP
1	B	384	ILE
1	B	385	ASP
1	B	125	ARG
1	B	157	LEU
1	B	221	GLY
1	B	279	THR
1	B	379	VAL
1	A	361	GLN
1	B	309	GLY
1	B	331	ARG
1	A	358	CYS
1	B	171	ARG
1	B	262	TYR
1	B	367	GLU
1	B	90	PRO
1	B	156	PRO
1	B	369	GLU
1	B	394	ASN
1	B	216	ALA
1	B	271	LEU
1	B	376	ALA
1	A	64	VAL
1	B	36	PRO
1	B	118	PRO
1	B	294	PRO

### 5.3.2 Protein sidechains [i](#)

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	320/334 (96%)	314 (98%)	6 (2%)	57	84
1	B	319/334 (96%)	312 (98%)	7 (2%)	52	81
All	All	639/668 (96%)	626 (98%)	13 (2%)	55	83

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	166	LEU
1	A	199	SER
1	A	236	ASP
1	A	250	LEU
1	A	298	PHE
1	A	397	ARG
1	B	75	CYS
1	B	159	VAL
1	B	301	THR
1	B	306	VAL
1	B	323	LEU
1	B	341	ILE
1	B	395	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	20	HIS
1	A	168	GLN
1	A	203	GLN
1	A	209	GLN
1	A	211	GLN
1	A	214	ASN
1	A	272	GLN
1	A	283	ASN
1	B	69	ASN
1	B	105	HIS
1	B	209	GLN
1	B	211	GLN
1	B	225	HIS
1	B	255	ASN
1	B	272	GLN
1	B	283	ASN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 10 ligands modelled in this entry, 1 is monoatomic - leaving 9 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
3	FIV	A	503	-	12,14,14	1.34	1 (8%)	16,19,19	0.62	0
4	SO4	B	503	-	4,4,4	0.15	0	6,6,6	0.22	0
2	HEM	B	501	1	27,50,50	2.14	6 (22%)	17,82,82	1.35	3 (17%)
3	FIV	A	502	-	12,14,14	1.65	1 (8%)	16,19,19	1.48	4 (25%)
3	FIV	B	502	-	12,14,14	1.45	1 (8%)	16,19,19	1.66	4 (25%)
4	SO4	A	505	-	4,4,4	0.11	0	6,6,6	0.49	0
4	SO4	A	504	-	4,4,4	0.18	0	6,6,6	0.28	0
5	GOL	A	506	-	5,5,5	0.38	0	5,5,5	0.62	0
2	HEM	A	501	1	27,50,50	2.09	5 (18%)	17,82,82	1.92	6 (35%)

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	B	501	1	-	0/6/54/54	-
3	FIV	A	502	-	-	0/0/4/4	0/2/2/2
3	FIV	B	502	-	-	0/0/4/4	0/2/2/2
3	FIV	A	503	-	-	0/0/4/4	0/2/2/2
5	GOL	A	506	-	-	2/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	HEM	A	501	1	-	0/6/54/54	-

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	HEM	C3D-C2D	5.44	1.53	1.37
2	B	501	HEM	C3C-C2C	-4.77	1.33	1.40
2	A	501	HEM	C3D-C2D	4.73	1.51	1.37
2	B	501	HEM	C3B-C2B	-4.57	1.34	1.40
2	A	501	HEM	C3C-C2C	-4.44	1.34	1.40
2	A	501	HEM	C3B-C2B	-4.43	1.34	1.40
3	A	502	FIV	C7-C11	4.20	1.51	1.47
2	A	501	HEM	C3B-CAB	3.96	1.56	1.47
2	A	501	HEM	C3C-CAC	3.58	1.55	1.47
3	B	502	FIV	C7-C11	3.53	1.50	1.47
3	A	503	FIV	C7-C11	3.32	1.50	1.47
2	B	501	HEM	C3B-CAB	3.27	1.54	1.47
2	B	501	HEM	C3C-CAC	3.20	1.54	1.47
2	B	501	HEM	CAA-C2A	2.03	1.55	1.52

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	HEM	CBD-CAD-C3D	-5.02	103.22	112.48
3	B	502	FIV	C5-C7-C11	3.48	124.94	120.36
3	B	502	FIV	C8-C7-C11	-3.28	115.96	120.37
3	A	502	FIV	C5-C7-C11	3.27	124.66	120.36
3	A	502	FIV	C8-C7-C11	-2.94	116.42	120.37
2	B	501	HEM	CAA-CBA-CGA	-2.63	108.26	112.67
2	A	501	HEM	C1D-C2D-C3D	-2.59	105.19	107.00
2	A	501	HEM	C4C-C3C-C2C	2.55	108.68	106.90
2	A	501	HEM	CMC-C2C-C3C	2.47	129.30	124.68
2	B	501	HEM	CMC-C2C-C3C	2.30	128.99	124.68
2	B	501	HEM	C4C-C3C-C2C	2.29	108.50	106.90
3	B	502	FIV	C4-C1-C2	2.23	122.91	118.92
3	B	502	FIV	C4-C1-C3	-2.12	118.23	123.19
3	A	502	FIV	C4-C1-C3	-2.07	118.36	123.19
2	A	501	HEM	CMA-C3A-C4A	-2.07	125.29	128.46
2	A	501	HEM	CAA-CBA-CGA	-2.01	109.29	112.67
3	A	502	FIV	C4-C1-C2	2.00	122.51	118.92

There are no chirality outliers.

All (2) torsion outliers are listed below:

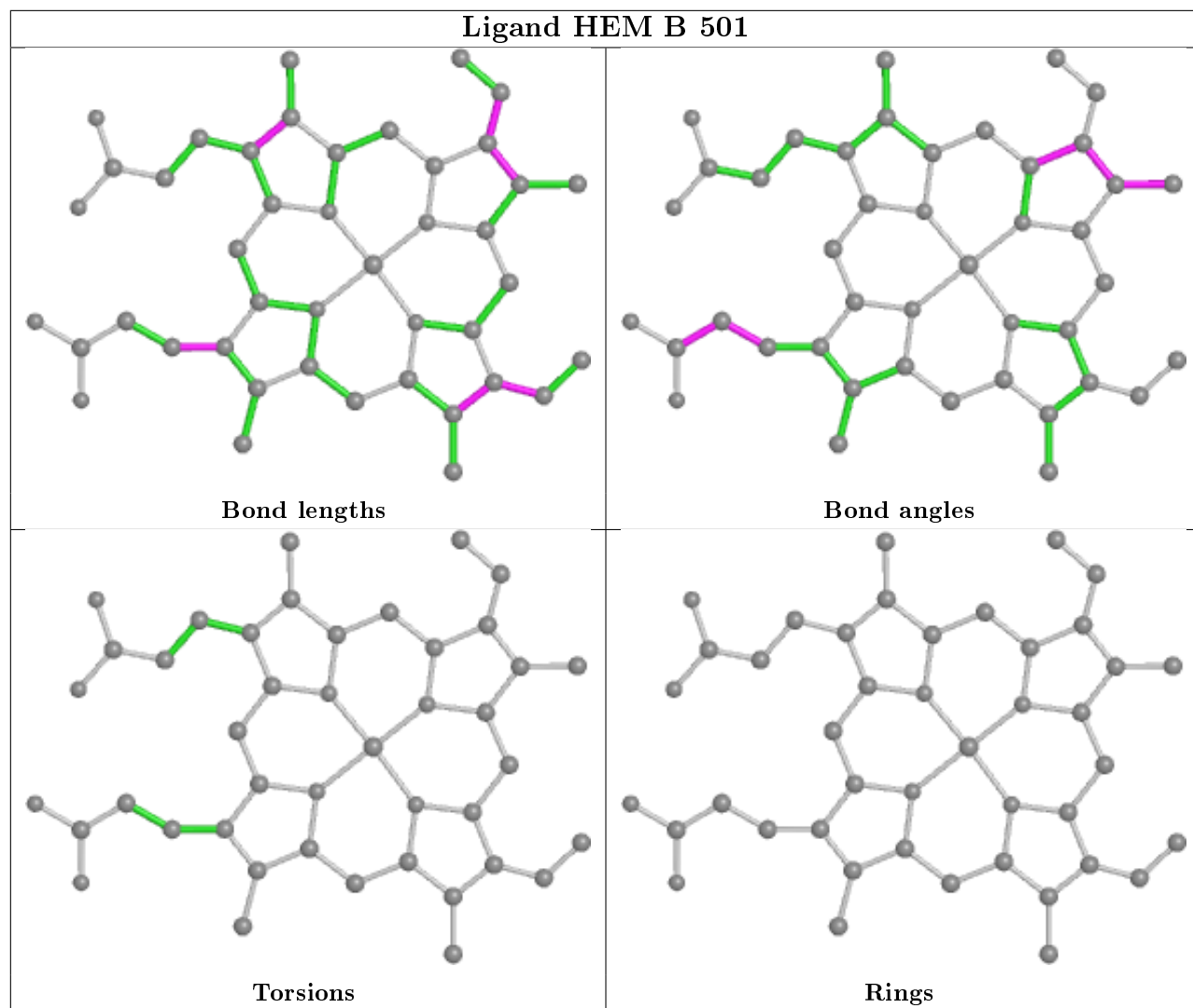
Mol	Chain	Res	Type	Atoms
5	A	506	GOL	O1-C1-C2-C3
5	A	506	GOL	O1-C1-C2-O2

There are no ring outliers.

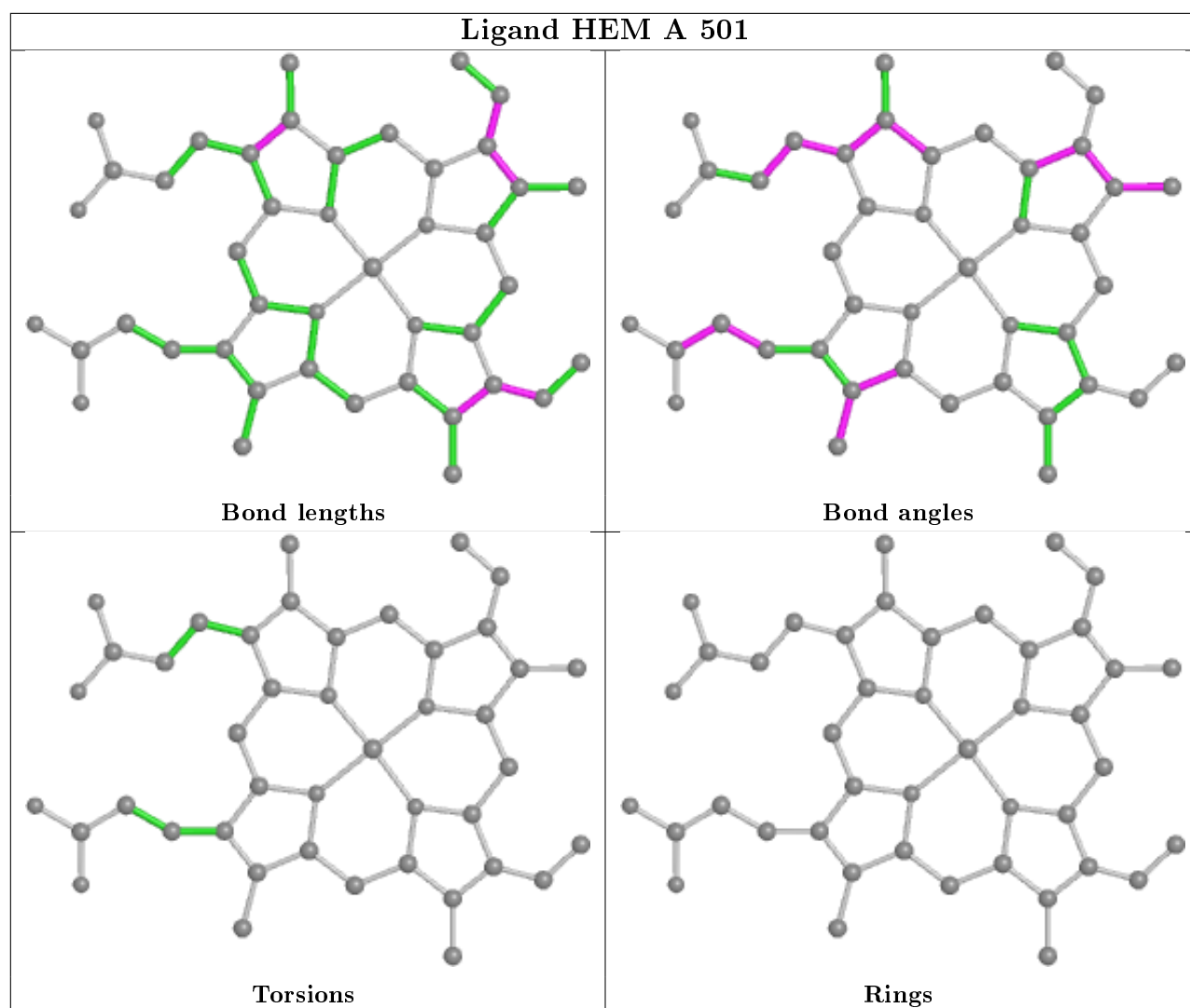
5 monomers are involved in 21 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	503	FIV	1	0
2	B	501	HEM	8	0
3	A	502	FIV	6	0
3	B	502	FIV	4	0
2	A	501	HEM	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.







## 5.7 Other polymers [i](#)

There are no such residues in this entry.

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

There are no chain breaks in this entry.

## 6 Fit of model and data [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	393/410 (95%)	-0.34	1 (0%) 94 84	34, 51, 69, 80	0
1	B	393/410 (95%)	0.02	7 (1%) 68 40	30, 69, 100, 110	0
All	All	786/820 (95%)	-0.16	8 (1%) 82 59	30, 57, 97, 110	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	198	ARG	3.1
1	B	199	SER	2.8
1	B	280	LEU	2.8
1	A	278	PRO	2.7
1	B	210	CYS	2.5
1	B	196	ILE	2.4
1	B	262	TYR	2.1
1	B	275	ARG	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.

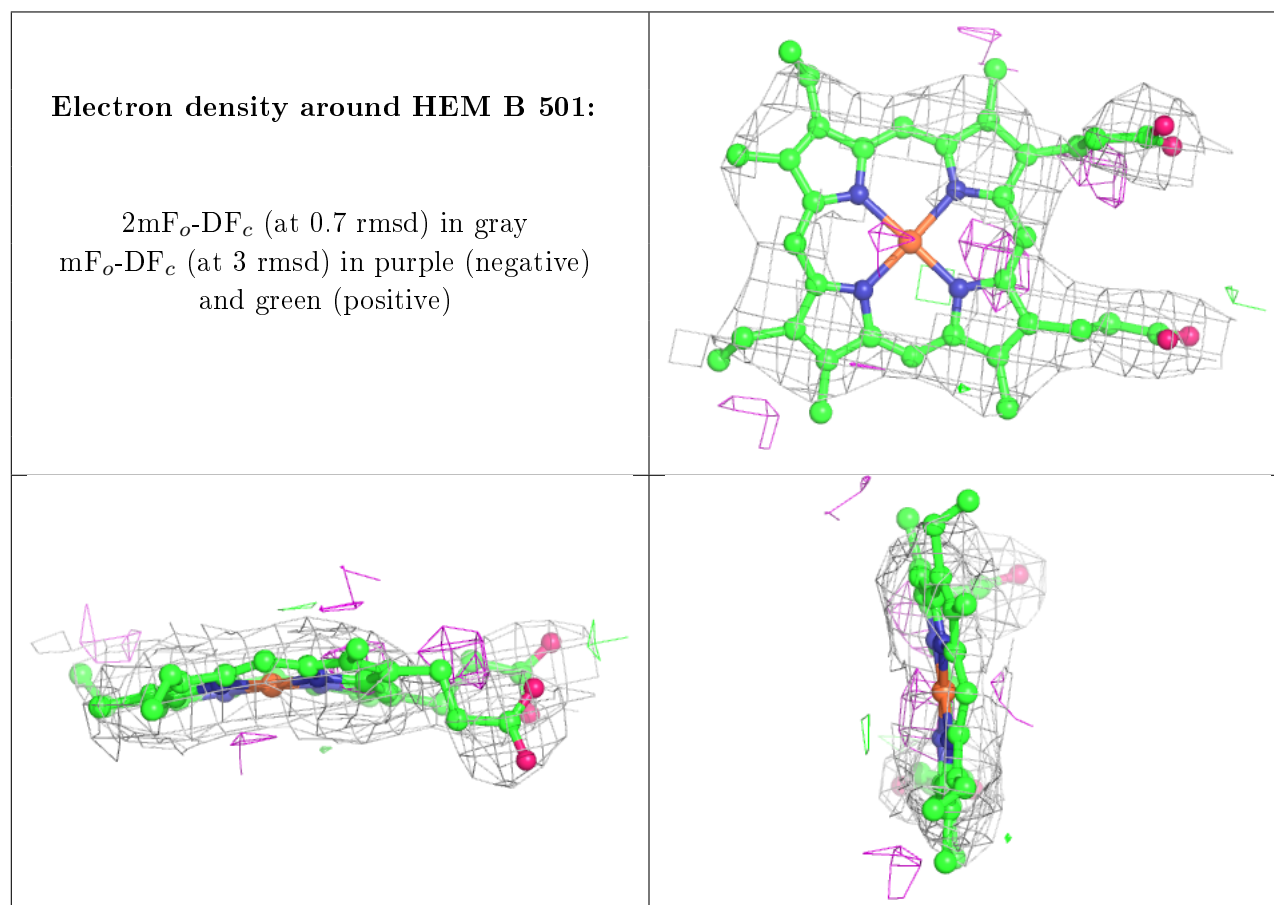
### 6.4 Ligands [i](#)

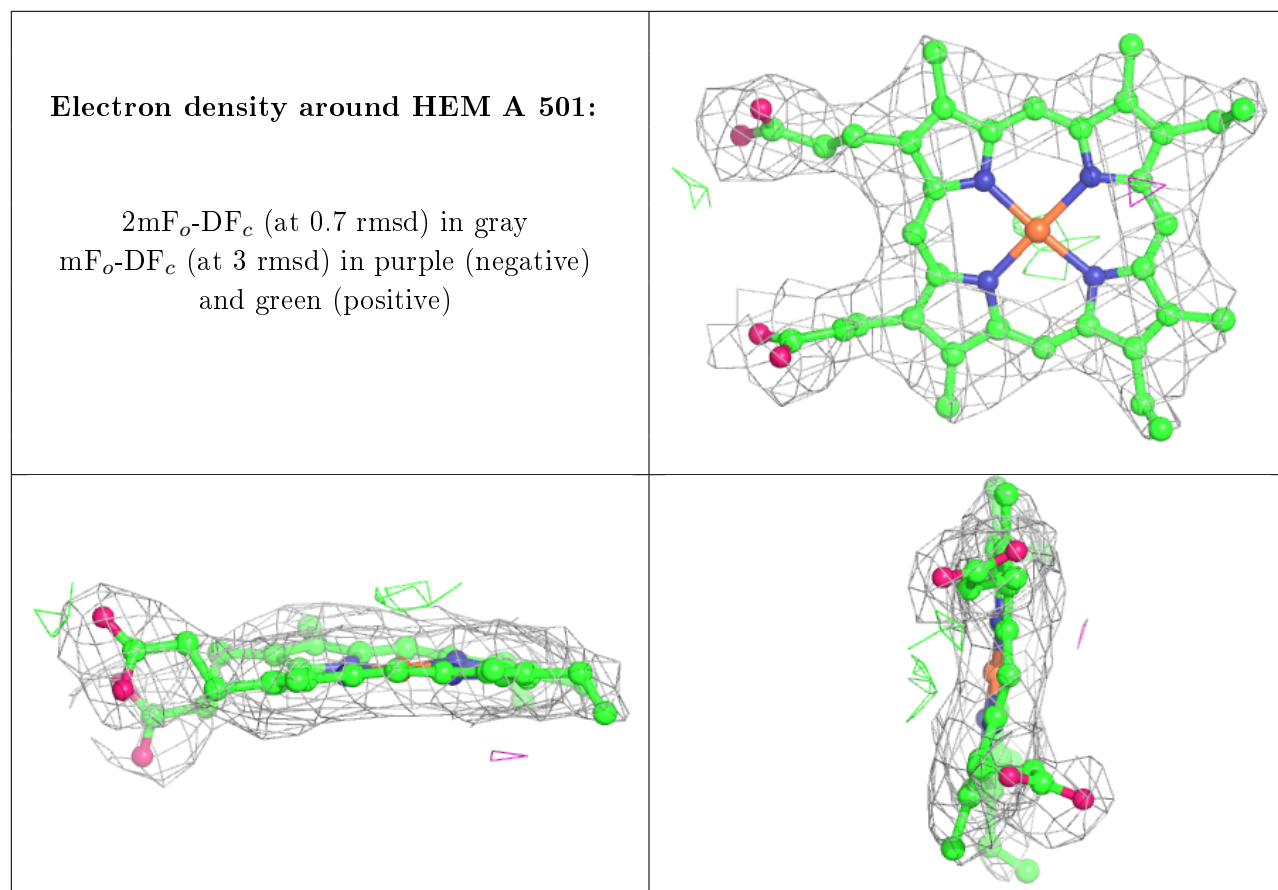
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
3	FIV	A	503	13/13	0.75	0.36	100,103,105,106	0
6	CL	A	507	1/1	0.86	0.17	71,71,71,71	0
3	FIV	B	502	13/13	0.88	0.40	54,56,58,58	0
5	GOL	A	506	6/6	0.94	0.21	58,61,62,65	0
4	SO4	A	505	5/5	0.95	0.16	77,79,79,79	0
4	SO4	A	504	5/5	0.95	0.12	90,90,90,90	0
4	SO4	B	503	5/5	0.95	0.14	80,80,80,81	0
2	HEM	B	501	43/43	0.96	0.26	49,56,59,60	0
3	FIV	A	502	13/13	0.96	0.23	45,48,55,56	0
2	HEM	A	501	43/43	0.98	0.19	31,36,41,43	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.





## 6.5 Other polymers [i](#)

There are no such residues in this entry.