



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 5, 2026 – 03:25 PM EDT

PDB ID : 3QM4 / pdb_00003qm4
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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.015 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.50

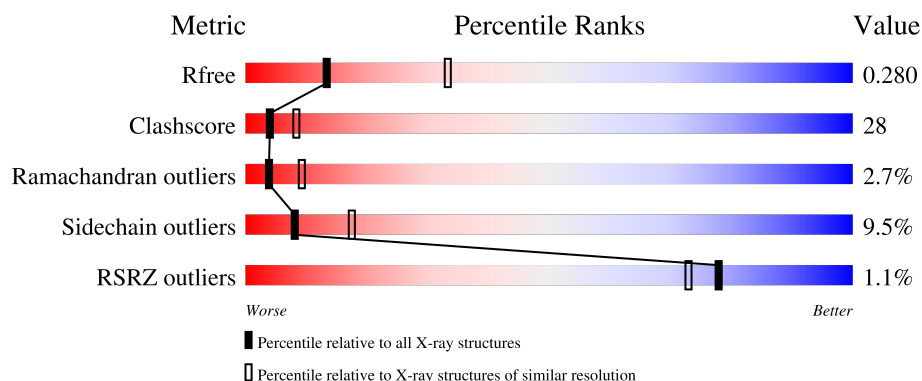
1 Overall quality at a glance

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}	180053	1407 (2.88-2.84)
Clashscore	190562	1446 (2.88-2.84)
Ramachandran outliers	187476	1406 (2.88-2.84)
Sidechain outliers	187428	1407 (2.88-2.84)
RSRZ outliers	180081	1408 (2.88-2.84)

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 of 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	% 49% 39% 9% • •
1	B	479	% 48% 39% 7% • 5%

2 Entry composition [i](#)

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:

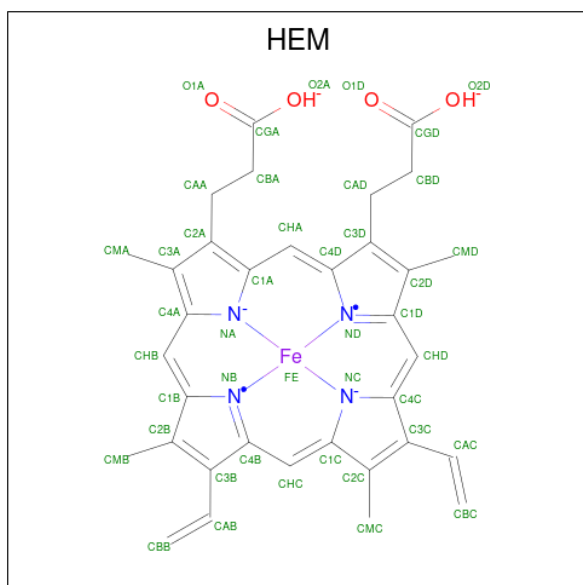
Chain	Residue	Modelled	Actual	Comment	Reference
A	23	MET	-	expression tag	UNP P10635
A	24	ALA	-	expression tag	UNP P10635
A	25	LYS	-	expression tag	UNP P10635
A	26	LYS	-	expression tag	UNP P10635
A	27	THR	-	expression tag	UNP P10635
A	28	SER	-	expression tag	UNP P10635
A	29	SER	-	expression tag	UNP P10635
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

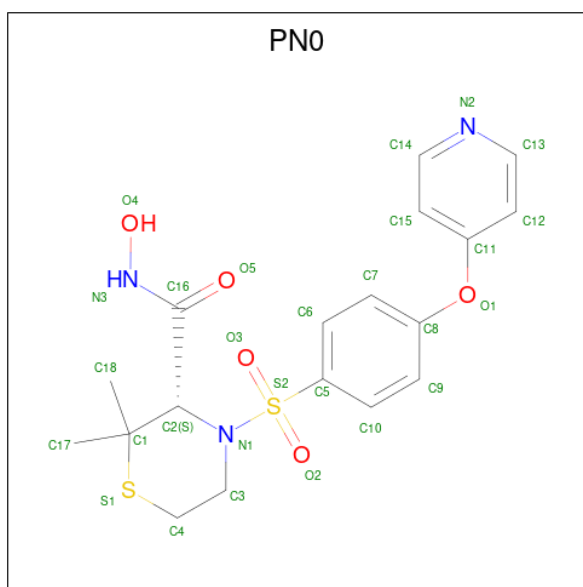
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Chain	Residue	Modelled	Actual	Comment	Reference
B	33	LEU	-	expression tag	UNP P10635
B	498	HIS	-	expression tag	UNP P10635
B	499	HIS	-	expression tag	UNP P10635
B	500	HIS	-	expression tag	UNP P10635
B	501	HIS	-	expression tag	UNP P10635

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





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 (CCD ID: 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		

Q250	L251	D252	R259	W262	R269	D270	L271	T272	E273	A274	F275	L276	A277	E278	M279	E280	K281	A282	K283	F290	L295	R296	I297	L302	F303	M307	V308	T309	T310	S311	T312	W316	L319	L320	M321	I322	L323	H324	Q328	R329	R330	V331	Q332	Q333	E334	I335	R343
M353	P354	Y355	T356	V359	I360	H361	E362	V363	D368	I369	V370	P371	L372	G373	V374	T375	H376	M377	T378	S379	E383	V384	F387	K391	G392	T393	T394	L395	S400	S401	V402	L403	E406	E410	K411	P412	P417	E418	H419	V428	R429	P430	E431	A432	F436	R440	R441
A442	C443	L444	G445	E446	P447	R450	M451	F454	L455	F456	F457	T458	S459	L460	L461	Q462	H463	P464	S465	F466	T470	G471	R474	P475	L484	V485	S486	P487	S488	P489	Y490	E491	L492	C493	A494	V495	P496	R497	HIS	HIS	HIS	HIS					

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 41.00 – 2.85	Depositor EDS
% Data completeness (in resolution range)	98.8 (41.00-2.85) 98.8 (41.00-2.85)	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 0.236 , 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)	0.30 , 38.9	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: HEM, PN0, 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	0.54	0/3777	1.06	23/5139 (0.4%)
1	B	0.52	0/3715	1.06	26/5051 (0.5%)
All	All	0.53	0/7492	1.06	49/10190 (0.5%)

There are no bond length outliers.

All (49) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	369	ILE	N-CA-C	8.39	119.19	110.72
1	B	369	ILE	N-CA-C	7.94	118.74	110.72
1	B	152	TRP	N-CA-C	-7.88	102.31	112.23
1	A	229	VAL	N-CA-C	-7.72	105.04	111.91
1	B	474	ARG	CA-C-N	7.61	127.66	119.90
1	B	474	ARG	C-N-CA	7.61	127.66	119.90
1	B	411	LYS	CA-C-N	7.16	127.48	119.32
1	B	411	LYS	C-N-CA	7.16	127.48	119.32
1	B	353	MET	CA-C-N	7.14	126.64	118.85
1	B	353	MET	C-N-CA	7.14	126.64	118.85
1	B	199	ASP	CA-C-N	7.14	126.39	118.97
1	B	199	ASP	C-N-CA	7.14	126.39	118.97
1	A	152	TRP	N-CA-C	-7.00	103.41	112.23
1	A	283	LYS	N-CA-C	-6.78	103.96	111.82
1	A	353	MET	CA-C-N	6.44	126.02	119.19
1	A	353	MET	C-N-CA	6.44	126.02	119.19
1	A	370	VAL	CA-C-N	6.27	127.68	119.84
1	A	370	VAL	C-N-CA	6.27	127.68	119.84
1	A	76	THR	N-CA-C	6.12	120.30	109.58
1	B	370	VAL	CA-C-N	6.12	127.50	119.84
1	B	370	VAL	C-N-CA	6.12	127.50	119.84
1	B	76	THR	N-CA-C	5.95	117.39	109.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	244	GLN	N-CA-C	-5.83	104.51	111.69
1	A	292	ASP	N-CA-C	5.74	118.28	111.33
1	B	108	GLN	N-CA-C	-5.69	105.15	111.36
1	B	172	PHE	N-CA-C	5.65	115.85	107.88
1	A	36	GLY	CA-C-N	5.65	126.90	119.84
1	A	36	GLY	C-N-CA	5.65	126.90	119.84
1	A	339	ILE	N-CA-C	5.60	115.80	110.53
1	A	155	GLU	N-CA-C	-5.53	105.16	111.07
1	B	495	VAL	CA-C-N	5.49	125.47	120.03
1	B	495	VAL	C-N-CA	5.49	125.47	120.03
1	A	495	VAL	CA-C-N	5.47	125.45	120.03
1	A	495	VAL	C-N-CA	5.47	125.45	120.03
1	A	481	PHE	N-CA-C	5.36	118.36	109.46
1	B	429	LYS	CA-C-N	5.29	125.27	120.03
1	B	429	LYS	C-N-CA	5.29	125.27	120.03
1	B	62	ARG	N-CA-C	-5.23	106.73	113.01
1	A	108	GLN	N-CA-C	-5.22	105.59	111.28
1	B	104	VAL	CA-C-N	5.20	126.34	119.84
1	B	104	VAL	C-N-CA	5.20	126.34	119.84
1	A	104	VAL	CA-C-N	5.19	126.33	119.84
1	A	104	VAL	C-N-CA	5.19	126.33	119.84
1	A	91	LEU	CA-C-N	-5.19	117.89	122.97
1	A	91	LEU	C-N-CA	-5.19	117.89	122.97
1	B	240	VAL	N-CA-C	5.15	120.05	109.34
1	A	172	PHE	N-CA-C	5.13	115.31	108.38
1	B	412	PRO	N-CA-C	5.01	121.22	113.75
1	B	48	HIS	N-CA-C	5.01	118.78	109.56

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	211	0
1	B	3618	0	3594	211	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
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	414	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 28.

All (414) 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	1.16	0.92
1:A:79:VAL:HG11	1:A:389:ILE:HD11	1.54	0.89
1:B:474:ARG:HG2	1:B:474:ARG:HH11	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:B:303:PHE:HE1	1:B:307:MET:HE3	1.41	0.86
1:A:444:LEU:HD23	1:A:444:LEU:C	2.03	0.83
1:A:47:LEU:HD23	1:A:47:LEU:H	1.42	0.83
1:A:321:MET:HE3	1:A:328:GLN:HA	1.59	0.82
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	1.64	0.80
1:A:276:LEU:HD23	1:A:279:MET:CE	2.12	0.80
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:HG2	1:B:474:ARG:NH1	1.99	0.73
1:B:102:PRO:HG3	1:B:394:THR:HG23	1.70	0.73
1:A:129:ARG:HH11	1:A:129:ARG:HG2	1.53	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:429:LYS:NZ	1:B:450:ARG:HH12	1.85	0.73
1:A:276:LEU:HD23	1:A:279:MET:HE1	1.69	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:444:LEU:C	1:B:444:LEU:HD23	2.14	0.71
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:HB2	1:B:129:ARG:CZ	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
1:B:89:GLU:OE1	1:B:384:VAL:HA	1.91	0.70
1:A:380:ARG:HB2	1:A:380:ARG:NH1	2.04	0.70
1:A:34:PRO:HA	1:A:387:PHE:CD2	2.26	0.70
1:A:321:MET:CE	1:A:328:GLN:HA	2.21	0.70
1:B:321:MET:HE2	1:B:457:PHE:HZ	1.55	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: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:75:TRP:CD1	1:A:75:TRP:H	2.14	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:276:LEU:HA	1:B:279:MET:HE3	1.79	0.64
1:B:151:GLN:HA	1:B:154:THR:HG23	1.81	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:206:LEU:HD23	1:B:307:MET:HE2	1.81	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
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:303:PHE:HE1	1:B:307:MET:CE	2.12	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:276:LEU:HD23	1:B:279:MET:HE1	1.83	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:B:124:TYR:HE1	1:B:440:ARG:HE	1.46	0.60
1:B:410:GLU:C	1:B:412:PRO:HD3	2.27	0.60
1:B:429:LYS:NZ	1:B:450:ARG:NH1	2.49	0.60
1:B:129:ARG:HB2	1:B:129:ARG:NH1	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:365:ARG:HH22	1:A:404:LYS:HG2	1.67	0.59
1:A:440:ARG:HH21	1:A:440:ARG:HG3	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:259:ARG:HD3	1:B:276:LEU:HD21	1.83	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:165:ALA:C	1:B:167:HIS:H	2.11	0.58
1:B:460:LEU:O	1:B:464:PHE:HB2	2.04	0.58
1:A:84:LEU:HD12	1:A:84:LEU:C	2.29	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:278:GLU:OE2	1:B:278:GLU:HA	2.03	0.57
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:191:CYS:HG	1:A:303:PHE:HE2	1.54	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:66:GLY:O	1:B:68:VAL:N	2.38	0.56
1:B:132:ARG:O	1:B:136:VAL:HG12	2.06	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:417:PRO:C	1:B:419:HIS:H	2.14	0.55
1:B:141:ASN:C	1:B:143:GLY:H	2.14	0.55
1:B:227:VAL:HG12	1:B:227:VAL:O	2.07	0.55
1:A:383:GLU:O	1:A:383:GLU:HG3	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:84:LEU:HD22	1:B:432:ALA:HB2	1.89	0.54
1:B:142:LEU:HD11	1:B:271:LEU:N	2.22	0.54
1:A:88:ARG:HG2	1:A:88:ARG:HH11	1.71	0.54
1:A:463:HIS:CG	1:A:463:HIS:O	2.61	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:A:36:GLY:C	1:A:65:PHE:CE2	2.86	0.53
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: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:41:PRO:HB3	1:B:42:GLY:H	1.73	0.53
1:A:232:HIS:O	1:A:233:ILE:HD13	2.08	0.53
1:A:470:THR:HG23	1:A:471:GLY:H	1.74	0.53
1:A:165:ALA:C	1:A:167:HIS:H	2.17	0.53
1:A:80:VAL:C	1:A:81:LEU:HD23	2.34	0.53
1:A:215:GLU:OE1	1:A:241:LEU:HA	2.08	0.53
1:A:233:ILE:HG21	1:A:236:LEU:HD12	1.90	0.53
1:A:313:THR:OG1	1:A:369:ILE:HD12	2.09	0.53
1:A:332:GLN:NE2	1:A:497:ARG:HD2	2.24	0.53
1:A:369:ILE:HG22	1:A:370:VAL:CG2	2.37	0.53
1:A:151:GLN:HA	1:A:154:THR:OG1	2.08	0.52
1:A:321:MET:HE3	1:A:328:GLN:CA	2.35	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: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:437:SER:OG	1:A:438:ALA:N	2.41	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:A:39:PRO:HG3	1:A:45:ASN:CG	2.35	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:123:ARG:NH2	1:B:377:MET:HE1	2.25	0.51
1:B:330:ARG:CB	1:B:355:TYR:CE2	2.94	0.51
1:A:232:HIS:C	1:A:234:PRO:HD3	2.36	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:O	1:A:428:VAL:HG23	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:A:141:ASN:C	1:A:143:GLY:H	2.19	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:303:PHE:CE1	1:B:307:MET:HE3	2.33	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:HD23	1:B:444:LEU:O	2.12	0.50
1:B:319:LEU:O	1:B:322:ILE:HG12	2.11	0.50
1:B:141:ASN:O	1:B:141:ASN:ND2	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
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:HG3	1:B:383:GLU:O	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:B:73:LEU:O	1:B:74:ALA:HB3	2.12	0.49
1:B:321:MET:HE2	1:B:457:PHE:CZ	2.42	0.49
1:A:39:PRO:HB3	1:A:44:GLY:C	2.38	0.48
1:A:231:LEU:O	1:A:232:HIS:HB2	2.12	0.48
1:A:440:ARG:HG3	1:A:440:ARG:NH2	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:476:SER:OG	1:A:478:HIS:ND1	2.47	0.48
1:A:490:TYR:CD1	1:A:490:TYR:C	2.91	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:89:GLU:HA	1:A:93:THR:OG1	2.13	0.48
1:A:444:LEU:HD23	1:A:445:GLY:N	2.28	0.48
1:B:428:VAL:O	1:B:428:VAL:HG23	2.14	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:B:41:PRO:O	1:B:42:GLY:C	2.57	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:279:MET:O	1:B:280:GLU:C	2.56	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:HB2	1.93	0.47
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:B:488:SER:O	1:B:489:PRO:C	2.58	0.47
1:A:433:PHE:O	1:A:434:LEU:HD23	2.15	0.46
1:B:173:ARG:NH1	1:B:490:TYR:H	2.05	0.46
1:B:335:ILE:HD13	1:B:462:GLN:HB2	1.96	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:75:TRP:CZ3	1:B:61:LEU:HD21	2.49	0.46
1:A:231:LEU:CD2	1:B:52:GLN:HB2	2.45	0.46
1:A:477:HIS:ND1	1:A:477:HIS:N	2.61	0.46
1:B:129:ARG:NH1	1:B:129:ARG:CB	2.79	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:410:GLU:C	1:A:412:PRO:HD3	2.41	0.46
1:A:417:PRO:C	1:A:419:HIS:H	2.23	0.46
1:A:466:PHE:HA	1:A:493:CYS:O	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:356:THR:HG21	1:B:458:THR:HG22	1.97	0.46
1:A:84:LEU:HD22	1:A:432:ALA:HA	1.97	0.46
1:A:106:ILE:HG12	1:A:106:ILE:O	2.16	0.46
1:A:75:TRP:H	1:A:75:TRP:HD1	1.63	0.46
1:A:413:PHE:CD1	1:A:413:PHE:N	2.84	0.46
1:B:221:ARG:O	1:B:225:ASN:HB2	2.16	0.46
1:B:224:LEU:HD12	1:B:240:VAL:HG21	1.98	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:HD23	1:B:224:LEU:H	1.81	0.45
1:B:165:ALA:C	1:B:167:HIS:N	2.75	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:B:417:PRO:C	1:B:419:HIS:N	2.74	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:HH11	1:B:115:ARG:HG2	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:74:ALA:HB3	1:B:226:ALA:HB1	1.99	0.45
1:B:440:ARG:HG2	1:B:440:ARG:NH2	2.31	0.45
1:B:444:LEU:C	1:B:444:LEU:CD2	2.86	0.44
1:A:39:PRO:CG	1:A:45:ASN:CB	2.94	0.44
1:A:88:ARG:HH11	1:A:88:ARG:CG	2.31	0.44
1:A:40:LEU:C	1:A:40:LEU:HD12	2.42	0.44
1:B:323:LEU:HD11	1:B:475:PRO:O	2.17	0.44
1:A:40:LEU:HB2	1:A:41:PRO:HD2	2.00	0.44
1:A:210:GLN:O	1:A:214:LYS:HG3	2.18	0.44
1:B:45:ASN:HA	1:B:48:HIS:CE1	2.53	0.44
1:B:320:LEU:O	1:B:324:HIS:HB2	2.18	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:A:357:THR:HG22	1:A:427:PHE:CE1	2.53	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:220:LEU:HD13	1:A:220:LEU:C	2.43	0.44
1:B:101:ARG:NH1	1:B:441:ARG:HH11	2.16	0.44
1:A:308:VAL:HG23	1:A:308:VAL:O	2.18	0.44
1:B:43:LEU:O	1:B:74:ALA:HA	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:370:VAL:CG1	1:B:373:GLY:HA2	2.48	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:276:LEU:HA	1:A:279:MET:HE2	1.99	0.43
1:A:444:LEU:HD22	2:A:502:HEM:C2D	2.53	0.43
1:A:444:LEU:CD2	2:A:502:HEM:HHD	2.48	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:43:LEU:HG	1:B:46:LEU:HD22	1.99	0.43
1:B:68:VAL:O	1:B:68:VAL:HG12	2.17	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:A:102:PRO:HB2	1:B:64:ARG:HH22	1.83	0.43
1:B:112:PHE:CD1	1:B:297:ILE:HD13	2.52	0.43
1:B:206:LEU:CD2	1:B:307:MET:HE2	2.48	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: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:B:333:GLN:O	1:B:334:GLU:C	2.61	0.43
1:A:109:ILE:O	1:A:109:ILE:HG22	2.19	0.43
1:A:131:GLN:OE1	1:A:131:GLN:HA	2.19	0.43
1:B:33:LEU:N	1:B:34:PRO:CD	2.82	0.43
1:B:33:LEU:N	1:B:34:PRO:HD3	2.34	0.43
1:B:241:LEU:CD1	1:B:241:LEU:N	2.81	0.43
1:A:450:ARG:HH11	1:A:450:ARG:HG3	1.84	0.42
1:B:281:LYS:C	1:B:283:LYS:H	2.26	0.42
1:B:379:SER:O	1:B:391:LYS:HE3	2.19	0.42
1:A:174:PRO:O	1:A:175:ASN:C	2.62	0.42
1:A:470:THR:CG2	1:A:471:GLY:N	2.82	0.42
1:A:305:ALA:C	1:A:307:MET:H	2.26	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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:487:PRO:O	1:A:488:SER:C	2.59	0.42
1:B:226:ALA:O	1:B:227:VAL:HG23	2.20	0.42
1:B:316:TRP:CE2	1:B:487:PRO:HG3	2.53	0.42
1:B:406:GLU:H	1:B:406:GLU:HG2	1.67	0.42
1:A:54:THR:HB	1:A:55:PRO:HD3	2.02	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:HE3	1:B:152:TRP:O	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:150:GLU:HB2	1:B:451:MET:HE2	2.01	0.42
1:A:65:PHE:CD1	1:A:65:PHE:N	2.88	0.42
1:A:205:LEU:HD12	1:A:205:LEU:HA	1.78	0.42
1:B:45:ASN:C	1:B:47:LEU:H	2.28	0.42
1:A:119:VAL:O	1:A:120:PHE:C	2.63	0.41
1:A:157:ALA:HA	1:A:456:PHE:HE1	1.85	0.41
1:A:165:ALA:C	1:A:167:HIS:N	2.77	0.41
1:B:279:MET:HG2	1:B:290:PHE:C	2.45	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:166:ASN:O	1:A:166:ASN:CG	2.62	0.41
1:B:123:ARG:HE	1:B:123:ARG:HB3	1.59	0.41
1:B:429:LYS:HZ2	1:B:450:ARG:NH1	1.96	0.41
1:A:220:LEU:HA	1:A:220:LEU:HD22	1.87	0.41
1:A:266:GLN:O	1:A:267:PRO:C	2.62	0.41
1:B:51:PHE:CZ	1:B:222:GLU:HG3	2.55	0.41
1:B:269:ARG:HG2	1:B:269:ARG:HH11	1.85	0.41
1:B:329:ARG:NH1	1:B:329:ARG:HG3	2.36	0.41
1:A:81:LEU:HD23	1:A:81:LEU:N	2.35	0.41
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:B:164:PHE:CE2	1:B:460:LEU:HD22	2.56	0.41
1:B:251:LEU:O	1:B:252:ASP:C	2.64	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:78:VAL:HG12	1:B:79:VAL:N	2.36	0.41
1:B:271:LEU:HD22	1:B:302:LEU:HD12	2.03	0.41
1:B:40:LEU:O	1:B:41:PRO:C	2.62	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:79:VAL:CG2	1:B:393:THR:HG21	2.51	0.41
1:B:226:ALA:C	1:B:227:VAL:HG23	2.46	0.41
1:B:417:PRO:O	1:B:419:HIS:N	2.54	0.41
1:A:47:LEU:CD2	1:A:47:LEU:N	2.76	0.41
1:A:121:LEU:HD23	1:A:121:LEU:HA	1.95	0.41
1:A:147:LYS:HA	1:A:151:GLN:HE21	1.85	0.41
1:A:442:ALA:O	1:A:443:CYS:C	2.63	0.41
1:B:88:ARG:O	1:B:92:VAL:HB	2.21	0.41
1:B:170:ARG:HA	1:B:170:ARG:HD2	1.96	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:A:481:PHE:CD2	1:A:481:PHE:C	2.98	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:B:303:PHE:CE1	1:B:307:MET:CE	2.98	0.40
1:A:35:PRO:HD3	1:A:387:PHE:CE2	2.57	0.40
1:A:469:PRO:O	1:A:470:THR:C	2.64	0.40
1:A:362:GLU:HG3	1:A:415:PHE:CE1	2.56	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	463/479 (97%)	396 (86%)	58 (12%)	9 (2%)	6	14
1	B	453/479 (95%)	382 (84%)	55 (12%)	16 (4%)	3	6
All	All	916/958 (96%)	778 (85%)	113 (12%)	25 (3%)	4	9

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%)	356 (90%)	41 (10%)	7	14
1	B	391/409 (96%)	357 (91%)	34 (9%)	9	21
All	All	788/818 (96%)	713 (90%)	75 (10%)	8	17

All (75) 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	81	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	220	LEU
1	A	227	VAL
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	308	VAL
1	A	321	MET
1	A	329	ARG
1	A	343	ARG
1	A	369	ILE
1	A	375	THR
1	A	383	GLU
1	A	395	LEU
1	A	403	LEU
1	A	444	LEU
1	A	451	MET
1	A	468	VAL
1	A	484	LEU
1	A	485	VAL
1	A	486	SER
1	A	488	SER
1	A	492	LEU
1	A	495	VAL
1	B	33	LEU
1	B	47	LEU

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Mol	Chain	Res	Type
1	B	48	HIS
1	B	71	LEU
1	B	72	GLN
1	B	75	TRP
1	B	93	THR
1	B	104	VAL
1	B	136	VAL
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	224	LEU
1	B	240	VAL
1	B	249	THR
1	B	250	GLN
1	B	323	LEU
1	B	329	ARG
1	B	331	VAL
1	B	335	ILE
1	B	369	ILE
1	B	379	SER
1	B	383	GLU
1	B	395	LEU
1	B	403	LEU
1	B	406	GLU
1	B	444	LEU
1	B	474	ARG
1	B	484	LEU
1	B	485	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	72	GLN
1	A	94	HIS
1	A	141	ASN
1	A	151	GLN
1	A	244	GLN
1	A	250	GLN

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Mol	Chain	Res	Type
1	A	426	HIS
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	244	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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	$\# Z > 2$
2	HEM	A	502	1,3	50,50,50	2.32	24 (48%)	67,82,82	1.36	10 (14%)
3	PN0	A	503	2	28,30,30	1.78	8 (28%)	37,44,44	1.60	9 (24%)
2	HEM	B	502	1,3	50,50,50	2.36	23 (46%)	67,82,82	1.38	12 (17%)
3	PN0	B	503	2	28,30,30	1.84	6 (21%)	37,44,44	1.72	9 (24%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	HEM	A	502	1,3	-	4/14/54/54	-
3	PN0	A	503	2	-	9/21/39/39	0/3/3/3
2	HEM	B	502	1,3	-	5/14/54/54	-
3	PN0	B	503	2	-	9/21/39/39	0/3/3/3

All (61) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	503	PN0	C1-S1	-4.58	1.80	1.84
2	A	502	HEM	C1C-NC	-4.50	1.31	1.39
2	A	502	HEM	C4A-NA	-4.23	1.31	1.39
2	B	502	HEM	C1C-NC	-4.21	1.31	1.39
2	B	502	HEM	FE-ND	4.19	2.07	1.94
2	B	502	HEM	FE-NB	4.16	2.07	1.94
2	B	502	HEM	CAC-C3C	-4.05	1.36	1.47
2	B	502	HEM	C4B-NB	-4.04	1.30	1.38
2	B	502	HEM	C1D-ND	-4.03	1.30	1.38
2	B	502	HEM	C4D-ND	-4.00	1.33	1.40
2	A	502	HEM	C1D-ND	-3.93	1.31	1.38
3	A	503	PN0	C1-S1	-3.93	1.80	1.84
3	B	503	PN0	C15-C11	3.89	1.46	1.38
2	A	502	HEM	C4B-NB	-3.87	1.31	1.38
2	A	502	HEM	FE-ND	3.85	2.06	1.94
2	B	502	HEM	C4A-NA	-3.71	1.32	1.39
2	A	502	HEM	CBB-CAB	3.68	1.48	1.30
2	A	502	HEM	C4D-ND	-3.68	1.33	1.40
2	A	502	HEM	CAC-C3C	-3.67	1.37	1.47
2	B	502	HEM	CBB-CAB	3.63	1.47	1.30
2	A	502	HEM	CMA-C3A	3.56	1.58	1.50
2	B	502	HEM	FE-NA	3.53	2.06	1.95
2	A	502	HEM	FE-NB	3.42	2.05	1.94
2	A	502	HEM	CMC-C2C	3.29	1.57	1.50
2	B	502	HEM	CMA-C3A	3.25	1.57	1.50
3	A	503	PN0	C15-C11	3.23	1.44	1.38
2	B	502	HEM	C1B-NB	-3.20	1.34	1.40
2	A	502	HEM	FE-NA	3.18	2.05	1.95
2	B	502	HEM	C1C-C2C	-3.16	1.39	1.45
3	A	503	PN0	C5-S2	-3.15	1.72	1.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	502	HEM	C1B-NB	-3.12	1.34	1.40
2	B	502	HEM	CMC-C2C	3.07	1.57	1.50
3	A	503	PN0	C2-C16	2.97	1.58	1.53
2	A	502	HEM	C1C-C2C	-2.91	1.39	1.45
3	B	503	PN0	C9-C8	2.91	1.44	1.38
3	B	503	PN0	C5-S2	-2.81	1.72	1.76
2	A	502	HEM	FE-NC	2.76	2.04	1.95
3	A	503	PN0	C6-C5	2.63	1.43	1.38
2	A	502	HEM	C3B-C2B	2.59	1.42	1.37
2	B	502	HEM	C3B-C2B	2.57	1.42	1.37
2	A	502	HEM	CBC-CAC	2.57	1.42	1.30
3	B	503	PN0	C6-C5	2.52	1.42	1.38
3	A	503	PN0	C9-C8	2.51	1.43	1.38
2	B	502	HEM	O2A-CGA	-2.50	1.22	1.30
2	A	502	HEM	C1A-NA	-2.49	1.34	1.39
2	A	502	HEM	C3C-C4C	-2.48	1.41	1.46
2	A	502	HEM	CAB-C3B	-2.38	1.41	1.47
2	A	502	HEM	C4C-NC	-2.35	1.35	1.39
3	B	503	PN0	C10-C5	2.34	1.42	1.38
2	B	502	HEM	CBC-CAC	2.34	1.41	1.30
2	B	502	HEM	FE-NC	2.34	2.02	1.95
3	A	503	PN0	C7-C8	2.28	1.43	1.38
2	B	502	HEM	C4D-C3D	-2.22	1.41	1.45
2	B	502	HEM	C1D-C2D	-2.22	1.40	1.44
3	A	503	PN0	O3-S2	2.21	1.45	1.43
2	A	502	HEM	C1D-C2D	-2.16	1.40	1.44
2	A	502	HEM	O2A-CGA	-2.12	1.23	1.30
2	A	502	HEM	CHD-C1D	-2.09	1.34	1.39
2	B	502	HEM	C3C-C2C	2.08	1.41	1.37
2	B	502	HEM	O2D-CGD	-2.08	1.23	1.30
2	B	502	HEM	CAB-C3B	-2.03	1.42	1.47

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	503	PN0	C11-O1-C8	4.13	128.21	118.78
3	A	503	PN0	C18-C1-C17	-3.87	105.15	109.64
3	B	503	PN0	C18-C1-C17	-3.81	105.22	109.64
3	A	503	PN0	O2-S2-C5	-3.67	103.54	108.10
2	A	502	HEM	CHD-C4C-NC	-3.55	120.59	124.45
3	A	503	PN0	C11-O1-C8	3.21	126.12	118.78
2	A	502	HEM	CHB-C1B-NB	-3.09	120.55	124.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	502	HEM	C3B-C2B-C1B	-2.99	104.17	106.41
2	B	502	HEM	C3B-C4B-NB	2.97	111.60	109.47
2	B	502	HEM	CHD-C4C-NC	-2.95	121.24	124.45
3	B	503	PN0	C6-C5-S2	-2.91	116.84	119.73
3	B	503	PN0	O3-S2-O2	-2.90	115.06	119.59
2	B	502	HEM	CHC-C4B-NB	-2.74	121.48	124.42
3	B	503	PN0	O2-S2-C5	-2.72	104.72	108.10
2	A	502	HEM	C1B-NB-C4B	2.70	108.41	105.21
2	A	502	HEM	CHC-C4B-NB	-2.68	121.54	124.42
3	B	503	PN0	O1-C11-C15	2.68	128.68	119.36
2	B	502	HEM	C4C-C3C-C2C	-2.61	104.55	106.81
2	A	502	HEM	C4B-C3B-C2B	-2.61	104.88	107.28
3	B	503	PN0	C10-C5-S2	2.59	122.31	119.73
2	B	502	HEM	CHB-C1B-NB	-2.59	121.17	124.37
2	B	502	HEM	C1C-CHC-C4B	2.56	131.46	126.02
3	B	503	PN0	C3-N1-C2	2.51	122.92	116.20
2	B	502	HEM	C4D-ND-C1D	2.46	108.12	105.21
2	A	502	HEM	C4D-ND-C1D	2.45	108.11	105.21
3	A	503	PN0	O1-C11-C15	2.41	127.75	119.36
3	B	503	PN0	O1-C11-C12	-2.40	110.99	119.36
3	A	503	PN0	C3-N1-C2	2.40	122.62	116.20
2	A	502	HEM	C4C-C3C-C2C	-2.40	104.73	106.81
2	B	502	HEM	C1B-NB-C4B	2.34	107.98	105.21
2	B	502	HEM	C4B-C3B-C2B	-2.32	105.15	107.28
3	A	503	PN0	O3-S2-O2	-2.29	116.01	119.59
3	A	503	PN0	C3-C4-S1	2.27	115.10	112.47
2	A	502	HEM	CHC-C1C-NC	-2.27	121.98	124.45
2	A	502	HEM	C1C-CHC-C4B	2.20	130.69	126.02
2	B	502	HEM	C4A-CHB-C1B	2.09	131.17	126.25
2	A	502	HEM	C3B-C2B-C1B	-2.08	104.85	106.41
3	A	503	PN0	O1-C11-C12	-2.07	112.15	119.36
2	B	502	HEM	C4C-CHD-C1D	2.02	130.32	126.02
3	A	503	PN0	C6-C5-S2	-2.01	117.74	119.73

There are no chirality outliers.

All (27) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	502	HEM	C2C-C3C-CAC-CBC
2	B	502	HEM	C4C-C3C-CAC-CBC
2	A	502	HEM	C2C-C3C-CAC-CBC
3	A	503	PN0	C3-N1-S2-O3

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Mol	Chain	Res	Type	Atoms
3	B	503	PN0	C3-N1-S2-O3
3	A	503	PN0	N3-C16-C2-N1
2	A	502	HEM	C4C-C3C-CAC-CBC
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
3	A	503	PN0	C10-C5-S2-O2
3	B	503	PN0	C10-C5-S2-O2
3	B	503	PN0	N3-C16-C2-N1
2	A	502	HEM	CAA-CBA-CGA-O2A
2	B	502	HEM	CAA-CBA-CGA-O2A
2	B	502	HEM	CAA-CBA-CGA-O1A
2	A	502	HEM	CAA-CBA-CGA-O1A
3	B	503	PN0	O5-C16-C2-N1
3	A	503	PN0	C6-C5-S2-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
2	B	502	HEM	C2B-C3B-CAB-CBB
3	A	503	PN0	C3-N1-S2-C5
3	B	503	PN0	C3-N1-S2-C5

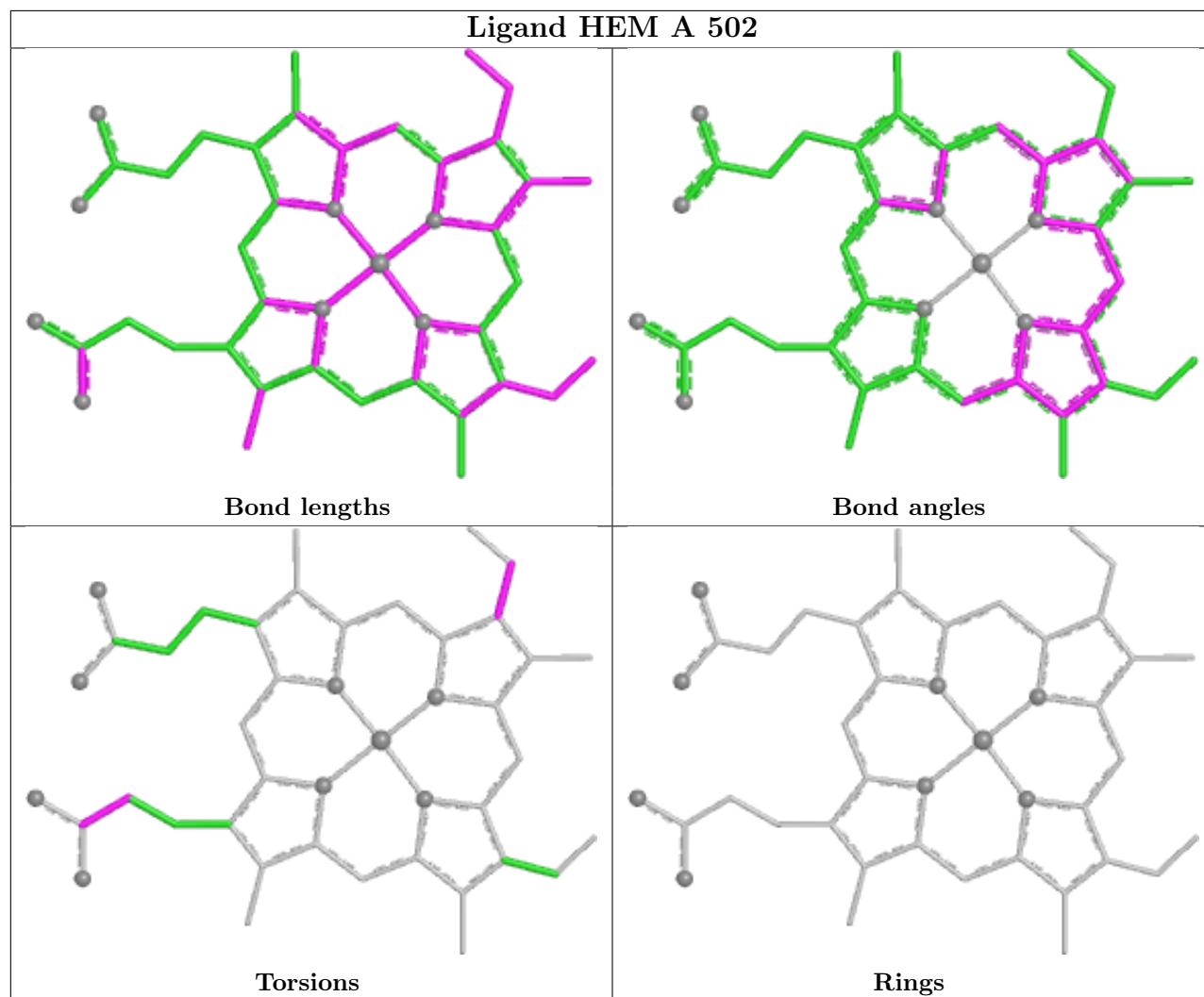
There are no ring outliers.

4 monomers are involved in 12 short contacts:

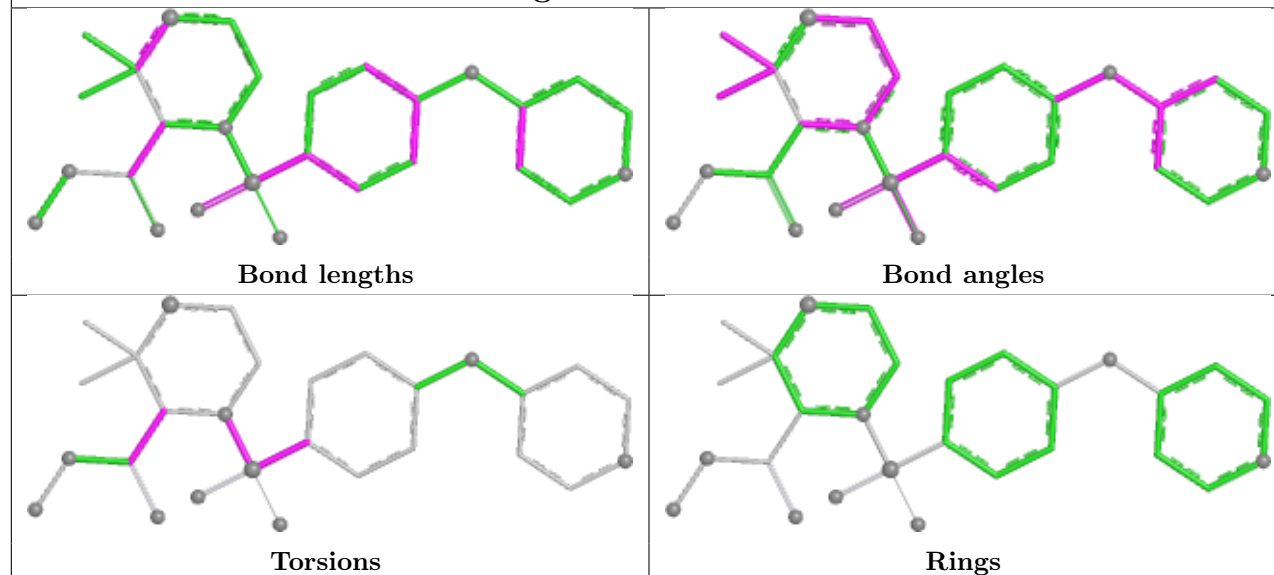
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	502	HEM	6	0
3	A	503	PN0	2	0
2	B	502	HEM	3	0
3	B	503	PN0	1	0

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

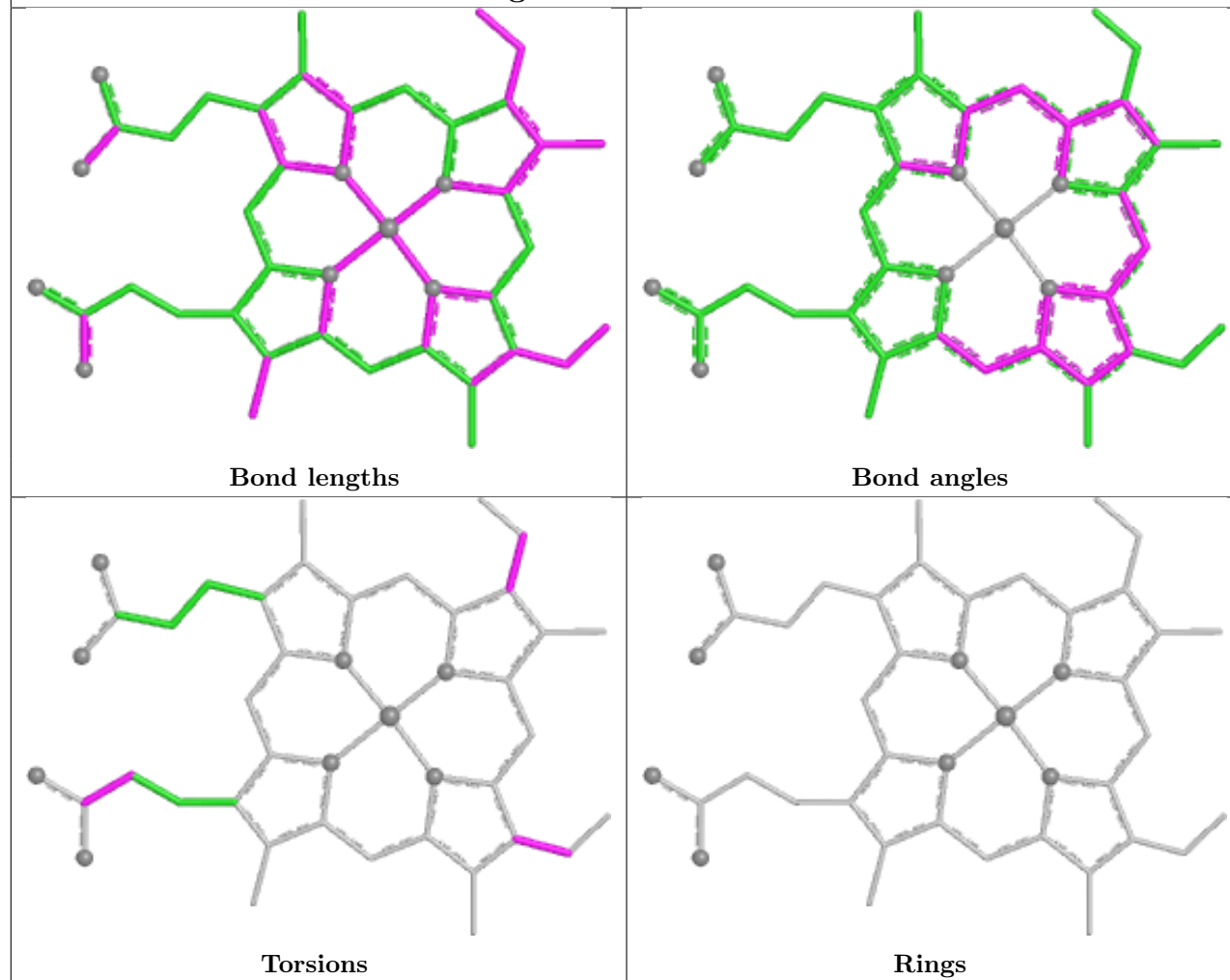
any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

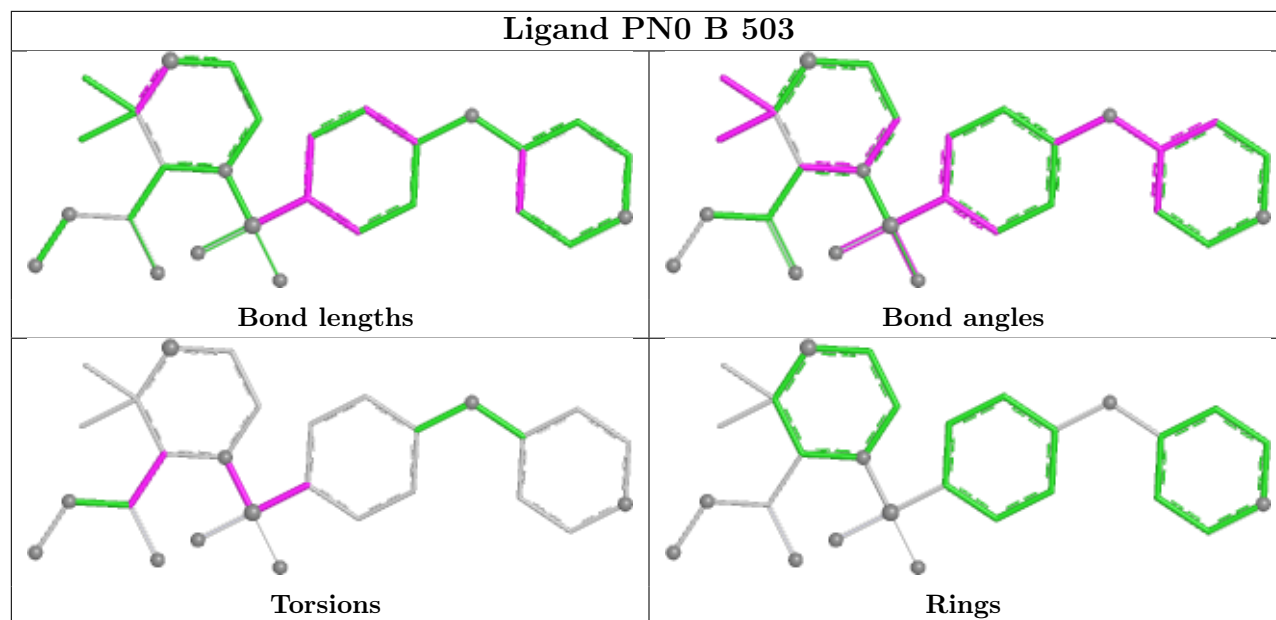


Ligand PN0 A 503



Ligand HEM B 502





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

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

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

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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%)	0.09	6 (1%) 75 68	43, 77, 101, 111	0
1	B	457/479 (95%)	0.14	4 (0%) 81 76	53, 79, 99, 111	0
All	All	922/958 (96%)	0.12	10 (1%) 78 73	43, 78, 100, 111	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	415	PHE	3.2
1	B	229	VAL	3.1
1	B	33	LEU	3.0
1	A	494	ALA	2.5
1	B	238	GLY	2.5
1	A	229	VAL	2.3
1	A	33	LEU	2.3
1	A	228	PRO	2.2
1	A	366	PHE	2.2
1	B	58	PHE	2.2

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 oligosaccharides in this entry.

6.4 Ligands [i](#)

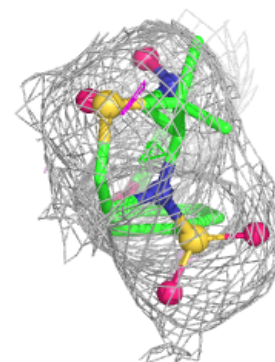
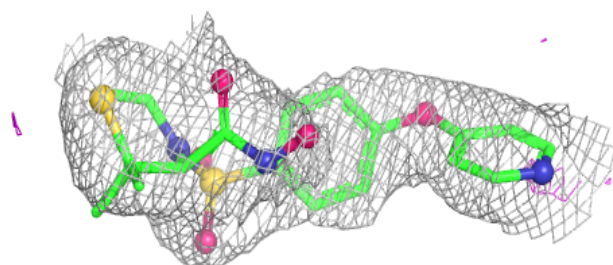
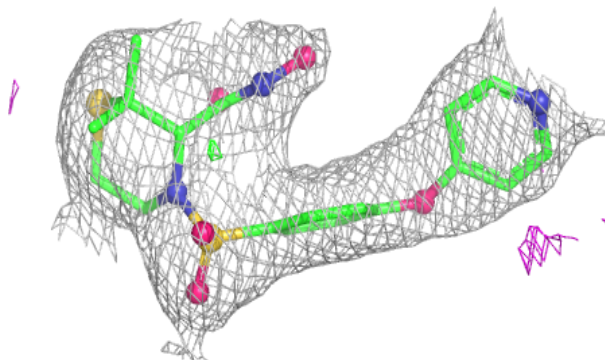
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	PN0	B	503	28/28	0.95	0.10	46,72,83,83	0
3	PN0	A	503	28/28	0.96	0.08	57,70,80,84	0
2	HEM	B	502	43/43	0.97	0.09	52,59,61,62	0
2	HEM	A	502	43/43	0.99	0.06	33,41,45,50	0
4	NI	B	600	1/1	0.99	0.02	73,73,73,73	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.

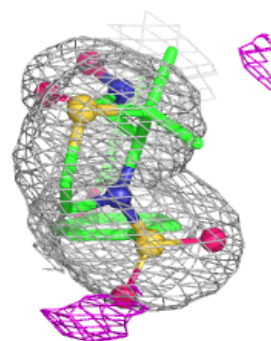
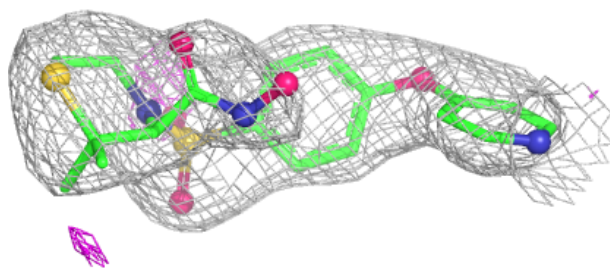
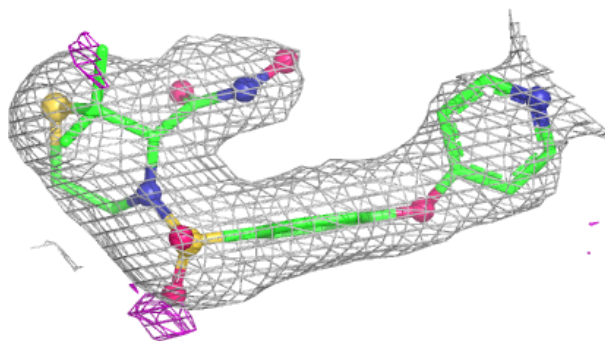
Electron density around PN0 B 503:

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



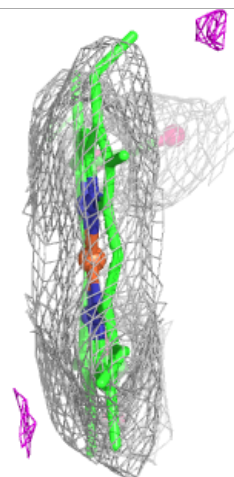
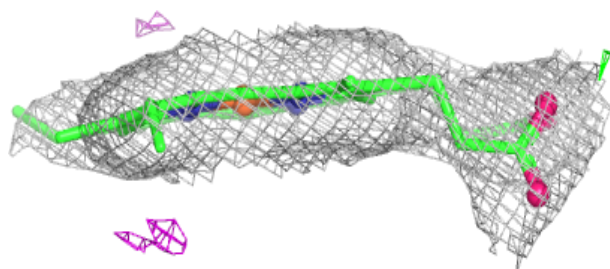
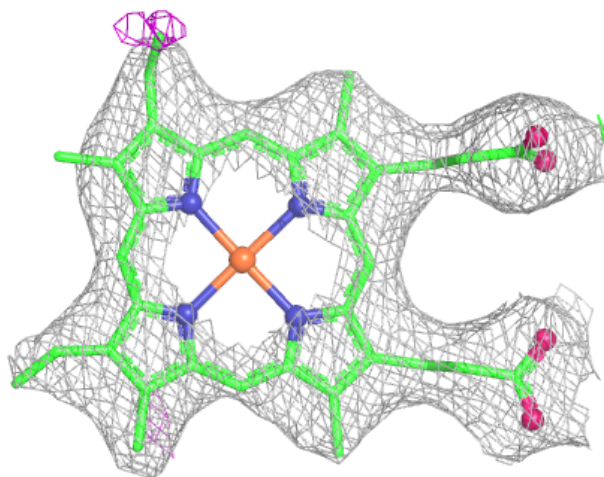
Electron density around PN0 A 503:

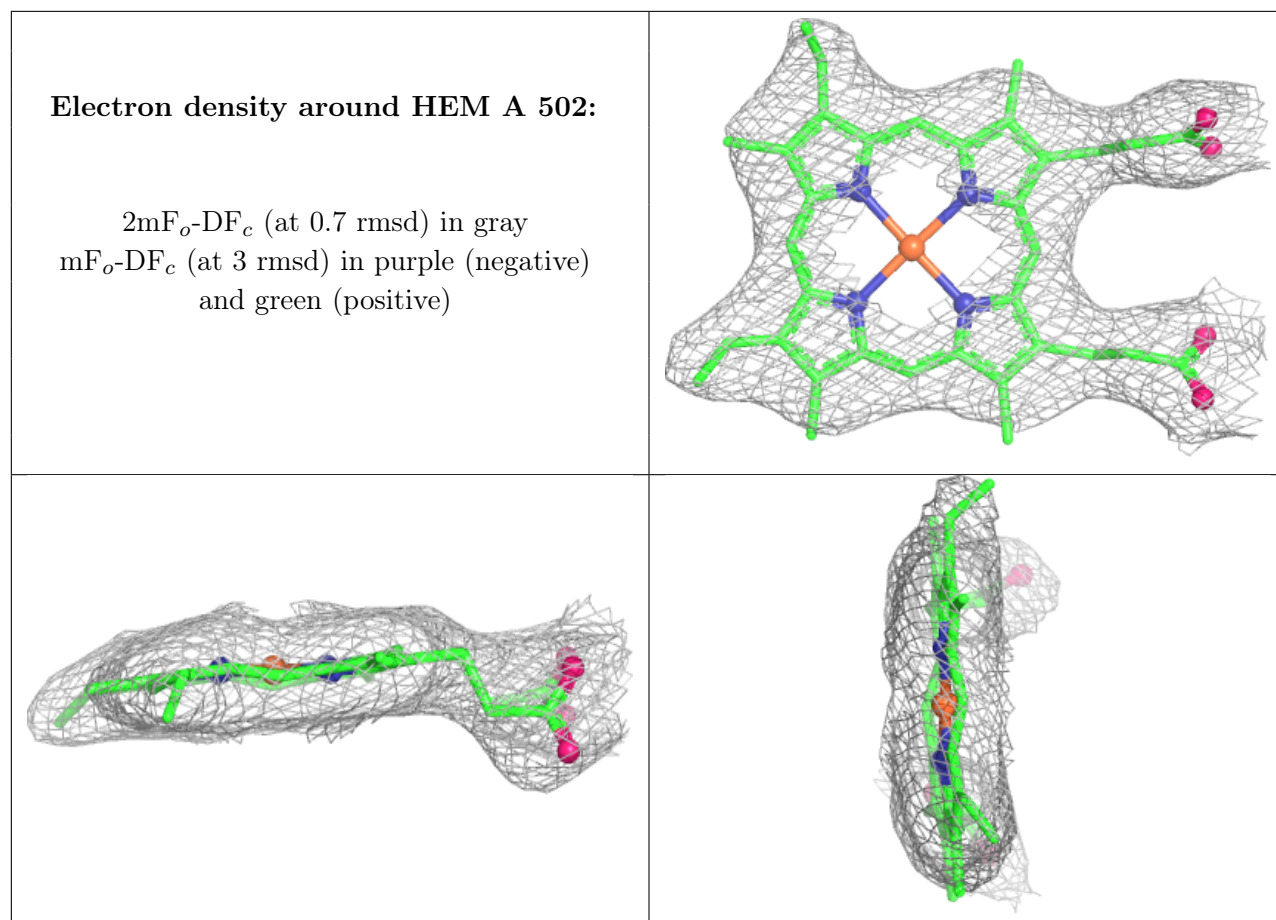
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



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.