



# Full wwPDB X-ray Structure Validation Report ⓘ

May 21, 2020 – 02:10 pm BST

PDB ID : 4GQS  
Title : Structure of Human Microsomal Cytochrome P450 (CYP) 2C19  
Authors : Reynald, R.L.; Sansen, S.; Stout, C.D.; Johnson, E.F.  
Deposited on : 2012-08-23  
Resolution : 2.87 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

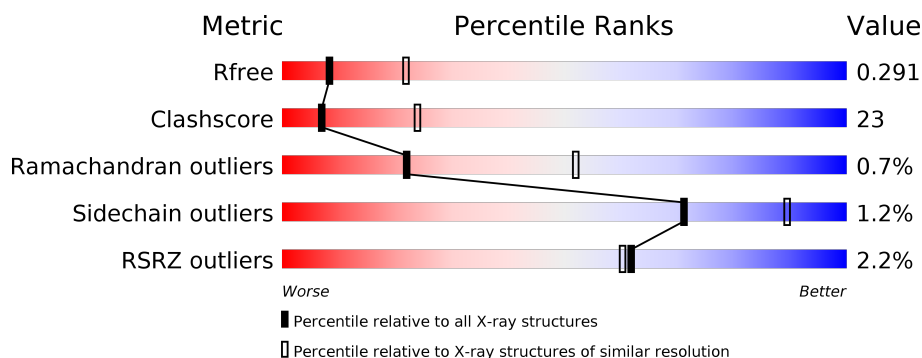
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.87 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	2691 (2.90-2.86)
Clashscore	141614	2947 (2.90-2.86)
Ramachandran outliers	138981	2868 (2.90-2.86)
Sidechain outliers	138945	2871 (2.90-2.86)
RSRZ outliers	127900	2629 (2.90-2.86)

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	477	<div> <div style="width: 61%; background-color: green;"></div> <div style="width: 35%; background-color: yellow;"></div> <div style="width: 4%; background-color: orange;"></div> <div style="width: 2%; background-color: red;"></div> <div style="width: 2%; background-color: grey;"></div> </div> <div>61% 35% . .</div>
1	B	477	<div> <div style="width: 51%; background-color: green;"></div> <div style="width: 42%; background-color: yellow;"></div> <div style="width: 4%; background-color: orange;"></div> <div style="width: 2%; background-color: red;"></div> <div style="width: 2%; background-color: grey;"></div> </div> <div>2% 51% 42% . .</div>
1	C	477	<div> <div style="width: 54%; background-color: green;"></div> <div style="width: 42%; background-color: yellow;"></div> <div style="width: 4%; background-color: orange;"></div> <div style="width: 2%; background-color: red;"></div> <div style="width: 2%; background-color: grey;"></div> </div> <div>2% 54% 42% . .</div>
1	D	477	<div> <div style="width: 53%; background-color: green;"></div> <div style="width: 42%; background-color: yellow;"></div> <div style="width: 5%; background-color: orange;"></div> <div style="width: 2%; background-color: red;"></div> <div style="width: 2%; background-color: grey;"></div> </div> <div>5% 53% 42% 5%</div>

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 15043 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 2C19.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	464	Total	C	N	O	S	0	0	0
			3735	2400	635	677	23			
1	B	456	Total	C	N	O	S	0	0	0
			3666	2354	623	666	23			
1	C	462	Total	C	N	O	S	0	0	0
			3707	2384	627	673	23			
1	D	452	Total	C	N	O	S	0	0	0
			3611	2319	611	659	22			

There are 40 discrepancies between the modelled and reference sequences:

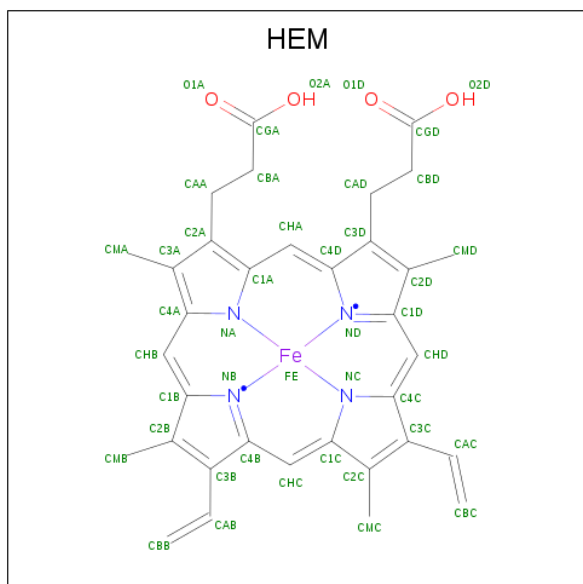
Chain	Residue	Modelled	Actual	Comment	Reference
A	18	MET	-	EXPRESSION TAG	UNP P33261
A	19	ALA	-	EXPRESSION TAG	UNP P33261
A	20	LYS	-	EXPRESSION TAG	UNP P33261
A	21	LYS	-	EXPRESSION TAG	UNP P33261
A	22	THR	-	EXPRESSION TAG	UNP P33261
A	490	ILE	-	EXPRESSION TAG	UNP P33261
A	491	HIS	-	EXPRESSION TAG	UNP P33261
A	492	HIS	-	EXPRESSION TAG	UNP P33261
A	493	HIS	-	EXPRESSION TAG	UNP P33261
A	494	HIS	-	EXPRESSION TAG	UNP P33261
B	18	MET	-	EXPRESSION TAG	UNP P33261
B	19	ALA	-	EXPRESSION TAG	UNP P33261
B	20	LYS	-	EXPRESSION TAG	UNP P33261
B	21	LYS	-	EXPRESSION TAG	UNP P33261
B	22	THR	-	EXPRESSION TAG	UNP P33261
B	490	ILE	-	EXPRESSION TAG	UNP P33261
B	491	HIS	-	EXPRESSION TAG	UNP P33261
B	492	HIS	-	EXPRESSION TAG	UNP P33261
B	493	HIS	-	EXPRESSION TAG	UNP P33261
B	494	HIS	-	EXPRESSION TAG	UNP P33261
C	18	MET	-	EXPRESSION TAG	UNP P33261

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Chain	Residue	Modelled	Actual	Comment	Reference
C	19	ALA	-	EXPRESSION TAG	UNP P33261
C	20	LYS	-	EXPRESSION TAG	UNP P33261
C	21	LYS	-	EXPRESSION TAG	UNP P33261
C	22	THR	-	EXPRESSION TAG	UNP P33261
C	490	ILE	-	EXPRESSION TAG	UNP P33261
C	491	HIS	-	EXPRESSION TAG	UNP P33261
C	492	HIS	-	EXPRESSION TAG	UNP P33261
C	493	HIS	-	EXPRESSION TAG	UNP P33261
C	494	HIS	-	EXPRESSION TAG	UNP P33261
D	18	MET	-	EXPRESSION TAG	UNP P33261
D	19	ALA	-	EXPRESSION TAG	UNP P33261
D	20	LYS	-	EXPRESSION TAG	UNP P33261
D	21	LYS	-	EXPRESSION TAG	UNP P33261
D	22	THR	-	EXPRESSION TAG	UNP P33261
D	490	ILE	-	EXPRESSION TAG	UNP P33261
D	491	HIS	-	EXPRESSION TAG	UNP P33261
D	492	HIS	-	EXPRESSION TAG	UNP P33261
D	493	HIS	-	EXPRESSION TAG	UNP P33261
D	494	HIS	-	EXPRESSION TAG	UNP P33261

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



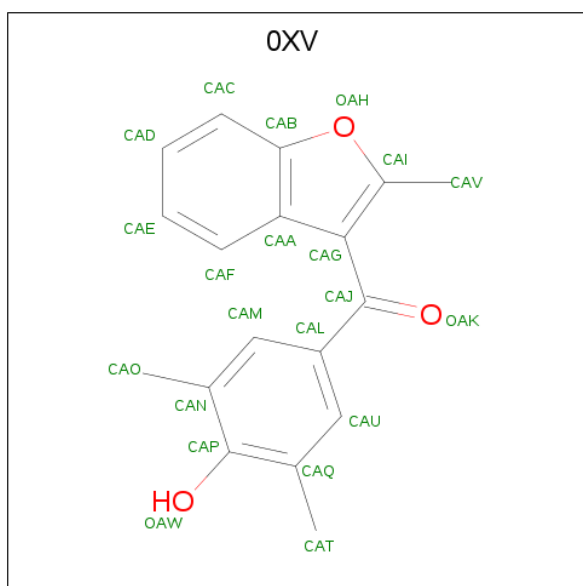
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

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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		
2	C	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	D	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- Molecule 3 is (4-hydroxy-3,5-dimethylphenyl)(2-methyl-1-benzofuran-3-yl)methanone (three-letter code: 0XV) (formula: C<sub>18</sub>H<sub>16</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			21	18	3		
3	B	1	Total	C	O	0	0
			21	18	3		
3	C	1	Total	C	O	0	0
			21	18	3		
3	D	1	Total	C	O	0	0
			21	18	3		

- Molecule 4 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
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		

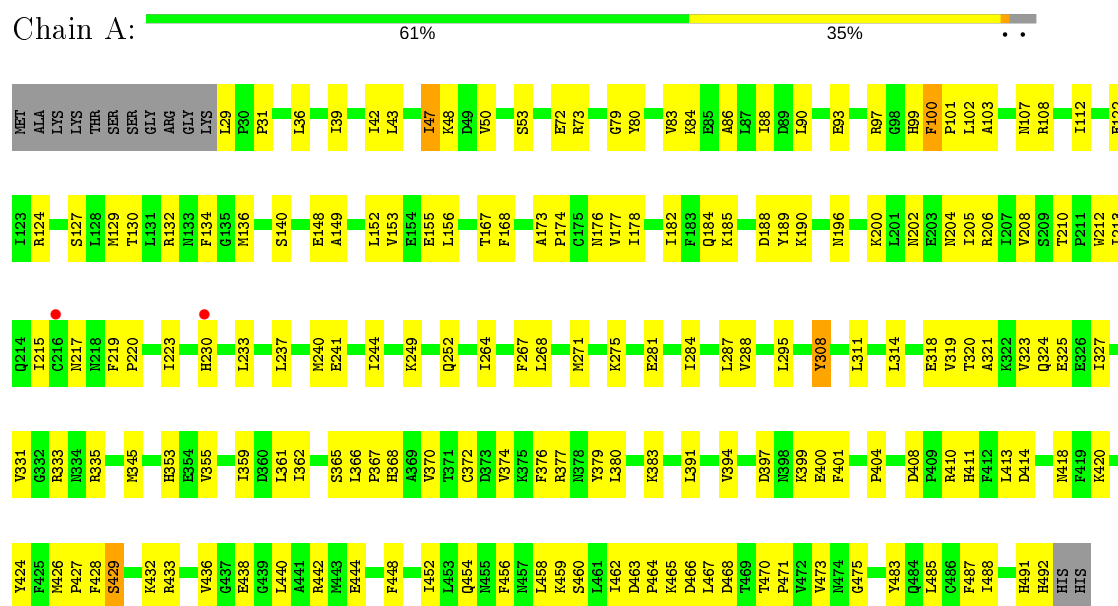
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	18	Total	O	0	0
			18	18		
5	B	13	Total	O	0	0
			13	13		
5	C	12	Total	O	0	0
			12	12		
5	D	7	Total	O	0	0
			7	7		

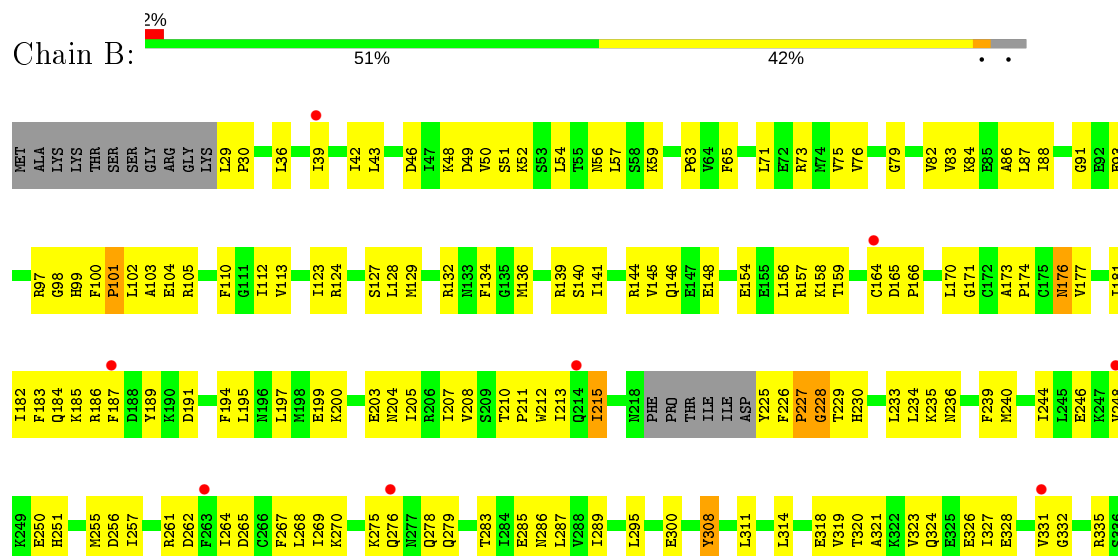
### 3 Residue-property plots

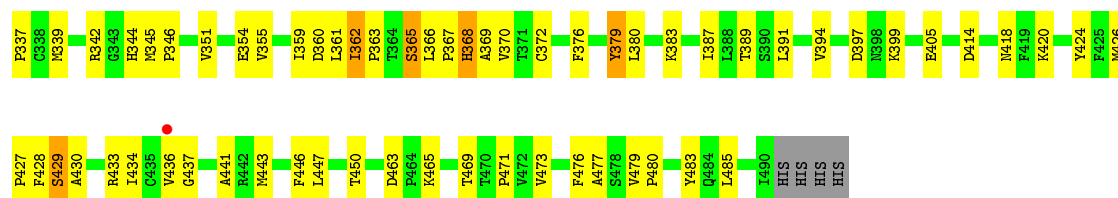
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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 2C19

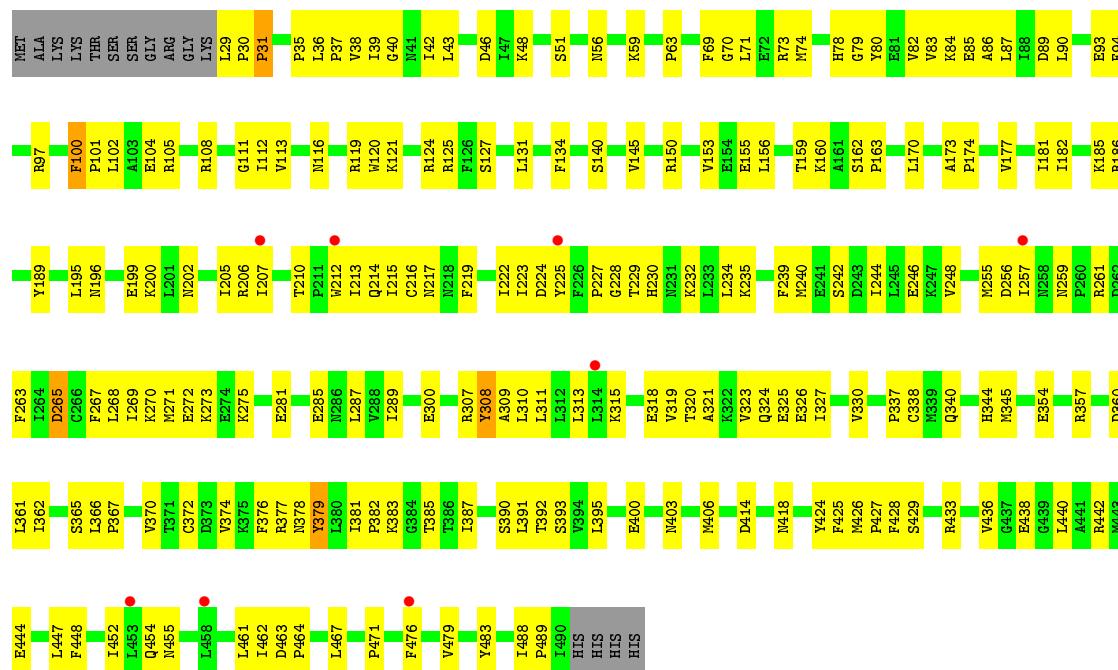


#### • Molecule 1: Cytochrome P450 2C19

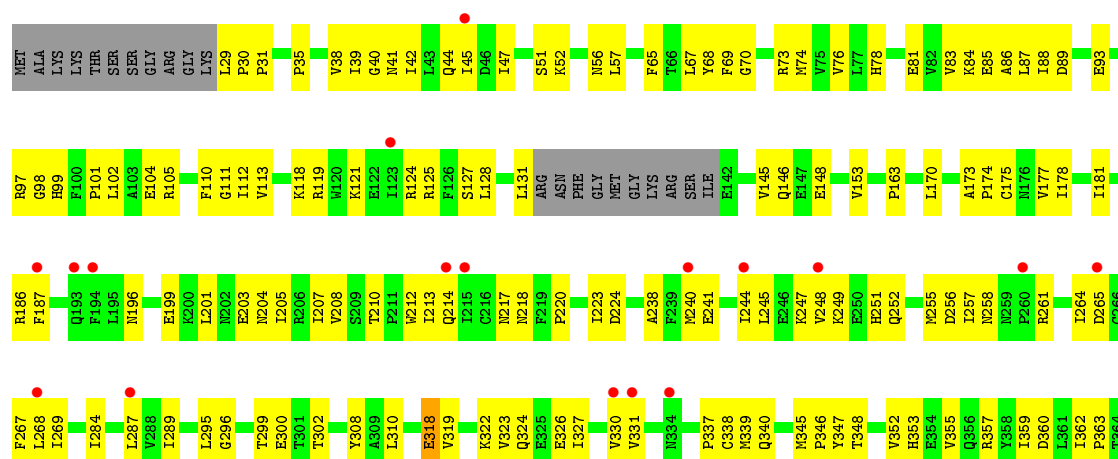




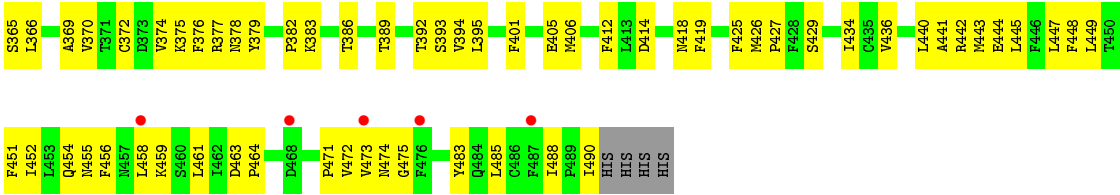
• Molecule 1: Cytochrome P450 2C19



• Molecule 1: Cytochrome P450 2C19







## 4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	159.19Å 159.19Å 450.03Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	79.60 – 2.87 79.59 – 2.87	Depositor EDS
% Data completeness (in resolution range)	99.7 (79.60-2.87) 99.9 (79.59-2.87)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.77 (at 2.86Å)	Xtriage
Refinement program	CNS 1.3	Depositor
R, $R_{free}$	0.250 , 0.296 0.245 , 0.291	Depositor DCC
$R_{free}$ test set	2574 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	82.6	Xtriage
Anisotropy	0.036	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 74.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	15043	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	85.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 17.95% 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, 0XV, HEM

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	0.26	0/3822	0.48	0/5172
1	B	0.26	0/3748	0.46	0/5068
1	C	0.27	0/3792	0.46	0/5132
1	D	0.26	0/3692	0.46	0/4999
All	All	0.26	0/15054	0.46	0/20371

There are no bond length outliers.

There are no bond angle outliers.

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	3735	0	3761	158	0
1	B	3666	0	3697	193	0
1	C	3707	0	3735	171	0
1	D	3611	0	3622	169	0
2	A	43	0	30	1	0
2	B	43	0	30	4	0
2	C	43	0	30	0	0
2	D	43	0	30	3	0
3	A	21	0	16	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	21	0	16	2	0
3	C	21	0	16	1	0
3	D	21	0	16	1	0
4	A	18	0	24	3	0
5	A	18	0	0	2	0
5	B	13	0	0	2	0
5	C	12	0	0	3	0
5	D	7	0	0	1	0
All	All	15043	0	15023	688	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 23.

All (688) 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:84:LYS:HG3	1:A:88:ILE:HD12	1.36	1.06
1:D:163:PRO:HB3	1:D:461:LEU:HD21	1.42	0.98
1:C:63:PRO:HB2	1:C:82:VAL:HG21	1.48	0.95
1:B:52:LYS:HE2	5:B:613:HOH:O	1.68	0.92
1:B:200:LYS:HG2	1:B:236:ASN:HD22	1.36	0.90
1:D:87:LEU:HD21	1:D:389:THR:HG21	1.56	0.88
1:A:173:ALA:HB3	1:A:174:PRO:HD3	1.59	0.84
1:C:269:ILE:HG22	1:C:273:LYS:HE3	1.61	0.82
1:C:462:ILE:HG13	1:C:463:ASP:H	1.45	0.82
1:D:45:ILE:HD11	1:D:57:LEU:HD11	1.62	0.81
1:B:173:ALA:HB3	1:B:174:PRO:HD3	1.60	0.81
1:C:310:LEU:HA	1:C:313:LEU:HD12	1.63	0.80
1:A:467:LEU:HD12	1:A:467:LEU:O	1.82	0.80
1:B:257:ILE:HD12	1:B:257:ILE:H	1.47	0.80
1:D:112:ILE:HD11	1:D:436:VAL:HG21	1.64	0.79
1:C:214:GLN:NE2	1:C:476:PHE:H	1.81	0.78
1:A:101:PRO:HG2	1:A:217:ASN:O	1.84	0.78
1:C:170:LEU:HD11	1:C:310:LEU:HD12	1.64	0.78
1:D:324:GLN:HA	1:D:327:ILE:HD12	1.65	0.78
1:D:39:ILE:HB	1:D:42:ILE:HG12	1.65	0.78
1:A:268:LEU:HD23	1:A:271:MET:HE3	1.64	0.77
1:C:46:ASP:OD2	1:C:48:LYS:HB3	1.86	0.76
1:B:285:GLU:HB2	5:B:611:HOH:O	1.87	0.75
1:C:112:ILE:HD11	1:C:436:VAL:HG21	1.69	0.75
1:D:455:ASN:HA	1:D:490:ILE:HG12	1.68	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:255:MET:HG2	1:B:256:ASP:H	1.52	0.75
1:D:148:GLU:HG2	1:D:177:VAL:HG23	1.70	0.74
1:D:39:ILE:O	1:D:70:GLY:HA2	1.89	0.73
1:A:268:LEU:HA	1:A:271:MET:HE2	1.68	0.73
1:B:251:HIS:HA	1:B:261:ARG:HH22	1.54	0.73
1:C:448:PHE:O	1:C:452:ILE:HG12	1.88	0.72
1:A:314:LEU:HD13	1:A:460:SER:HB3	1.70	0.72
1:B:166:PRO:HB2	1:B:170:LEU:HD12	1.71	0.72
1:B:36:LEU:HD22	1:B:39:ILE:HD11	1.70	0.72
1:B:101:PRO:O	1:B:104:GLU:HG2	1.90	0.72
1:C:185:LYS:HG2	1:C:186:ARG:H	1.55	0.72
1:C:479:VAL:CG1	5:C:612:HOH:O	2.37	0.72
1:C:121:LYS:O	1:C:125:ARG:HG2	1.90	0.72
1:D:173:ALA:HB3	1:D:174:PRO:HD3	1.73	0.71
1:C:101:PRO:HG2	1:C:217:ASN:O	1.92	0.70
1:B:283:THR:HG23	1:B:286:ASN:H	1.56	0.70
1:B:71:LEU:HD23	1:B:71:LEU:H	1.57	0.70
1:C:173:ALA:HB3	1:C:174:PRO:HD3	1.74	0.70
1:C:268:LEU:HD23	1:C:271:MET:HE3	1.73	0.70
1:C:56:ASN:HA	1:C:59:LYS:HD2	1.74	0.70
1:B:39:ILE:HD12	1:B:43:LEU:HD13	1.74	0.69
1:C:257:ILE:H	1:C:257:ILE:HD12	1.55	0.69
1:D:47:ILE:HD12	1:D:47:ILE:H	1.57	0.69
1:A:129:MET:HA	1:A:132:ARG:CD	2.22	0.68
1:A:468:ASP:OD2	1:A:470:THR:HB	1.93	0.68
1:A:397:ASP:OD2	1:A:399:LYS:HB3	1.94	0.68
1:C:272:GLU:HA	1:C:275:LYS:HE3	1.76	0.68
1:B:267:PHE:CE2	1:B:287:LEU:HB2	2.29	0.67
1:B:98:GLY:HA2	1:B:369:ALA:HB2	1.75	0.67
1:C:370:VAL:HG13	1:C:387:ILE:HD11	1.75	0.67
1:B:200:LYS:HG2	1:B:236:ASN:ND2	2.08	0.67
1:A:212:TRP:HA	1:A:215:ILE:HD12	1.76	0.67
1:C:270:LYS:HA	1:C:273:LYS:HD2	1.77	0.67
1:A:100:PHE:HE1	1:A:367:PRO:HG2	1.60	0.66
1:A:491:HIS:O	1:A:492:HIS:HB2	1.95	0.66
1:D:251:HIS:HA	1:D:261:ARG:HH21	1.61	0.66
1:A:90:LEU:O	1:A:93:GLU:HG2	1.95	0.66
1:A:215:ILE:HA	4:A:505:GOL:O2	1.95	0.66
1:B:366:LEU:HG	3:B:502:OXV:H2	1.78	0.66
1:D:366:LEU:HG	3:D:502:OXV:H2	1.77	0.65
1:B:311:LEU:HG	1:B:469:THR:HG22	1.76	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:223:ILE:HD11	1:A:230:HIS:HB2	1.78	0.65
1:C:479:VAL:HG13	5:C:612:HOH:O	1.95	0.65
1:D:146:GLN:HG2	1:D:447:LEU:HD13	1.79	0.65
1:B:197:LEU:HD11	1:B:295:LEU:HD21	1.77	0.65
1:D:110:PHE:HA	1:D:119:ARG:NH1	2.11	0.65
1:D:39:ILE:HD12	1:D:42:ILE:HD11	1.78	0.65
1:B:204:ASN:O	1:B:207:ILE:HG22	1.96	0.65
1:D:173:ALA:O	1:D:177:VAL:HG23	1.97	0.64
1:A:365:SER:HB2	1:A:391:LEU:HD11	1.80	0.64
1:D:178:ILE:HG21	1:D:295:LEU:HA	1.80	0.64
1:D:326:GLU:O	1:D:330:VAL:HG22	1.98	0.64
1:D:426:MET:N	1:D:427:PRO:HD3	2.12	0.64
1:C:202:ASN:O	1:C:206:ARG:HG3	1.98	0.64
1:C:102:LEU:HD11	1:C:213:ILE:HG23	1.80	0.63
1:C:29:LEU:HD11	1:C:73:ARG:CZ	2.28	0.63
1:B:351:VAL:O	1:B:355:VAL:HG23	1.99	0.63
1:A:473:VAL:HG12	1:A:475:GLY:H	1.62	0.63
1:D:207:ILE:HG22	1:D:213:ILE:HD12	1.80	0.63
1:D:459:LYS:NZ	1:D:459:LYS:HA	2.13	0.63
1:B:46:ASP:OD2	1:B:48:LYS:HG2	1.99	0.63
1:C:426:MET:N	1:C:427:PRO:HD3	2.13	0.63
1:D:214:GLN:HE22	1:D:475:GLY:HA2	1.62	0.63
1:C:214:GLN:HE22	1:C:476:PHE:H	1.45	0.63
1:D:319:VAL:O	1:D:323:VAL:HG23	1.99	0.63
1:C:267:PHE:CE2	1:C:287:LEU:HB2	2.34	0.63
1:D:203:GLU:O	1:D:207:ILE:HG12	1.99	0.63
1:D:98:GLY:HA2	1:D:369:ALA:HB2	1.81	0.62
1:A:93:GLU:O	1:A:370:VAL:HA	1.98	0.62
1:B:50:VAL:HG13	1:B:51:SER:H	1.64	0.62
1:C:39:ILE:HB	1:C:42:ILE:HG12	1.81	0.62
1:D:323:VAL:O	1:D:327:ILE:HG13	1.99	0.62
1:D:458:LEU:HB3	1:D:485:LEU:HD12	1.81	0.62
1:B:157:ARG:HH11	1:B:157:ARG:HG2	1.63	0.62
1:D:148:GLU:HA	1:D:148:GLU:OE2	1.98	0.62
1:D:256:ASP:OD1	1:D:258:ASN:HB2	1.99	0.62
1:B:211:PRO:O	1:B:215:ILE:HG12	1.98	0.62
1:A:458:LEU:HD22	1:A:485:LEU:HD12	1.80	0.62
1:A:361:LEU:C	1:A:362:ILE:HD13	2.21	0.62
1:A:99:HIS:CE1	5:A:614:HOH:O	2.51	0.62
1:B:141:ILE:O	1:B:145:VAL:HG23	1.99	0.62
1:D:267:PHE:CG	1:D:287:LEU:HD13	2.35	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:188:ASP:OD2	1:A:190:LYS:HB2	2.00	0.61
1:B:323:VAL:O	1:B:327:ILE:HG13	2.00	0.61
1:A:155:GLU:HG3	1:A:189:TYR:CE1	2.35	0.61
1:B:204:ASN:OD1	1:B:233:LEU:HD12	2.00	0.61
1:B:255:MET:HG2	1:B:256:ASP:N	2.14	0.61
1:A:333:ARG:HG3	1:A:333:ARG:HH11	1.66	0.61
1:B:397:ASP:OD2	1:B:399:LYS:HB2	2.01	0.61
1:B:63:PRO:HB2	1:B:82:VAL:HG11	1.81	0.61
1:A:102:LEU:HD11	1:A:233:LEU:CD2	2.31	0.61
1:A:112:ILE:HD11	1:A:127:SER:OG	2.00	0.61
1:C:224:ASP:O	1:C:227:PRO:HD3	2.01	0.61
1:D:87:LEU:HD21	1:D:389:THR:CG2	2.29	0.61
1:D:348:THR:O	1:D:352:VAL:HG23	2.01	0.60
1:D:472:VAL:HG12	1:D:474:ASN:H	1.65	0.60
1:A:102:LEU:HD11	1:A:233:LEU:HD23	1.84	0.60
1:D:318:GLU:OE2	1:D:319:VAL:HG23	1.99	0.60
1:B:251:HIS:HA	1:B:261:ARG:NH2	2.16	0.60
1:A:36:LEU:HD21	1:C:43:LEU:HD22	1.84	0.60
1:C:155:GLU:HG3	1:C:189:TYR:CD1	2.37	0.60
1:A:134:PHE:CE2	1:A:140:SER:HB3	2.36	0.60
1:B:148:GLU:OE2	1:B:186:ARG:HD2	2.01	0.60
1:B:362:ILE:HD12	1:B:362:ILE:H	1.67	0.60
1:C:185:LYS:HG2	1:C:186:ARG:N	2.16	0.60
1:A:426:MET:N	1:A:427:PRO:HD3	2.17	0.59
1:D:121:LYS:HD3	1:D:125:ARG:NH2	2.17	0.59
1:D:39:ILE:HB	1:D:42:ILE:CG1	2.31	0.59
1:A:267:PHE:CG	1:A:287:LEU:HD13	2.37	0.59
1:A:308:TYR:OH	1:A:471:PRO:HG3	2.02	0.59
1:B:207:ILE:HG21	1:B:233:LEU:HB2	1.85	0.59
1:C:464:PRO:HA	1:C:467:LEU:HD13	1.85	0.59
1:D:101:PRO:HG2	1:D:217:ASN:O	2.03	0.59
1:B:139:ARG:HH21	1:B:144:ARG:HA	1.68	0.59
1:B:128:LEU:HD12	1:B:436:VAL:HG13	1.85	0.59
1:B:87:LEU:HD11	1:B:368:HIS:CE1	2.38	0.59
1:A:167:THR:HG23	1:A:168:PHE:N	2.18	0.59
1:C:79:GLY:O	1:C:83:VAL:HG23	2.03	0.59
1:C:462:ILE:HG13	1:C:463:ASP:N	2.15	0.58
1:D:105:ARG:HH22	1:D:224:ASP:N	2.01	0.58
1:A:90:LEU:HB3	1:A:93:GLU:CG	2.34	0.58
1:C:156:LEU:HD11	1:C:452:ILE:HD12	1.84	0.58
1:C:210:THR:O	1:C:214:GLN:HG3	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:321:ALA:O	1:A:325:GLU:HG3	2.04	0.58
1:B:244:ILE:O	1:B:248:VAL:HG23	2.02	0.58
1:C:100:PHE:O	1:C:104:GLU:HG2	2.02	0.58
1:C:156:LEU:O	1:C:159:THR:HG22	2.03	0.58
1:B:171:GLY:O	1:B:174:PRO:HD2	2.03	0.58
1:C:155:GLU:HG3	1:C:189:TYR:CE1	2.38	0.58
1:A:323:VAL:O	1:A:327:ILE:HG13	2.04	0.58
1:C:36:LEU:HB3	1:C:37:PRO:HD2	1.85	0.58
1:C:223:ILE:HD11	1:C:230:HIS:ND1	2.19	0.58
1:A:129:MET:HA	1:A:132:ARG:HD2	1.84	0.57
1:D:148:GLU:HG2	1:D:177:VAL:CG2	2.33	0.57
1:A:36:LEU:CD2	1:A:39:ILE:HD11	2.33	0.57
1:D:39:ILE:O	1:D:70:GLY:CA	2.51	0.57
1:D:339:MET:SD	1:D:443:MET:HG3	2.44	0.57
1:D:97:ARG:NH1	1:D:366:LEU:HB3	2.19	0.57
1:B:365:SER:HB2	1:B:391:LEU:CD1	2.34	0.57
1:C:101:PRO:HD2	1:C:217:ASN:ND2	2.20	0.57
1:D:459:LYS:HA	1:D:459:LYS:HZ3	1.68	0.57
1:D:74:MET:HB3	1:D:386:THR:HB	1.86	0.57
1:B:191:ASP:O	1:B:195:LEU:HG	2.05	0.57
1:C:244:ILE:O	1:C:248:VAL:HG23	2.04	0.57
1:C:381:ILE:HD12	1:C:381:ILE:N	2.20	0.57
1:A:414:ASP:OD2	1:A:418:ASN:HB2	2.05	0.56
1:C:173:ALA:O	1:C:177:VAL:HG23	2.04	0.56
1:C:36:LEU:HB3	1:C:37:PRO:CD	2.35	0.56
1:A:249:LYS:HA	1:A:252:GLN:OE1	2.05	0.56
1:C:267:PHE:CG	1:C:287:LEU:HD13	2.40	0.56
1:D:127:SER:O	1:D:131:LEU:HD13	2.06	0.56
1:B:227:PRO:HG2	1:B:228:GLY:H	1.70	0.56
1:B:36:LEU:CD2	1:B:39:ILE:HD11	2.36	0.56
1:D:170:LEU:HD11	1:D:310:LEU:HD12	1.87	0.56
1:A:100:PHE:CE1	1:A:367:PRO:HG2	2.39	0.56
1:D:454:GLN:O	1:D:490:ILE:HD13	2.06	0.56
1:C:177:VAL:O	1:C:181:ILE:HG12	2.05	0.56
1:C:182:ILE:HA	1:C:263:PHE:HB3	1.87	0.56
1:A:240:MET:O	1:A:244:ILE:HG13	2.06	0.56
1:A:206:ARG:O	1:A:210:THR:HG23	2.05	0.56
1:B:189:TYR:O	1:B:195:LEU:HD21	2.06	0.56
1:C:196:ASN:HA	1:C:199:GLU:OE1	2.06	0.56
1:A:129:MET:O	1:A:132:ARG:HD3	2.06	0.56
1:B:365:SER:HB2	1:B:391:LEU:HD11	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:127:SER:O	1:C:131:LEU:HG	2.06	0.56
1:D:84:LYS:CG	1:D:88:ILE:HD12	2.36	0.56
1:B:104:GLU:HG3	1:B:105:ARG:N	2.21	0.56
1:D:441:ALA:O	1:D:445:LEU:HD23	2.06	0.56
1:B:187:PHE:CD1	1:B:194:PHE:HB2	2.41	0.55
1:B:240:MET:O	1:B:244:ILE:HG13	2.06	0.55
1:C:80:TYR:CZ	1:C:84:LYS:HD2	2.41	0.55
1:D:377:ARG:O	1:D:378:ASN:HB3	2.06	0.55
1:B:166:PRO:HB2	1:B:170:LEU:CD1	2.36	0.55
1:B:318:GLU:HG2	1:B:319:VAL:N	2.21	0.55
1:D:267:PHE:CE2	1:D:287:LEU:HB2	2.41	0.55
1:B:63:PRO:HB2	1:B:82:VAL:CG1	2.37	0.55
1:A:223:ILE:HD13	1:A:230:HIS:ND1	2.21	0.55
1:B:314:LEU:HD11	1:B:485:LEU:HD13	1.88	0.55
1:C:112:ILE:HG23	1:C:113:VAL:N	2.21	0.55
1:A:284:ILE:O	1:A:288:VAL:HG23	2.07	0.55
1:B:100:PHE:HB3	1:B:103:ALA:HB3	1.89	0.55
1:D:252:GLN:HA	1:D:268:LEU:HD21	1.88	0.55
1:D:173:ALA:HB1	1:D:448:PHE:CE2	2.42	0.55
1:D:39:ILE:HD12	1:D:42:ILE:CD1	2.36	0.55
1:D:73:ARG:HH11	1:D:73:ARG:HB3	1.72	0.55
1:C:150:ARG:O	1:C:153:VAL:HG12	2.07	0.55
1:C:200:LYS:HE3	1:C:240:MET:HB2	1.89	0.55
1:B:134:PHE:CZ	1:B:140:SER:HB2	2.42	0.54
1:B:331:VAL:HG21	1:B:337:PRO:HB3	1.89	0.54
1:B:405:GLU:OE2	1:B:405:GLU:HA	2.08	0.54
1:D:175:CYS:HB2	1:D:299:THR:HG23	1.89	0.54
1:A:130:THR:HG22	1:A:136:MET:HE2	1.89	0.54
1:A:353:HIS:HE1	1:A:442:ARG:HH22	1.55	0.54
1:B:372:CYS:HA	1:B:383:LYS:HB2	1.88	0.54
1:D:44:GLN:NE2	5:D:607:HOH:O	2.41	0.54
1:B:124:ARG:O	1:B:128:LEU:HD13	2.08	0.54
1:B:187:PHE:HB3	1:B:194:PHE:HB2	1.90	0.54
1:A:413:LEU:O	1:A:420:LYS:HE2	2.08	0.54
1:A:90:LEU:HB3	1:A:93:GLU:HG2	1.90	0.54
1:C:255:MET:HE1	1:C:257:ILE:HG13	1.90	0.54
1:D:452:ILE:O	1:D:456:PHE:HB2	2.07	0.54
1:A:267:PHE:CE2	1:A:287:LEU:HB2	2.42	0.54
1:B:362:ILE:HD12	1:B:362:ILE:N	2.23	0.54
1:B:427:PRO:HG2	1:B:428:PHE:CD2	2.42	0.54
1:C:71:LEU:O	1:C:71:LEU:HD23	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:436:VAL:O	1:D:436:VAL:HG12	2.07	0.54
1:B:368:HIS:NE2	1:B:433:ARG:HB2	2.23	0.54
1:C:414:ASP:OD2	1:C:418:ASN:HB2	2.08	0.54
1:A:408:ASP:OD1	1:A:410:ARG:HD2	2.08	0.53
1:D:153:VAL:CG1	1:D:455:ASN:HD21	2.21	0.53
1:A:366:LEU:HG	3:A:502:OXV:H2	1.90	0.53
1:C:361:LEU:HB3	1:C:362:ILE:HD12	1.89	0.53
1:D:29:LEU:HD11	1:D:73:ARG:HE	1.73	0.53
1:A:122:GLU:HG3	1:A:281:GLU:HG3	1.91	0.53
1:B:157:ARG:HG2	1:B:157:ARG:NH1	2.24	0.53
1:C:265:ASP:O	1:C:269:ILE:HG13	2.07	0.53
1:D:186:ARG:HG2	1:D:187:PHE:N	2.23	0.53
1:A:394:VAL:HG21	1:A:426:MET:HE1	1.90	0.53
1:B:49:ASP:OD1	1:B:473:VAL:HG11	2.09	0.53
1:C:338:CYS:SG	1:C:340:GLN:HB2	2.48	0.53
1:C:376:PHE:O	1:C:377:ARG:HG3	2.09	0.53
1:C:51:SER:OG	1:C:392:THR:HG21	2.09	0.53
1:D:196:ASN:O	1:D:199:GLU:HB3	2.09	0.53
1:A:372:CYS:HA	1:A:383:LYS:HG3	1.89	0.53
1:A:353:HIS:CE1	1:A:442:ARG:HH22	2.26	0.53
1:B:355:VAL:O	1:B:359:ILE:HG12	2.09	0.53
1:C:131:LEU:HD12	1:C:436:VAL:HG12	1.90	0.53
1:C:156:LEU:CD1	1:C:452:ILE:HD12	2.39	0.53
1:A:361:LEU:O	1:A:362:ILE:HD13	2.09	0.52
1:A:79:GLY:O	1:A:83:VAL:HG23	2.09	0.52
1:D:355:VAL:O	1:D:359:ILE:HG23	2.08	0.52
1:A:167:THR:CG2	1:A:168:PHE:N	2.73	0.52
1:B:84:LYS:O	1:B:88:ILE:HB	2.09	0.52
1:C:365:SER:HB2	1:C:391:LEU:HD11	1.91	0.52
1:D:442:ARG:HH11	1:D:442:ARG:HG2	1.75	0.52
1:B:476:PHE:CG	1:B:477:ALA:N	2.78	0.52
1:C:134:PHE:CE2	1:C:140:SER:HB3	2.45	0.52
1:A:428:PHE:CD2	1:A:438:GLU:HG3	2.45	0.52
1:A:318:GLU:HG2	1:A:319:VAL:N	2.25	0.52
1:C:285:GLU:O	1:C:289:ILE:HD13	2.09	0.52
1:D:330:VAL:HG23	1:D:331:VAL:HG23	1.92	0.52
1:D:153:VAL:HG13	1:D:456:PHE:HE2	1.75	0.52
1:A:129:MET:SD	1:B:328:GLU:OE1	2.67	0.52
1:B:394:VAL:HG11	1:B:426:MET:CE	2.39	0.52
1:B:65:PHE:CE2	1:B:76:VAL:HB	2.45	0.52
1:C:365:SER:OG	1:C:366:LEU:N	2.42	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:261:ARG:HG3	1:C:261:ARG:HH11	1.75	0.52
1:D:245:LEU:HD11	1:D:284:ILE:HD11	1.91	0.52
1:A:129:MET:HA	1:A:132:ARG:HD3	1.91	0.52
1:B:446:PHE:HD1	1:B:447:LEU:HD23	1.75	0.52
1:D:372:CYS:O	1:D:374:VAL:HG23	2.10	0.52
1:B:134:PHE:CE2	1:B:140:SER:HB2	2.45	0.51
1:B:82:VAL:HG23	1:B:83:VAL:N	2.25	0.51
1:C:100:PHE:CD2	1:C:367:PRO:HG2	2.46	0.51
1:A:220:PRO:O	1:A:223:ILE:HG22	2.11	0.51
1:C:100:PHE:HD2	1:C:367:PRO:HG2	1.75	0.51
1:B:199:GLU:O	1:B:203:GLU:HB2	2.11	0.51
1:A:100:PHE:HB3	1:A:103:ALA:HB3	1.92	0.51
1:A:462:ILE:HD12	1:A:463:ASP:O	2.10	0.51
1:C:86:ALA:HB2	1:C:376:PHE:CE1	2.45	0.51
1:D:463:ASP:OD1	1:D:464:PRO:HD2	2.10	0.51
1:B:139:ARG:NH2	1:B:144:ARG:HA	2.26	0.51
1:B:39:ILE:HB	1:B:42:ILE:CG2	2.41	0.51
1:B:50:VAL:HG13	1:B:51:SER:N	2.25	0.51
1:A:320:THR:O	1:A:324:GLN:HG3	2.10	0.51
1:B:257:ILE:H	1:B:257:ILE:CD1	2.20	0.51
1:B:257:ILE:HD12	1:B:257:ILE:N	2.22	0.51
1:B:361:LEU:HB3	1:B:362:ILE:HD12	1.93	0.51
1:C:223:ILE:HD11	1:C:230:HIS:CG	2.45	0.51
1:D:145:VAL:HG21	1:D:443:MET:SD	2.50	0.51
1:D:445:LEU:O	1:D:449:LEU:HB2	2.10	0.51
1:D:440:LEU:O	1:D:444:GLU:HG3	2.10	0.51
1:D:310:LEU:HB2	1:D:483:TYR:OH	2.11	0.51
1:C:326:GLU:HG2	1:C:345:MET:HG2	1.92	0.50
1:D:318:GLU:O	1:D:322:LYS:HG3	2.10	0.50
1:A:462:ILE:HD11	1:A:467:LEU:HB3	1.92	0.50
1:B:113:VAL:HG21	3:B:502:OXV:H10	1.93	0.50
1:D:395:LEU:HD21	1:D:426:MET:H	1.76	0.50
1:A:80:TYR:OH	1:A:84:LYS:HD3	2.11	0.50
1:B:361:LEU:O	1:B:363:PRO:HD3	2.11	0.50
1:D:330:VAL:HG21	1:D:345:MET:HG2	1.93	0.50
1:D:73:ARG:HH11	1:D:73:ARG:CB	2.24	0.50
1:A:208:VAL:HA	1:A:213:ILE:HG21	1.92	0.50
1:A:237:LEU:HB3	1:A:241:GLU:OE2	2.12	0.50
1:D:205:ILE:HG13	1:D:300:GLU:HG2	1.93	0.50
1:A:130:THR:HG22	1:A:136:MET:CE	2.42	0.50
1:A:202:ASN:O	1:A:205:ILE:HG22	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:275:LYS:HG2	1:B:276:GLN:N	2.26	0.50
1:C:327:ILE:HG12	1:C:345:MET:HE1	1.93	0.50
1:D:73:ARG:NH1	1:D:73:ARG:HB3	2.27	0.50
1:A:448:PHE:O	1:A:452:ILE:HG13	2.11	0.50
1:B:128:LEU:CD1	1:B:436:VAL:HG13	2.41	0.50
1:C:390:SER:OG	1:C:393:SER:HB3	2.12	0.50
1:D:257:ILE:HG13	1:D:258:ASN:N	2.27	0.50
1:A:36:LEU:HD23	1:A:39:ILE:HD11	1.93	0.50
1:D:81:GLU:O	1:D:85:GLU:HG3	2.11	0.50
1:D:84:LYS:HG3	1:D:88:ILE:HD12	1.94	0.50
1:C:195:LEU:O	1:C:199:GLU:HG3	2.12	0.49
1:D:414:ASP:OD1	1:D:418:ASN:N	2.40	0.49
1:B:87:LEU:HD21	1:B:389:THR:HG21	1.94	0.49
1:D:177:VAL:O	1:D:181:ILE:HG13	2.12	0.49
1:C:93:GLU:O	1:C:370:VAL:HA	2.12	0.49
1:B:327:ILE:HA	1:B:345:MET:HE3	1.94	0.49
1:B:29:LEU:HD23	1:B:380:LEU:O	2.13	0.49
1:C:268:LEU:HD23	1:C:271:MET:CE	2.42	0.49
1:C:39:ILE:O	1:C:42:ILE:HG12	2.13	0.49
1:C:207:ILE:O	1:C:210:THR:HG22	2.12	0.49
1:B:283:THR:HG22	1:B:286:ASN:HB2	1.95	0.49
1:A:365:SER:HB2	1:A:391:LEU:CD1	2.43	0.49
1:A:467:LEU:HD12	1:A:467:LEU:C	2.33	0.49
1:B:176:ASN:OD1	1:B:186:ARG:HG2	2.13	0.49
1:A:155:GLU:HG3	1:A:189:TYR:HE1	1.78	0.49
1:B:105:ARG:HH11	1:B:234:LEU:HD22	1.78	0.49
1:B:145:VAL:HG21	1:B:443:MET:SD	2.53	0.49
1:B:428:PHE:O	1:B:429:SER:HB3	2.13	0.49
1:B:93:GLU:N	1:B:93:GLU:OE2	2.46	0.49
1:D:363:PRO:HB3	1:D:392:THR:HG1	1.78	0.49
1:B:269:ILE:HG22	1:B:270:LYS:N	2.28	0.48
1:B:446:PHE:O	1:B:450:THR:HG23	2.13	0.48
1:D:338:CYS:SG	1:D:340:GLN:HB2	2.53	0.48
1:D:394:VAL:HG11	1:D:426:MET:SD	2.52	0.48
1:B:101:PRO:HG2	1:B:102:LEU:H	1.78	0.48
1:B:177:VAL:O	1:B:181:ILE:HG13	2.12	0.48
1:B:29:LEU:HD11	1:B:73:ARG:NH1	2.28	0.48
1:C:177:VAL:HG11	1:C:440:LEU:HD11	1.95	0.48
1:A:264:ILE:HG22	1:A:268:LEU:HD12	1.95	0.48
1:C:163:PRO:HB3	1:C:461:LEU:HD21	1.96	0.48
1:D:244:ILE:O	1:D:248:VAL:HG23	2.12	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:82:VAL:HG23	1:B:83:VAL:H	1.77	0.48
1:C:69:PHE:HD2	1:C:74:MET:SD	2.36	0.48
1:D:363:PRO:HB3	1:D:392:THR:OG1	2.14	0.48
1:D:395:LEU:CD2	1:D:426:MET:H	2.26	0.48
1:A:204:ASN:O	1:A:208:VAL:HG23	2.13	0.48
1:A:42:ILE:HD12	4:A:505:GOL:H11	1.93	0.48
1:B:184:GLN:OE1	1:B:261:ARG:HD2	2.13	0.48
1:C:374:VAL:HG12	1:C:376:PHE:N	2.28	0.48
1:A:97:ARG:HD3	1:A:367:PRO:O	2.14	0.48
1:A:428:PHE:O	1:A:429:SER:HB3	2.13	0.48
1:B:212:TRP:HZ3	1:B:229:THR:HG1	1.61	0.48
1:C:387:ILE:N	1:C:387:ILE:HD12	2.29	0.48
1:D:78:HIS:NE2	1:D:393:SER:HB3	2.28	0.48
1:C:370:VAL:HG13	1:C:387:ILE:CD1	2.40	0.48
1:D:39:ILE:O	1:D:70:GLY:N	2.47	0.48
1:A:365:SER:OG	1:A:366:LEU:N	2.46	0.48
1:B:210:THR:HG22	1:B:212:TRP:H	1.78	0.48
1:D:30:PRO:HG3	1:D:379:TYR:HB3	1.95	0.48
1:D:85:GLU:HA	1:D:89:ASP:OD2	2.13	0.47
1:A:132:ARG:HH22	1:B:321:ALA:HA	1.79	0.47
1:B:79:GLY:O	1:B:83:VAL:HG23	2.13	0.47
1:C:455:ASN:N	1:C:455:ASN:HD22	2.12	0.47
1:D:201:LEU:HD11	1:D:296:GLY:HA2	1.95	0.47
1:C:89:ASP:C	1:C:90:LEU:HD12	2.34	0.47
1:D:110:PHE:HA	1:D:119:ARG:HH12	1.76	0.47
1:D:78:HIS:CD2	1:D:393:SER:HB3	2.49	0.47
1:A:355:VAL:O	1:A:359:ILE:HG12	2.13	0.47
1:A:36:LEU:HD22	1:A:39:ILE:HD11	1.95	0.47
1:B:359:ILE:HG13	1:B:360:ASP:N	2.29	0.47
1:D:105:ARG:HH12	1:D:223:ILE:C	2.18	0.47
1:B:436:VAL:HG23	2:B:501:HEM:HBD2	1.97	0.47
1:D:146:GLN:NE2	1:D:338:CYS:HA	2.29	0.47
1:A:124:ARG:NH1	1:A:436:VAL:HG23	2.29	0.47
1:B:105:ARG:NH1	1:B:234:LEU:HD13	2.30	0.47
1:B:437:GLY:HA3	2:B:501:HEM:HBC2	1.97	0.47
1:C:225:TYR:N	1:C:225:TYR:CD2	2.81	0.47
1:D:124:ARG:HH22	1:D:434:ILE:HG22	1.79	0.47
1:B:75:VAL:HB	1:B:387:ILE:CD1	2.44	0.47
1:C:185:LYS:HG2	1:C:186:ARG:HG2	1.96	0.47
1:C:235:LYS:HE2	1:C:239:PHE:CE2	2.50	0.47
1:C:374:VAL:HG12	1:C:376:PHE:H	1.80	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:105:ARG:HD3	1:B:234:LEU:HD11	1.96	0.47
1:C:97:ARG:NH1	1:C:366:LEU:HB3	2.29	0.47
1:D:247:LYS:HD2	1:D:251:HIS:HE1	1.79	0.47
1:A:333:ARG:NH1	1:A:333:ARG:HG3	2.29	0.47
1:B:394:VAL:HG11	1:B:426:MET:HE3	1.97	0.47
1:D:337:PRO:HD2	1:D:451:PHE:CE1	2.50	0.47
1:D:426:MET:N	1:D:427:PRO:CD	2.77	0.47
1:A:178:ILE:HG22	1:A:182:ILE:HD12	1.96	0.47
1:A:394:VAL:HG21	1:A:426:MET:CE	2.45	0.47
1:B:205:ILE:HG21	1:B:300:GLU:HG2	1.96	0.47
1:B:264:ILE:O	1:B:268:LEU:HG	2.15	0.47
1:C:357:ARG:HA	1:C:395:LEU:HD22	1.96	0.47
1:C:442:ARG:HH11	1:C:442:ARG:HG2	1.79	0.47
1:D:265:ASP:O	1:D:269:ILE:HG13	2.15	0.47
1:A:323:VAL:HG12	1:A:327:ILE:HD11	1.96	0.47
1:B:112:ILE:HD11	1:B:127:SER:OG	2.14	0.47
1:B:265:ASP:O	1:B:269:ILE:HD13	2.15	0.47
1:D:425:PHE:C	1:D:427:PRO:HD3	2.35	0.47
1:B:183:PHE:HA	1:B:262:ASP:OD2	2.15	0.46
1:B:342:ARG:HA	1:B:446:PHE:CZ	2.50	0.46
1:D:251:HIS:HD1	1:D:251:HIS:H	1.63	0.46
1:D:35:PRO:HG3	1:D:68:TYR:CD1	2.50	0.46
1:A:39:ILE:HB	1:A:42:ILE:CG2	2.45	0.46
1:A:462:ILE:HD13	1:A:466:ASP:HB2	1.98	0.46
1:B:105:ARG:NH1	1:B:234:LEU:HD22	2.30	0.46
1:B:186:ARG:HG2	1:B:187:PHE:H	1.80	0.46
1:B:146:GLN:HG2	1:B:447:LEU:HD13	1.97	0.46
1:A:155:GLU:HG3	1:A:189:TYR:CD1	2.50	0.46
1:B:394:VAL:HG21	1:B:426:MET:HE3	1.97	0.46
1:B:363:PRO:HD2	1:B:476:PHE:O	2.16	0.46
1:D:101:PRO:HD2	1:D:217:ASN:ND2	2.31	0.46
1:D:251:HIS:HA	1:D:261:ARG:NH2	2.29	0.46
1:D:264:ILE:O	1:D:268:LEU:HB2	2.16	0.46
1:D:41:ASN:HA	1:D:44:GLN:OE1	2.15	0.46
1:B:344:HIS:C	1:B:346:PRO:HD3	2.36	0.46
1:B:39:ILE:HB	1:B:42:ILE:HG22	1.98	0.46
1:C:170:LEU:HD13	1:C:307:ARG:HA	1.97	0.46
1:C:320:THR:O	1:C:324:GLN:HG3	2.15	0.46
1:A:324:GLN:HA	1:A:327:ILE:HD12	1.98	0.46
1:A:327:ILE:HG12	1:A:345:MET:HE1	1.98	0.46
1:A:83:VAL:HG11	1:A:426:MET:HE1	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:30:PRO:HA	1:B:379:TYR:CD2	2.51	0.46
1:B:331:VAL:HA	1:B:335:ARG:NH2	2.31	0.46
1:A:408:ASP:HB3	1:A:411:HIS:CE1	2.50	0.46
1:D:57:LEU:HD12	1:D:67:LEU:HD11	1.98	0.46
1:B:394:VAL:HG21	1:B:426:MET:CE	2.46	0.46
1:C:337:PRO:HD3	1:C:454:GLN:OE1	2.15	0.46
1:D:84:LYS:HG2	1:D:88:ILE:HD12	1.98	0.46
1:C:318:GLU:HG3	1:C:319:VAL:N	2.30	0.46
1:D:69:PHE:HB3	1:D:218:ASN:OD1	2.15	0.46
1:D:255:MET:HE2	1:D:268:LEU:HD23	1.98	0.46
1:D:429:SER:HB3	2:D:501:HEM:HBA1	1.98	0.46
1:A:196:ASN:O	1:A:200:LYS:HB2	2.16	0.46
1:B:146:GLN:HG2	1:B:447:LEU:CD1	2.45	0.46
1:D:52:LYS:O	1:D:56:ASN:ND2	2.49	0.46
1:A:420:LYS:HG3	1:A:420:LYS:O	2.16	0.45
1:B:165:ASP:OD2	1:B:166:PRO:HD2	2.16	0.45
1:B:75:VAL:HB	1:B:387:ILE:HD13	1.98	0.45
1:C:111:GLY:CA	1:C:289:ILE:HG22	2.45	0.45
1:A:314:LEU:HD11	1:A:485:LEU:CB	2.46	0.45
1:C:326:GLU:O	1:C:330:VAL:HG23	2.16	0.45
1:A:176:ASN:HD22	1:A:189:TYR:HE2	1.65	0.45
1:A:31:PRO:HD3	1:A:379:TYR:CE2	2.52	0.45
1:D:245:LEU:O	1:D:249:LYS:HG2	2.17	0.45
1:D:124:ARG:HG3	1:D:128:LEU:CD1	2.47	0.45
1:D:85:GLU:O	1:D:89:ASP:HB2	2.17	0.45
1:A:215:ILE:HG12	4:A:505:GOL:H32	1.97	0.45
1:B:441:ALA:HB1	2:B:501:HEM:HAB	1.98	0.45
1:C:362:ILE:HG22	1:C:365:SER:N	2.31	0.45
1:A:374:VAL:O	1:A:380:LEU:HD12	2.17	0.45
1:B:56:ASN:O	1:B:59:LYS:HB2	2.17	0.45
1:C:102:LEU:HD11	1:C:213:ILE:CG2	2.45	0.45
1:C:318:GLU:HG3	1:C:319:VAL:HG23	1.98	0.45
1:C:97:ARG:NH2	1:C:433:ARG:HD2	2.32	0.45
1:A:50:VAL:O	1:A:53:SER:HB2	2.17	0.45
1:B:394:VAL:HG12	1:B:424:TYR:O	2.16	0.45
1:C:321:ALA:O	1:C:325:GLU:HG3	2.17	0.45
1:D:170:LEU:HD11	1:D:310:LEU:CD1	2.46	0.45
1:A:458:LEU:HD22	1:A:485:LEU:CD1	2.47	0.45
1:A:156:LEU:HD22	1:A:487:PHE:CE2	2.52	0.45
1:B:327:ILE:HG12	1:B:345:MET:HE1	1.99	0.45
1:B:365:SER:HB2	1:B:391:LEU:HG	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:210:THR:HG22	1:D:212:TRP:H	1.82	0.45
1:B:328:GLU:O	1:B:332:GLY:HA2	2.17	0.44
1:C:309:ALA:O	1:C:313:LEU:HG	2.17	0.44
1:B:207:ILE:O	1:B:213:ILE:HG13	2.17	0.44
1:B:420:LYS:HG3	1:B:420:LYS:O	2.16	0.44
1:C:242:SER:O	1:C:246:GLU:HG2	2.17	0.44
1:A:401:PHE:HB3	1:A:411:HIS:CD2	2.52	0.44
1:C:112:ILE:CD1	1:C:436:VAL:HG21	2.42	0.44
1:C:97:ARG:HB3	1:C:116:ASN:HD21	1.83	0.44
1:B:204:ASN:O	1:B:208:VAL:HG23	2.17	0.44
1:B:87:LEU:HD12	1:B:430:ALA:O	2.16	0.44
1:B:124:ARG:NH2	1:B:434:ILE:O	2.50	0.44
1:C:78:HIS:CE1	1:C:393:SER:HB2	2.52	0.44
1:D:324:GLN:HA	1:D:327:ILE:CD1	2.42	0.44
1:D:73:ARG:HH21	1:D:382:PRO:HG2	1.82	0.44
1:A:29:LEU:HD11	1:A:73:ARG:CZ	2.48	0.44
1:B:164:CYS:SG	1:B:165:ASP:N	2.90	0.44
1:B:285:GLU:O	1:B:289:ILE:HG12	2.18	0.44
1:B:414:ASP:N	1:B:418:ASN:O	2.37	0.44
1:C:205:ILE:HG12	1:C:300:GLU:HG2	1.99	0.44
1:C:162:SER:HB3	1:C:163:PRO:HD2	1.99	0.44
1:D:124:ARG:NH2	1:D:434:ILE:HG22	2.33	0.44
1:D:261:ARG:HB2	1:D:265:ASP:OD2	2.18	0.44
1:D:372:CYS:HA	1:D:383:LYS:HB2	1.99	0.44
1:C:261:ARG:HG3	1:C:261:ARG:NH1	2.33	0.44
1:C:372:CYS:HA	1:C:383:LYS:HB2	2.00	0.44
1:A:107:ASN:O	1:A:108:ARG:HB2	2.18	0.44
1:A:153:VAL:HG13	1:A:456:PHE:CE2	2.53	0.44
1:A:470:THR:HA	1:A:471:PRO:HD3	1.89	0.44
1:B:139:ARG:NE	1:B:144:ARG:HG2	2.33	0.44
1:D:240:MET:O	1:D:244:ILE:HG13	2.17	0.44
1:D:353:HIS:CE1	1:D:419:PHE:HZ	2.36	0.44
1:A:153:VAL:HG13	1:A:456:PHE:HE2	1.83	0.44
1:A:47:ILE:HA	1:A:47:ILE:HD12	1.89	0.44
1:A:86:ALA:HB2	1:A:376:PHE:CE1	2.53	0.44
1:B:344:HIS:O	1:B:346:PRO:HD3	2.17	0.44
1:B:426:MET:HG2	1:B:430:ALA:HB2	2.00	0.44
1:C:315:LYS:HD3	1:C:467:LEU:O	2.18	0.44
1:D:128:LEU:HD12	1:D:128:LEU:H	1.83	0.44
1:D:30:PRO:HA	1:D:379:TYR:CD2	2.53	0.44
1:A:327:ILE:O	1:A:331:VAL:HB	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:136:MET:SD	1:B:181:ILE:HG23	2.57	0.43
1:C:440:LEU:O	1:C:444:GLU:HG3	2.18	0.43
1:D:330:VAL:HG21	1:D:345:MET:CG	2.47	0.43
1:B:414:ASP:HB3	1:B:418:ASN:H	1.83	0.43
1:C:319:VAL:O	1:C:323:VAL:HG23	2.17	0.43
1:C:330:VAL:HG11	1:C:344:HIS:HB2	2.00	0.43
1:C:70:GLY:O	1:C:71:LEU:HB3	2.18	0.43
1:D:365:SER:OG	1:D:366:LEU:N	2.49	0.43
1:A:454:GLN:O	1:A:454:GLN:HG2	2.19	0.43
1:B:129:MET:O	1:B:132:ARG:HG3	2.17	0.43
1:B:246:GLU:O	1:B:250:GLU:HG3	2.18	0.43
1:C:119:ARG:HG2	1:C:281:GLU:CD	2.39	0.43
1:A:331:VAL:HA	1:A:335:ARG:NH2	2.34	0.43
1:B:52:LYS:HD3	1:B:52:LYS:O	2.19	0.43
1:C:104:GLU:O	1:C:108:ARG:HA	2.18	0.43
1:C:229:THR:O	1:C:232:LYS:HB3	2.17	0.43
1:C:308:TYR:OH	1:C:471:PRO:HG3	2.18	0.43
1:C:488:ILE:N	1:C:488:ILE:HD12	2.33	0.43
1:A:429:SER:CB	2:A:501:HEM:HBA1	2.49	0.43
1:C:30:PRO:HA	1:C:379:TYR:CD2	2.53	0.43
1:D:251:HIS:O	1:D:255:MET:HB2	2.18	0.43
1:D:205:ILE:HG13	1:D:300:GLU:CG	2.49	0.43
1:A:223:ILE:CD1	1:A:230:HIS:ND1	2.82	0.43
1:A:432:LYS:HE3	1:A:432:LYS:HB3	1.79	0.43
1:B:283:THR:CG2	1:B:286:ASN:HB2	2.48	0.43
1:C:366:LEU:HG	3:C:502:OXV:H2	2.01	0.43
1:D:218:ASN:O	1:D:220:PRO:HD3	2.19	0.43
1:D:376:PHE:CD2	1:D:377:ARG:HG3	2.54	0.43
1:D:414:ASP:OD2	1:D:418:ASN:HB2	2.19	0.43
1:B:154:GLU:O	1:B:158:LYS:HG2	2.18	0.43
1:B:311:LEU:HB2	1:B:483:TYR:CE2	2.54	0.43
1:C:145:VAL:HG11	1:C:447:LEU:HD12	2.01	0.43
1:A:440:LEU:O	1:A:444:GLU:HG3	2.19	0.43
1:C:267:PHE:CD2	1:C:287:LEU:HD22	2.52	0.43
1:C:360:ASP:OD2	1:C:392:THR:HB	2.19	0.43
1:D:111:GLY:CA	1:D:289:ILE:HG22	2.49	0.43
1:B:235:LYS:O	1:B:239:PHE:HD1	2.02	0.43
1:D:93:GLU:O	1:D:370:VAL:HA	2.18	0.43
1:B:354:GLU:HA	1:B:354:GLU:OE1	2.19	0.42
1:B:433:ARG:HB3	2:B:501:HEM:O2A	2.18	0.42
1:C:256:ASP:HB3	1:C:259:ASN:HB3	1.99	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:272:GLU:HA	1:C:275:LYS:HB2	2.01	0.42
1:D:238:ALA:O	1:D:241:GLU:HB2	2.19	0.42
1:A:314:LEU:HD11	1:A:485:LEU:HB3	1.99	0.42
1:A:401:PHE:O	1:A:404:PRO:HD3	2.18	0.42
1:A:48:LYS:HE3	1:A:48:LYS:HB2	1.83	0.42
1:A:72:GLU:HG2	5:A:614:HOH:O	2.19	0.42
1:C:31:PRO:HD3	1:C:379:TYR:CE2	2.54	0.42
1:B:355:VAL:O	1:B:359:ILE:HG23	2.19	0.42
1:B:49:ASP:HA	1:B:473:VAL:HG11	2.02	0.42
1:A:90:LEU:HB3	1:A:93:GLU:HG3	2.01	0.42
1:B:105:ARG:CZ	1:B:234:LEU:HD13	2.49	0.42
1:C:36:LEU:HD12	1:C:43:LEU:HD12	2.01	0.42
1:A:442:ARG:HG2	1:A:442:ARG:HH11	1.84	0.42
1:C:105:ARG:HE	1:C:234:LEU:CD1	2.31	0.42
1:D:113:VAL:HG22	2:D:501:HEM:HAD1	2.01	0.42
1:D:346:PRO:HG2	1:D:347:TYR:H	1.83	0.42
1:A:149:ALA:O	1:A:153:VAL:HG23	2.20	0.42
1:B:368:HIS:NE2	1:B:433:ARG:CB	2.82	0.42
1:C:85:GLU:O	1:C:89:ASP:HB2	2.20	0.42
1:A:178:ILE:HG21	1:A:295:LEU:HA	2.02	0.42
1:A:42:ILE:HG23	1:A:43:LEU:N	2.32	0.42
1:B:139:ARG:HD3	1:B:144:ARG:NH1	2.34	0.42
1:B:267:PHE:CD2	1:B:287:LEU:HB2	2.55	0.42
1:C:307:ARG:NH1	5:C:612:HOH:O	2.47	0.42
1:C:330:VAL:O	1:C:330:VAL:HG12	2.19	0.42
1:C:425:PHE:C	1:C:427:PRO:HD3	2.40	0.42
1:D:38:VAL:HG12	1:D:39:ILE:N	2.35	0.42
1:D:412:PHE:HB3	1:D:419:PHE:HE2	1.85	0.42
1:D:83:VAL:HG13	1:D:87:LEU:HD23	2.01	0.42
1:A:152:LEU:O	1:A:156:LEU:HG	2.20	0.42
1:B:463:ASP:OD2	1:B:465:LYS:HD2	2.19	0.42
1:C:85:GLU:O	1:C:90:LEU:HD13	2.20	0.42
1:D:30:PRO:HG3	1:D:379:TYR:CB	2.50	0.42
1:A:84:LYS:CG	1:A:88:ILE:HD12	2.26	0.42
1:B:320:THR:O	1:B:324:GLN:HG3	2.19	0.42
1:D:445:LEU:O	1:D:449:LEU:HD12	2.20	0.42
1:B:275:LYS:HG2	1:B:276:GLN:HG3	2.01	0.42
1:B:308:TYR:OH	1:B:471:PRO:HG3	2.20	0.42
1:C:212:TRP:O	1:C:215:ILE:HG22	2.20	0.42
1:C:426:MET:N	1:C:427:PRO:CD	2.82	0.42
1:A:173:ALA:HB3	1:A:174:PRO:CD	2.42	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:156:LEU:O	1:B:159:THR:HG22	2.20	0.41
1:B:86:ALA:HB2	1:B:376:PHE:CE1	2.54	0.41
1:C:428:PHE:O	1:C:429:SER:HB3	2.20	0.41
1:D:118:LYS:O	1:D:119:ARG:C	2.58	0.41
1:B:339:MET:O	1:B:342:ARG:HB3	2.20	0.41
1:C:403:ASN:HB3	1:C:406:MET:HG3	2.00	0.41
1:D:436:VAL:HG23	2:D:501:HEM:HBD2	2.02	0.41
1:A:184:GLN:HG3	1:A:185:LYS:HG2	2.01	0.41
1:B:182:ILE:HG22	1:B:183:PHE:CE1	2.55	0.41
1:B:123:ILE:HG13	1:B:286:ASN:ND2	2.35	0.41
1:B:365:SER:HB2	1:B:391:LEU:CG	2.51	0.41
1:B:49:ASP:CG	1:B:52:LYS:HB2	2.41	0.41
1:C:455:ASN:O	1:C:489:PRO:HA	2.20	0.41
1:A:252:GLN:HG2	1:A:268:LEU:CD2	2.50	0.41
1:A:444:GLU:HB3	1:A:448:PHE:CE2	2.55	0.41
1:B:370:VAL:HG23	1:B:370:VAL:O	2.20	0.41
1:C:35:PRO:HB2	1:C:40:GLY:HA2	2.02	0.41
1:C:370:VAL:HG21	1:C:381:ILE:HG22	2.01	0.41
1:C:377:ARG:O	1:C:378:ASN:C	2.58	0.41
1:D:204:ASN:O	1:D:208:VAL:HG23	2.20	0.41
1:D:471:PRO:HB2	1:D:473:VAL:HG13	2.03	0.41
1:A:252:GLN:HG2	1:A:268:LEU:HD21	2.03	0.41
1:B:49:ASP:HA	1:B:473:VAL:CG1	2.51	0.41
1:C:311:LEU:HB2	1:C:483:TYR:CE2	2.55	0.41
1:A:178:ILE:HG22	1:A:182:ILE:CD1	2.51	0.41
1:A:223:ILE:O	1:A:223:ILE:HD12	2.21	0.41
1:A:426:MET:N	1:A:427:PRO:CD	2.84	0.41
1:B:187:PHE:CG	1:B:194:PHE:HB2	2.55	0.41
1:C:181:ILE:O	1:C:263:PHE:HB2	2.21	0.41
1:C:382:PRO:HG2	1:C:385:THR:OG1	2.20	0.41
1:D:86:ALA:HB2	1:D:376:PHE:CE1	2.55	0.41
1:A:173:ALA:O	1:A:177:VAL:HG23	2.19	0.41
1:B:110:PHE:HB2	1:B:289:ILE:HG13	2.03	0.41
1:C:39:ILE:HB	1:C:42:ILE:CG1	2.49	0.41
1:A:148:GLU:OE2	1:A:176:ASN:HB3	2.21	0.41
1:A:400:GLU:HG2	1:A:424:TYR:CG	2.56	0.41
1:B:278:GLN:HG3	1:B:279:GLN:NE2	2.36	0.41
1:B:430:ALA:HA	1:B:434:ILE:HD12	2.01	0.41
1:B:479:VAL:HG13	1:B:480:PRO:HD2	2.02	0.41
1:C:318:GLU:CG	1:C:319:VAL:N	2.84	0.41
1:C:85:GLU:CD	1:C:377:ARG:HH21	2.24	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:38:VAL:HG12	1:C:38:VAL:O	2.21	0.41
1:D:51:SER:OG	1:D:363:PRO:HB3	2.20	0.41
1:D:444:GLU:O	1:D:448:PHE:HB2	2.20	0.41
1:D:488:ILE:N	1:D:488:ILE:HD12	2.36	0.41
1:A:205:ILE:HA	1:A:205:ILE:HD12	1.72	0.41
1:A:42:ILE:CG2	1:A:43:LEU:N	2.84	0.41
1:C:36:LEU:HD23	1:C:36:LEU:HA	1.94	0.41
1:C:428:PHE:CG	1:C:438:GLU:HG2	2.56	0.41
1:D:102:LEU:HB2	1:D:217:ASN:OD1	2.21	0.41
1:D:153:VAL:HG13	1:D:455:ASN:HD21	1.86	0.41
1:B:230:HIS:O	1:B:233:LEU:HB3	2.21	0.41
1:B:124:ARG:NH1	1:B:433:ARG:HA	2.36	0.41
1:B:97:ARG:HD3	1:B:367:PRO:O	2.21	0.41
1:C:219:PHE:O	1:C:222:ILE:HG22	2.21	0.41
1:C:225:TYR:HD2	1:C:225:TYR:N	2.19	0.41
1:D:101:PRO:O	1:D:104:GLU:HB3	2.21	0.41
1:D:65:PHE:CE1	1:D:76:VAL:HB	2.56	0.41
1:A:368:HIS:HE1	1:A:433:ARG:HB2	1.86	0.41
1:B:173:ALA:CB	1:B:174:PRO:HD3	2.41	0.41
1:D:357:ARG:HD2	1:D:401:PHE:CD1	2.56	0.41
1:D:405:GLU:HG2	1:D:406:MET:N	2.36	0.41
1:A:213:ILE:HD13	1:A:213:ILE:HA	1.92	0.40
1:A:463:ASP:OD2	1:A:465:LYS:HB2	2.21	0.40
1:B:257:ILE:O	1:B:257:ILE:HG22	2.20	0.40
1:B:54:LEU:O	1:B:57:LEU:HB2	2.21	0.40
1:C:120:TRP:CZ2	1:C:124:ARG:HD2	2.56	0.40
1:C:30:PRO:HA	1:C:379:TYR:CG	2.56	0.40
1:C:400:GLU:HG2	1:C:424:TYR:HB2	2.04	0.40
1:B:110:PHE:N	1:B:110:PHE:CD1	2.89	0.40
1:B:225:TYR:CG	1:B:226:PHE:N	2.89	0.40
1:C:182:ILE:HA	1:C:263:PHE:CB	2.51	0.40
1:C:69:PHE:HB2	1:C:74:MET:CE	2.51	0.40
1:D:148:GLU:OE1	1:D:173:ALA:HA	2.21	0.40
1:D:39:ILE:HG13	1:D:40:GLY:N	2.36	0.40
1:A:275:LYS:O	1:A:275:LYS:HG2	2.21	0.40
1:A:311:LEU:HB2	1:A:483:TYR:CE2	2.56	0.40
1:A:376:PHE:O	1:A:377:ARG:HD3	2.22	0.40
1:C:354:GLU:HA	1:C:354:GLU:OE2	2.21	0.40
1:D:248:VAL:HA	1:D:264:ILE:HG21	2.03	0.40
1:A:463:ASP:OD1	1:A:464:PRO:HD2	2.20	0.40
1:B:184:GLN:HG3	1:B:185:LYS:N	2.36	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:159:THR:O	1:C:160:LYS:HB2	2.21	0.40
1:C:327:ILE:HA	1:C:345:MET:HE1	2.04	0.40
1:A:459:LYS:HB3	1:A:488:ILE:CD1	2.51	0.40
1:B:326:GLU:HG2	1:B:345:MET:HE2	2.04	0.40
1:C:86:ALA:HB1	1:C:94:PHE:CZ	2.57	0.40
1:D:359:ILE:HG13	1:D:360:ASP:N	2.35	0.40
1:D:374:VAL:HG12	1:D:375:LYS:N	2.35	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	462/477 (97%)	418 (90%)	42 (9%)	2 (0%)	34 64
1	B	452/477 (95%)	411 (91%)	34 (8%)	7 (2%)	10 32
1	C	460/477 (96%)	415 (90%)	43 (9%)	2 (0%)	34 64
1	D	448/477 (94%)	396 (88%)	51 (11%)	1 (0%)	47 76
All	All	1822/1908 (96%)	1640 (90%)	170 (9%)	12 (1%)	22 52

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	228	GLY
1	B	365	SER
1	A	429	SER
1	D	31	PRO
1	B	429	SER
1	C	228	GLY
1	B	91	GLY
1	B	215	ILE

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Mol	Chain	Res	Type
1	B	227	PRO
1	C	31	PRO
1	A	47	ILE
1	B	101	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	421/431 (98%)	418 (99%)	3 (1%)	84	94
1	B	413/431 (96%)	407 (98%)	6 (2%)	65	86
1	C	417/431 (97%)	411 (99%)	6 (1%)	67	87
1	D	405/431 (94%)	400 (99%)	5 (1%)	71	89
All	All	1656/1724 (96%)	1636 (99%)	20 (1%)	71	89

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	100	PHE
1	A	219	PHE
1	A	308	TYR
1	B	99	HIS
1	B	176	ASN
1	B	308	TYR
1	B	362	ILE
1	B	368	HIS
1	B	379	TYR
1	C	87	LEU
1	C	100	PHE
1	C	216	CYS
1	C	265	ASP
1	C	308	TYR
1	C	379	TYR
1	D	99	HIS
1	D	302	THR

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Mol	Chain	Res	Type
1	D	308	TYR
1	D	318	GLU
1	D	362	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	192	GLN
1	A	457	ASN
1	A	474	ASN
1	B	146	GLN
1	B	231	ASN
1	B	236	ASN
1	C	196	ASN
1	C	214	GLN
1	C	218	ASN
1	C	378	ASN
1	D	56	ASN
1	D	176	ASN
1	D	192	GLN
1	D	214	GLN
1	D	231	ASN
1	D	277	ASN
1	D	403	ASN
1	D	454	GLN
1	D	455	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

### 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry

11 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
4	GOL	A	504	-	5,5,5	0.35	0	5,5,5	0.25	0
2	HEM	D	501	1	27,50,50	1.97	10 (37%)	17,82,82	1.69	4 (23%)
4	GOL	A	503	-	5,5,5	0.30	0	5,5,5	0.24	0
2	HEM	B	501	1	27,50,50	1.99	10 (37%)	17,82,82	1.77	4 (23%)
3	0XV	C	502	-	18,23,23	2.26	9 (50%)	25,34,34	1.40	4 (16%)
2	HEM	C	501	1	27,50,50	1.99	11 (40%)	17,82,82	1.77	4 (23%)
3	0XV	A	502	-	18,23,23	2.20	8 (44%)	25,34,34	1.40	4 (16%)
3	0XV	B	502	-	18,23,23	2.21	9 (50%)	25,34,34	1.37	4 (16%)
4	GOL	A	505	-	5,5,5	0.32	0	5,5,5	0.26	0
3	0XV	D	502	-	18,23,23	2.24	9 (50%)	25,34,34	1.39	4 (16%)
2	HEM	A	501	1,5	27,50,50	1.98	10 (37%)	17,82,82	1.78	4 (23%)

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	GOL	A	504	-	-	2/4/4/4	-
2	HEM	D	501	1	-	0/6/54/54	-
4	GOL	A	503	-	-	2/4/4/4	-
2	HEM	B	501	1	-	0/6/54/54	-
3	0XV	C	502	-	-	0/5/8/8	0/3/3/3
2	HEM	C	501	1	-	0/6/54/54	-
3	0XV	A	502	-	-	0/5/8/8	0/3/3/3
3	0XV	B	502	-	-	0/5/8/8	0/3/3/3
4	GOL	A	505	-	-	0/4/4/4	-
3	0XV	D	502	-	-	0/5/8/8	0/3/3/3
2	HEM	A	501	1,5	-	0/6/54/54	-



All (76) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	HEM	C3C-CAC	-5.01	1.37	1.47
2	B	501	HEM	C3C-CAC	-5.01	1.37	1.47
2	D	501	HEM	C3C-CAC	-5.00	1.37	1.47
2	C	501	HEM	C3C-CAC	-4.97	1.37	1.47
3	C	502	0XV	CAL-CAJ	4.17	1.56	1.49
3	A	502	0XV	CAL-CAJ	4.14	1.56	1.49
3	D	502	0XV	CAL-CAJ	4.12	1.56	1.49
3	B	502	0XV	CAL-CAJ	4.05	1.56	1.49
3	A	502	0XV	CAG-CAJ	3.72	1.55	1.50
3	C	502	0XV	CAV-CAI	3.65	1.53	1.48
3	B	502	0XV	CAG-CAJ	3.65	1.55	1.50
3	C	502	0XV	CAG-CAJ	3.65	1.55	1.50
3	D	502	0XV	CAG-CAJ	3.64	1.55	1.50
3	D	502	0XV	CAV-CAI	3.56	1.53	1.48
3	A	502	0XV	CAV-CAI	3.40	1.53	1.48
3	B	502	0XV	CAV-CAI	3.36	1.53	1.48
2	B	501	HEM	CMA-C3A	3.19	1.58	1.51
2	A	501	HEM	CAD-C3D	3.17	1.57	1.52
2	D	501	HEM	C1D-ND	-3.16	1.29	1.36
2	C	501	HEM	CMA-C3A	3.15	1.58	1.51
2	D	501	HEM	CMA-C3A	3.13	1.58	1.51
2	B	501	HEM	CAD-C3D	3.12	1.57	1.52
2	C	501	HEM	CAD-C3D	3.11	1.57	1.52
2	A	501	HEM	CMA-C3A	3.11	1.58	1.51
2	B	501	HEM	C1D-ND	-3.09	1.29	1.36
2	A	501	HEM	C1D-ND	-3.06	1.29	1.36
2	C	501	HEM	C1D-ND	-2.97	1.30	1.36
2	A	501	HEM	CBB-CAB	2.95	1.48	1.29
2	D	501	HEM	CAD-C3D	2.93	1.57	1.52
2	D	501	HEM	CBB-CAB	2.92	1.48	1.29
2	C	501	HEM	CBB-CAB	2.90	1.48	1.29
2	B	501	HEM	CBB-CAB	2.90	1.48	1.29
2	D	501	HEM	C3B-CAB	-2.75	1.42	1.47
2	C	501	HEM	C3B-CAB	-2.64	1.42	1.47
2	B	501	HEM	C3B-CAB	-2.60	1.42	1.47
3	C	502	0XV	CAU-CAL	2.58	1.43	1.39
3	A	502	0XV	CAU-CAL	2.54	1.43	1.39
3	D	502	0XV	CAE-CAF	2.52	1.42	1.36
3	B	502	0XV	CAU-CAL	2.51	1.43	1.39
2	A	501	HEM	C3B-CAB	-2.51	1.42	1.47
3	D	502	0XV	CAU-CAL	2.49	1.43	1.39
3	C	502	0XV	CAD-CAC	2.48	1.42	1.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	502	0XV	CAD-CAC	2.48	1.42	1.36
3	B	502	0XV	CAE-CAF	2.47	1.42	1.36
3	C	502	0XV	CAE-CAF	2.45	1.42	1.36
3	B	502	0XV	CAD-CAC	2.45	1.42	1.36
3	A	502	0XV	CAE-CAF	2.35	1.42	1.36
3	D	502	0XV	CAM-CAL	2.32	1.42	1.39
3	A	502	0XV	CAD-CAC	2.29	1.41	1.36
3	B	502	0XV	CAM-CAL	2.27	1.42	1.39
2	A	501	HEM	CMC-C2C	2.26	1.57	1.51
2	B	501	HEM	C4A-NA	-2.26	1.31	1.36
3	A	502	0XV	CAM-CAL	2.26	1.42	1.39
2	B	501	HEM	CMC-C2C	2.25	1.56	1.51
2	C	501	HEM	CAA-C2A	2.25	1.55	1.52
2	C	501	HEM	C4A-NA	-2.23	1.31	1.36
2	C	501	HEM	CMC-C2C	2.22	1.56	1.51
3	C	502	0XV	CAP-CAN	2.22	1.43	1.40
2	D	501	HEM	CAA-C2A	2.21	1.55	1.52
2	D	501	HEM	CMC-C2C	2.21	1.56	1.51
2	A	501	HEM	CAA-C2A	2.21	1.55	1.52
3	A	502	0XV	CAP-CAN	2.20	1.43	1.40
3	C	502	0XV	CAM-CAL	2.20	1.42	1.39
3	B	502	0XV	CAP-CAN	2.19	1.43	1.40
3	D	502	0XV	CAP-CAN	2.17	1.42	1.40
2	A	501	HEM	C4A-NA	-2.16	1.31	1.36
2	B	501	HEM	CAA-C2A	2.14	1.55	1.52
2	D	501	HEM	C4B-NB	-2.12	1.31	1.36
3	C	502	0XV	CAP-CAQ	2.11	1.42	1.40
2	D	501	HEM	C4A-NA	-2.08	1.31	1.36
2	B	501	HEM	C4B-NB	-2.07	1.31	1.36
3	B	502	0XV	CAP-CAQ	2.06	1.42	1.40
2	C	501	HEM	C4B-NB	-2.04	1.32	1.36
2	A	501	HEM	C4B-NB	-2.04	1.32	1.36
3	D	502	0XV	CAP-CAQ	2.03	1.42	1.40
2	C	501	HEM	CBC-CAC	2.01	1.42	1.29

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	502	0XV	CAQ-CAP-CAN	-3.83	118.92	121.65
3	A	502	0XV	CAQ-CAP-CAN	-3.77	118.96	121.65
3	D	502	0XV	CAQ-CAP-CAN	-3.74	118.98	121.65
3	B	502	0XV	CAQ-CAP-CAN	-3.73	118.99	121.65

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	HEM	CAD-CBD-CGD	3.71	118.90	112.67
2	C	501	HEM	CAD-CBD-CGD	3.62	118.75	112.67
2	B	501	HEM	CAD-CBD-CGD	3.55	118.63	112.67
2	D	501	HEM	CAD-CBD-CGD	3.36	118.31	112.67
3	C	502	0XV	CAM-CAN-CAP	3.25	120.73	118.08
3	A	502	0XV	CAM-CAN-CAP	3.22	120.71	118.08
2	B	501	HEM	C4C-C3C-C2C	-3.22	104.65	106.90
3	D	502	0XV	CAM-CAN-CAP	3.18	120.68	118.08
3	B	502	0XV	CAM-CAN-CAP	3.18	120.68	118.08
2	C	501	HEM	C4C-C3C-C2C	-3.16	104.69	106.90
2	D	501	HEM	C4C-C3C-C2C	-3.11	104.73	106.90
2	A	501	HEM	C4C-C3C-C2C	-3.09	104.74	106.90
2	B	501	HEM	CMB-C2B-C3B	2.94	130.18	124.68
2	A	501	HEM	CMB-C2B-C3B	2.93	130.15	124.68
3	A	502	0XV	CAU-CAQ-CAP	2.92	120.46	118.08
2	C	501	HEM	CMB-C2B-C3B	2.91	130.12	124.68
3	D	502	0XV	CAU-CAQ-CAP	2.89	120.44	118.08
3	C	502	0XV	CAU-CAQ-CAP	2.88	120.43	118.08
3	B	502	0XV	CAU-CAQ-CAP	2.84	120.40	118.08
2	D	501	HEM	CMB-C2B-C3B	2.78	129.88	124.68
3	D	502	0XV	CAG-CAJ-CAL	2.65	123.23	119.53
3	A	502	0XV	CAG-CAJ-CAL	2.62	123.19	119.53
3	C	502	0XV	CAG-CAJ-CAL	2.61	123.17	119.53
3	B	502	0XV	CAG-CAJ-CAL	2.45	122.95	119.53
2	B	501	HEM	CBA-CAA-C2A	2.42	116.95	112.49
2	C	501	HEM	CBA-CAA-C2A	2.42	116.94	112.49
2	D	501	HEM	CBA-CAA-C2A	2.37	116.85	112.49
2	A	501	HEM	CBA-CAA-C2A	2.35	116.83	112.49

There are no chirality outliers.

All (4) torsion outliers are listed below:

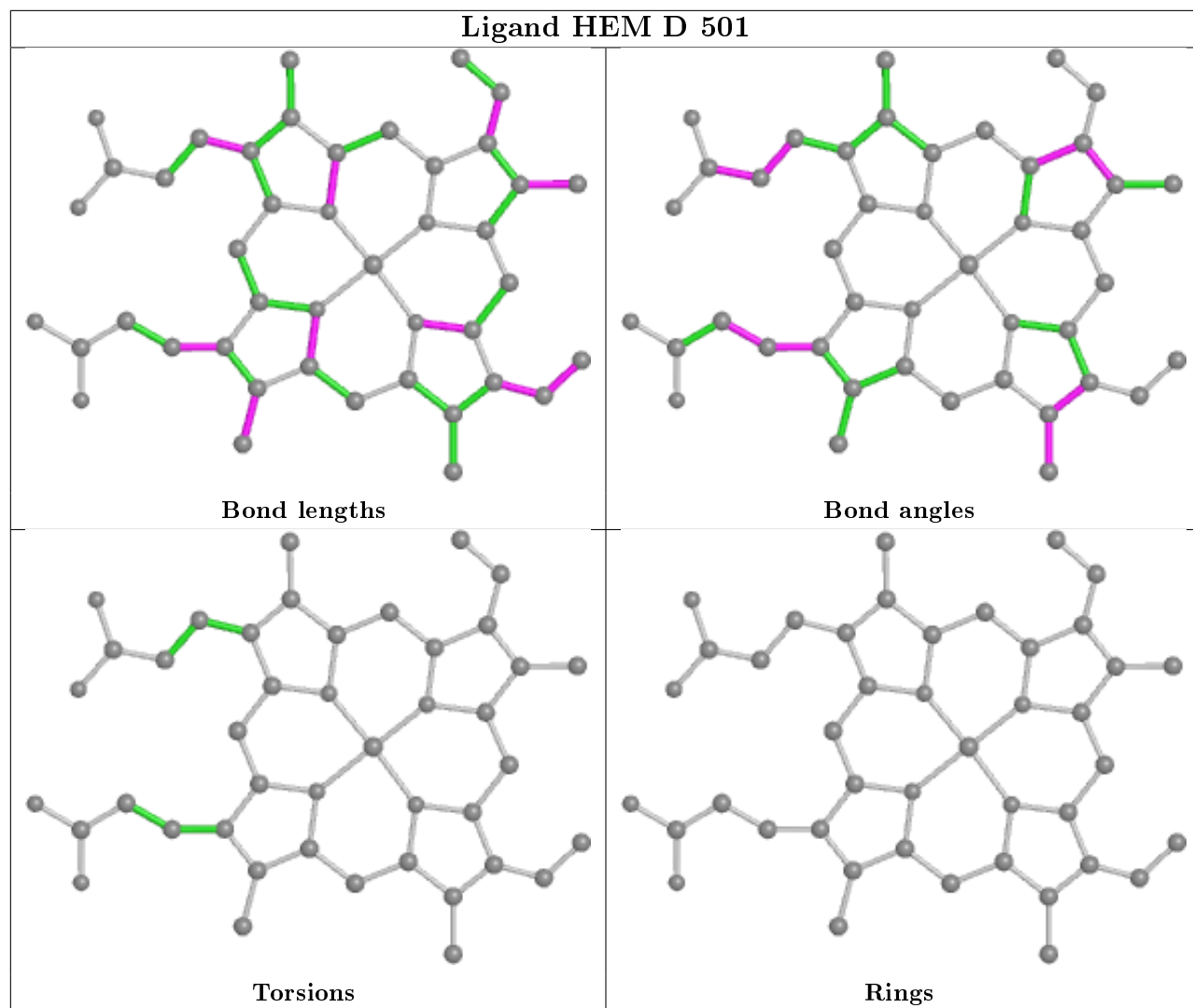
Mol	Chain	Res	Type	Atoms
4	A	504	GOL	O1-C1-C2-C3
4	A	503	GOL	C1-C2-C3-O3
4	A	504	GOL	O1-C1-C2-O2
4	A	503	GOL	O2-C2-C3-O3

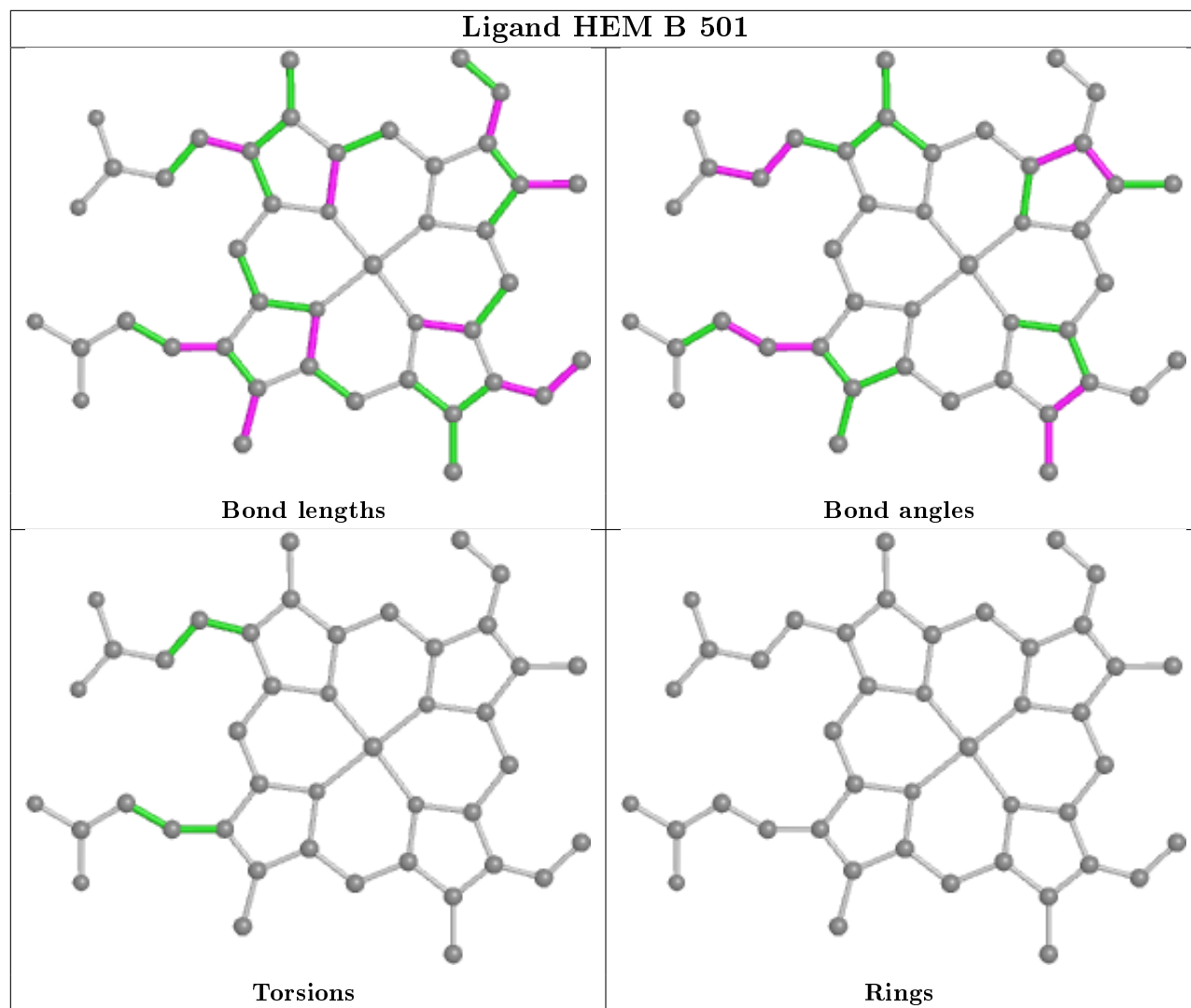
There are no ring outliers.

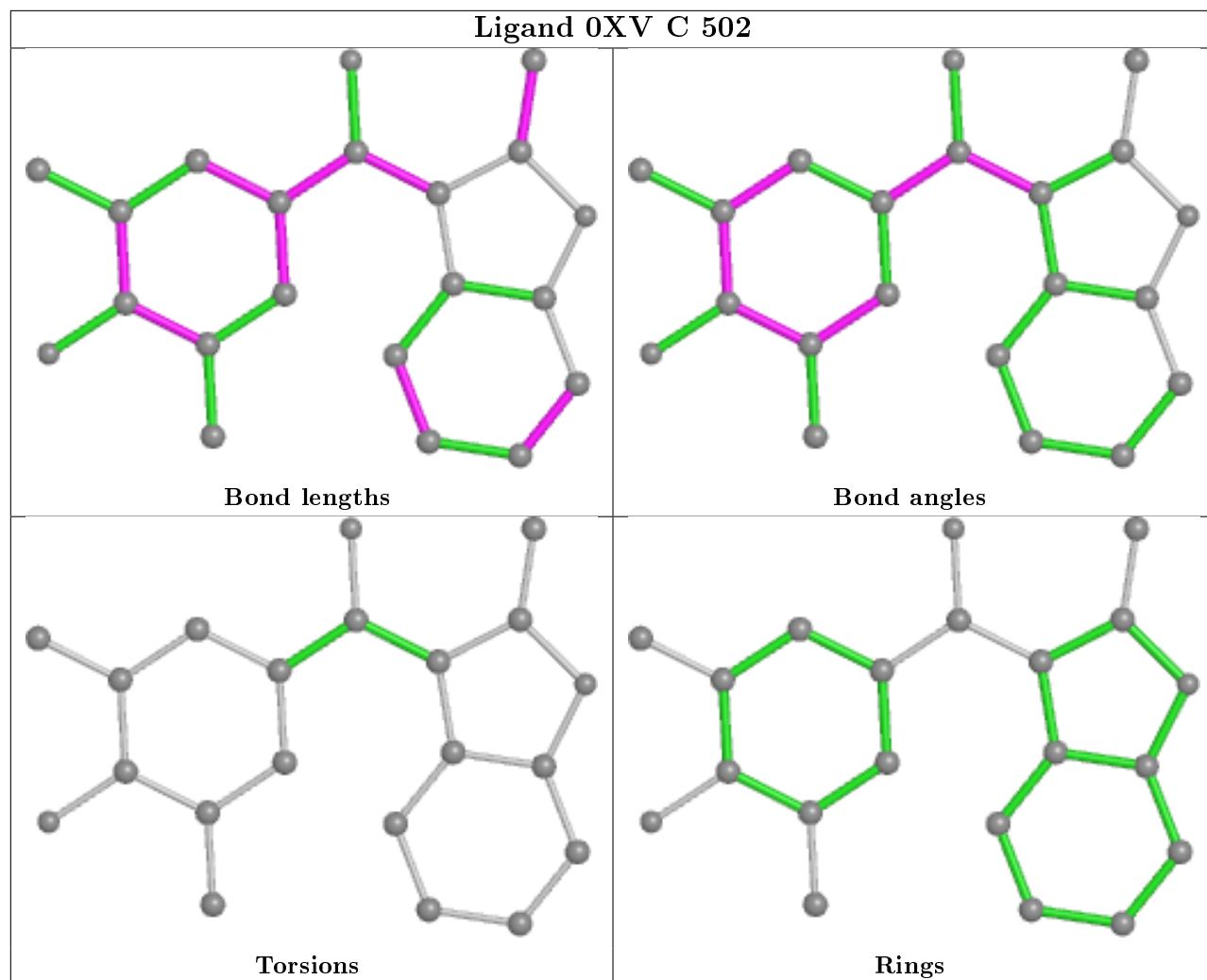
8 monomers are involved in 16 short contacts:

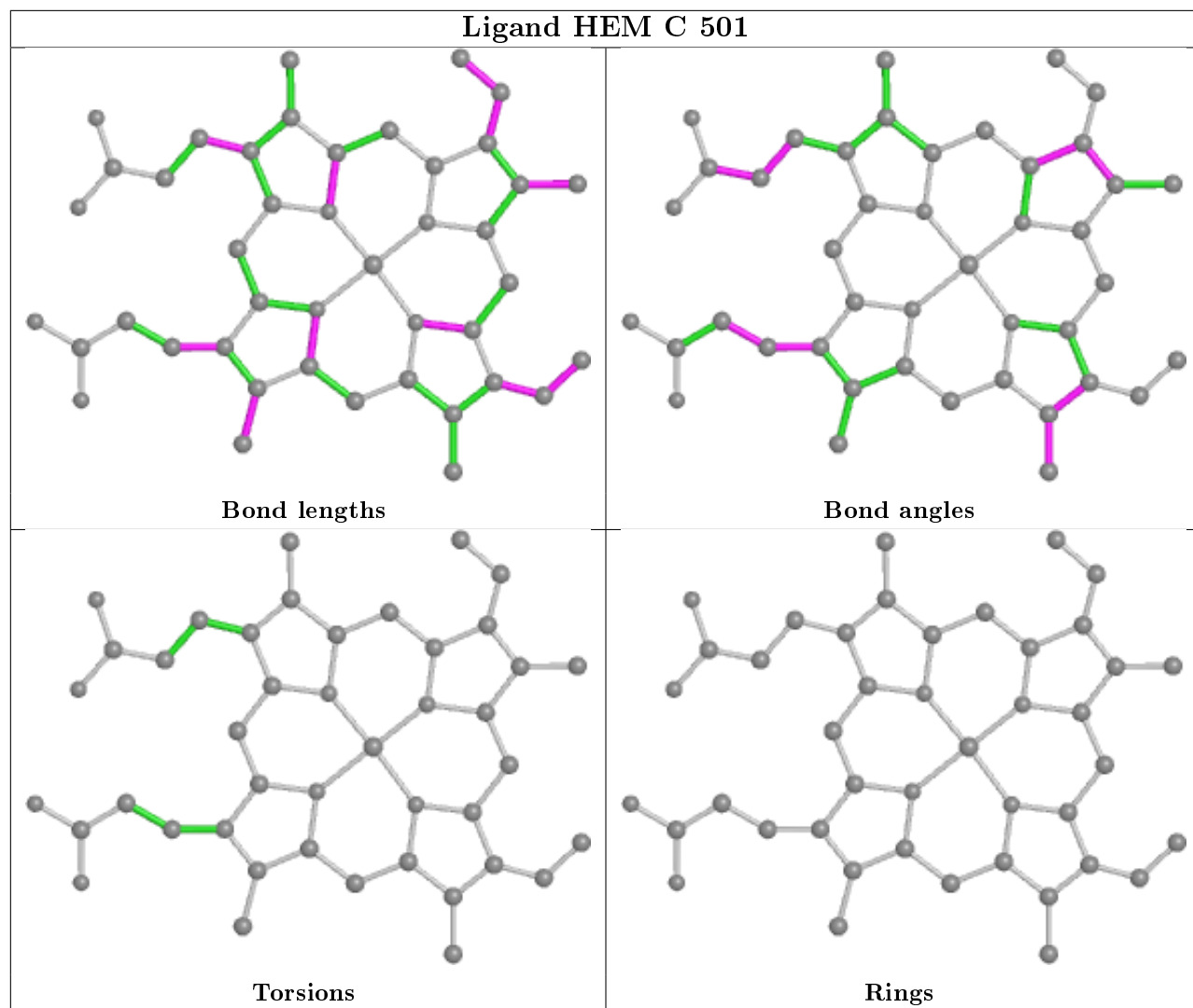
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	501	HEM	3	0
2	B	501	HEM	4	0
3	C	502	0XV	1	0
3	A	502	0XV	1	0
3	B	502	0XV	2	0
4	A	505	GOL	3	0
3	D	502	0XV	1	0
2	A	501	HEM	1	0

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.

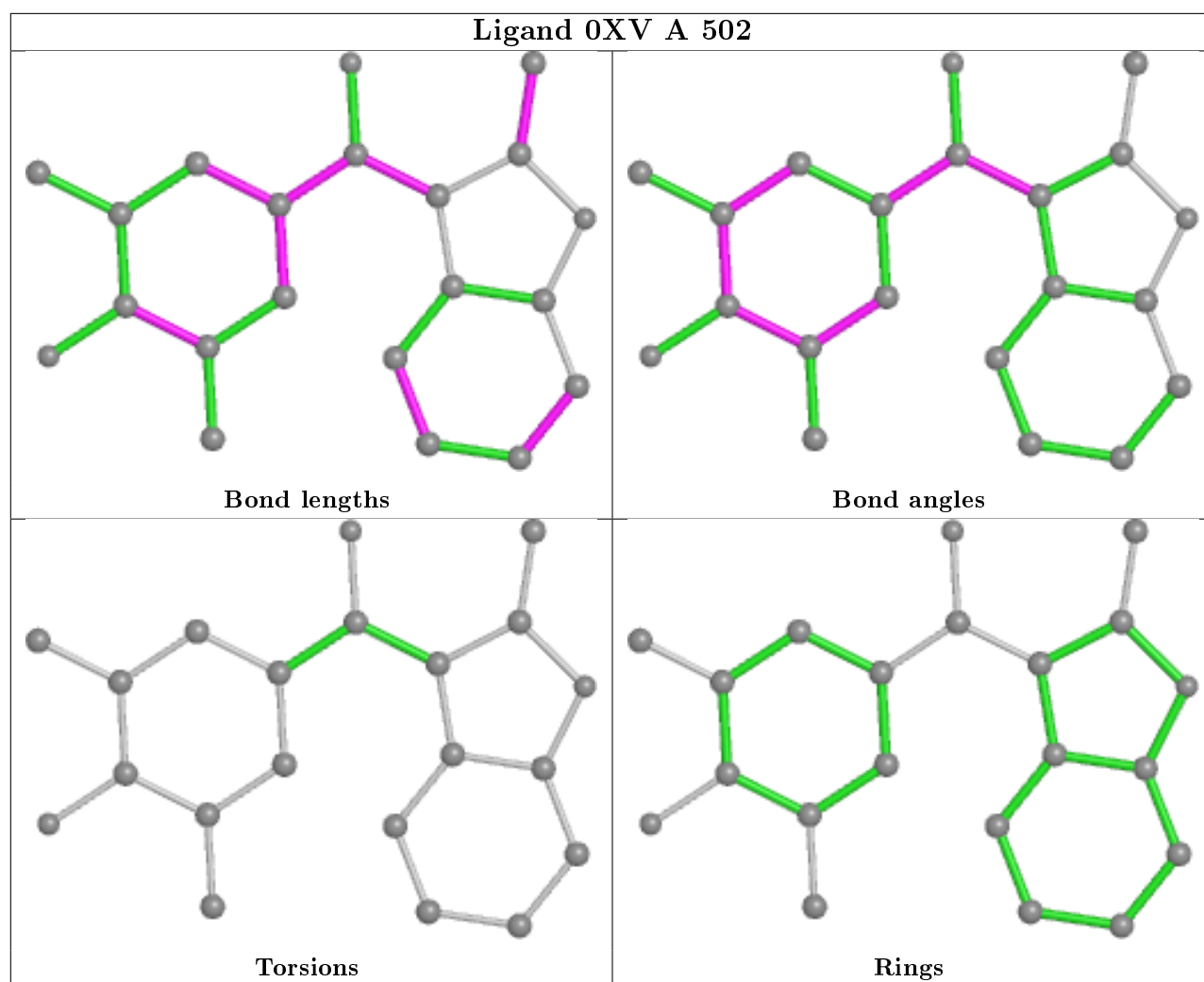


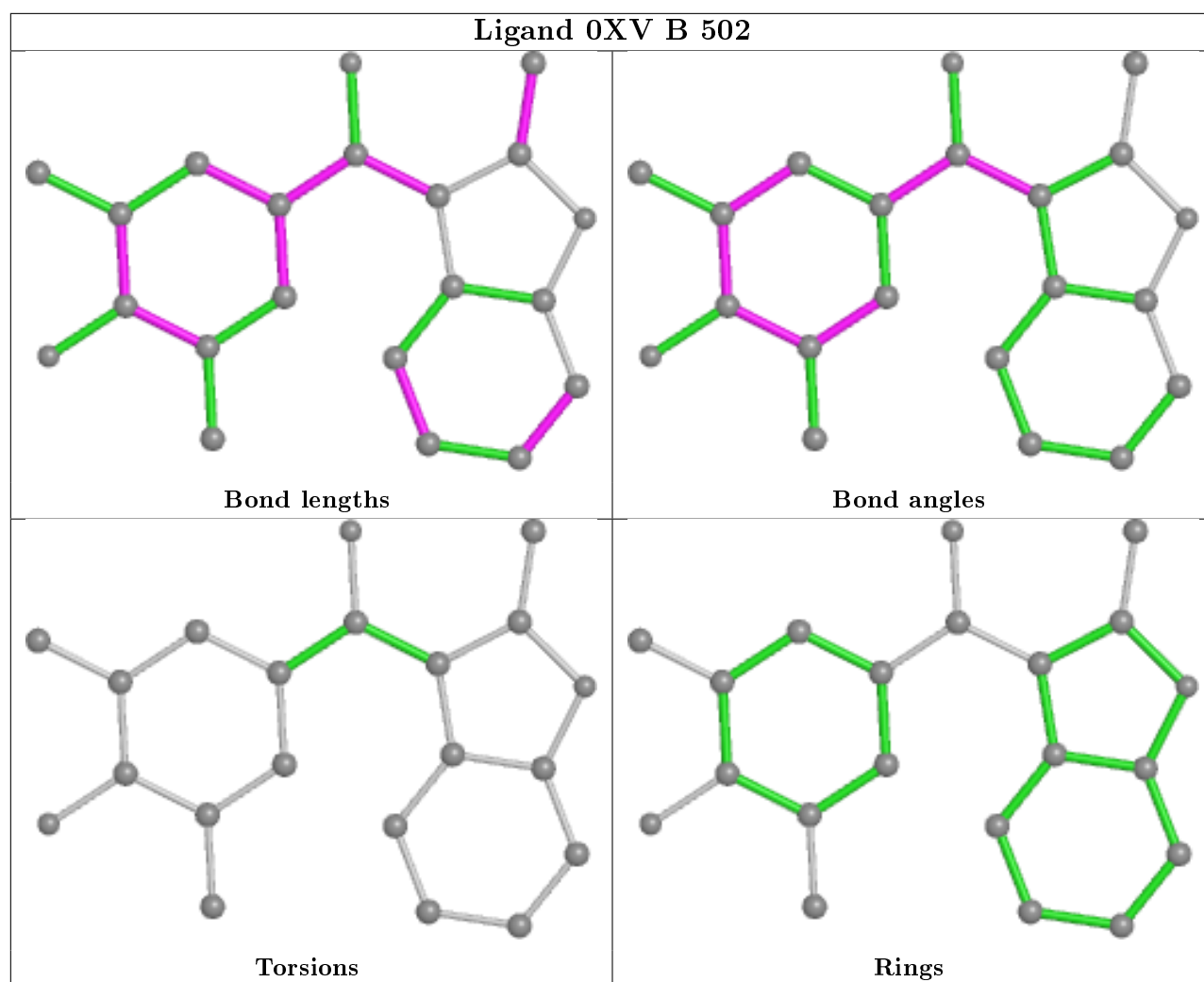


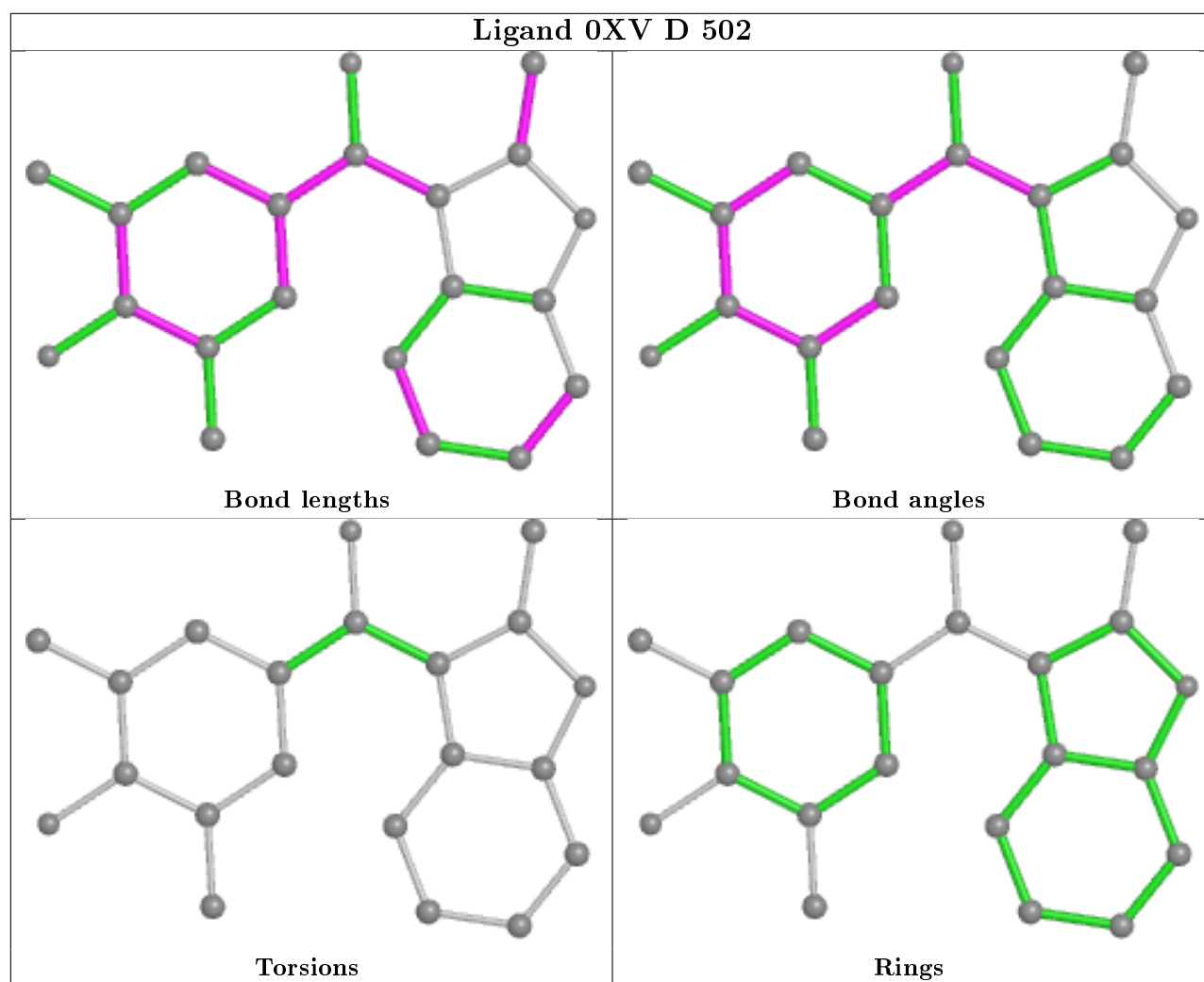


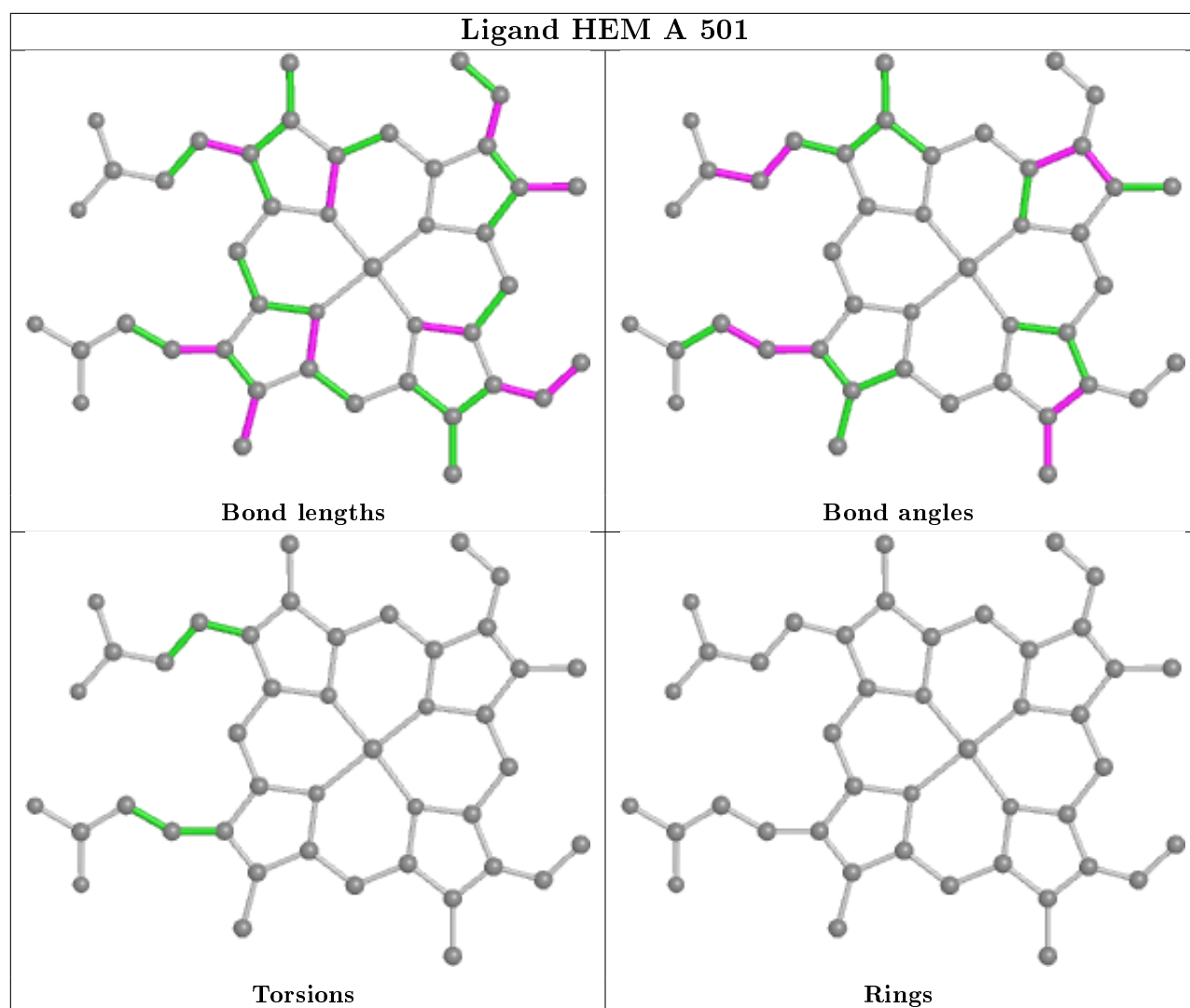












## 5.7 Other polymers [i](#)

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

### 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, 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	464/477 (97%)	0.10	2 (0%) 92 92	47, 68, 92, 113	0
1	B	456/477 (95%)	0.34	9 (1%) 65 63	55, 87, 117, 125	0
1	C	462/477 (96%)	0.21	8 (1%) 70 70	57, 86, 107, 127	0
1	D	452/477 (94%)	0.41	22 (4%) 29 26	65, 102, 128, 133	0
All	All	1834/1908 (96%)	0.26	41 (2%) 62 60	47, 85, 119, 133	0

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	214	GLN	3.8
1	D	473	VAL	3.7
1	D	214	GLN	3.3
1	C	458	LEU	3.2
1	D	260	PRO	3.1
1	D	476	PHE	2.9
1	D	487	PHE	2.9
1	C	476	PHE	2.8
1	C	225	TYR	2.8
1	B	164	CYS	2.7
1	C	257	ILE	2.5
1	A	216	CYS	2.5
1	C	212	TRP	2.5
1	D	123	ILE	2.5
1	D	265	ASP	2.4
1	D	194	PHE	2.4
1	D	45	ILE	2.4
1	D	215	ILE	2.4
1	D	244	ILE	2.3
1	D	248	VAL	2.3
1	B	248	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	331	VAL	2.3
1	C	314	LEU	2.3
1	B	276	GLN	2.3
1	D	193	GLN	2.3
1	D	268	LEU	2.2
1	C	207	ILE	2.2
1	D	468	ASP	2.2
1	B	39	ILE	2.2
1	D	240	MET	2.1
1	D	187	PHE	2.1
1	D	330	VAL	2.1
1	B	436	VAL	2.1
1	B	187	PHE	2.1
1	D	287	LEU	2.1
1	A	230	HIS	2.0
1	B	263	PHE	2.0
1	D	331	VAL	2.0
1	D	334	ASN	2.0
1	C	453	LEU	2.0
1	D	458	LEU	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( $\text{\AA}^2$ )	Q<0.9
4	GOL	A	504	6/6	0.87	0.17	75,75,75,75	0
4	GOL	A	505	6/6	0.87	0.18	89,90,90,90	0
3	0XV	B	502	21/21	0.88	0.35	77,84,86,87	0

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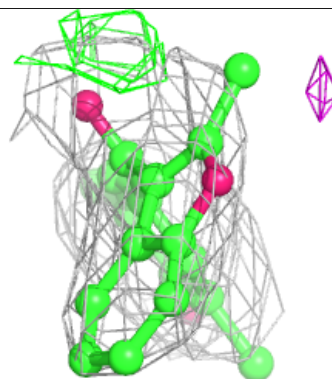
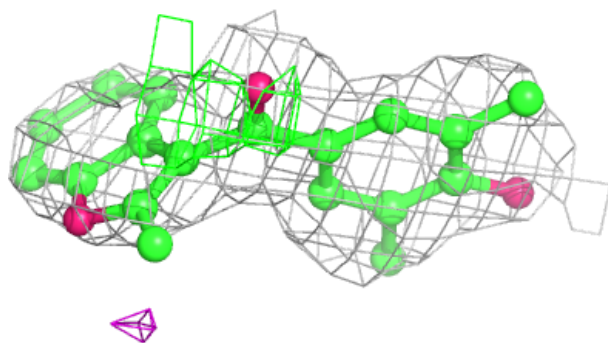
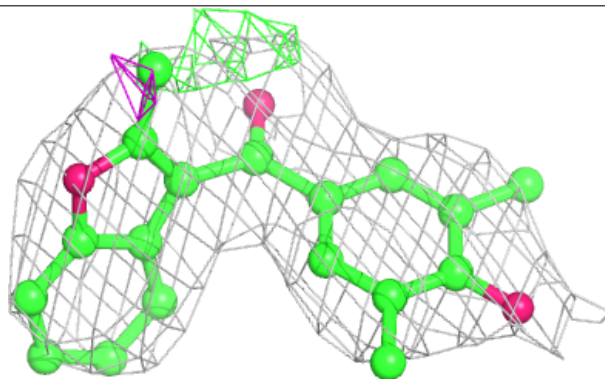
*Continued from previous page...*

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	0XV	D	502	21/21	0.88	0.34	95,97,99,99	0
3	0XV	C	502	21/21	0.89	0.39	78,81,83,83	0
4	GOL	A	503	6/6	0.89	0.18	68,70,71,72	0
3	0XV	A	502	21/21	0.95	0.28	52,54,58,58	0
2	HEM	B	501	43/43	0.97	0.28	77,80,88,90	0
2	HEM	D	501	43/43	0.97	0.34	91,95,98,99	0
2	HEM	C	501	43/43	0.98	0.22	55,59,61,66	0
2	HEM	A	501	43/43	0.98	0.19	37,44,54,57	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.

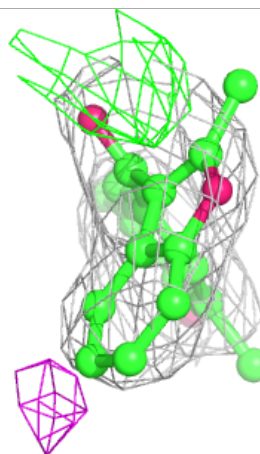
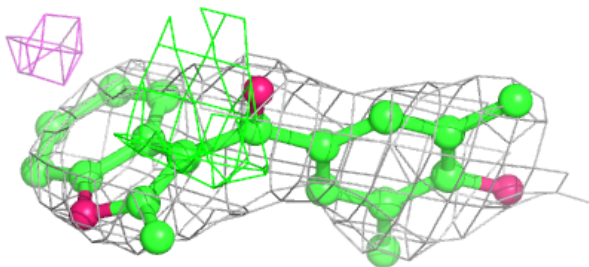
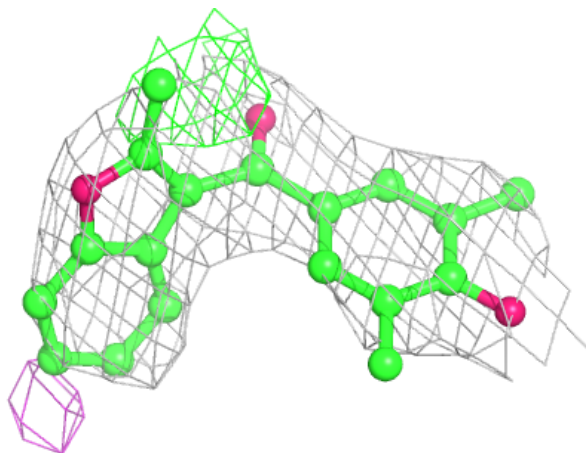
**Electron density around 0XV B 502:**

2mF<sub>o</sub>-DF<sub>c</sub> (at 0.7 rmsd) in gray  
mF<sub>o</sub>-DF<sub>c</sub> (at 3 rmsd) in purple (negative)  
and green (positive)



**Electron density around 0XV D 502:**

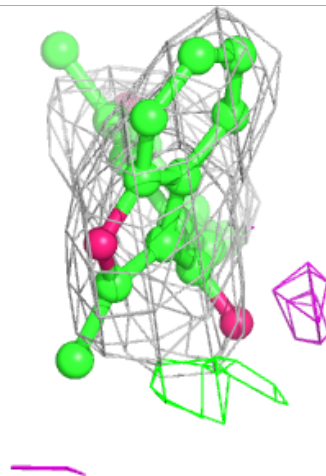
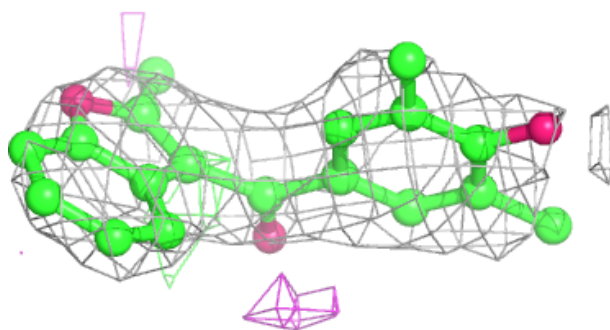
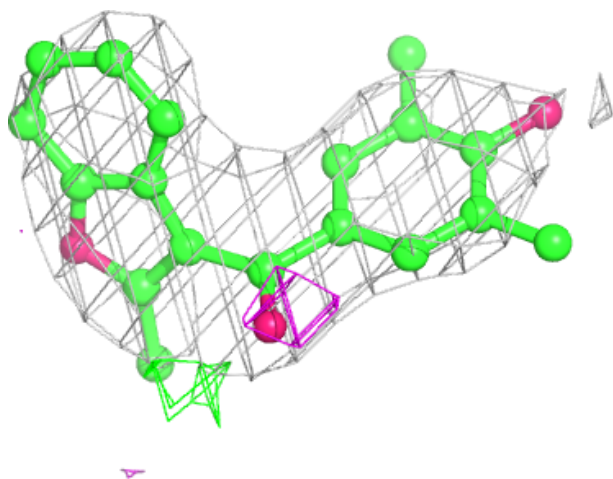
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





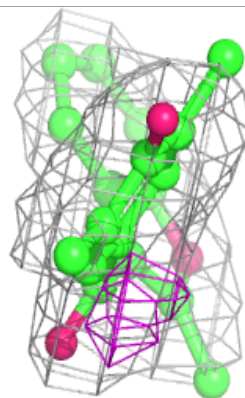
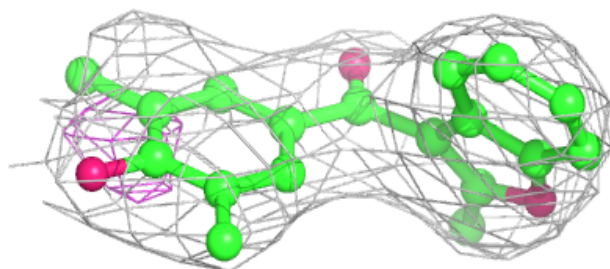
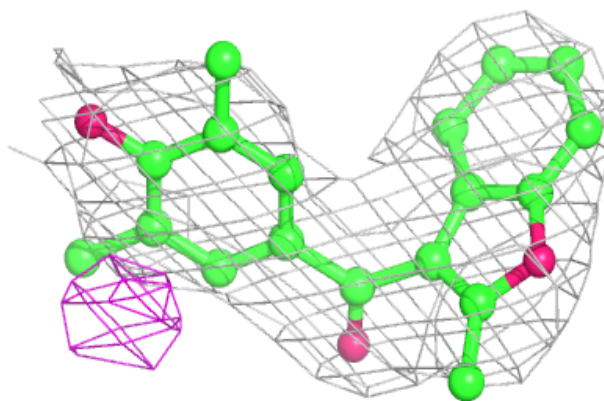
**Electron density around 0XV C 502:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

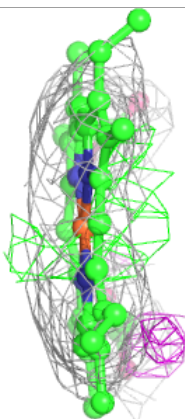
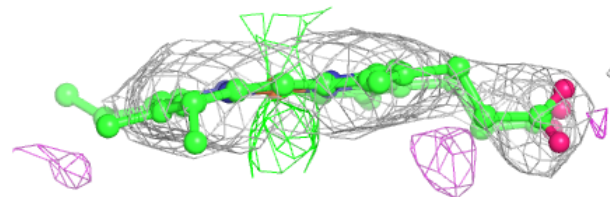
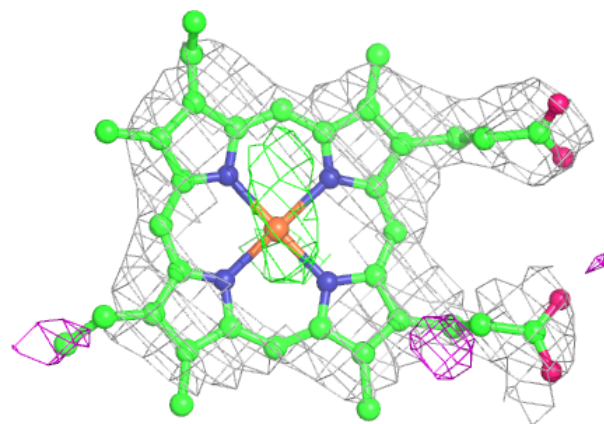


**Electron density around 0XV A 502:**

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 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

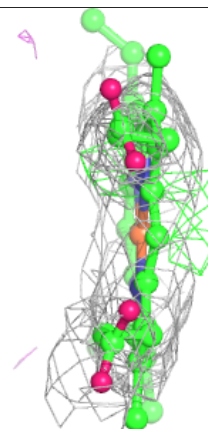
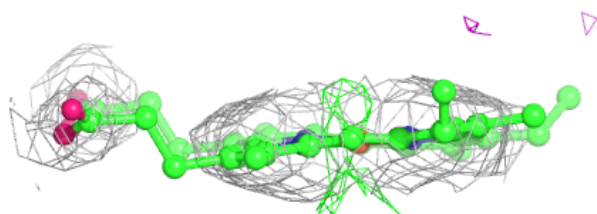
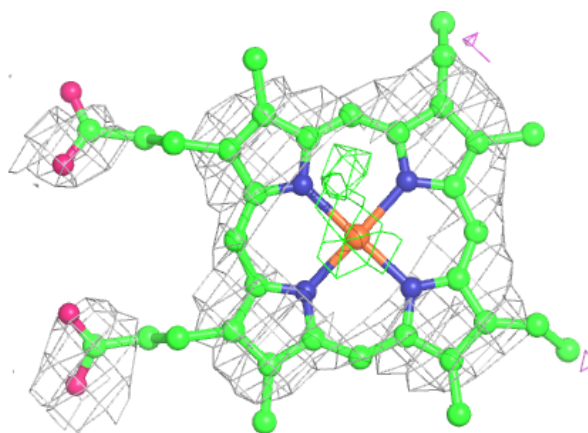
**Electron density around HEM B 501:**

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 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



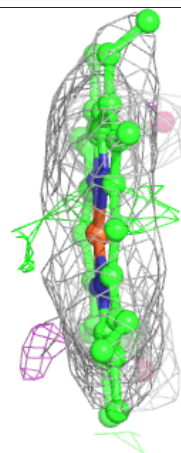
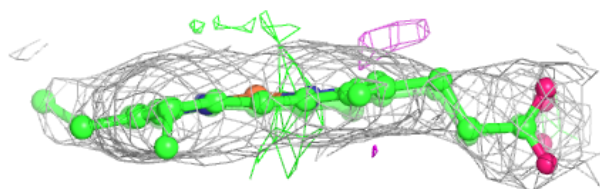
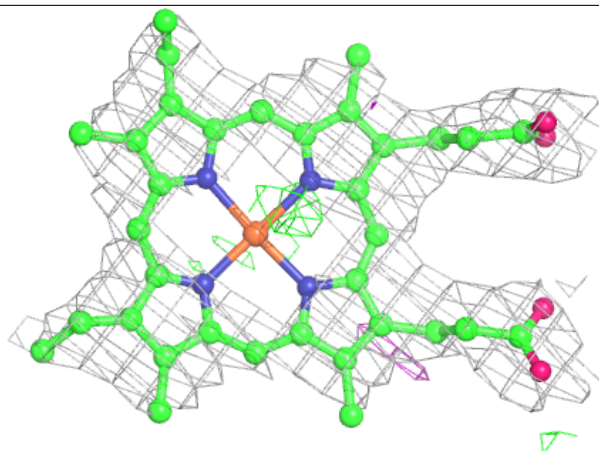
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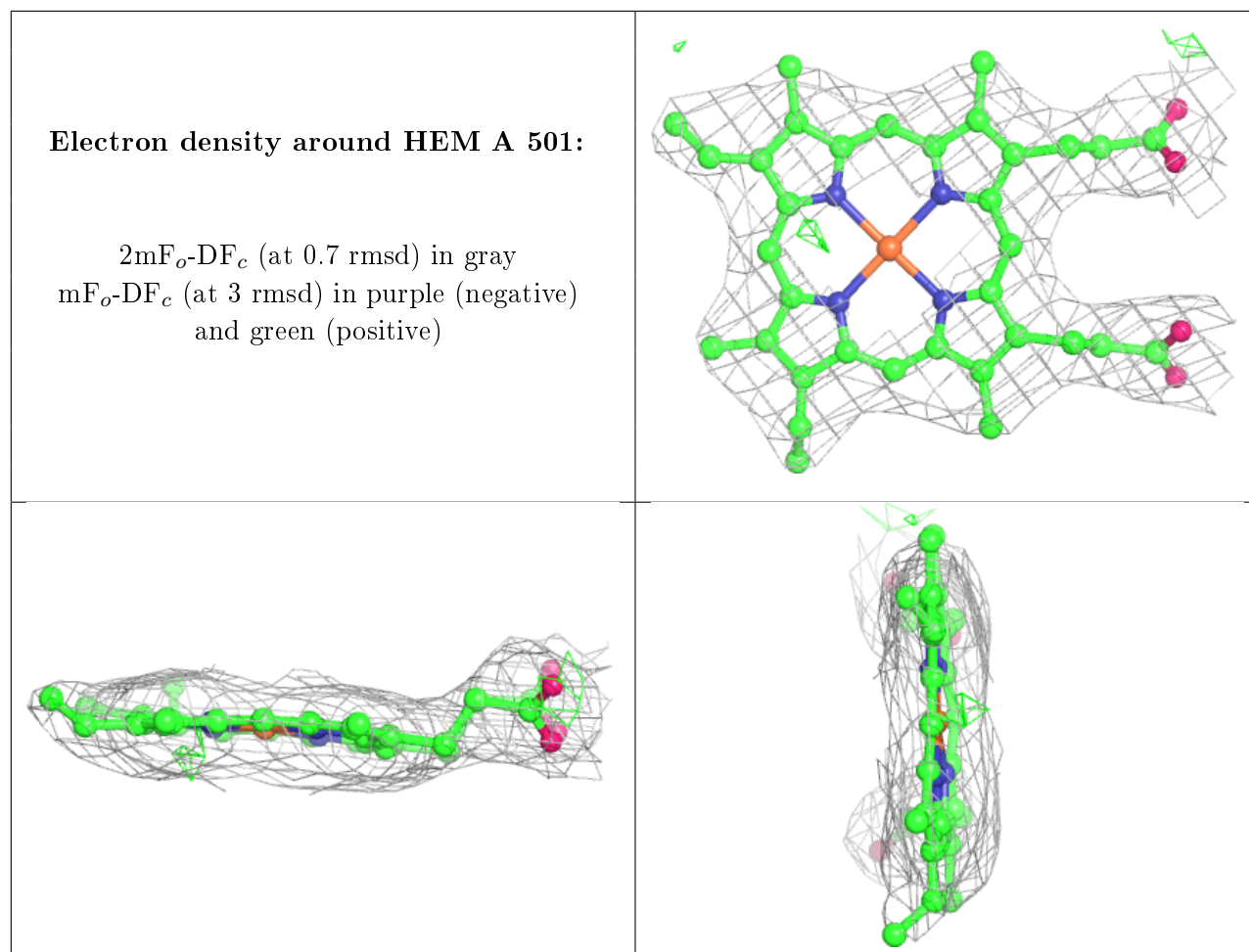
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



**Electron density around HEM C 501:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





## 6.5 Other polymers [i](#)

There are no such residues in this entry.