



Full wwPDB X-ray Structure Validation Report ⓘ

May 24, 2020 – 07:49 am BST

PDB ID : 3QM4
Title : Human Cytochrome P450 (CYP) 2D6 - Prinomastat Complex
Authors : Wang, A.; Stout, C.D.; Johnson, E.F.
Deposited on : 2011-02-03
Resolution : 2.85 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (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

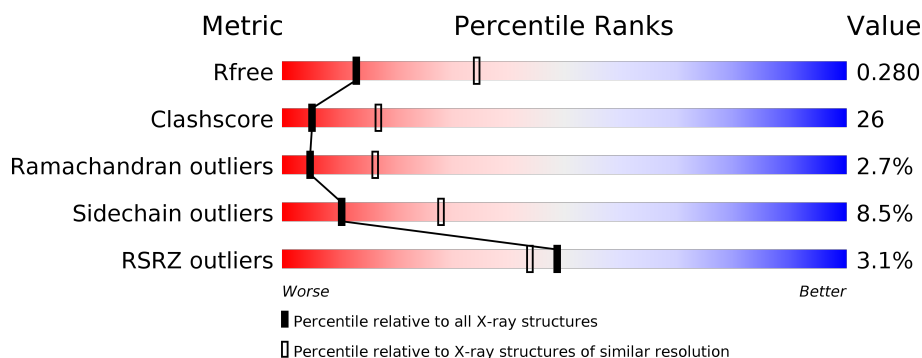
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.85 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	3168 (2.90-2.82)
Clashscore	141614	3438 (2.90-2.82)
Ramachandran outliers	138981	3348 (2.90-2.82)
Sidechain outliers	138945	3351 (2.90-2.82)
RSRZ outliers	127900	3103 (2.90-2.82)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria 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	479	<div> <div>4%</div> <div> <div></div> <div>52%</div> <div>39%</div> <div>6%</div> <div></div> </div> </div>
1	B	479	<div> <div>3%</div> <div> <div></div> <div>51%</div> <div>39%</div> <div>6%</div> <div>5%</div> </div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 7445 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 2D6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	465	Total	C	N	O	S	0	0	0
			3677	2359	651	653	14			
1	B	457	Total	C	N	O	S	0	0	0
			3618	2318	641	645	14			

There are 30 discrepancies between the modelled and reference sequences:

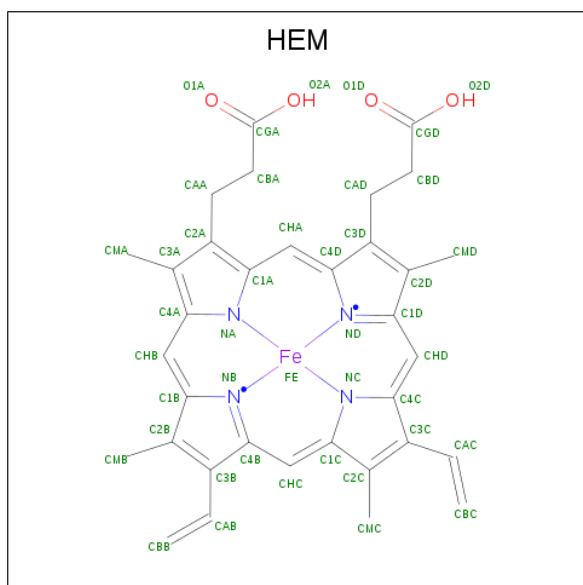
Chain	Residue	Modelled	Actual	Comment	Reference
A	23	MET	-	EXPRESSION TAG	UNP P10635
A	24	ALA	-	EXPRESSION TAG	UNP P10635
A	25	LYS	-	EXPRESSION TAG	UNP P10635
A	26	LYS	-	EXPRESSION TAG	UNP P10635
A	27	THR	-	EXPRESSION TAG	UNP P10635
A	28	SER	-	EXPRESSION TAG	UNP P10635
A	29	SER	-	EXPRESSION TAG	UNP P10635
A	30	LYS	-	EXPRESSION TAG	UNP P10635
A	31	GLY	-	EXPRESSION TAG	UNP P10635
A	32	LYS	-	EXPRESSION TAG	UNP P10635
A	33	LEU	-	EXPRESSION TAG	UNP P10635
A	498	HIS	-	EXPRESSION TAG	UNP P10635
A	499	HIS	-	EXPRESSION TAG	UNP P10635
A	500	HIS	-	EXPRESSION TAG	UNP P10635
A	501	HIS	-	EXPRESSION TAG	UNP P10635
B	23	MET	-	EXPRESSION TAG	UNP P10635
B	24	ALA	-	EXPRESSION TAG	UNP P10635
B	25	LYS	-	EXPRESSION TAG	UNP P10635
B	26	LYS	-	EXPRESSION TAG	UNP P10635
B	27	THR	-	EXPRESSION TAG	UNP P10635
B	28	SER	-	EXPRESSION TAG	UNP P10635
B	29	SER	-	EXPRESSION TAG	UNP P10635
B	30	LYS	-	EXPRESSION TAG	UNP P10635
B	31	GLY	-	EXPRESSION TAG	UNP P10635
B	32	LYS	-	EXPRESSION TAG	UNP P10635

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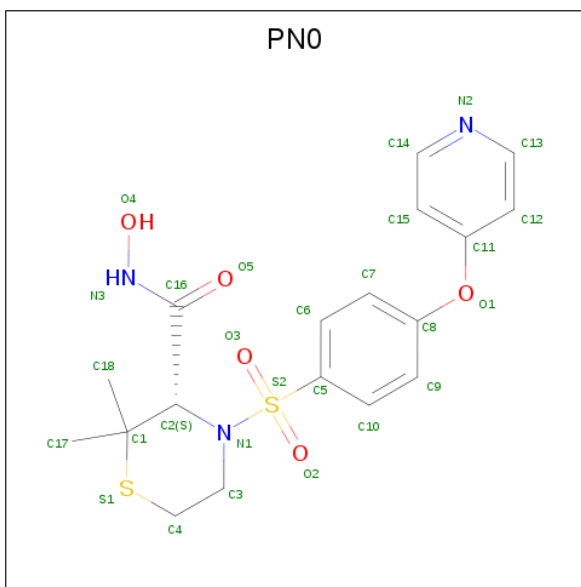
Chain	Residue	Modelled	Actual	Comment	Reference
B	33	LEU	-	EXPRESSION TAG	UNP P10635
B	498	HIS	-	EXPRESSION TAG	UNP P10635
B	499	HIS	-	EXPRESSION TAG	UNP P10635
B	500	HIS	-	EXPRESSION TAG	UNP P10635
B	501	HIS	-	EXPRESSION TAG	UNP P10635

- 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		
2	B	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- Molecule 3 is Prinomastat (three-letter code: PN0) (formula: $C_{18}H_{21}N_3O_5S_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			28	18	3	5	2		
3	B	1	Total	C	N	O	S	0	0
			28	18	3	5	2		

- Molecule 4 is NICKEL (II) ION (three-letter code: NI) (formula: Ni).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Ni	0	0
			1	1		

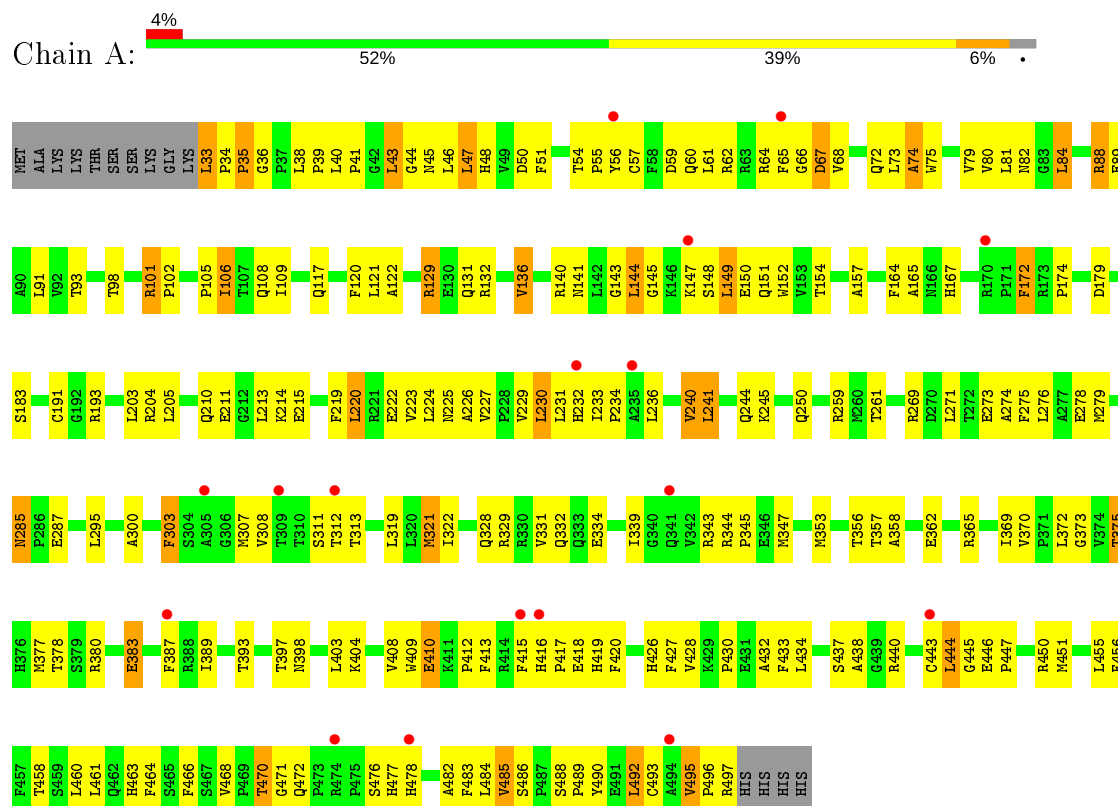
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	3	Total	O	0	0
			3	3		
5	B	4	Total	O	0	0
			4	4		

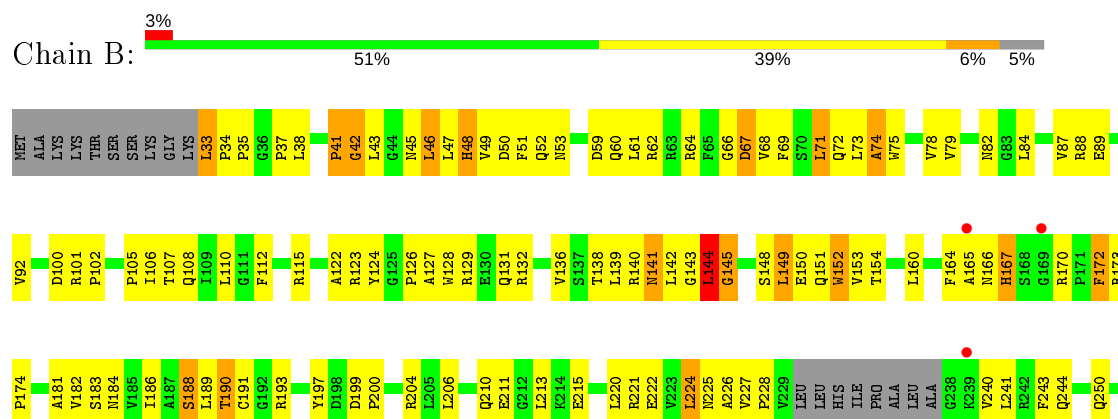
3 Residue-property plots [i](#)

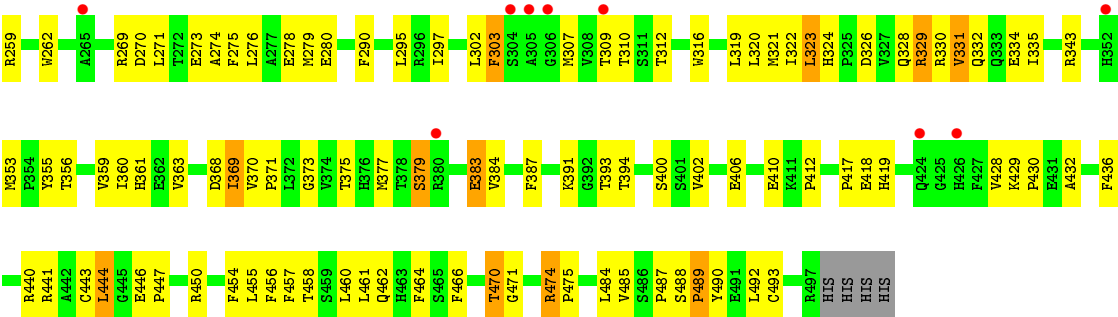
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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 ($\text{RSRZ} > 2$). 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 2D6



• Molecule 1: Cytochrome P450 2D6





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	194.54Å 55.07Å 145.87Å 90.00° 134.97° 90.00°	Depositor
Resolution (Å)	41.00 – 2.85 36.41 – 2.70	Depositor EDS
% Data completeness (in resolution range)	98.8 (41.00-2.85) 98.8 (36.41-2.70)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.28 (at 2.68Å)	Xtriage
Refinement program	CNS 1.3	Depositor
R, R_{free}	0.243 , 0.288 0.237 , 0.280	Depositor DCC
R_{free} test set	1314 reflections (4.36%)	wwPDB-VP
Wilson B-factor (Å ²)	63.9	Xtriage
Anisotropy	0.553	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 45.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7445	wwPDB-VP
Average B, all atoms (Å ²)	77.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.45% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PN0, HEM, NI

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	0.45	0/3777	0.67	0/5139
1	B	0.42	0/3715	0.65	0/5051
All	All	0.44	0/7492	0.66	0/10190

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

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	3677	0	3663	199	0
1	B	3618	0	3594	201	0
2	A	43	0	30	6	0
2	B	43	0	30	3	0
3	A	28	0	21	2	0
3	B	28	0	21	1	0
4	B	1	0	0	0	0
5	A	3	0	0	0	0
5	B	4	0	0	0	0
All	All	7445	0	7359	392	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 26.

All (392) 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:72:GLN:HG2	1:B:38:LEU:HD22	1.36	1.06
1:A:40:LEU:HD21	1:A:47:LEU:HD11	1.49	0.93
1:B:429:LYS:HZ2	1:B:450:ARG:HH12	1.17	0.92
1:A:79:VAL:HG11	1:A:389:ILE:HD11	1.54	0.89
1:B:474:ARG:HH11	1:B:474:ARG:HG2	1.37	0.88
1:A:132:ARG:O	1:A:136:VAL:HG12	1.73	0.87
1:B:312:THR:HB	1:B:369:ILE:HD11	1.55	0.87
1:A:47:LEU:H	1:A:47:LEU:HD23	1.42	0.83
1:B:110:LEU:HD12	1:B:112:PHE:HE2	1.43	0.82
1:A:380:ARG:HH11	1:A:380:ARG:HB2	1.44	0.81
1:A:149:LEU:HB3	1:A:451:MET:HE1	1.63	0.80
1:A:276:LEU:HD23	1:A:279:MET:CE	2.12	0.80
1:A:444:LEU:C	1:A:444:LEU:HD23	2.03	0.79
1:B:410:GLU:O	1:B:412:PRO:HD3	1.84	0.78
1:A:369:ILE:HG22	1:A:370:VAL:HG23	1.66	0.77
1:A:213:LEU:HD21	1:A:484:LEU:HD22	1.67	0.77
1:B:173:ARG:NH1	1:B:489:PRO:HA	1.99	0.76
1:B:107:THR:HB	1:B:112:PHE:CD2	2.21	0.75
1:B:173:ARG:HH11	1:B:490:TYR:H	1.32	0.75
1:A:105:PRO:HG3	1:B:60:GLN:OE1	1.86	0.75
1:B:211:GLU:HB3	1:B:243:PHE:HD2	1.52	0.74
1:B:474:ARG:NH1	1:B:474:ARG:HG2	1.99	0.73
1:B:102:PRO:HG3	1:B:394:THR:HG23	1.70	0.73
1:A:129:ARG:HG2	1:A:129:ARG:HH11	1.53	0.73
1:A:276:LEU:HD23	1:A:279:MET:HE1	1.69	0.73
1:B:429:LYS:NZ	1:B:450:ARG:HH12	1.85	0.73
1:B:220:LEU:HD13	1:B:220:LEU:O	1.88	0.73
1:A:105:PRO:O	1:A:108:GLN:HG3	1.90	0.72
1:B:359:VAL:O	1:B:363:VAL:HG23	1.88	0.72
1:B:67:ASP:HA	1:B:82:ASN:HB2	1.71	0.72
1:B:174:PRO:HD3	1:B:492:LEU:HD22	1.72	0.72
1:B:186:ILE:O	1:B:190:THR:HG23	1.89	0.72
1:B:211:GLU:CB	1:B:243:PHE:HD2	2.04	0.71
1:A:410:GLU:O	1:A:412:PRO:HD3	1.91	0.71
1:B:129:ARG:CZ	1:B:129:ARG:HB2	2.20	0.71
1:B:312:THR:CB	1:B:369:ILE:HD11	2.21	0.71
1:A:412:PRO:HB2	1:A:413:PHE:CD1	2.25	0.71
1:A:62:ARG:O	1:A:66:GLY:HA2	1.91	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:89:GLU:OE1	1:B:384:VAL:HA	1.91	0.70
1:A:380:ARG:NH1	1:A:380:ARG:HB2	2.04	0.70
1:A:321:MET:CE	1:A:328:GLN:HA	2.21	0.70
1:A:34:PRO:HA	1:A:387:PHE:CD2	2.26	0.70
1:B:321:MET:HE2	1:B:457:PHE:HZ	1.56	0.70
1:B:211:GLU:HB3	1:B:243:PHE:CD2	2.27	0.70
1:B:215:GLU:OE2	1:B:244:GLN:HG3	1.91	0.70
1:A:39:PRO:CG	1:A:45:ASN:HB2	2.20	0.70
1:A:43:LEU:O	1:A:46:LEU:HB2	1.93	0.69
1:B:173:ARG:HD2	1:B:490:TYR:O	1.93	0.69
1:B:220:LEU:HD13	1:B:224:LEU:HG	1.74	0.69
1:A:79:VAL:HG11	1:A:389:ILE:CD1	2.22	0.68
1:B:62:ARG:O	1:B:66:GLY:HA2	1.94	0.68
1:B:87:VAL:HG21	1:B:402:VAL:CG2	2.23	0.68
1:A:75:TRP:HE1	1:A:226:ALA:HA	1.59	0.68
1:A:179:ASP:OD1	1:A:311:SER:HB2	1.94	0.68
1:A:312:THR:HG22	1:A:369:ILE:HD11	1.76	0.68
1:B:279:MET:HG2	1:B:290:PHE:O	1.95	0.67
1:A:244:GLN:NE2	3:A:503:PN0:H3	2.10	0.67
1:B:444:LEU:C	1:B:444:LEU:HD23	2.14	0.67
1:A:339:ILE:HD11	1:A:353:MET:SD	2.35	0.67
1:A:47:LEU:H	1:A:47:LEU:CD2	2.06	0.66
1:B:182:VAL:HG11	1:B:310:THR:HB	1.77	0.66
1:A:33:LEU:O	1:A:387:PHE:HD2	1.77	0.66
1:A:285:ASN:HD22	1:A:285:ASN:H	1.44	0.66
1:A:408:VAL:HG11	1:A:432:ALA:CB	2.26	0.66
1:A:275:PHE:CD2	1:A:295:LEU:HD13	2.31	0.65
1:A:319:LEU:HB2	1:A:490:TYR:CE2	2.31	0.65
1:A:132:ARG:O	1:A:136:VAL:CG1	2.43	0.65
1:A:59:ASP:HA	1:A:62:ARG:HG2	1.79	0.64
1:B:150:GLU:O	1:B:154:THR:HG23	1.97	0.64
1:B:441:ARG:HD2	2:B:502:HEM:O2D	1.96	0.64
1:B:151:GLN:HA	1:B:154:THR:HG23	1.81	0.63
1:B:303:PHE:HE1	1:B:307:MET:CE	2.12	0.63
1:A:203:LEU:HD23	1:A:203:LEU:O	1.99	0.63
1:A:370:VAL:HG12	1:A:373:GLY:HA2	1.80	0.63
1:A:213:LEU:HD22	1:A:308:VAL:HG21	1.80	0.63
1:B:369:ILE:HG22	1:B:370:VAL:HG23	1.80	0.63
1:B:174:PRO:HD3	1:B:492:LEU:CD2	2.28	0.62
1:B:213:LEU:HD21	1:B:484:LEU:HD22	1.80	0.62
1:A:39:PRO:HG2	1:A:45:ASN:HB2	1.82	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:470:THR:HG23	1:A:471:GLY:N	2.15	0.62
1:A:91:LEU:HD11	1:A:397:THR:HG21	1.79	0.62
1:B:206:LEU:O	1:B:210:GLN:HG2	2.00	0.61
1:B:204:ARG:HG2	1:B:204:ARG:HH11	1.66	0.61
1:B:331:VAL:HG11	1:B:461:LEU:HD12	1.83	0.61
1:A:73:LEU:HD12	1:A:73:LEU:O	2.01	0.60
1:B:149:LEU:HD11	1:B:189:LEU:HD11	1.81	0.60
1:B:454:PHE:HD1	1:B:455:LEU:HD23	1.65	0.60
1:A:274:ALA:O	1:A:278:GLU:HG2	2.00	0.60
1:B:316:TRP:CD2	1:B:487:PRO:HG3	2.37	0.60
1:B:122:ALA:O	1:B:441:ARG:NH2	2.25	0.60
1:A:321:MET:HE3	1:A:328:GLN:HA	1.83	0.60
1:B:124:TYR:HE1	1:B:440:ARG:HE	1.46	0.60
1:B:206:LEU:HD23	1:B:307:MET:HE2	1.83	0.60
1:B:429:LYS:NZ	1:B:450:ARG:NH1	2.49	0.60
1:B:129:ARG:NH1	1:B:129:ARG:HB2	2.17	0.60
1:B:312:THR:CG2	1:B:369:ILE:HD11	2.32	0.60
1:B:49:VAL:HG23	1:B:49:VAL:O	2.01	0.59
1:A:428:VAL:O	1:A:430:PRO:HD3	2.02	0.59
1:A:321:MET:HE1	1:A:328:GLN:HA	1.84	0.59
1:A:365:ARG:HH22	1:A:404:LYS:HG2	1.67	0.59
1:A:440:ARG:HG3	1:A:440:ARG:HH21	1.67	0.59
1:A:236:LEU:O	1:A:240:VAL:HG23	2.03	0.59
1:A:443:CYS:HB2	2:A:502:HEM:C4A	2.37	0.59
1:A:84:LEU:HD12	1:A:84:LEU:O	2.02	0.59
1:B:259:ARG:HD3	1:B:276:LEU:HD21	1.83	0.59
1:A:191:CYS:HG	1:A:303:PHE:HE2	1.51	0.59
1:B:128:TRP:NE1	1:B:441:ARG:NE	2.51	0.59
1:A:446:GLU:HB3	1:A:447:PRO:HD3	1.84	0.58
1:B:35:PRO:HD3	1:B:387:PHE:CE2	2.38	0.58
1:B:460:LEU:O	1:B:464:PHE:HB2	2.04	0.58
1:B:278:GLU:OE2	1:B:278:GLU:HA	2.03	0.58
1:A:380:ARG:NH1	1:A:380:ARG:CB	2.67	0.58
1:B:470:THR:HG23	1:B:471:GLY:H	1.68	0.58
1:A:67:ASP:HA	1:A:82:ASN:HB2	1.85	0.57
1:B:167:HIS:ND1	1:B:172:PHE:CD1	2.71	0.57
1:B:330:ARG:HB2	1:B:355:TYR:CE2	2.38	0.57
1:A:444:LEU:HD22	2:A:502:HEM:HMD3	1.85	0.57
1:B:140:ARG:HA	1:B:144:LEU:HB3	1.86	0.57
1:A:211:GLU:O	1:A:215:GLU:HG3	2.04	0.57
1:B:303:PHE:HE1	1:B:307:MET:HE3	1.68	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:328:GLN:O	1:A:332:GLN:HG3	2.05	0.57
1:B:440:ARG:HG2	1:B:440:ARG:HH21	1.68	0.57
1:A:334:GLU:OE2	1:A:353:MET:HB3	2.05	0.57
1:B:446:GLU:O	1:B:450:ARG:HG2	2.05	0.57
1:A:129:ARG:HH11	1:A:129:ARG:CG	2.17	0.56
1:B:101:ARG:HD2	1:B:375:THR:O	2.05	0.56
1:A:259:ARG:HD3	1:A:276:LEU:CD2	2.35	0.56
1:A:285:ASN:OD1	1:A:287:GLU:HB2	2.05	0.56
1:B:361:HIS:NE2	1:B:450:ARG:NH1	2.52	0.56
1:A:117:GLN:HB3	1:A:122:ALA:HA	1.88	0.56
1:A:233:ILE:N	1:A:234:PRO:HD3	2.20	0.56
1:A:229:VAL:HG22	1:A:229:VAL:O	2.05	0.56
1:B:276:LEU:HD23	1:B:279:MET:CE	2.35	0.56
1:B:132:ARG:O	1:B:136:VAL:HG12	2.06	0.56
1:B:66:GLY:O	1:B:68:VAL:N	2.38	0.56
1:A:72:GLN:CG	1:B:38:LEU:HD22	2.25	0.56
1:B:79:VAL:HG23	1:B:393:THR:HG21	1.86	0.55
1:B:410:GLU:C	1:B:412:PRO:HD3	2.27	0.55
1:B:227:VAL:HG12	1:B:227:VAL:O	2.07	0.55
1:A:383:GLU:HG3	1:A:383:GLU:O	2.07	0.55
1:A:41:PRO:CB	1:B:42:GLY:H	2.20	0.55
1:B:138:THR:CG2	1:B:271:LEU:HD12	2.37	0.55
1:A:174:PRO:HD3	1:A:492:LEU:HD22	1.88	0.54
1:B:269:ARG:HB2	1:B:273:GLU:OE2	2.07	0.54
1:A:219:PHE:O	1:A:223:VAL:HG23	2.07	0.54
1:A:66:GLY:O	1:A:68:VAL:N	2.40	0.54
1:B:142:LEU:HD11	1:B:271:LEU:N	2.22	0.54
1:B:84:LEU:HD22	1:B:432:ALA:HB2	1.89	0.54
1:A:463:HIS:CG	1:A:463:HIS:O	2.61	0.54
1:A:88:ARG:HG2	1:A:88:ARG:HH11	1.71	0.54
1:B:126:PRO:HG2	1:B:127:ALA:H	1.73	0.54
1:B:276:LEU:HA	1:B:279:MET:CE	2.37	0.54
1:B:224:LEU:N	1:B:224:LEU:HD23	2.23	0.54
1:A:408:VAL:HG11	1:A:432:ALA:HB3	1.90	0.54
1:B:165:ALA:C	1:B:167:HIS:H	2.11	0.54
1:A:39:PRO:HG3	1:A:45:ASN:OD1	2.08	0.53
1:B:275:PHE:CD2	1:B:295:LEU:HD13	2.43	0.53
1:A:84:LEU:HD12	1:A:84:LEU:C	2.29	0.53
1:B:183:SER:O	1:B:186:ILE:HG22	2.08	0.53
1:B:456:PHE:O	1:B:460:LEU:HG	2.08	0.53
1:A:232:HIS:O	1:A:233:ILE:HD13	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:41:PRO:HB3	1:B:42:GLY:H	1.73	0.53
1:A:470:THR:HG23	1:A:471:GLY:H	1.74	0.53
1:A:233:ILE:HG21	1:A:236:LEU:HD12	1.90	0.53
1:A:215:GLU:OE1	1:A:241:LEU:HA	2.08	0.53
1:A:313:THR:OG1	1:A:369:ILE:HD12	2.09	0.53
1:A:369:ILE:HG22	1:A:370:VAL:CG2	2.37	0.53
1:A:332:GLN:NE2	1:A:497:ARG:HD2	2.24	0.53
1:A:151:GLN:HA	1:A:154:THR:OG1	2.08	0.52
1:A:416:HIS:CE1	1:A:418:GLU:HB2	2.45	0.52
1:B:274:ALA:O	1:B:278:GLU:HG2	2.09	0.52
1:A:38:LEU:N	1:A:39:PRO:HD3	2.24	0.52
1:B:138:THR:HG21	1:B:271:LEU:HD12	1.90	0.52
1:A:51:PHE:CD2	1:A:222:GLU:HG3	2.45	0.52
1:B:370:VAL:HG12	1:B:373:GLY:HA2	1.91	0.52
1:A:437:SER:OG	1:A:438:ALA:N	2.41	0.52
1:B:271:LEU:CD2	1:B:302:LEU:HD12	2.38	0.52
1:B:446:GLU:HB3	1:B:447:PRO:HD3	1.92	0.52
1:A:147:LYS:HA	1:A:151:GLN:NE2	2.25	0.51
1:A:224:LEU:HD12	1:A:240:VAL:HG21	1.92	0.51
1:A:443:CYS:HB2	2:A:502:HEM:NA	2.25	0.51
1:B:330:ARG:CB	1:B:355:TYR:CE2	2.94	0.51
1:A:140:ARG:HA	1:A:144:LEU:HB3	1.91	0.51
1:B:123:ARG:NH2	1:B:377:MET:CE	2.73	0.51
1:A:428:VAL:HG23	1:A:428:VAL:O	2.11	0.51
1:A:64:ARG:HD2	1:A:65:PHE:CE1	2.46	0.51
1:B:184:ASN:ND2	1:B:197:TYR:HE2	2.09	0.51
1:B:141:ASN:C	1:B:143:GLY:H	2.14	0.51
1:A:412:PRO:HB2	1:A:413:PHE:HD1	1.71	0.51
1:B:369:ILE:O	1:B:371:PRO:HD3	2.10	0.51
1:A:259:ARG:HD3	1:A:276:LEU:HD22	1.91	0.50
1:B:417:PRO:C	1:B:419:HIS:H	2.14	0.50
1:B:356:THR:O	1:B:360:ILE:HG13	2.11	0.50
1:A:51:PHE:CE2	1:A:222:GLU:HG3	2.45	0.50
1:B:100:ASP:HA	1:B:124:TYR:HB2	1.94	0.50
1:B:172:PHE:N	1:B:172:PHE:CD2	2.80	0.50
1:B:191:CYS:HG	1:B:303:PHE:HE2	1.58	0.50
1:B:444:LEU:O	1:B:444:LEU:HD23	2.12	0.50
1:B:319:LEU:O	1:B:322:ILE:HG12	2.11	0.50
1:B:141:ASN:ND2	1:B:141:ASN:O	2.45	0.50
1:B:244:GLN:NE2	3:B:503:PN0:H3	2.27	0.50
1:B:332:GLN:NE2	1:B:464:PHE:O	2.45	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:241:LEU:O	1:A:245:LYS:HD3	2.13	0.49
1:A:332:GLN:HE21	1:A:497:ARG:HH11	1.59	0.49
1:B:59:ASP:HA	1:B:62:ARG:HG2	1.94	0.49
1:B:383:GLU:O	1:B:383:GLU:HG3	2.12	0.49
1:B:33:LEU:O	1:B:387:PHE:HD2	1.96	0.49
1:A:227:VAL:O	1:A:229:VAL:N	2.39	0.49
1:B:139:LEU:HB3	1:B:144:LEU:HD12	1.94	0.49
1:A:120:PHE:HD2	1:A:121:LEU:HG	1.78	0.49
1:A:36:GLY:C	1:A:65:PHE:CE2	2.86	0.49
1:B:73:LEU:O	1:B:74:ALA:HB3	2.12	0.49
1:A:231:LEU:O	1:A:232:HIS:HB2	2.12	0.48
1:A:440:ARG:NH2	1:A:440:ARG:HG3	2.28	0.48
1:B:34:PRO:HA	1:B:387:PHE:CD2	2.47	0.48
1:A:220:LEU:HD12	1:A:240:VAL:CG1	2.42	0.48
1:A:444:LEU:HD22	2:A:502:HEM:CMD	2.43	0.48
1:A:106:ILE:HG22	1:A:225:ASN:OD1	2.14	0.48
1:A:150:GLU:OE1	1:A:347:MET:N	2.44	0.48
1:A:165:ALA:C	1:A:167:HIS:H	2.17	0.48
1:A:476:SER:OG	1:A:478:HIS:ND1	2.47	0.48
1:B:428:VAL:O	1:B:430:PRO:HD3	2.14	0.48
1:A:276:LEU:HA	1:A:279:MET:CE	2.44	0.48
1:A:444:LEU:HD23	1:A:445:GLY:N	2.28	0.48
1:A:89:GLU:HA	1:A:93:THR:OG1	2.13	0.48
1:B:428:VAL:O	1:B:428:VAL:HG23	2.14	0.48
1:A:80:VAL:C	1:A:81:LEU:HD23	2.34	0.48
1:B:128:TRP:CE2	1:B:441:ARG:NE	2.81	0.48
1:B:334:GLU:OE2	1:B:353:MET:HB3	2.14	0.48
1:B:87:VAL:HG21	1:B:402:VAL:HG21	1.93	0.47
1:A:357:THR:HG22	1:A:427:PHE:CD1	2.50	0.47
1:B:172:PHE:N	1:B:172:PHE:HD2	2.11	0.47
1:A:74:ALA:CB	1:A:226:ALA:HB1	2.44	0.47
1:A:230:LEU:HD23	1:A:231:LEU:N	2.29	0.47
1:A:370:VAL:HG12	1:A:373:GLY:CA	2.45	0.47
1:B:164:PHE:O	1:B:167:HIS:HB2	2.14	0.47
1:B:123:ARG:NH2	1:B:377:MET:HE1	2.30	0.47
1:B:331:VAL:CG1	1:B:461:LEU:HD12	2.44	0.47
1:A:362:GLU:OE1	1:A:420:PHE:HE1	1.98	0.47
1:B:131:GLN:OE1	1:B:131:GLN:HA	2.13	0.47
1:B:328:GLN:O	1:B:332:GLN:HG3	2.15	0.47
1:A:39:PRO:HG3	1:A:45:ASN:CG	2.35	0.47
1:A:39:PRO:HG3	1:A:45:ASN:HB2	1.93	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:48:HIS:NE2	1:A:61:LEU:HD21	2.30	0.47
1:B:262:TRP:CE3	1:B:276:LEU:HD13	2.50	0.47
1:A:232:HIS:C	1:A:234:PRO:HD3	2.36	0.46
1:A:433:PHE:O	1:A:434:LEU:HD23	2.15	0.46
1:B:335:ILE:HD13	1:B:462:GLN:HB2	1.96	0.46
1:B:173:ARG:NH1	1:B:490:TYR:H	2.05	0.46
1:A:39:PRO:HG3	1:A:45:ASN:CB	2.45	0.46
1:A:75:TRP:CD1	1:A:75:TRP:N	2.82	0.46
1:B:343:ARG:O	1:B:343:ARG:HG3	2.15	0.46
1:A:477:HIS:N	1:A:477:HIS:ND1	2.61	0.46
1:B:129:ARG:NH1	1:B:129:ARG:CB	2.79	0.46
1:A:231:LEU:CD2	1:B:52:GLN:HB2	2.45	0.46
1:A:75:TRP:CZ3	1:B:61:LEU:HD21	2.49	0.46
1:B:276:LEU:HA	1:B:279:MET:HE2	1.95	0.46
1:A:485:VAL:O	1:A:485:VAL:HG22	2.16	0.46
1:B:220:LEU:CD1	1:B:224:LEU:HG	2.43	0.46
1:A:466:PHE:HA	1:A:493:CYS:O	2.16	0.46
1:B:356:THR:HG21	1:B:458:THR:HG22	1.97	0.46
1:A:106:ILE:HG12	1:A:106:ILE:O	2.16	0.46
1:A:84:LEU:HD22	1:A:432:ALA:HA	1.97	0.46
1:A:141:ASN:C	1:A:143:GLY:H	2.19	0.46
1:A:413:PHE:N	1:A:413:PHE:CD1	2.84	0.46
1:A:75:TRP:HD1	1:A:75:TRP:H	1.63	0.46
1:B:224:LEU:HD12	1:B:240:VAL:HG21	1.98	0.46
1:B:221:ARG:O	1:B:225:ASN:HB2	2.16	0.46
1:A:358:ALA:HB1	1:A:420:PHE:HB2	1.98	0.45
1:A:482:ALA:O	1:A:483:PHE:HB3	2.15	0.45
1:B:224:LEU:H	1:B:224:LEU:HD23	1.81	0.45
1:B:276:LEU:HD23	1:B:279:MET:HE3	1.98	0.45
1:A:269:ARG:HG2	1:A:273:GLU:OE2	2.14	0.45
1:B:126:PRO:HG2	1:B:127:ALA:N	2.31	0.45
1:A:461:LEU:HD23	1:A:466:PHE:HE1	1.82	0.45
1:B:444:LEU:HD22	2:B:502:HEM:HMD2	1.99	0.45
1:B:115:ARG:HG2	1:B:115:ARG:HH11	1.82	0.45
1:B:443:CYS:HB2	2:B:502:HEM:NA	2.31	0.45
1:B:470:THR:HG23	1:B:471:GLY:N	2.32	0.45
1:A:120:PHE:CD2	1:A:121:LEU:HG	2.52	0.45
1:A:331:VAL:HG11	1:A:461:LEU:HD12	1.99	0.45
1:B:440:ARG:NH2	1:B:440:ARG:HG2	2.31	0.45
1:B:74:ALA:HB3	1:B:226:ALA:HB1	1.99	0.45
1:A:39:PRO:CG	1:A:45:ASN:CB	2.94	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:88:ARG:CG	1:A:88:ARG:HH11	2.31	0.44
1:B:323:LEU:HD11	1:B:475:PRO:O	2.17	0.44
1:A:210:GLN:O	1:A:214:LYS:HG3	2.18	0.44
1:A:40:LEU:HB2	1:A:41:PRO:HD2	2.00	0.44
1:A:490:TYR:CD1	1:A:490:TYR:C	2.91	0.44
1:B:320:LEU:O	1:B:324:HIS:HB2	2.18	0.44
1:B:45:ASN:HA	1:B:48:HIS:CE1	2.53	0.44
1:A:357:THR:HG22	1:A:427:PHE:CE1	2.53	0.44
1:A:39:PRO:HB3	1:A:45:ASN:CA	2.48	0.44
1:A:56:TYR:O	1:A:60:GLN:HG2	2.17	0.44
1:B:370:VAL:HG12	1:B:373:GLY:CA	2.48	0.44
1:B:466:PHE:HA	1:B:493:CYS:O	2.18	0.44
1:A:39:PRO:HB3	1:A:44:GLY:C	2.38	0.44
1:B:101:ARG:NH1	1:B:441:ARG:HH11	2.16	0.44
1:A:308:VAL:O	1:A:308:VAL:HG23	2.18	0.44
1:B:370:VAL:CG1	1:B:373:GLY:HA2	2.48	0.44
1:B:68:VAL:O	1:B:68:VAL:HG12	2.17	0.44
1:B:43:LEU:O	1:B:74:ALA:HA	2.18	0.44
1:A:33:LEU:HD12	1:A:387:PHE:HA	2.00	0.43
1:A:365:ARG:HD2	1:A:409:TRP:CE2	2.52	0.43
1:B:188:SER:OG	1:B:189:LEU:N	2.50	0.43
1:B:270:ASP:OD2	1:B:273:GLU:HG3	2.18	0.43
1:A:444:LEU:CD2	2:A:502:HEM:HHD	2.48	0.43
1:A:444:LEU:HD22	2:A:502:HEM:C2D	2.53	0.43
1:A:172:PHE:N	1:A:172:PHE:CD2	2.85	0.43
1:A:172:PHE:N	1:A:172:PHE:HD2	2.16	0.43
1:A:451:MET:HG2	1:A:455:LEU:HD12	2.01	0.43
1:B:160:LEU:HD13	1:B:181:ALA:HB2	2.01	0.43
1:B:262:TRP:CZ3	1:B:276:LEU:HB3	2.52	0.43
1:B:43:LEU:HG	1:B:46:LEU:HD22	1.99	0.43
1:B:112:PHE:CD1	1:B:297:ILE:HD13	2.52	0.43
1:A:102:PRO:HB2	1:B:64:ARG:HH22	1.83	0.43
1:B:316:TRP:CG	1:B:487:PRO:HG3	2.54	0.43
1:A:460:LEU:O	1:A:464:PHE:HB2	2.17	0.43
1:B:276:LEU:HA	1:B:279:MET:HE3	2.00	0.43
1:B:45:ASN:HA	1:B:48:HIS:HE1	1.83	0.43
1:A:300:ALA:HB1	3:A:503:PN0:H17	2.01	0.43
1:B:143:GLY:O	1:B:145:GLY:N	2.51	0.43
1:B:368:ASP:OD1	1:B:400:SER:HA	2.18	0.43
1:B:69:PHE:CE1	1:B:71:LEU:HD21	2.53	0.43
1:A:109:ILE:HG22	1:A:109:ILE:O	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:131:GLN:HA	1:A:131:GLN:OE1	2.19	0.43
1:B:33:LEU:N	1:B:34:PRO:CD	2.82	0.43
1:B:41:PRO:O	1:B:42:GLY:C	2.57	0.43
1:B:241:LEU:CD1	1:B:241:LEU:N	2.81	0.43
1:B:33:LEU:N	1:B:34:PRO:HD3	2.34	0.43
1:A:276:LEU:HA	1:A:279:MET:HE2	2.00	0.42
1:A:450:ARG:HH11	1:A:450:ARG:HG3	1.84	0.42
1:B:379:SER:O	1:B:391:LYS:HE3	2.19	0.42
1:A:470:THR:CG2	1:A:471:GLY:N	2.82	0.42
1:B:276:LEU:HD23	1:B:279:MET:HE1	2.00	0.42
1:A:495:VAL:HA	1:A:496:PRO:HD3	1.72	0.42
1:B:105:PRO:O	1:B:108:GLN:HG3	2.19	0.42
1:B:279:MET:O	1:B:280:GLU:C	2.56	0.42
1:A:50:ASP:HB3	1:A:57:CYS:SG	2.59	0.42
1:A:98:THR:HG22	1:A:378:THR:HG22	2.00	0.42
1:B:226:ALA:O	1:B:227:VAL:HG23	2.20	0.42
1:B:406:GLU:HG2	1:B:406:GLU:H	1.67	0.42
1:B:316:TRP:CE2	1:B:487:PRO:HG3	2.53	0.42
1:A:54:THR:HB	1:A:55:PRO:HD3	2.02	0.42
1:A:417:PRO:C	1:A:419:HIS:H	2.23	0.42
1:B:436:PHE:CE1	1:B:446:GLU:HG3	2.54	0.42
1:A:344:ARG:HA	1:A:345:PRO:HD3	1.93	0.42
1:B:152:TRP:O	1:B:152:TRP:HE3	2.03	0.42
1:A:102:PRO:HA	1:A:377:MET:SD	2.60	0.42
1:A:372:LEU:HA	1:A:398:ASN:HA	2.02	0.42
1:B:309:THR:OG1	1:B:310:THR:N	2.52	0.42
1:B:488:SER:O	1:B:489:PRO:C	2.58	0.42
1:B:321:MET:HE2	1:B:457:PHE:CZ	2.45	0.42
1:A:205:LEU:HA	1:A:205:LEU:HD12	1.78	0.42
1:A:65:PHE:N	1:A:65:PHE:CD1	2.88	0.42
1:A:157:ALA:HA	1:A:456:PHE:HE1	1.85	0.41
1:A:233:ILE:N	1:A:234:PRO:CD	2.80	0.41
1:A:389:ILE:HG23	1:A:393:THR:HG21	2.01	0.41
1:B:149:LEU:O	1:B:153:VAL:HG23	2.20	0.41
1:B:199:ASP:HA	1:B:200:PRO:HD3	1.91	0.41
1:A:410:GLU:C	1:A:412:PRO:HD3	2.41	0.41
1:B:123:ARG:HE	1:B:123:ARG:HB3	1.59	0.41
1:A:220:LEU:HD22	1:A:220:LEU:HA	1.87	0.41
1:B:269:ARG:HH11	1:B:269:ARG:HG2	1.85	0.41
1:B:329:ARG:NH1	1:B:329:ARG:HG3	2.36	0.41
1:B:51:PHE:CZ	1:B:222:GLU:HG3	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:183:SER:OG	1:A:307:MET:HG3	2.21	0.41
1:A:370:VAL:CG1	1:A:373:GLY:HA2	2.48	0.41
1:A:81:LEU:HD23	1:A:81:LEU:N	2.35	0.41
1:B:164:PHE:CE2	1:B:460:LEU:HD22	2.56	0.41
1:A:356:THR:HG21	1:A:458:THR:CG2	2.50	0.41
1:A:483:PHE:CD2	1:A:484:LEU:HD13	2.56	0.41
1:B:271:LEU:HD22	1:B:302:LEU:HD12	2.03	0.41
1:B:78:VAL:HG12	1:B:79:VAL:N	2.36	0.41
1:B:303:PHE:CE1	1:B:307:MET:HE1	2.55	0.41
1:B:79:VAL:CG2	1:B:393:THR:HG21	2.51	0.41
1:B:417:PRO:O	1:B:419:HIS:N	2.54	0.41
1:A:121:LEU:HA	1:A:121:LEU:HD23	1.95	0.41
1:B:170:ARG:HA	1:B:170:ARG:HD2	1.96	0.41
1:B:417:PRO:C	1:B:419:HIS:N	2.74	0.41
1:B:88:ARG:O	1:B:92:VAL:HB	2.21	0.41
1:A:147:LYS:HA	1:A:151:GLN:HE21	1.85	0.41
1:A:47:LEU:N	1:A:47:LEU:CD2	2.76	0.41
1:A:319:LEU:O	1:A:322:ILE:HG12	2.20	0.41
1:B:107:THR:O	1:B:112:PHE:HD2	2.04	0.41
1:A:426:HIS:CD2	1:A:426:HIS:N	2.89	0.40
1:B:204:ARG:HG2	1:B:204:ARG:NH1	2.35	0.40
1:A:101:ARG:HD2	1:A:375:THR:O	2.21	0.40
1:A:35:PRO:HD3	1:A:387:PHE:CE2	2.57	0.40
1:A:40:LEU:HD12	1:A:40:LEU:C	2.42	0.40
1:A:362:GLU:HG3	1:A:415:PHE:CE1	2.56	0.40
1:B:444:LEU:C	1:B:444:LEU:CD2	2.86	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	463/479 (97%)	396 (86%)	58 (12%)	9 (2%)	8	24
1	B	453/479 (95%)	382 (84%)	55 (12%)	16 (4%)	3	12
All	All	916/958 (96%)	778 (85%)	113 (12%)	25 (3%)	5	16

All (25) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	67	ASP
1	A	470	THR
1	B	67	ASP
1	B	470	THR
1	B	42	GLY
1	A	74	ALA
1	B	37	PRO
1	B	144	LEU
1	B	418	GLU
1	B	489	PRO
1	A	35	PRO
1	B	74	ALA
1	A	148	SER
1	B	46	LEU
1	B	50	ASP
1	B	148	SER
1	B	166	ASN
1	A	410	GLU
1	B	41	PRO
1	B	106	ILE
1	B	228	PRO
1	A	106	ILE
1	A	145	GLY
1	B	145	GLY
1	A	489	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	397/409 (97%)	359 (90%)	38 (10%)	8	22
1	B	391/409 (96%)	362 (93%)	29 (7%)	13	35
All	All	788/818 (96%)	721 (92%)	67 (8%)	10	28

All (67) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	33	LEU
1	A	43	LEU
1	A	47	LEU
1	A	84	LEU
1	A	88	ARG
1	A	101	ARG
1	A	129	ARG
1	A	136	VAL
1	A	144	LEU
1	A	149	LEU
1	A	152	TRP
1	A	164	PHE
1	A	172	PHE
1	A	193	ARG
1	A	204	ARG
1	A	220	LEU
1	A	230	LEU
1	A	240	VAL
1	A	241	LEU
1	A	250	GLN
1	A	261	THR
1	A	271	LEU
1	A	285	ASN
1	A	303	PHE
1	A	321	MET
1	A	329	ARG
1	A	343	ARG
1	A	375	THR
1	A	383	GLU
1	A	403	LEU
1	A	444	LEU
1	A	468	VAL
1	A	472	GLN
1	A	485	VAL
1	A	486	SER

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Mol	Chain	Res	Type
1	A	488	SER
1	A	492	LEU
1	A	495	VAL
1	B	33	LEU
1	B	47	LEU
1	B	48	HIS
1	B	53	ASN
1	B	71	LEU
1	B	72	GLN
1	B	75	TRP
1	B	141	ASN
1	B	144	LEU
1	B	149	LEU
1	B	152	TRP
1	B	167	HIS
1	B	172	PHE
1	B	188	SER
1	B	190	THR
1	B	193	ARG
1	B	224	LEU
1	B	250	GLN
1	B	303	PHE
1	B	323	LEU
1	B	326	ASP
1	B	329	ARG
1	B	331	VAL
1	B	369	ILE
1	B	379	SER
1	B	383	GLU
1	B	444	LEU
1	B	474	ARG
1	B	485	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	53	ASN
1	A	141	ASN
1	A	151	GLN
1	A	250	GLN
1	A	332	GLN
1	A	426	HIS

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Mol	Chain	Res	Type
1	B	48	HIS
1	B	53	ASN
1	B	72	GLN
1	B	141	ASN
1	B	151	GLN
1	B	210	GLN
1	B	332	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 5 ligands modelled in this entry, 1 is monoatomic - leaving 4 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	PN0	B	503	2	28,30,30	1.80	6 (21%)	36,44,44	1.73	9 (25%)
2	HEM	A	502	1,3	27,50,50	1.87	10 (37%)	17,82,82	1.54	3 (17%)
2	HEM	B	502	1,3	27,50,50	1.92	9 (33%)	17,82,82	1.83	3 (17%)
3	PN0	A	503	2	28,30,30	1.77	8 (28%)	36,44,44	1.60	8 (22%)

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
3	PN0	B	503	2	-	8/21/39/39	0/3/3/3
2	HEM	A	502	1,3	-	0/6/54/54	-
2	HEM	B	502	1,3	-	0/6/54/54	-
3	PN0	A	503	2	-	8/21/39/39	0/3/3/3

All (33) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	502	HEM	C3C-CAC	-5.47	1.36	1.47
2	A	502	HEM	C3C-CAC	-4.98	1.37	1.47
3	B	503	PN0	C1-S1	-4.52	1.80	1.84
3	A	503	PN0	C1-S1	-3.88	1.80	1.84
3	B	503	PN0	C15-C11	3.68	1.46	1.38
2	A	502	HEM	C3B-CAB	-3.39	1.41	1.47
2	A	502	HEM	CMA-C3A	3.09	1.58	1.51
3	A	503	PN0	C15-C11	3.07	1.44	1.38
3	A	503	PN0	C5-S2	-3.04	1.72	1.76
3	A	503	PN0	C2-C16	3.03	1.58	1.53
2	B	502	HEM	C3B-CAB	-2.92	1.42	1.47
2	A	502	HEM	CBB-CAB	2.84	1.48	1.29
2	B	502	HEM	CBB-CAB	2.80	1.47	1.29
2	B	502	HEM	CMA-C3A	2.79	1.57	1.51
3	B	503	PN0	C9-C8	2.77	1.44	1.38
3	B	503	PN0	C5-S2	-2.72	1.72	1.76
3	A	503	PN0	C6-C5	2.64	1.43	1.38
2	B	502	HEM	C4B-NB	-2.54	1.30	1.36
2	B	502	HEM	C1D-ND	-2.54	1.30	1.36
3	B	503	PN0	C6-C5	2.53	1.42	1.38
2	A	502	HEM	CMC-C2C	2.50	1.57	1.51
2	A	502	HEM	C1D-ND	-2.44	1.31	1.36
3	A	503	PN0	C9-C8	2.40	1.43	1.38
2	A	502	HEM	C4B-NB	-2.38	1.31	1.36
2	A	502	HEM	C1D-CHD	-2.36	1.34	1.41
3	B	503	PN0	C10-C5	2.35	1.42	1.38
2	B	502	HEM	CMC-C2C	2.31	1.57	1.51
2	B	502	HEM	CAD-C3D	2.29	1.56	1.52
2	A	502	HEM	C4A-NA	-2.19	1.31	1.36
3	A	503	PN0	C7-C8	2.18	1.43	1.38
3	A	503	PN0	O3-S2	2.10	1.45	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	502	HEM	C1D-CHD	-2.07	1.35	1.41
2	A	502	HEM	CBC-CAC	2.03	1.42	1.29

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	503	PN0	C11-O1-C8	4.02	128.21	118.80
3	A	503	PN0	C18-C1-C17	-3.83	105.15	109.63
3	B	503	PN0	C18-C1-C17	-3.77	105.22	109.63
2	B	502	HEM	CAD-CBD-CGD	3.66	118.81	112.67
3	A	503	PN0	O2-S2-C5	-3.56	103.54	108.05
2	B	502	HEM	C4C-C3C-C2C	-3.37	104.55	106.90
3	B	503	PN0	C3-N1-C2	3.20	122.92	115.70
2	A	502	HEM	CMB-C2B-C3B	3.17	130.61	124.68
3	A	503	PN0	C11-O1-C8	3.13	126.12	118.80
2	A	502	HEM	C4C-C3C-C2C	-3.10	104.73	106.90
2	B	502	HEM	CMB-C2B-C3B	3.10	130.47	124.68
3	A	503	PN0	C3-N1-C2	3.07	122.62	115.70
3	B	503	PN0	C6-C5-S2	-2.77	116.84	119.76
3	B	503	PN0	O3-S2-O2	-2.75	115.06	119.52
3	B	503	PN0	O1-C11-C15	2.64	128.68	119.38
2	A	502	HEM	CAD-CBD-CGD	2.64	117.10	112.67
3	B	503	PN0	O2-S2-C5	-2.62	104.72	108.05
3	B	503	PN0	C10-C5-S2	2.42	122.31	119.76
3	B	503	PN0	O1-C11-C12	-2.38	110.99	119.38
3	A	503	PN0	O1-C11-C15	2.38	127.75	119.38
3	A	503	PN0	O3-S2-O2	-2.17	116.01	119.52
3	A	503	PN0	C3-C4-S1	2.06	115.10	112.52
3	A	503	PN0	O1-C11-C12	-2.06	112.15	119.38

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	503	PN0	C3-N1-S2-O3
3	B	503	PN0	C3-N1-S2-O3
3	A	503	PN0	N3-C16-C2-N1
3	A	503	PN0	C3-N1-S2-O2
3	A	503	PN0	O5-C16-C2-N1
3	B	503	PN0	C3-N1-S2-O2
3	A	503	PN0	C6-C5-S2-O2
3	B	503	PN0	C6-C5-S2-O2

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Mol	Chain	Res	Type	Atoms
3	A	503	PN0	C10-C5-S2-O2
3	B	503	PN0	C10-C5-S2-O2
3	B	503	PN0	N3-C16-C2-N1
3	A	503	PN0	C6-C5-S2-N1
3	B	503	PN0	O5-C16-C2-N1
3	B	503	PN0	C6-C5-S2-N1
3	A	503	PN0	C10-C5-S2-N1
3	B	503	PN0	C10-C5-S2-N1

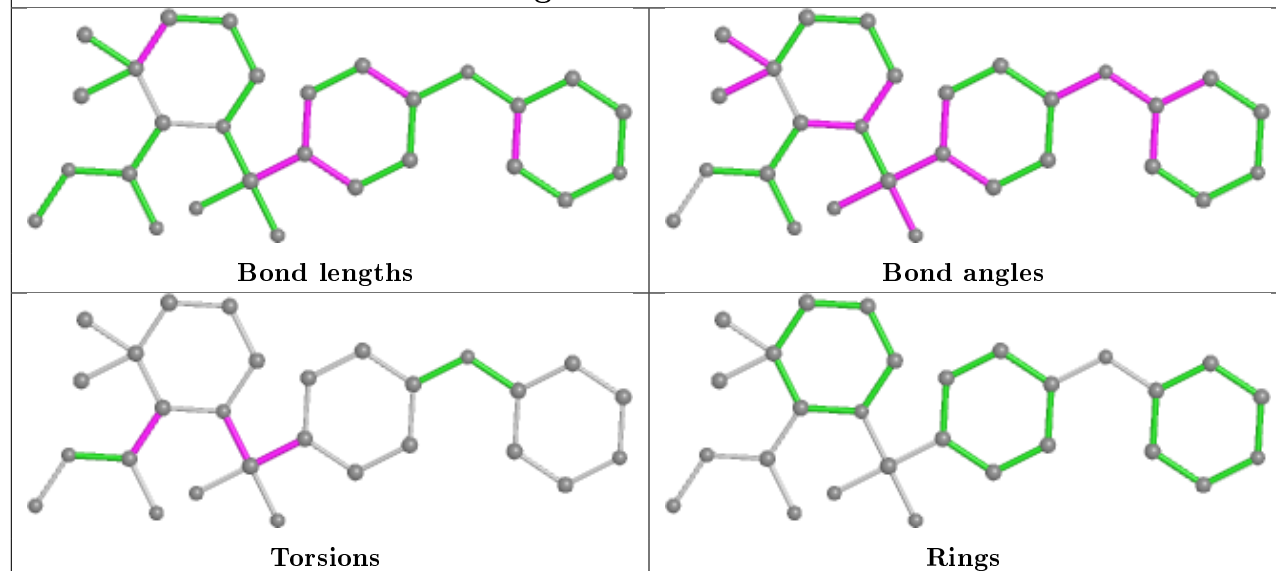
There are no ring outliers.

4 monomers are involved in 12 short contacts:

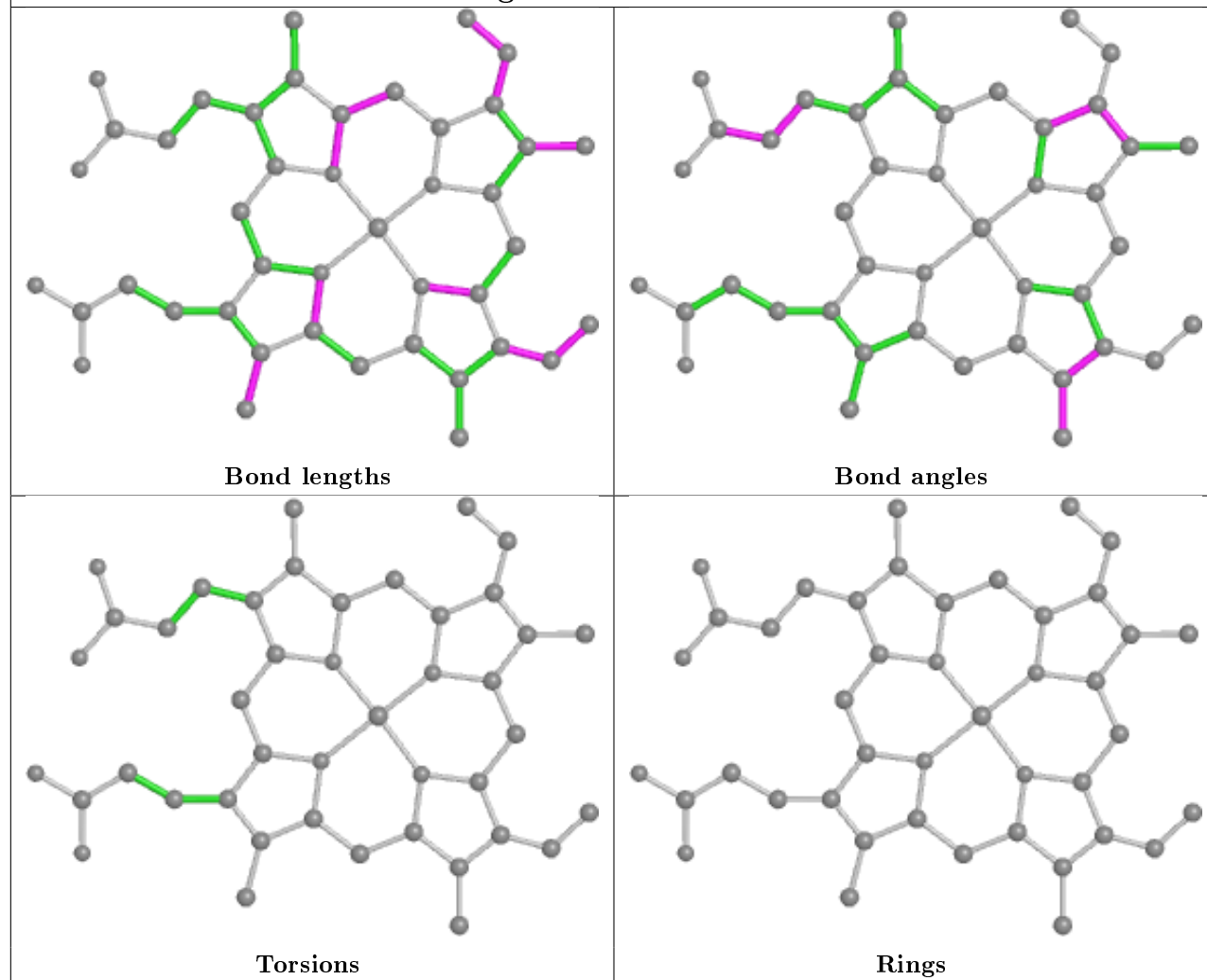
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	503	PN0	1	0
2	A	502	HEM	6	0
2	B	502	HEM	3	0
3	A	503	PN0	2	0

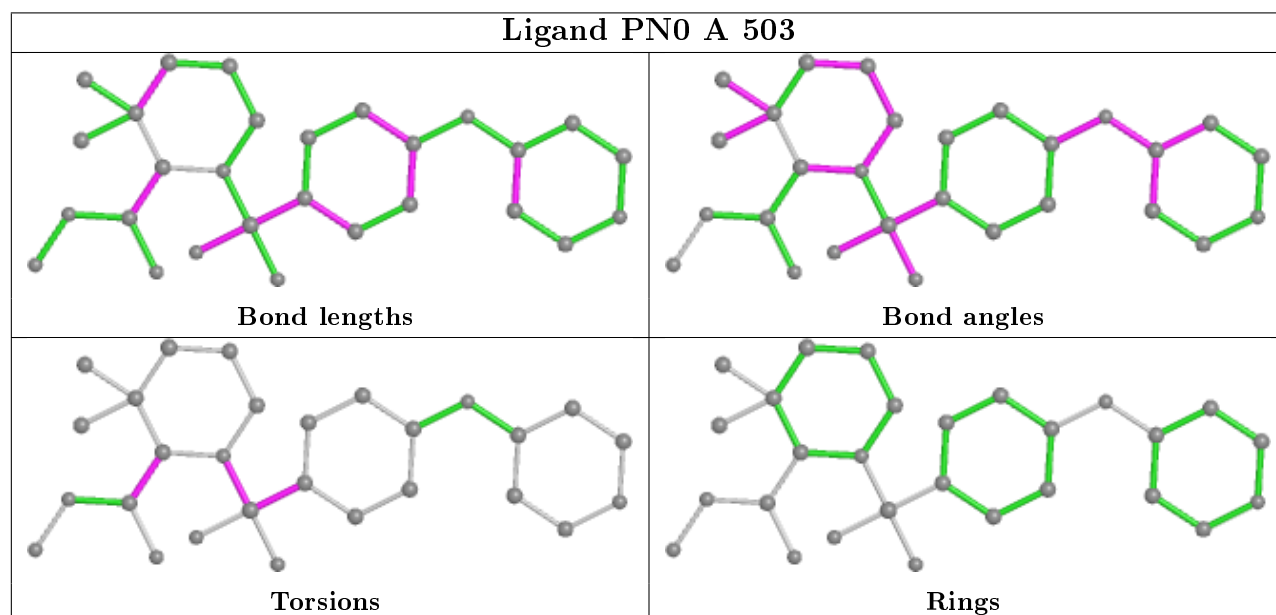
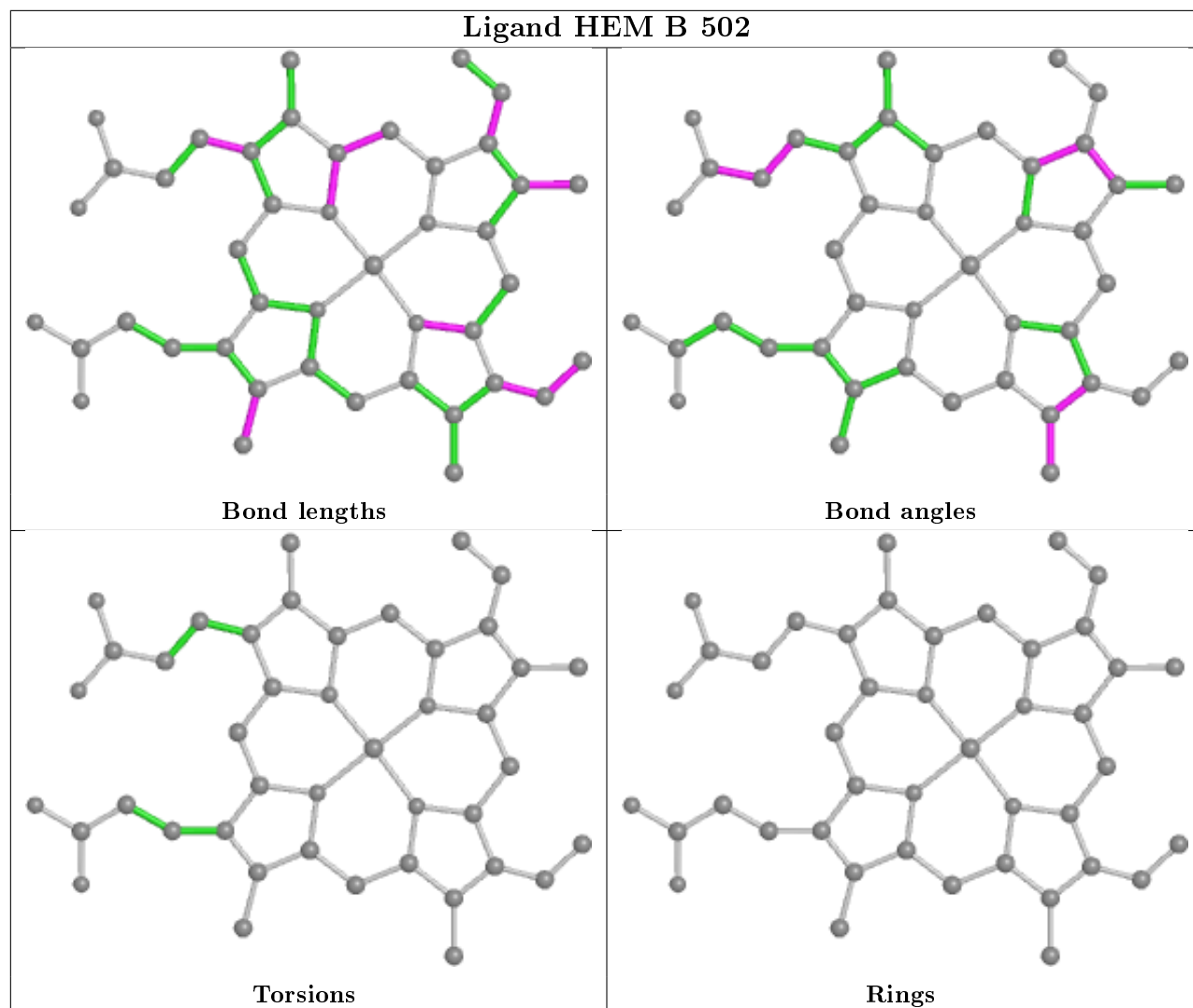
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

Ligand PN0 B 503



Ligand HEM A 502





5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	465/479 (97%)	0.07	17 (3%) 41 36	43, 77, 101, 111	0
1	B	457/479 (95%)	0.09	12 (2%) 56 52	53, 79, 99, 111	0
All	All	922/958 (96%)	0.08	29 (3%) 49 44	43, 78, 100, 111	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	232	HIS	4.7
1	B	309	THR	3.1
1	B	352	HIS	3.1
1	A	341	GLN	3.0
1	A	474	ARG	2.9
1	A	305	ALA	2.7
1	A	309	THR	2.7
1	B	239	LYS	2.6
1	A	170	ARG	2.6
1	B	306	GLY	2.5
1	B	380	ARG	2.5
1	B	169	GLY	2.5
1	A	65	PHE	2.4
1	B	305	ALA	2.4
1	A	416	HIS	2.4
1	A	478	HIS	2.3
1	B	426	HIS	2.3
1	B	165	ALA	2.3
1	A	147	LYS	2.2
1	B	304	SER	2.2
1	B	424	GLN	2.2
1	A	387	PHE	2.1
1	A	443	CYS	2.1
1	A	235	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	494	ALA	2.1
1	B	265	ALA	2.1
1	A	415	PHE	2.1
1	A	56	TYR	2.0
1	A	312	THR	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](#)

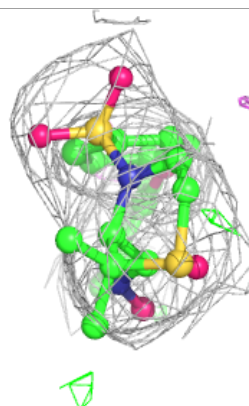
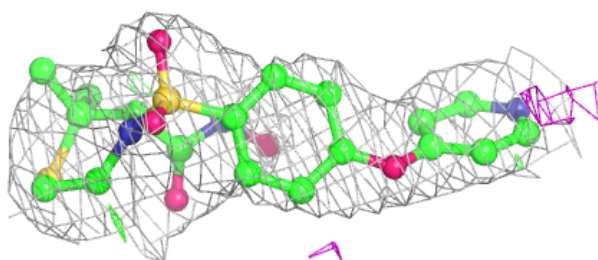
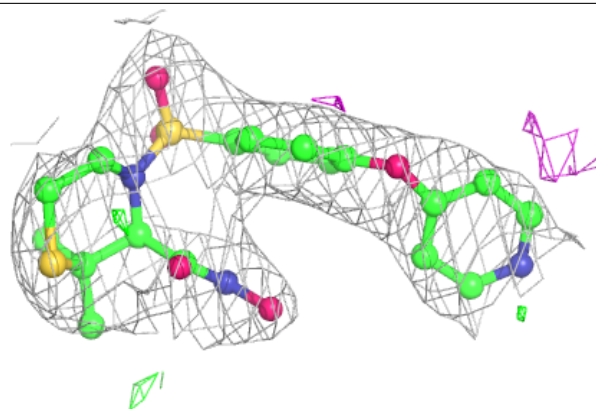
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	PN0	B	503	28/28	0.95	0.26	46,72,83,83	0
3	PN0	A	503	28/28	0.95	0.23	57,70,80,84	0
2	HEM	B	502	43/43	0.97	0.28	52,59,61,62	0
4	NI	B	600	1/1	0.98	0.07	73,73,73,73	0
2	HEM	A	502	43/43	0.98	0.23	33,41,45,50	0

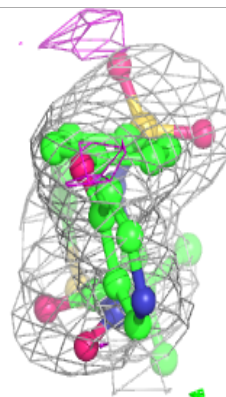
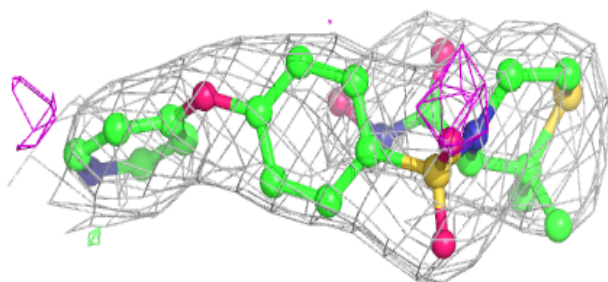
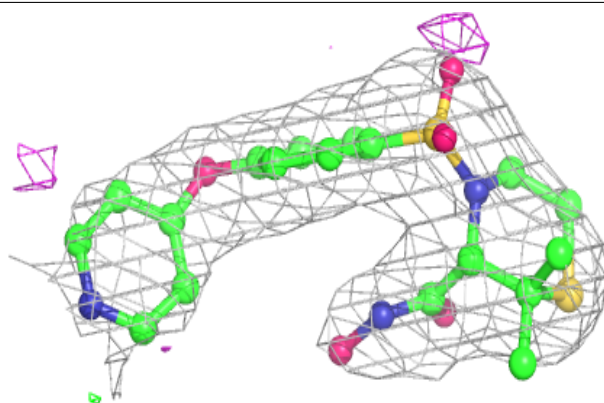
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around PN0 B 503:

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

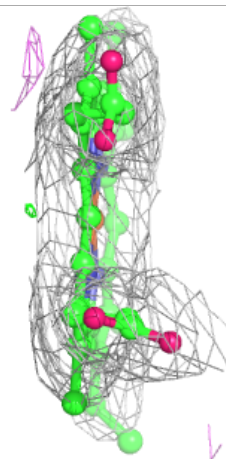
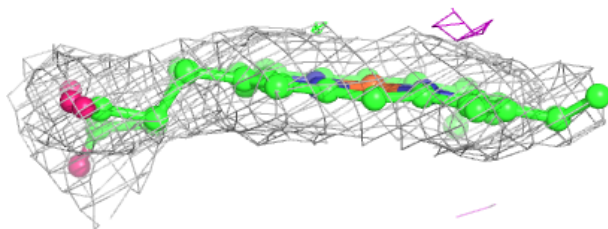
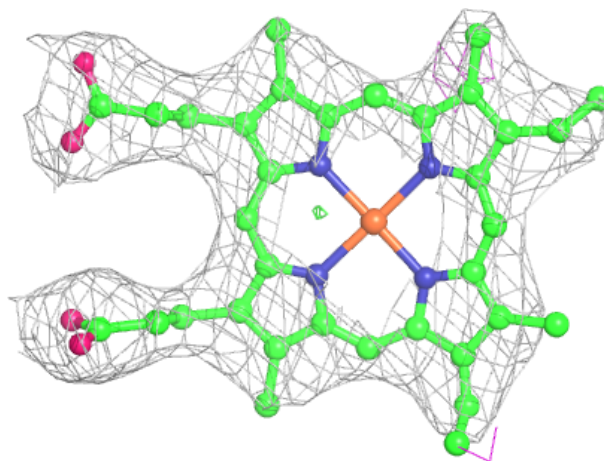
**Electron density around PN0 A 503:**

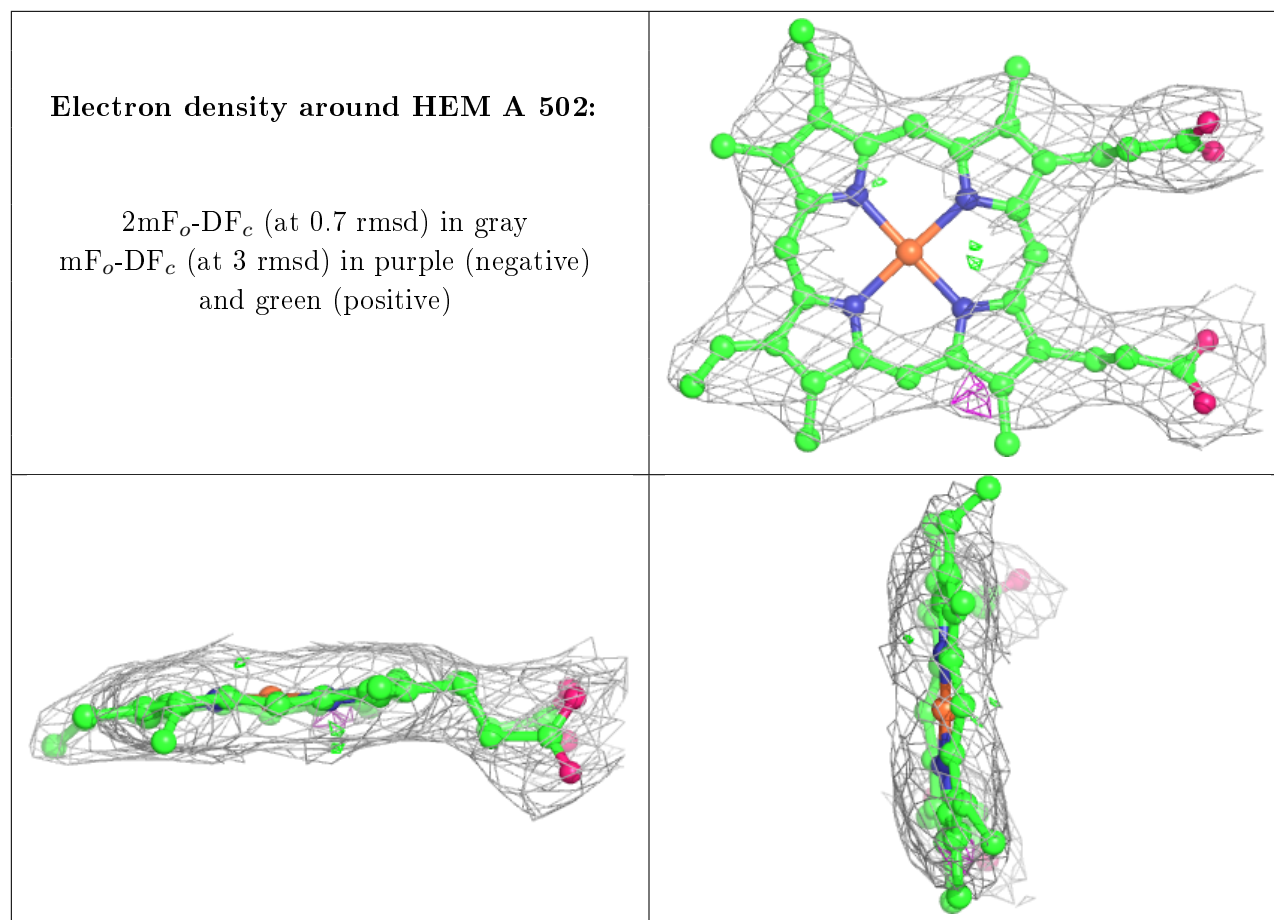
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around HEM B 502:

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





6.5 Other polymers ⓘ

There are no such residues in this entry.