



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 07:22 PM UTC

PDB ID : 3BE0 / pdb_00003be0
Title : The Role of Asn 242 in P450cin
Authors : Mehareenna, Y.T.; Poulos, T.L.
Deposited on : 2007-11-15
Resolution : 3.05 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 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

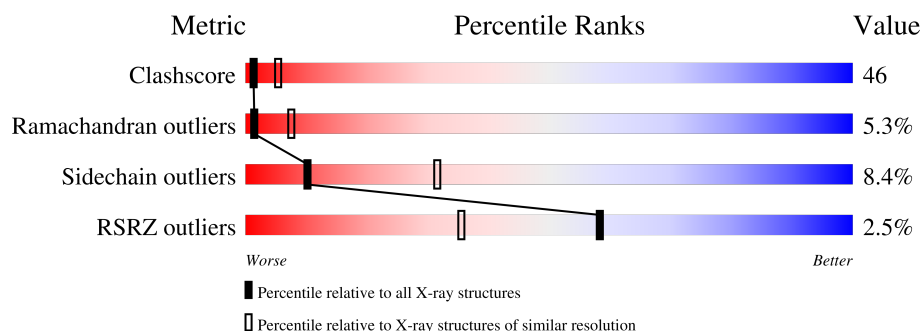
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.05 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	2569 (3.10-3.02)
Ramachandran outliers	187476	2424 (3.10-3.02)
Sidechain outliers	187428	2423 (3.10-3.02)
RSRZ outliers	180081	2469 (3.10-3.02)

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	397	2% 36% 53% 11%
1	B	397	3% 33% 55% 11%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6489 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 P450cin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	397	Total	C	N	O	S	0	0	0
			3146	2006	548	581	11			
1	B	397	Total	C	N	O	S	0	0	0
			3146	2006	548	581	11			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	242	ALA	ASN	engineered mutation	UNP Q8VQF6
B	242	ALA	ASN	engineered mutation	UNP Q8VQF6

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



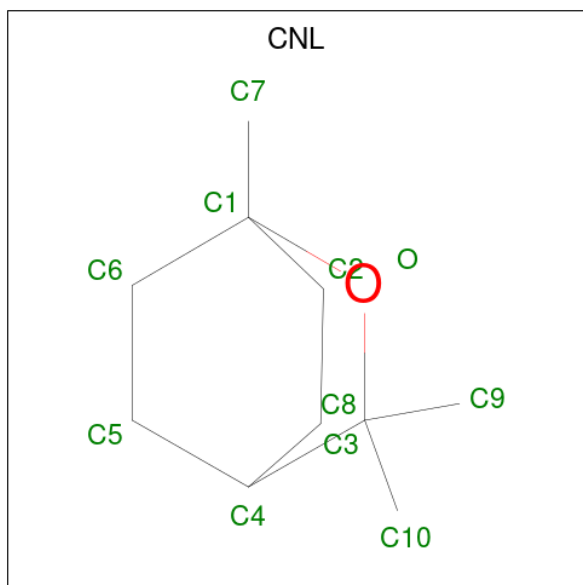
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

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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		

- Molecule 3 is 1,3,3-TRIMETHYL-2-OXABICYCLO[2.2.2]OCTANE (CCD ID: CNL) (formula: C₁₀H₁₈O).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			11	10	1		
3	B	1	Total	C	O	0	0
			11	10	1		

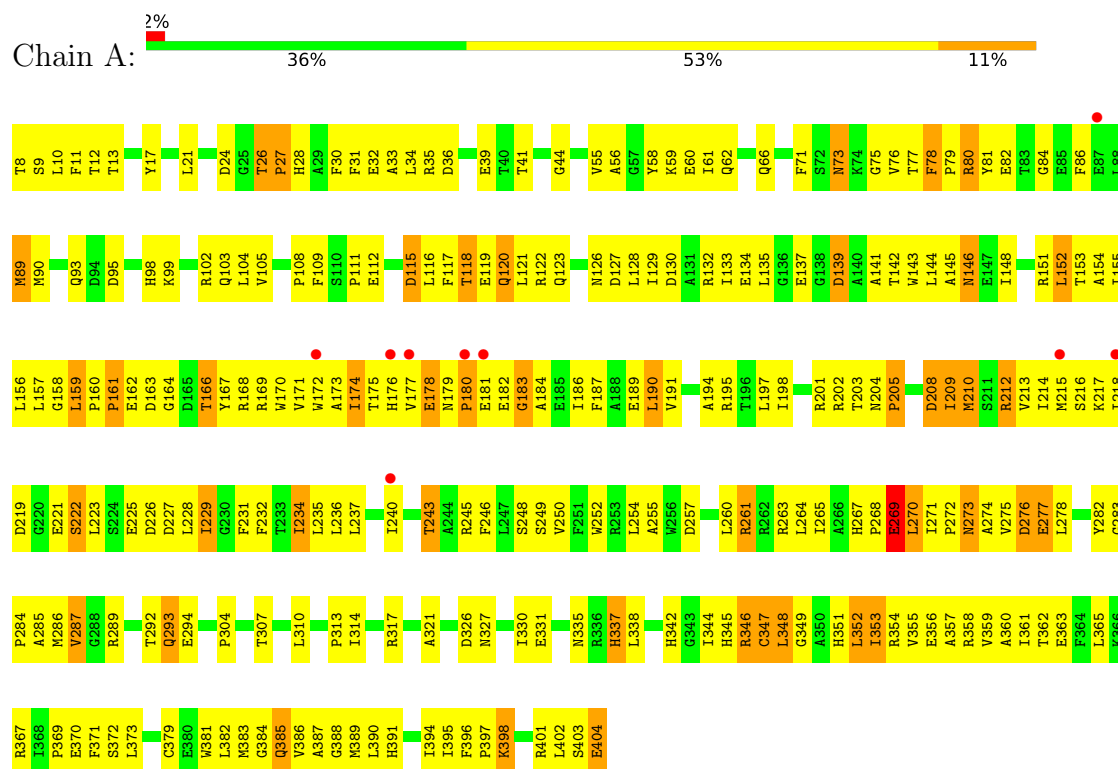
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	56	Total	O	0	0
			56	56		
4	B	33	Total	O	0	0
			33	33		

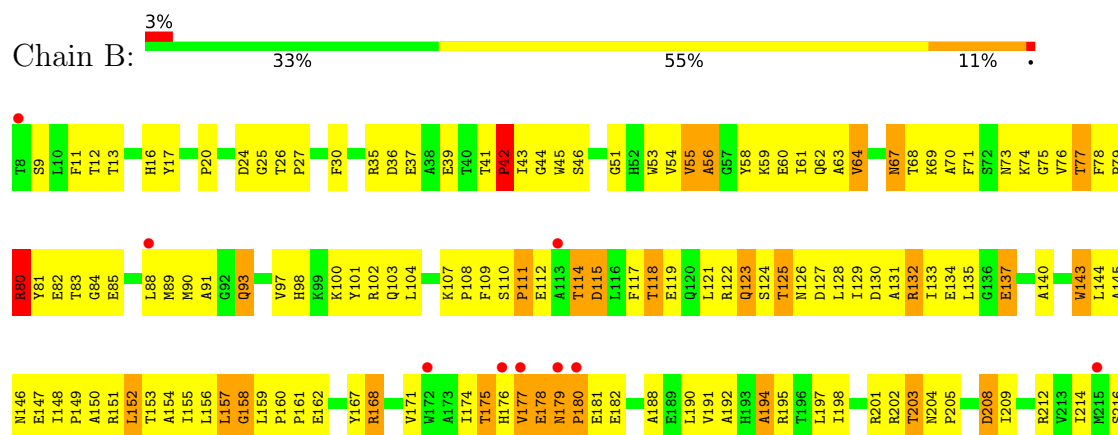
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: P450cin



• Molecule 1: P450cin



T362	K217	Q293	H267	A350	L402
E363	I218	E294	P268	H351	S403
F364	D219	V295	E269	L352	E404
L365	G220	T296	L270	I353	
K366	E221	V297	I271	R354	
R367	G222	G298	P272	V355	
T368	L223	T299	N273	E356	
F369	S224	I300	A274	A357	
E370	E225	T301	D276	R358	
F371		M302	L279	V359	
S372	L228	K303	R280	A360	
L373	I229	P304	P284	I361	
D374	G230	G305	A285		
F375	F231	Q306	N286		
N376		T307	V287		
K377	I234		G288		
E378	L235	L310	R289		
C379	L236	F311	L290		
	L237	F312	V291		
		P313	T292		
		I314			
		A315			
		A321			
		F322			
		D323			
		S324			
		F325			
		D326			
		N327			
		I328			
		V329			
		I330			
		E331			
		R336			
		H337			
		L338			
		S339			
		L340			
		G341			
		H342			
		G343			
		I344			
		H345			
		R346			
		C347			
		A350			
		H351			
		L352			
		I353			
		R354			
		V355			
		E356			
		A357			
		R358			
		V359			
		A360			
		I361			

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	59.40Å 128.68Å 69.02Å 90.00° 97.84° 90.00°	Depositor
Resolution (Å)	47.95 – 3.05 47.95 – 3.05	Depositor EDS
% Data completeness (in resolution range)	91.4 (47.95-3.05) 91.4 (47.95-3.05)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	0.15	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.67 (at 3.07Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.236 , 0.322 0.244 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	54.7	Xtriage
Anisotropy	0.630	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 69.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	6489	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 18.82% 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, CNL

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	0.51	0/3226	1.07	25/4386 (0.6%)
1	B	0.50	0/3226	1.04	19/4386 (0.4%)
All	All	0.51	0/6452	1.05	44/8772 (0.5%)

There are no bond length outliers.

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	182	GLU	N-CA-C	-8.40	102.58	113.17
1	A	204	ASN	CA-C-N	8.31	130.22	119.84
1	A	204	ASN	C-N-CA	8.31	130.22	119.84
1	B	157	LEU	N-CA-C	-7.82	103.47	113.16
1	A	146	ASN	N-CA-C	7.77	119.53	111.14
1	A	283	GLY	CA-C-N	7.71	127.89	119.87
1	A	283	GLY	C-N-CA	7.71	127.89	119.87
1	A	17	TYR	N-CA-C	-7.59	104.04	113.38
1	A	120	GLN	N-CA-C	-7.28	104.65	113.97
1	B	125	THR	N-CA-C	-7.21	103.11	110.97
1	B	287	VAL	N-CA-C	7.17	119.14	108.96
1	A	174	ILE	N-CA-C	-6.95	104.00	110.53
1	B	17	TYR	N-CA-C	-6.60	104.80	112.92
1	A	183	GLY	N-CA-C	-6.30	104.41	113.86
1	A	115	ASP	N-CA-C	-6.24	105.69	113.55
1	B	248	SER	N-CA-C	-6.23	104.40	111.07
1	B	147	GLU	N-CA-C	6.18	120.08	112.54
1	A	269	GLU	N-CA-C	-6.14	104.64	111.71
1	B	115	ASP	N-CA-C	-6.12	105.52	114.39
1	A	128	LEU	N-CA-C	-6.08	105.91	113.38
1	B	267	HIS	CA-C-N	6.00	126.11	119.87
1	B	267	HIS	C-N-CA	6.00	126.11	119.87

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	242	ALA	N-CA-C	-5.93	104.15	111.33
1	A	385	GLN	N-CA-C	-5.89	105.94	113.01
1	A	229	ILE	N-CA-C	-5.83	105.42	113.00
1	A	190	LEU	N-CA-C	-5.80	104.95	111.28
1	A	347	CYS	CA-C-N	-5.74	112.14	120.29
1	A	347	CYS	C-N-CA	-5.74	112.14	120.29
1	A	227	ASP	N-CA-C	-5.66	106.73	113.97
1	A	162	GLU	N-CA-C	5.61	117.62	109.71
1	B	16	HIS	N-CA-C	5.45	119.78	113.18
1	A	352	LEU	N-CA-C	5.43	117.20	111.28
1	A	159	LEU	CA-C-N	5.29	125.83	120.38
1	A	159	LEU	C-N-CA	5.29	125.83	120.38
1	B	378	GLU	N-CA-C	5.21	118.22	110.14
1	B	194	ALA	N-CA-C	-5.18	105.81	111.82
1	B	114	THR	N-CA-C	-5.17	107.48	112.97
1	B	224	SER	N-CA-C	5.12	118.10	110.52
1	A	287	VAL	N-CA-C	5.11	116.25	108.80
1	B	295	VAL	N-CA-C	5.11	114.59	108.06
1	B	56	ALA	N-CA-C	5.06	118.56	112.38
1	B	77	THR	N-CA-C	-5.04	102.24	110.20
1	A	148	ILE	CB-CA-C	-5.04	109.01	114.35
1	A	277	GLU	N-CA-C	-5.00	106.02	111.82

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	3146	0	3098	295	0
1	B	3146	0	3098	294	1
2	A	43	0	30	7	0
2	B	43	0	30	5	0
3	A	11	0	18	0	0
3	B	11	0	18	0	0
4	A	56	0	0	23	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	33	0	0	8	0
All	All	6489	0	6292	583	1

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 46.

All (583) 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:326:ASP:HB3	1:B:263:ARG:NH2	1.76	0.99
1:B:122:ARG:HH22	1:B:362:THR:HG21	1.28	0.97
1:B:398:LYS:NZ	1:B:398:LYS:HB3	1.82	0.95
1:B:133:ILE:HG13	1:B:368:ILE:HD11	1.49	0.92
1:A:385:GLN:OE1	4:A:538:HOH:O	1.89	0.91
1:B:80:ARG:HD3	1:B:81:TYR:N	1.88	0.89
1:B:158:GLY:HA3	1:B:201:ARG:HH22	1.37	0.89
1:A:370:GLU:HG2	4:A:549:HOH:O	1.74	0.88
1:A:370:GLU:CG	4:A:549:HOH:O	2.21	0.87
1:A:122:ARG:HG3	1:A:363:GLU:HG3	1.57	0.86
1:A:327:ASN:HA	4:B:531:HOH:O	1.74	0.86
1:B:157:LEU:HB2	1:B:159:LEU:CD1	2.06	0.85
1:B:90:MET:HE3	1:B:102:ARG:HA	1.58	0.84
1:B:157:LEU:HB2	1:B:159:LEU:HD11	1.58	0.84
1:B:375:PRO:HD3	1:B:395:ILE:HD13	1.60	0.83
1:B:354:ARG:HB3	1:B:358:ARG:HH22	1.43	0.83
1:A:357:ALA:O	1:A:361:ILE:HG22	1.78	0.82
1:A:287:VAL:HA	4:A:522:HOH:O	1.81	0.80
1:B:179:ASN:C	1:B:181:GLU:H	1.89	0.80
1:A:212:ARG:NH1	1:A:212:ARG:HB2	1.96	0.79
1:B:272:PRO:HG3	1:B:358:ARG:HD2	1.63	0.79
1:A:194:ALA:O	1:A:198:ILE:HG13	1.83	0.79
1:A:270:LEU:HD12	4:A:524:HOH:O	1.83	0.77
1:A:217:LYS:HG2	1:A:222:SER:HA	1.67	0.77
1:B:74:LYS:HG3	1:B:290:LEU:HD23	1.66	0.77
1:A:130:ASP:HB3	1:A:401:ARG:CZ	2.16	0.77
1:A:353:ILE:HD13	1:A:353:ILE:H	1.48	0.77
1:A:151:ARG:O	1:A:155:ILE:HG12	1.85	0.76
1:A:367:ARG:HH22	1:A:401:ARG:HA	1.51	0.76
1:B:398:LYS:HB3	1:B:398:LYS:HZ2	1.51	0.75
1:A:264:LEU:HD13	1:A:361:ILE:HD11	1.67	0.75
1:B:146:ASN:HD22	1:B:168:ARG:HH21	1.35	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:151:ARG:O	1:B:155:ILE:HD13	1.87	0.74
1:A:396:PHE:HB2	1:A:397:PRO:HD2	1.69	0.74
1:B:152:LEU:HD21	1:B:352:LEU:HD11	1.68	0.74
1:A:327:ASN:HB3	1:B:327:ASN:HD22	1.51	0.74
1:A:172:TRP:CD1	1:A:177:VAL:HG21	2.23	0.73
1:A:225:GLU:HA	1:A:228:LEU:HG	1.69	0.73
1:A:212:ARG:HB2	1:A:212:ARG:HH11	1.52	0.73
1:A:61:ILE:HG23	1:A:310:LEU:HD22	1.70	0.73
1:A:143:TRP:O	1:A:144:LEU:HD23	1.89	0.73
1:A:156:LEU:HB3	1:A:209:ILE:HD11	1.72	0.72
1:B:133:ILE:HG13	1:B:368:ILE:CD1	2.21	0.71
1:A:156:LEU:O	1:A:208:ASP:HB2	1.91	0.71
1:B:137:GLU:HB3	1:B:395:ILE:HA	1.73	0.69
1:A:157:LEU:HB2	1:A:159:LEU:HD12	1.74	0.69
1:B:194:ALA:O	1:B:198:ILE:HG13	1.92	0.69
1:B:235:LEU:C	1:B:237:LEU:H	1.99	0.69
1:B:111:PRO:HB3	1:B:351:HIS:CD2	2.28	0.69
1:B:122:ARG:HA	1:B:359:VAL:HG21	1.75	0.69
1:A:358:ARG:NH2	4:A:545:HOH:O	2.24	0.68
1:B:67:ASN:HD21	1:B:70:ALA:HB3	1.57	0.68
1:A:159:LEU:O	1:A:161:PRO:HD3	1.94	0.68
1:A:401:ARG:HH21	1:A:403:SER:HB3	1.59	0.68
1:B:67:ASN:ND2	1:B:70:ALA:HB3	2.09	0.68
1:B:236:LEU:O	1:B:236:LEU:HG	1.94	0.67
1:A:121:LEU:HD21	1:A:152:LEU:HD22	1.77	0.67
1:A:146:ASN:HB2	4:A:539:HOH:O	1.94	0.67
1:B:131:ALA:O	1:B:132:ARG:HG2	1.95	0.67
1:B:337:HIS:HD2	1:B:339:SER:HB2	1.58	0.67
1:B:123:GLN:HG3	1:B:127:ASP:OD1	1.94	0.66
1:B:398:LYS:HB3	1:B:398:LYS:HZ3	1.57	0.66
1:A:111:PRO:HA	1:A:351:HIS:CD2	2.29	0.66
1:A:401:ARG:HH21	1:A:403:SER:CB	2.08	0.66
1:A:358:ARG:NE	4:A:545:HOH:O	2.24	0.66
1:A:209:ILE:O	1:A:212:ARG:HG2	1.95	0.66
1:A:55:VAL:O	1:A:61:ILE:HD11	1.97	0.65
1:B:346:ARG:HD3	1:B:347:CYS:N	2.11	0.65
1:B:179:ASN:O	1:B:181:GLU:N	2.29	0.65
1:A:126:ASN:ND2	1:A:363:GLU:OE2	2.27	0.65
1:B:267:HIS:HB2	1:B:270:LEU:HD12	1.79	0.65
1:A:213:VAL:O	1:A:216:SER:HB2	1.97	0.65
1:A:152:LEU:HD12	1:A:152:LEU:O	1.96	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:146:ASN:HD22	1:A:168:ARG:HD3	1.62	0.65
1:B:370:GLU:O	1:B:397:PRO:HD2	1.97	0.64
1:B:245:ARG:HA	1:B:245:ARG:NE	2.12	0.64
1:A:337:HIS:O	1:A:338:LEU:HD23	1.97	0.64
1:B:125:THR:HG23	1:B:148:ILE:HD11	1.80	0.64
1:B:35:ARG:HD2	1:B:36:ASP:N	2.13	0.64
1:A:130:ASP:OD2	1:A:403:SER:HB2	1.98	0.63
1:B:203:THR:C	1:B:205:PRO:HD3	2.23	0.63
1:B:80:ARG:HD3	1:B:81:TYR:H	1.64	0.63
1:B:243:THR:HG23	1:B:353:ILE:HD13	1.80	0.63
1:A:367:ARG:HH12	1:A:402:LEU:HG	1.63	0.63
1:A:348:LEU:O	1:A:348:LEU:HD22	1.99	0.63
1:B:237:LEU:C	1:B:237:LEU:HD23	2.24	0.62
1:A:35:ARG:HD2	1:A:35:ARG:C	2.24	0.62
1:A:89:MET:HB3	1:A:90:MET:HE2	1.80	0.62
1:A:250:VAL:O	1:A:254:LEU:HG	1.98	0.62
1:A:284:PRO:HB3	1:A:387:ALA:O	1.99	0.62
1:A:104:LEU:HD11	1:A:218:ILE:HD11	1.80	0.62
1:B:93:GLN:HE21	1:B:93:GLN:N	1.97	0.62
1:A:367:ARG:NH1	1:A:402:LEU:HG	2.14	0.62
1:A:234:ILE:HD12	2:A:450:HEM:CMD	2.30	0.62
1:B:76:VAL:HG13	1:B:77:THR:N	2.15	0.61
1:A:164:GLY:HA3	1:A:168:ARG:NH2	2.15	0.61
1:B:175:THR:HG22	1:B:175:THR:O	2.01	0.61
1:A:217:LYS:HG2	1:A:222:SER:CA	2.29	0.61
1:A:268:PRO:HA	1:A:271:ILE:HD12	1.82	0.61
1:B:24:ASP:OD1	1:B:26:THR:HG23	2.01	0.61
1:A:55:VAL:HB	1:A:310:LEU:HD23	1.83	0.61
1:A:129:ILE:O	1:A:133:ILE:HB	2.00	0.61
1:B:382:LEU:O	1:B:387:ALA:HB1	2.01	0.60
1:B:107:LYS:HD3	1:B:212:ARG:CZ	2.31	0.60
1:A:172:TRP:O	1:A:177:VAL:HG23	2.01	0.60
4:A:511:HOH:O	1:B:327:ASN:HA	2.02	0.60
1:A:12:THR:HG23	1:A:13:THR:N	2.16	0.60
1:A:90:MET:O	2:A:450:HEM:HBD2	2.02	0.60
1:A:122:ARG:HD3	1:A:363:GLU:OE2	2.02	0.60
1:A:11:PHE:CD2	1:A:44:GLY:HA3	2.37	0.60
1:B:132:ARG:HB2	1:B:143:TRP:CZ2	2.37	0.60
1:A:33:ALA:O	1:A:36:ASP:HB2	2.01	0.59
1:A:177:VAL:HG12	1:A:177:VAL:O	2.00	0.59
1:B:35:ARG:HD2	1:B:35:ARG:C	2.27	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:62:GLN:C	1:B:64:VAL:H	2.10	0.59
1:B:217:LYS:HG2	1:B:222:SER:HA	1.83	0.59
1:A:31:PHE:CD1	1:A:314:ILE:HD11	2.38	0.59
1:B:345:HIS:HA	2:B:450:HEM:O1D	2.02	0.59
1:B:268:PRO:HA	1:B:271:ILE:HD13	1.85	0.59
1:B:158:GLY:CA	1:B:201:ARG:HH22	2.13	0.59
1:A:129:ILE:HD13	1:A:367:ARG:HE	1.67	0.59
1:A:218:ILE:O	1:A:219:ASP:HB2	2.03	0.59
1:A:270:LEU:C	1:A:272:PRO:HD2	2.28	0.58
1:B:76:VAL:HG13	1:B:77:THR:H	1.67	0.58
1:B:146:ASN:HB3	1:B:168:ARG:NH2	2.18	0.58
1:B:253:ARG:CZ	1:B:328:ILE:HD12	2.33	0.58
1:A:12:THR:HA	4:A:505:HOH:O	2.02	0.58
1:B:111:PRO:HG2	1:B:112:GLU:OE1	2.03	0.58
1:A:152:LEU:HD12	1:A:152:LEU:C	2.28	0.58
1:A:234:ILE:HD12	2:A:450:HEM:HMD1	1.85	0.58
1:A:179:ASN:CG	1:A:180:PRO:HD2	2.29	0.58
1:B:231:PHE:HA	1:B:234:ILE:HG22	1.84	0.58
1:A:395:ILE:O	1:A:395:ILE:HG13	2.04	0.58
1:B:174:ILE:C	1:B:176:HIS:H	2.11	0.58
1:A:359:VAL:O	1:A:360:ALA:C	2.47	0.57
1:A:28:HIS:HE1	1:A:282:TYR:CE1	2.21	0.57
1:A:36:ASP:O	1:A:39:GLU:HG3	2.03	0.57
1:B:42:PRO:HB3	1:B:60:GLU:HG2	1.85	0.57
1:A:181:GLU:O	1:A:184:ALA:N	2.35	0.57
1:A:111:PRO:HA	1:A:351:HIS:HD2	1.67	0.57
1:A:121:LEU:HD21	1:A:152:LEU:CD2	2.34	0.57
1:B:121:LEU:C	1:B:123:GLN:H	2.12	0.57
1:A:252:TRP:HD1	1:A:373:LEU:HD21	1.69	0.57
1:B:46:SER:O	1:B:51:GLY:HA2	2.05	0.57
1:B:134:GLU:HG2	1:B:399:GLY:O	2.04	0.57
1:A:271:ILE:N	1:A:272:PRO:HD2	2.20	0.57
1:B:384:GLY:O	4:B:525:HOH:O	2.18	0.57
1:A:102:ARG:C	1:A:104:LEU:H	2.13	0.57
1:B:336:ARG:HG2	4:B:526:HOH:O	2.04	0.56
1:B:74:LYS:CG	1:B:290:LEU:HD23	2.35	0.56
1:B:228:LEU:C	1:B:230:GLY:H	2.13	0.56
1:A:243:THR:HG23	1:A:353:ILE:HG22	1.88	0.56
1:B:109:PHE:CZ	2:B:450:HEM:HBC1	2.40	0.56
1:B:122:ARG:NH2	1:B:362:THR:HG21	2.10	0.56
1:B:359:VAL:O	1:B:360:ALA:C	2.48	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:398:LYS:HB2	1:A:398:LYS:NZ	2.21	0.56
1:A:118:THR:C	1:A:120:GLN:H	2.14	0.56
1:A:252:TRP:CD1	1:A:373:LEU:HD21	2.41	0.56
1:A:285:ALA:HB3	4:A:501:HOH:O	2.05	0.56
1:A:358:ARG:CZ	4:A:545:HOH:O	2.54	0.56
1:A:123:GLN:O	1:A:127:ASP:HB2	2.06	0.56
1:B:230:GLY:O	1:B:234:ILE:HG22	2.05	0.56
1:A:172:TRP:NE1	1:A:177:VAL:HG21	2.21	0.56
1:B:150:ALA:HA	1:B:240:ILE:CD1	2.37	0.56
1:B:68:THR:HG22	1:B:344:ILE:HD11	1.87	0.55
1:A:129:ILE:HD12	1:A:130:ASP:N	2.21	0.55
1:B:235:LEU:C	1:B:237:LEU:N	2.65	0.55
1:A:153:THR:O	1:A:154:ALA:C	2.50	0.55
1:A:179:ASN:CG	1:A:182:GLU:HB2	2.31	0.55
1:B:64:VAL:HG11	1:B:310:LEU:HD21	1.89	0.55
1:B:379:CYS:HB3	1:B:389:MET:HE3	1.88	0.55
1:A:175:THR:O	1:A:176:HIS:HB3	2.07	0.55
1:B:102:ARG:C	1:B:104:LEU:H	2.14	0.55
1:B:252:TRP:HB2	1:B:392:VAL:HG11	1.88	0.55
1:A:246:PHE:O	1:A:249:SER:HB2	2.07	0.55
1:B:71:PHE:HB3	1:B:289:ARG:HD2	1.88	0.55
1:A:172:TRP:C	1:A:175:THR:HG22	2.32	0.55
1:A:214:ILE:HB	1:A:215:MET:HE2	1.87	0.55
1:A:277:GLU:HA	1:A:335:ASN:HD21	1.72	0.55
1:B:100:LYS:HD3	1:B:218:ILE:HG21	1.87	0.55
1:B:117:PHE:O	1:B:118:THR:C	2.49	0.55
1:B:176:HIS:O	1:B:177:VAL:C	2.50	0.55
1:B:292:THR:O	1:B:304:PRO:HG3	2.07	0.55
1:A:326:ASP:HB3	1:B:263:ARG:CZ	2.36	0.55
1:A:349:GLY:O	1:A:353:ILE:HD13	2.07	0.54
1:A:186:ILE:HA	1:A:189:GLU:OE1	2.07	0.54
1:A:203:THR:C	1:A:205:PRO:HD3	2.32	0.54
1:A:292:THR:C	1:A:304:PRO:HB3	2.33	0.54
1:B:152:LEU:HD12	1:B:152:LEU:O	2.07	0.54
1:A:8:THR:N	4:A:515:HOH:O	2.41	0.54
1:A:62:GLN:O	1:A:66:GLN:HG3	2.07	0.54
1:B:133:ILE:CG1	1:B:368:ILE:HD11	2.31	0.54
1:A:95:ASP:OD1	1:A:99:LYS:N	2.37	0.54
1:B:398:LYS:NZ	1:B:398:LYS:CB	2.64	0.54
1:A:129:ILE:HD12	1:A:129:ILE:C	2.33	0.54
1:A:179:ASN:ND2	1:A:182:GLU:HB2	2.23	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:162:GLU:CD	1:B:162:GLU:H	2.16	0.54
1:B:204:ASN:N	1:B:205:PRO:HD3	2.23	0.54
1:B:243:THR:HA	2:B:450:HEM:HBB1	1.90	0.54
1:A:213:VAL:HA	1:A:216:SER:OG	2.08	0.53
1:A:89:MET:HE2	1:A:93:GLN:HE22	1.72	0.53
1:B:216:SER:C	1:B:217:LYS:HG3	2.32	0.53
1:B:123:GLN:O	1:B:126:ASN:N	2.41	0.53
1:B:231:PHE:HA	1:B:234:ILE:CG2	2.38	0.53
1:A:367:ARG:NH2	1:A:401:ARG:HA	2.22	0.53
1:A:401:ARG:HE	1:A:403:SER:HB3	1.74	0.53
1:B:126:ASN:OD1	1:B:363:GLU:HB3	2.08	0.53
1:A:76:VAL:HG13	1:A:77:THR:N	2.23	0.53
1:A:153:THR:O	1:A:156:LEU:N	2.36	0.53
1:A:154:ALA:HA	1:A:159:LEU:HB2	1.90	0.53
1:A:381:TRP:CZ3	1:A:388:GLY:HA2	2.44	0.53
1:A:163:ASP:HB2	1:A:167:TYR:CE2	2.44	0.53
1:A:289:ARG:O	1:A:307:THR:HG23	2.08	0.53
1:B:358:ARG:NH2	4:B:510:HOH:O	2.42	0.53
1:A:201:ARG:C	1:A:203:THR:H	2.16	0.52
1:B:107:LYS:HD3	1:B:212:ARG:NH1	2.23	0.52
1:B:350:ALA:O	1:B:353:ILE:HG22	2.09	0.52
1:B:67:ASN:HD21	1:B:70:ALA:CB	2.21	0.52
1:B:272:PRO:O	1:B:274:ALA:N	2.42	0.52
1:A:24:ASP:OD2	1:A:26:THR:HG23	2.08	0.52
1:A:235:LEU:C	1:A:237:LEU:H	2.16	0.52
1:B:241:ASP:O	1:B:242:ALA:C	2.51	0.52
1:B:351:HIS:CE1	4:B:515:HOH:O	2.61	0.52
1:A:163:ASP:O	1:A:164:GLY:C	2.53	0.52
1:A:357:ALA:O	1:A:358:ARG:C	2.53	0.52
1:B:76:VAL:HA	1:B:288:GLY:O	2.10	0.52
1:B:160:PRO:HB3	1:B:162:GLU:OE1	2.08	0.52
1:B:228:LEU:C	1:B:230:GLY:N	2.65	0.52
1:B:286:MET:HE2	1:B:313:PRO:CD	2.39	0.52
1:B:359:VAL:HG23	1:B:363:GLU:HG2	1.91	0.52
1:A:89:MET:HG2	1:A:231:PHE:CE1	2.45	0.52
1:A:146:ASN:CB	4:A:539:HOH:O	2.56	0.52
1:A:126:ASN:C	1:A:129:ILE:HG13	2.35	0.52
1:A:175:THR:O	1:A:176:HIS:CB	2.58	0.52
1:B:153:THR:O	1:B:154:ALA:C	2.53	0.52
1:B:218:ILE:C	1:B:220:GLY:H	2.18	0.52
1:B:272:PRO:C	1:B:274:ALA:H	2.18	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:284:PRO:O	1:B:386:VAL:HA	2.10	0.52
1:A:132:ARG:HA	1:A:135:LEU:HD13	1.92	0.51
1:B:280:ARG:HG3	1:B:337:HIS:HB3	1.90	0.51
1:A:41:THR:O	1:A:41:THR:HG22	2.11	0.51
1:A:270:LEU:CD1	4:A:524:HOH:O	2.47	0.51
1:B:269:GLU:O	1:B:272:PRO:HD2	2.10	0.51
1:A:272:PRO:HG2	1:A:273:ASN:H	1.76	0.51
1:A:276:ASP:OD1	1:A:354:ARG:NH1	2.43	0.51
1:B:110:SER:C	1:B:112:GLU:H	2.18	0.51
1:B:140:ALA:HB3	1:B:392:VAL:HG23	1.92	0.51
1:B:274:ALA:C	1:B:276:ASP:N	2.66	0.51
1:B:337:HIS:CD2	1:B:339:SER:HB2	2.44	0.51
1:A:370:GLU:CD	4:A:549:HOH:O	2.50	0.51
1:B:229:ILE:O	1:B:229:ILE:HG22	2.11	0.51
1:A:354:ARG:HG2	1:A:354:ARG:HH11	1.76	0.51
1:B:253:ARG:NH2	1:B:328:ILE:HD12	2.25	0.51
1:B:93:GLN:N	1:B:93:GLN:NE2	2.58	0.51
1:B:304:PRO:C	1:B:306:GLN:H	2.19	0.51
1:A:78:PHE:HA	1:A:79:PRO:C	2.35	0.51
1:A:84:GLY:HA3	4:A:534:HOH:O	2.11	0.51
1:A:179:ASN:OD1	1:A:182:GLU:HB2	2.11	0.51
1:A:267:HIS:N	1:A:268:PRO:HD3	2.26	0.51
1:B:42:PRO:HB3	1:B:60:GLU:CD	2.36	0.51
1:B:268:PRO:HA	1:B:271:ILE:CD1	2.41	0.51
1:B:352:LEU:O	1:B:356:GLU:HG3	2.11	0.51
1:B:336:ARG:NH1	1:B:336:ARG:HB3	2.26	0.51
1:A:170:TRP:HB3	1:A:190:LEU:HD12	1.93	0.50
1:A:396:PHE:HB2	1:A:397:PRO:CD	2.40	0.50
1:B:26:THR:HB	1:B:27:PRO:HD2	1.93	0.50
1:B:178:GLU:O	1:B:179:ASN:HB2	2.10	0.50
1:B:181:GLU:HA	1:B:181:GLU:OE1	2.11	0.50
1:A:326:ASP:HB3	1:B:263:ARG:HH21	1.72	0.50
1:A:76:VAL:HG13	1:A:77:THR:HG23	1.93	0.50
1:A:139:ASP:HB3	1:A:142:THR:OG1	2.11	0.50
1:B:43:ILE:HG12	1:B:55:VAL:HG13	1.92	0.50
1:B:148:ILE:N	1:B:149:PRO:HD2	2.26	0.50
1:B:359:VAL:HG23	1:B:363:GLU:CG	2.41	0.50
1:A:156:LEU:HB3	1:A:209:ILE:CD1	2.41	0.50
1:A:102:ARG:C	1:A:104:LEU:N	2.69	0.50
1:A:223:LEU:N	1:A:223:LEU:HD12	2.27	0.50
1:A:397:PRO:HG2	4:A:527:HOH:O	2.10	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:353:ILE:HG23	1:B:354:ARG:N	2.27	0.50
1:A:164:GLY:HA3	1:A:168:ARG:HH22	1.77	0.50
1:A:286:MET:HE2	1:A:313:PRO:HD2	1.94	0.50
1:B:102:ARG:O	1:B:104:LEU:N	2.42	0.50
1:B:243:THR:HG23	1:B:353:ILE:CD1	2.41	0.50
1:B:80:ARG:HD3	1:B:80:ARG:C	2.37	0.49
1:B:122:ARG:O	1:B:122:ARG:HG2	2.10	0.49
1:A:108:PRO:HG3	1:A:212:ARG:HD3	1.94	0.49
1:A:95:ASP:OD1	1:A:98:HIS:HB3	2.12	0.49
1:B:79:PRO:O	1:B:80:ARG:C	2.54	0.49
1:A:269:GLU:HG2	1:A:270:LEU:HD13	1.95	0.49
1:A:351:HIS:O	1:A:355:VAL:HG23	2.11	0.49
1:B:159:LEU:HB3	1:B:167:TYR:OH	2.13	0.49
1:B:237:LEU:HD23	1:B:237:LEU:O	2.12	0.49
1:A:86:PHE:HE2	1:A:226:ASP:HA	1.78	0.49
1:A:156:LEU:O	1:A:209:ILE:HG13	2.11	0.49
1:A:181:GLU:C	1:A:183:GLY:N	2.70	0.49
1:B:122:ARG:HD2	4:B:503:HOH:O	2.12	0.49
1:B:124:SER:O	1:B:128:LEU:HG	2.13	0.49
1:B:67:ASN:ND2	1:B:70:ALA:H	2.11	0.49
1:B:271:ILE:O	1:B:275:VAL:HG23	2.11	0.49
1:A:171:VAL:HA	1:A:174:ILE:HD12	1.95	0.49
1:B:375:PRO:CD	1:B:395:ILE:HD13	2.39	0.49
1:B:101:TYR:HA	1:B:104:LEU:HD12	1.95	0.48
1:B:174:ILE:O	1:B:176:HIS:N	2.46	0.48
1:B:377:LYS:HB3	1:B:393:PRO:HG3	1.94	0.48
1:A:198:ILE:HD11	1:A:229:ILE:HG12	1.94	0.48
1:A:235:LEU:C	1:A:237:LEU:N	2.68	0.48
1:A:245:ARG:NE	1:A:245:ARG:HA	2.28	0.48
1:B:56:ALA:HB1	1:B:314:ILE:HG21	1.95	0.48
1:B:89:MET:HG3	1:B:90:MET:HG3	1.95	0.48
1:B:109:PHE:HA	1:B:114:THR:CG2	2.42	0.48
1:B:158:GLY:C	1:B:159:LEU:HD12	2.39	0.48
1:A:79:PRO:O	1:A:80:ARG:C	2.57	0.48
1:A:10:LEU:HB2	1:A:34:LEU:HD13	1.94	0.48
1:A:102:ARG:O	1:A:104:LEU:N	2.47	0.48
1:A:130:ASP:OD1	1:A:401:ARG:HB2	2.12	0.48
1:A:181:GLU:C	1:A:183:GLY:H	2.21	0.48
1:A:181:GLU:C	1:A:181:GLU:CD	2.81	0.48
1:B:73:ASN:HB3	1:B:91:ALA:O	2.13	0.48
1:B:123:GLN:C	1:B:125:THR:N	2.69	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:129:ILE:O	1:B:131:ALA:N	2.43	0.48
1:B:329:VAL:CG1	1:B:331:GLU:HG2	2.44	0.48
1:B:270:LEU:O	1:B:271:ILE:C	2.56	0.48
1:B:110:SER:O	1:B:112:GLU:N	2.46	0.48
1:A:261:ARG:HG3	1:A:261:ARG:HH11	1.78	0.48
1:A:263:ARG:NH2	1:B:326:ASP:HB3	2.29	0.48
1:B:12:THR:HG23	1:B:13:THR:HG23	1.95	0.48
1:B:130:ASP:O	1:B:130:ASP:OD2	2.32	0.48
1:B:119:GLU:C	1:B:121:LEU:N	2.71	0.47
1:A:152:LEU:HD11	1:A:156:LEU:HD21	1.95	0.47
1:A:177:VAL:C	1:A:179:ASN:N	2.70	0.47
1:B:361:ILE:O	1:B:364:PHE:N	2.47	0.47
1:B:217:LYS:HE3	1:B:222:SER:OG	2.15	0.47
1:A:154:ALA:O	1:A:159:LEU:N	2.46	0.47
1:B:346:ARG:CZ	1:B:350:ALA:HB1	2.44	0.47
1:A:361:ILE:HG23	1:A:362:THR:N	2.29	0.47
1:A:365:LEU:O	1:A:369:PRO:HG3	2.14	0.47
1:B:43:ILE:HG22	1:B:44:GLY:N	2.30	0.47
1:B:20:PRO:HD2	1:B:30:PHE:CD2	2.50	0.47
1:A:11:PHE:CG	1:A:44:GLY:HA3	2.48	0.47
1:A:55:VAL:HG12	1:A:61:ILE:CG1	2.45	0.47
1:A:187:PHE:O	1:A:191:VAL:HG23	2.15	0.47
1:B:45:TRP:CZ2	1:B:51:GLY:HA3	2.50	0.47
1:B:195:ARG:HD3	1:B:225:GLU:OE2	2.14	0.47
1:B:234:ILE:HG13	2:B:450:HEM:HMD1	1.94	0.47
1:A:112:GLU:OE1	1:A:112:GLU:N	2.45	0.47
1:B:41:THR:N	1:B:42:PRO:HD3	2.30	0.47
1:B:271:ILE:HB	1:B:272:PRO:HD3	1.97	0.47
1:A:157:LEU:O	1:A:158:GLY:C	2.58	0.47
1:A:269:GLU:HG2	1:A:270:LEU:N	2.30	0.47
1:B:41:THR:HG22	1:B:41:THR:O	2.14	0.47
1:B:98:HIS:CE1	1:B:344:ILE:HB	2.50	0.47
1:B:156:LEU:O	1:B:208:ASP:HB2	2.15	0.47
1:A:56:ALA:HB1	1:A:314:ILE:HG21	1.96	0.46
1:B:354:ARG:HH11	1:B:354:ARG:HG2	1.78	0.46
1:A:345:HIS:O	2:A:450:HEM:HAA2	2.15	0.46
1:A:12:THR:CG2	1:A:13:THR:N	2.79	0.46
1:A:155:ILE:HD11	1:A:161:PRO:HG3	1.96	0.46
1:A:264:LEU:CD1	1:A:361:ILE:HD11	2.40	0.46
1:B:12:THR:HG23	1:B:13:THR:N	2.29	0.46
1:B:100:LYS:HB3	1:B:218:ILE:HD13	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:361:ILE:O	1:B:363:GLU:N	2.49	0.46
1:B:274:ALA:C	1:B:276:ASP:H	2.23	0.46
1:B:376:ASN:O	1:B:377:LYS:HG2	2.16	0.46
1:A:354:ARG:NH1	1:A:354:ARG:HG2	2.30	0.46
1:B:237:LEU:HD21	4:B:507:HOH:O	2.16	0.46
1:B:179:ASN:ND2	1:B:180:PRO:HD2	2.31	0.46
1:B:361:ILE:CG2	1:B:365:LEU:HD12	2.46	0.46
1:A:126:ASN:HA	1:A:129:ILE:HG12	1.98	0.46
1:A:177:VAL:O	1:A:179:ASN:N	2.49	0.46
1:B:201:ARG:O	1:B:214:ILE:HD12	2.16	0.46
1:A:212:ARG:HG3	1:A:213:VAL:N	2.31	0.46
1:A:278:LEU:HD23	1:A:278:LEU:HA	1.68	0.46
1:A:77:THR:HB	1:A:81:TYR:CD1	2.51	0.45
1:A:122:ARG:HD3	1:A:363:GLU:CD	2.40	0.45
1:A:176:HIS:CD2	1:A:178:GLU:H	2.34	0.45
1:A:123:GLN:O	1:A:127:ASP:N	2.49	0.45
1:A:129:ILE:HD13	1:A:367:ARG:NE	2.30	0.45
1:A:352:LEU:O	1:A:356:GLU:HG3	2.16	0.45
1:A:397:PRO:O	1:A:398:LYS:C	2.59	0.45
1:B:237:LEU:C	1:B:237:LEU:CD2	2.89	0.45
1:B:254:LEU:HB3	1:B:371:PHE:HZ	1.81	0.45
1:A:274:ALA:HB2	1:A:330:ILE:HB	1.98	0.45
1:B:78:PHE:HA	1:B:80:ARG:N	2.32	0.45
1:B:217:LYS:HE2	1:B:220:GLY:O	2.16	0.45
1:A:231:PHE:O	1:A:232:PHE:C	2.59	0.45
1:B:67:ASN:HD21	1:B:70:ALA:H	1.63	0.45
1:B:363:GLU:OE1	1:B:363:GLU:HA	2.16	0.45
1:B:384:GLY:O	1:B:385:GLN:C	2.60	0.45
1:B:403:SER:O	1:B:404:GLU:CB	2.63	0.45
1:A:121:LEU:HD11	1:A:152:LEU:CD2	2.47	0.45
1:A:345:HIS:HB3	2:A:450:HEM:CGA	2.47	0.45
1:B:272:PRO:C	1:B:274:ALA:N	2.74	0.45
1:A:141:ALA:HA	1:A:145:ALA:HB3	1.99	0.45
1:A:209:ILE:O	1:A:210:MET:C	2.59	0.45
1:B:109:PHE:HA	1:B:114:THR:HG21	1.98	0.45
1:B:129:ILE:C	1:B:131:ALA:H	2.24	0.45
1:A:157:LEU:HB2	1:A:159:LEU:CD1	2.43	0.45
1:A:402:LEU:O	4:A:528:HOH:O	2.21	0.45
1:B:88:LEU:O	1:B:89:MET:HB3	2.16	0.45
1:B:121:LEU:C	1:B:123:GLN:N	2.75	0.45
1:A:172:TRP:O	1:A:175:THR:HG22	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:111:PRO:O	1:B:115:ASP:HB2	2.15	0.45
1:B:359:VAL:HA	1:B:362:THR:HB	1.98	0.45
1:B:371:PHE:CB	1:B:396:PHE:HB3	2.47	0.45
1:A:109:PHE:CG	1:A:348:LEU:HD11	2.51	0.45
1:A:118:THR:O	1:A:120:GLN:N	2.50	0.45
1:A:130:ASP:HB3	1:A:401:ARG:NE	2.31	0.45
1:B:179:ASN:C	1:B:181:GLU:N	2.59	0.45
1:B:198:ILE:CG2	1:B:228:LEU:HD13	2.47	0.45
1:A:30:PHE:N	4:A:512:HOH:O	2.39	0.44
1:A:236:LEU:HD12	1:A:240:ILE:CD1	2.48	0.44
1:B:122:ARG:O	1:B:122:ARG:CG	2.65	0.44
1:B:202:ARG:O	1:B:202:ARG:HG3	2.16	0.44
1:B:337:HIS:HD2	1:B:339:SER:CB	2.28	0.44
1:A:132:ARG:C	1:A:134:GLU:H	2.24	0.44
1:A:89:MET:HE2	1:A:93:GLN:NE2	2.32	0.44
1:A:248:SER:CB	1:A:389:MET:O	2.65	0.44
1:B:42:PRO:HB3	1:B:60:GLU:CG	2.47	0.44
1:A:126:ASN:HD21	1:A:363:GLU:CD	2.23	0.44
1:B:303:LYS:HB3	1:B:304:PRO:CD	2.48	0.44
1:A:71:PHE:HB3	1:A:289:ARG:HB3	1.99	0.44
1:A:170:TRP:O	1:A:174:ILE:HG13	2.17	0.44
1:A:170:TRP:CE3	1:A:236:LEU:HD21	2.53	0.44
1:A:330:ILE:HG12	4:A:503:HOH:O	2.17	0.44
1:B:127:ASP:C	1:B:129:ILE:N	2.74	0.44
1:B:195:ARG:HG2	1:B:229:ILE:HD11	1.99	0.44
1:A:143:TRP:C	1:A:144:LEU:HD23	2.42	0.44
1:A:381:TRP:NE1	1:A:382:LEU:O	2.51	0.44
1:A:384:GLY:C	1:A:386:VAL:H	2.25	0.44
1:B:9:SER:HB2	1:B:37:GLU:OE2	2.17	0.44
1:B:133:ILE:HD13	1:B:367:ARG:NH2	2.33	0.44
1:A:144:LEU:O	1:A:145:ALA:C	2.59	0.44
1:B:108:PRO:HB2	1:B:209:ILE:HD11	2.00	0.44
1:B:174:ILE:C	1:B:176:HIS:N	2.76	0.44
1:B:221:GLU:OE2	1:B:221:GLU:HA	2.18	0.44
1:A:255:ALA:HA	1:A:371:PHE:CE1	2.53	0.44
1:B:132:ARG:HG3	1:B:143:TRP:NE1	2.33	0.44
1:B:217:LYS:HA	1:B:223:LEU:CD1	2.47	0.44
1:A:168:ARG:NH1	1:A:168:ARG:HG3	2.33	0.43
1:A:217:LYS:HG2	1:A:222:SER:N	2.32	0.43
1:B:361:ILE:O	1:B:362:THR:C	2.60	0.43
1:A:73:ASN:CG	1:A:345:HIS:NE2	2.77	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:115:ASP:C	1:A:117:PHE:H	2.26	0.43
1:A:170:TRP:HE3	1:A:236:LEU:HD21	1.82	0.43
1:B:80:ARG:HD2	1:B:81:TYR:O	2.17	0.43
1:B:123:GLN:O	1:B:124:SER:C	2.61	0.43
1:B:275:VAL:HG11	1:B:357:ALA:HB3	2.00	0.43
1:A:212:ARG:HH11	1:A:212:ARG:CB	2.26	0.43
1:A:381:TRP:CD1	1:A:382:LEU:O	2.71	0.43
1:B:37:GLU:C	1:B:39:GLU:H	2.26	0.43
1:B:130:ASP:O	1:B:130:ASP:CG	2.60	0.43
1:B:262:ARG:HG2	1:B:262:ARG:HH11	1.83	0.43
1:A:55:VAL:HG12	1:A:61:ILE:HG13	2.00	0.43
1:A:129:ILE:HD13	1:A:367:ARG:HH21	1.82	0.43
1:A:137:GLU:HB2	1:A:394:ILE:O	2.19	0.43
1:A:245:ARG:HD2	1:A:382:LEU:CD2	2.48	0.43
1:A:269:GLU:CG	1:A:270:LEU:HD13	2.48	0.43
1:B:155:ILE:HD11	1:B:161:PRO:HG3	1.99	0.43
1:B:404:GLU:C	1:B:404:GLU:OE1	2.61	0.43
1:A:9:SER:O	1:A:12:THR:HG22	2.19	0.43
1:A:166:THR:O	1:A:169:ARG:HB3	2.18	0.43
1:B:93:GLN:HG3	1:B:97:VAL:CG1	2.49	0.43
1:B:109:PHE:CE1	2:B:450:HEM:HBC1	2.53	0.43
1:B:338:LEU:C	1:B:340:LEU:H	2.27	0.43
1:A:214:ILE:CG2	1:A:215:MET:HE2	2.49	0.43
1:A:218:ILE:O	1:A:219:ASP:CB	2.66	0.43
1:B:11:PHE:HD2	1:B:54:VAL:HB	1.82	0.43
1:B:191:VAL:HG23	1:B:229:ILE:HG21	2.00	0.43
1:A:126:ASN:HA	1:A:129:ILE:CG1	2.48	0.43
1:B:379:CYS:HB3	1:B:389:MET:CE	2.49	0.43
1:A:143:TRP:NE1	1:A:144:LEU:HD21	2.34	0.43
1:B:218:ILE:C	1:B:220:GLY:N	2.73	0.43
1:A:98:HIS:HE1	1:A:344:ILE:O	2.02	0.43
1:A:132:ARG:C	1:A:134:GLU:N	2.74	0.43
1:B:122:ARG:CG	1:B:363:GLU:OE2	2.66	0.43
1:A:26:THR:OG1	1:A:28:HIS:HB2	2.19	0.42
1:A:104:LEU:HD11	1:A:218:ILE:CD1	2.46	0.42
1:A:221:GLU:O	1:A:222:SER:O	2.37	0.42
1:A:346:ARG:O	1:A:347:CYS:C	2.62	0.42
1:A:384:GLY:C	1:A:386:VAL:N	2.76	0.42
1:B:11:PHE:CD2	1:B:54:VAL:HB	2.54	0.42
1:B:42:PRO:CB	1:B:60:GLU:HG2	2.49	0.42
1:A:58:TYR:HB3	4:A:519:HOH:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:105:VAL:C	1:A:108:PRO:HD2	2.44	0.42
1:A:139:ASP:O	1:A:143:TRP:HB3	2.18	0.42
1:A:361:ILE:HG23	1:A:362:THR:H	1.84	0.42
1:A:382:LEU:HG	1:A:383:MET:N	2.34	0.42
1:B:25:GLY:HA3	4:B:519:HOH:O	2.19	0.42
1:B:54:VAL:HG13	1:B:311:TRP:CD1	2.55	0.42
1:B:82:GLU:O	1:B:84:GLY:N	2.46	0.42
1:B:108:PRO:O	1:B:114:THR:HG22	2.18	0.42
1:B:228:LEU:O	1:B:230:GLY:N	2.53	0.42
1:A:243:THR:HG21	1:A:356:GLU:OE1	2.19	0.42
1:B:102:ARG:C	1:B:104:LEU:N	2.77	0.42
1:A:126:ASN:O	1:A:129:ILE:HD11	2.19	0.42
1:B:132:ARG:HD2	1:B:135:LEU:HD12	2.00	0.42
1:A:26:THR:C	1:A:28:HIS:H	2.28	0.42
1:A:201:ARG:C	1:A:203:THR:N	2.77	0.42
1:A:225:GLU:O	1:A:229:ILE:HG13	2.19	0.42
1:B:122:ARG:HG3	1:B:363:GLU:OE2	2.19	0.42
1:B:145:ALA:O	1:B:149:PRO:CD	2.67	0.42
1:B:374:ASP:HA	1:B:375:PRO:HD3	1.75	0.42
1:A:73:ASN:C	1:A:75:GLY:H	2.27	0.42
1:B:69:LYS:HD2	1:B:293:GLN:OE1	2.20	0.42
1:A:109:PHE:CD1	1:A:348:LEU:HD11	2.55	0.42
1:A:202:ARG:NH1	1:A:214:ILE:HG23	2.34	0.42
1:B:61:ILE:O	1:B:64:VAL:HB	2.19	0.42
1:A:24:ASP:HB3	1:B:262:ARG:HD3	2.02	0.42
1:A:76:VAL:CG1	1:A:77:THR:N	2.82	0.42
1:B:43:ILE:HB	1:B:300:ILE:HD12	2.00	0.42
1:B:175:THR:O	1:B:176:HIS:HB3	2.20	0.42
1:B:188:ALA:O	1:B:191:VAL:HG12	2.20	0.42
1:B:276:ASP:O	1:B:279:LEU:HB2	2.20	0.42
1:B:144:LEU:O	1:B:145:ALA:C	2.63	0.42
1:B:372:SER:O	1:B:395:ILE:N	2.45	0.42
1:A:174:ILE:HD11	1:A:190:LEU:HD13	2.01	0.41
1:B:145:ALA:O	1:B:149:PRO:HD2	2.19	0.41
1:B:286:MET:HE3	1:B:286:MET:HB2	1.84	0.41
1:A:173:ALA:HB1	1:A:183:GLY:HA2	2.02	0.41
1:A:245:ARG:HD2	1:A:382:LEU:HD22	2.01	0.41
1:A:390:LEU:HG	1:A:391:HIS:CE1	2.55	0.41
1:B:58:TYR:HA	1:B:315:ALA:HB1	2.01	0.41
1:B:79:PRO:O	1:B:80:ARG:O	2.38	0.41
1:B:323:ASP:O	1:B:324:SER:C	2.63	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:170:TRP:CB	1:A:190:LEU:HD12	2.50	0.41
1:A:348:LEU:HD22	1:A:348:LEU:C	2.45	0.41
1:A:371:PHE:O	1:A:372:SER:HB3	2.20	0.41
1:B:53:TRP:CG	1:B:302:MET:HE2	2.56	0.41
1:B:60:GLU:OE1	1:B:60:GLU:N	2.51	0.41
1:B:76:VAL:CG1	1:B:77:THR:N	2.82	0.41
1:A:108:PRO:HG3	1:A:212:ARG:CD	2.49	0.41
1:B:354:ARG:HG2	1:B:354:ARG:NH1	2.36	0.41
1:A:313:PRO:O	1:A:317:ARG:NH1	2.52	0.41
1:B:78:PHE:CD2	1:B:307:THR:HG21	2.56	0.41
1:B:190:LEU:C	1:B:190:LEU:HD23	2.45	0.41
1:A:270:LEU:O	1:A:271:ILE:C	2.64	0.41
1:B:365:LEU:O	1:B:369:PRO:HD3	2.20	0.41
1:A:153:THR:C	1:A:155:ILE:N	2.77	0.41
1:A:236:LEU:O	1:A:240:ILE:HD12	2.21	0.41
1:A:257:ASP:O	1:A:260:LEU:N	2.53	0.41
1:B:73:ASN:C	1:B:75:GLY:H	2.27	0.41
1:A:59:LYS:HG3	1:A:60:GLU:CD	2.46	0.41
1:A:284:PRO:CB	1:A:387:ALA:O	2.68	0.41
1:B:68:THR:CG2	1:B:344:ILE:HD11	2.50	0.41
1:B:304:PRO:O	1:B:306:GLN:N	2.54	0.41
1:A:118:THR:C	1:A:120:GLN:N	2.79	0.41
1:A:159:LEU:HA	1:A:160:PRO:HD3	1.87	0.41
1:A:191:VAL:O	1:A:195:ARG:HG3	2.21	0.41
1:A:203:THR:O	1:A:205:PRO:HD3	2.21	0.41
1:A:273:ASN:HB2	4:A:517:HOH:O	2.20	0.41
1:A:293:GLN:NE2	1:A:294:GLU:O	2.53	0.41
1:B:61:ILE:HD12	1:B:315:ALA:HB2	2.03	0.41
1:B:89:MET:C	1:B:91:ALA:H	2.29	0.41
1:B:286:MET:HE2	1:B:313:PRO:CG	2.51	0.41
1:A:56:ALA:HB1	1:A:314:ILE:CG2	2.50	0.41
1:A:177:VAL:C	1:A:179:ASN:H	2.29	0.41
1:A:367:ARG:C	1:A:369:PRO:HD3	2.46	0.41
1:B:88:LEU:O	1:B:89:MET:CB	2.69	0.41
1:B:192:ALA:O	1:B:195:ARG:HB2	2.21	0.41
1:B:346:ARG:HG2	1:B:346:ARG:HH11	1.86	0.41
1:B:395:ILE:O	1:B:395:ILE:HG13	2.20	0.41
1:A:109:PHE:CE1	2:A:450:HEM:HBC2	2.56	0.40
1:A:132:ARG:HD3	1:A:143:TRP:CD2	2.56	0.40
1:A:401:ARG:NH2	1:A:403:SER:HB3	2.31	0.40
1:B:194:ALA:O	1:B:198:ILE:CG1	2.66	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:312:PHE:N	1:B:313:PRO:HD2	2.36	0.40
1:B:392:VAL:HG23	1:B:392:VAL:O	2.21	0.40
1:A:21:LEU:CD2	1:A:27:PRO:HB3	2.51	0.40
1:A:404:GLU:O	1:A:404:GLU:OE1	2.39	0.40
1:B:197:LEU:C	1:B:197:LEU:HD23	2.46	0.40
1:B:304:PRO:C	1:B:306:GLN:N	2.78	0.40
1:B:119:GLU:C	1:B:121:LEU:H	2.30	0.40
1:B:125:THR:HG23	1:B:148:ILE:CD1	2.47	0.40
1:B:132:ARG:HB2	1:B:143:TRP:CE2	2.57	0.40
1:B:154:ALA:O	1:B:158:GLY:N	2.55	0.40
1:A:254:LEU:O	1:A:261:ARG:NH1	2.55	0.40
1:A:353:ILE:HG21	2:A:450:HEM:CBB	2.52	0.40
1:B:77:THR:HG21	1:B:81:TYR:CD2	2.56	0.40
1:B:261:ARG:O	1:B:262:ARG:C	2.64	0.40
1:A:160:PRO:HB2	1:A:163:ASP:OD1	2.22	0.40
1:B:93:GLN:HG3	1:B:97:VAL:HG12	2.03	0.40
1:B:132:ARG:C	1:B:134:GLU:N	2.77	0.40
1:B:175:THR:C	1:B:177:VAL:H	2.29	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:366:LYS:NZ	4:A:530:HOH:O[1_556]	2.06	0.14

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	395/397 (100%)	313 (79%)	68 (17%)	14 (4%)	3 13
1	B	395/397 (100%)	298 (75%)	69 (18%)	28 (7%)	1 4

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	790/794 (100%)	611 (77%)	137 (17%)	42 (5%)	1 7

All (42) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	80	ARG
1	A	118	THR
1	A	180	PRO
1	A	205	PRO
1	A	222	SER
1	B	80	ARG
1	B	83	THR
1	B	118	THR
1	B	175	THR
1	B	178	GLU
1	B	362	THR
1	B	398	LYS
1	A	119	GLU
1	A	178	GLU
1	A	337	HIS
1	B	67	ASN
1	B	177	VAL
1	B	208	ASP
1	B	273	ASN
1	A	82	GLU
1	A	103	GLN
1	A	321	ALA
1	B	63	ALA
1	B	103	GLN
1	B	111	PRO
1	B	272	PRO
1	B	305	GLY
1	B	361	ILE
1	B	42	PRO
1	B	143	TRP
1	B	236	LEU
1	B	298	GLY
1	B	321	ALA
1	A	139	ASP
1	A	208	ASP
1	B	132	ARG
1	A	116	LEU

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Mol	Chain	Res	Type
1	B	158	GLY
1	B	180	PRO
1	B	339	SER
1	B	179	ASN
1	B	375	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	332/332 (100%)	301 (91%)	31 (9%)	8	28
1	B	332/332 (100%)	307 (92%)	25 (8%)	12	36
All	All	664/664 (100%)	608 (92%)	56 (8%)	10	31

All (56) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	26	THR
1	A	27	PRO
1	A	32	GLU
1	A	73	ASN
1	A	78	PHE
1	A	89	MET
1	A	152	LEU
1	A	161	PRO
1	A	166	THR
1	A	197	LEU
1	A	209	ILE
1	A	210	MET
1	A	212	ARG
1	A	234	ILE
1	A	243	THR
1	A	261	ARG
1	A	265	ILE
1	A	269	GLU

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Mol	Chain	Res	Type
1	A	270	LEU
1	A	273	ASN
1	A	275	VAL
1	A	276	ASP
1	A	293	GLN
1	A	331	GLU
1	A	342	HIS
1	A	346	ARG
1	A	348	LEU
1	A	353	ILE
1	A	379	CYS
1	A	398	LYS
1	A	404	GLU
1	B	42	PRO
1	B	55	VAL
1	B	59	LYS
1	B	64	VAL
1	B	80	ARG
1	B	85	GLU
1	B	93	GLN
1	B	123	GLN
1	B	137	GLU
1	B	152	LEU
1	B	168	ARG
1	B	171	VAL
1	B	203	THR
1	B	235	LEU
1	B	240	ILE
1	B	269	GLU
1	B	292	THR
1	B	297	VAL
1	B	342	HIS
1	B	344	ILE
1	B	347	CYS
1	B	359	VAL
1	B	371	PHE
1	B	398	LYS
1	B	404	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	28	HIS
1	A	66	GLN
1	A	103	GLN
1	A	126	ASN
1	A	146	ASN
1	A	176	HIS
1	A	293	GLN
1	A	335	ASN
1	A	351	HIS
1	A	391	HIS
1	B	62	GLN
1	B	66	GLN
1	B	67	ASN
1	B	179	ASN
1	B	267	HIS
1	B	327	ASN
1	B	337	HIS

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](#)

4 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
2	HEM	A	450	1	50,50,50	1.12	3 (6%)	67,82,82	1.10	5 (7%)
2	HEM	B	450	1	50,50,50	1.21	4 (8%)	67,82,82	1.36	6 (8%)
3	CNL	A	500	-	12,12,12	2.72	7 (58%)	19,20,20	1.25	2 (10%)
3	CNL	B	500	-	12,12,12	2.58	7 (58%)	19,20,20	1.25	2 (10%)

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	450	1	-	6/14/54/54	-
2	HEM	B	450	1	-	10/14/54/54	-
3	CNL	A	500	-	-	-	0/3/2/2
3	CNL	B	500	-	-	-	0/3/2/2

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	500	CNL	C2-C3	3.75	1.60	1.53
3	A	500	CNL	O-C1	3.73	1.51	1.45
3	A	500	CNL	C6-C5	3.67	1.60	1.53
3	B	500	CNL	C2-C3	3.51	1.60	1.53
3	B	500	CNL	O-C1	3.51	1.51	1.45
3	B	500	CNL	C6-C5	3.49	1.60	1.53
2	B	450	HEM	FE-NA	3.26	2.05	1.95
3	A	500	CNL	O-C8	3.24	1.51	1.46
3	A	500	CNL	C5-C4	3.16	1.60	1.53
2	A	450	HEM	CAC-C3C	-3.13	1.39	1.47
3	B	500	CNL	O-C8	2.97	1.51	1.46
3	B	500	CNL	C3-C4	2.93	1.60	1.53
3	B	500	CNL	C5-C4	2.89	1.60	1.53
3	A	500	CNL	C3-C4	2.73	1.59	1.53
2	A	450	HEM	CAB-C3B	-2.70	1.40	1.47
2	B	450	HEM	CAC-C3C	-2.61	1.40	1.47
2	B	450	HEM	C2A-C3A	-2.56	1.32	1.38
2	A	450	HEM	C2A-C3A	-2.45	1.32	1.38
2	B	450	HEM	CAB-C3B	-2.43	1.40	1.47
3	A	500	CNL	C6-C1	2.36	1.60	1.52
3	B	500	CNL	C2-C1	2.20	1.59	1.52

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	450	HEM	CAA-C2A-C1A	5.32	135.32	124.94
2	B	450	HEM	CBA-CAA-C2A	5.17	126.83	112.53
3	B	500	CNL	C10-C8-C9	-4.24	103.26	110.51
3	A	500	CNL	C10-C8-C9	-4.21	103.33	110.51
2	B	450	HEM	CAA-C2A-C3A	-3.98	118.13	127.07
2	A	450	HEM	C3B-C4B-NB	3.27	111.82	109.47
2	A	450	HEM	CAD-C3D-C4D	2.71	129.42	124.70
2	A	450	HEM	C3C-C2C-C1C	-2.61	104.58	107.05
2	B	450	HEM	CAD-C3D-C4D	2.50	129.05	124.70
2	B	450	HEM	CBD-CAD-C3D	2.31	118.92	112.53
2	B	450	HEM	C4A-C3A-C2A	2.19	109.33	106.82
3	A	500	CNL	O-C1-C7	2.18	112.39	105.44
3	B	500	CNL	O-C1-C7	2.11	112.18	105.44
2	A	450	HEM	CBA-CAA-C2A	2.10	118.33	112.53
2	A	450	HEM	CMC-C2C-C1C	2.01	128.27	124.73

There are no chirality outliers.

All (16) torsion outliers are listed below:

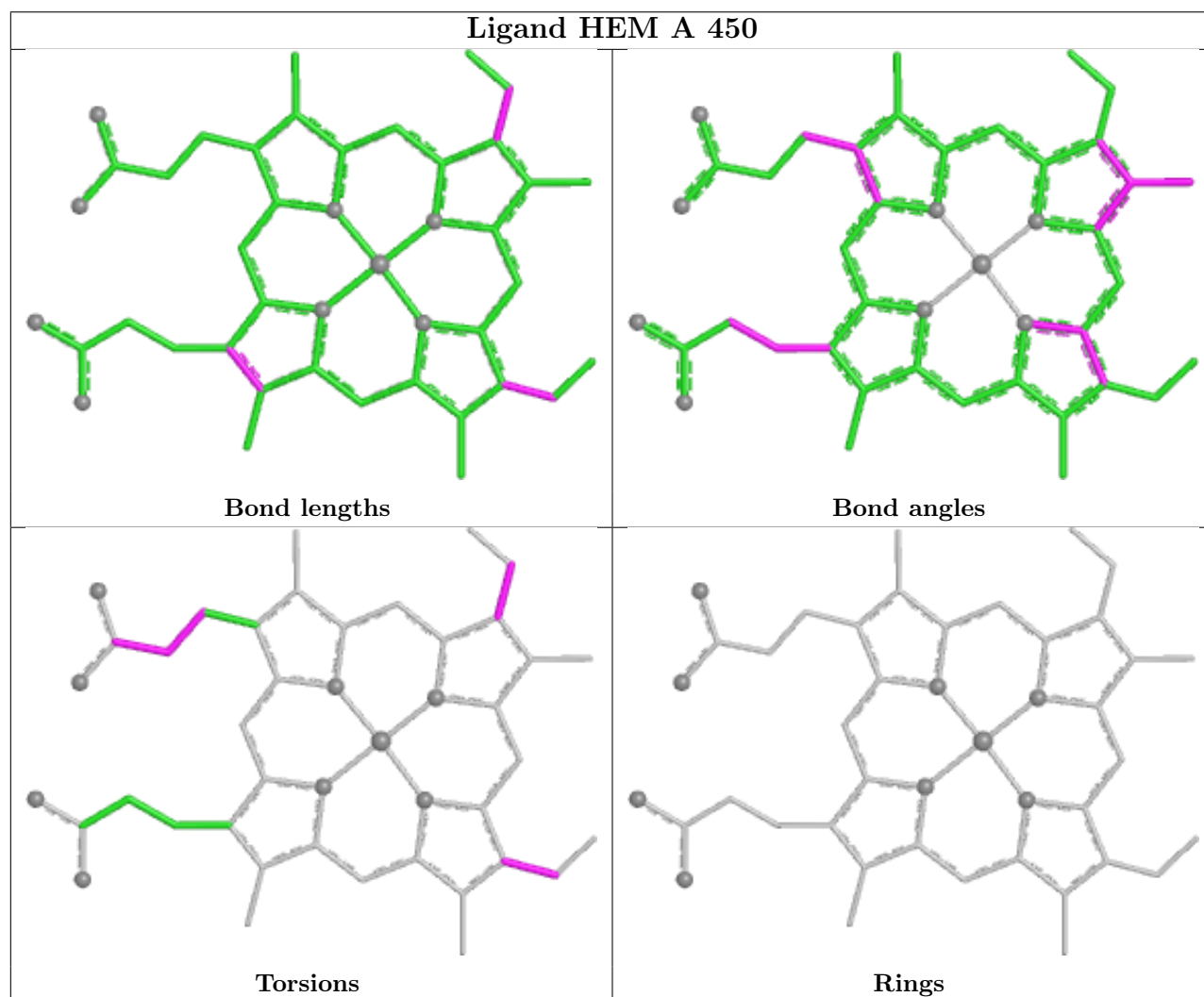
Mol	Chain	Res	Type	Atoms
2	A	450	HEM	C2C-C3C-CAC-CBC
2	A	450	HEM	C4C-C3C-CAC-CBC
2	B	450	HEM	C1A-C2A-CAA-CBA
2	B	450	HEM	C2C-C3C-CAC-CBC
2	B	450	HEM	C4C-C3C-CAC-CBC
2	B	450	HEM	C3A-C2A-CAA-CBA
2	B	450	HEM	C4D-C3D-CAD-CBD
2	B	450	HEM	C2D-C3D-CAD-CBD
2	A	450	HEM	C3D-CAD-CBD-CGD
2	B	450	HEM	C2B-C3B-CAB-CBB
2	A	450	HEM	CAD-CBD-CGD-O2D
2	A	450	HEM	C2B-C3B-CAB-CBB
2	A	450	HEM	CAD-CBD-CGD-O1D
2	B	450	HEM	CAD-CBD-CGD-O2D
2	B	450	HEM	CAD-CBD-CGD-O1D
2	B	450	HEM	CAA-CBA-CGA-O2A

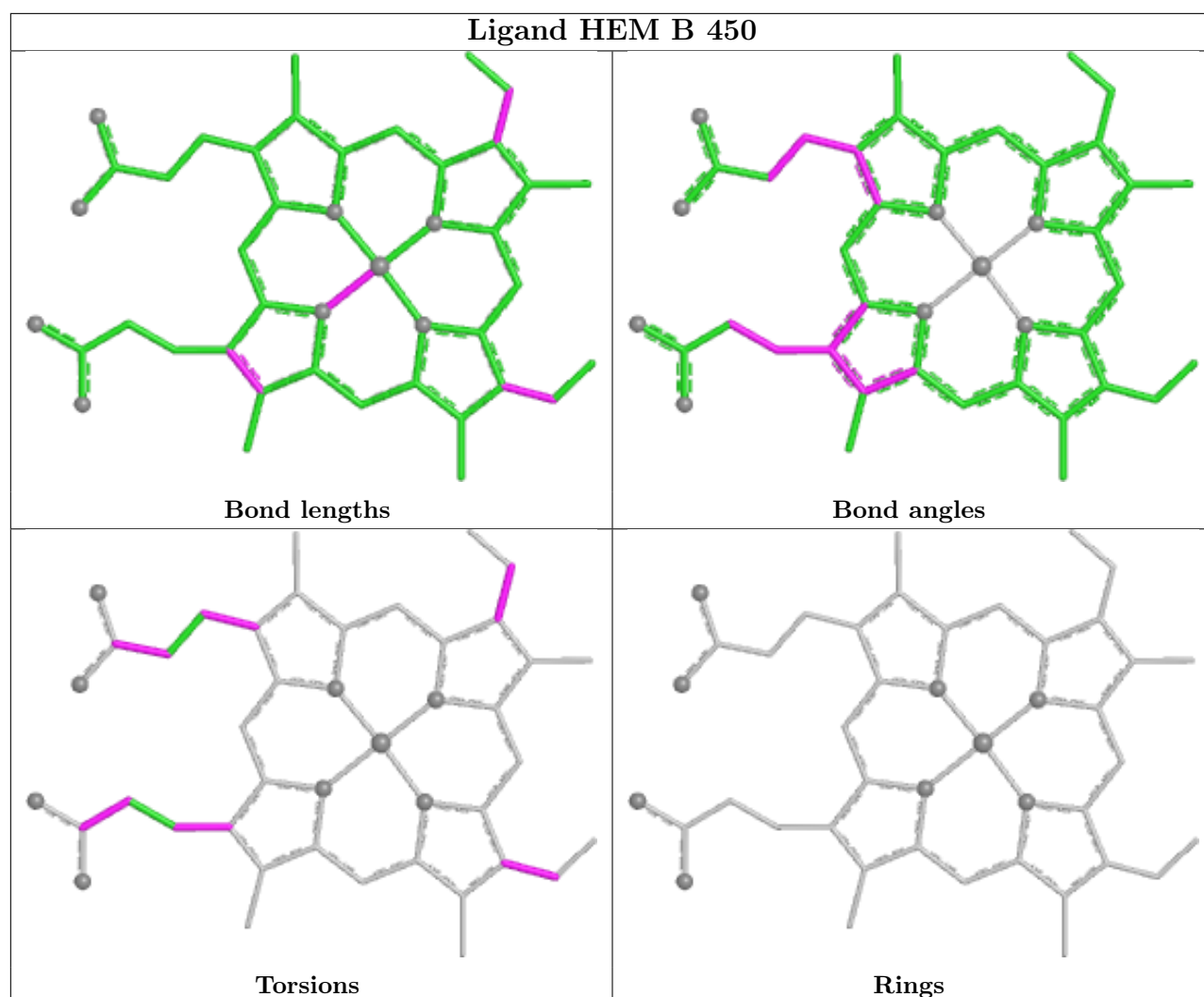
There are no ring outliers.

2 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	450	HEM	7	0
2	B	450	HEM	5	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, 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	397/397 (100%)	0.46	9 (2%) 61 38	25, 59, 90, 120	0
1	B	397/397 (100%)	0.47	11 (2%) 55 32	24, 58, 90, 117	0
All	All	794/794 (100%)	0.46	20 (2%) 58 35	24, 58, 90, 120	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	402	LEU	3.0
1	B	177	VAL	3.0
1	B	176	HIS	2.8
1	B	88	LEU	2.7
1	B	172	TRP	2.7
1	B	113	ALA	2.6
1	A	218	ILE	2.5
1	B	215	MET	2.5
1	A	215	MET	2.5
1	A	172	TRP	2.5
1	B	179	ASN	2.4
1	A	177	VAL	2.4
1	B	180	PRO	2.4
1	A	180	PRO	2.2
1	B	218	ILE	2.1
1	A	176	HIS	2.1
1	A	87	GLU	2.1
1	A	181	GLU	2.1
1	A	240	ILE	2.0
1	B	8	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

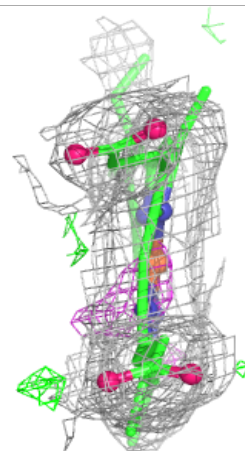
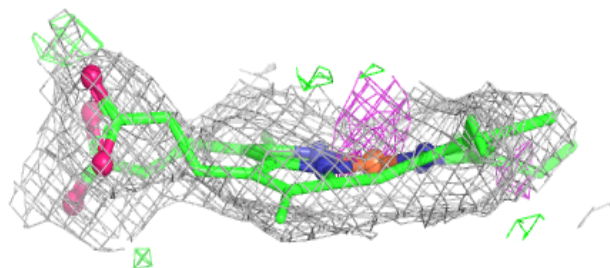
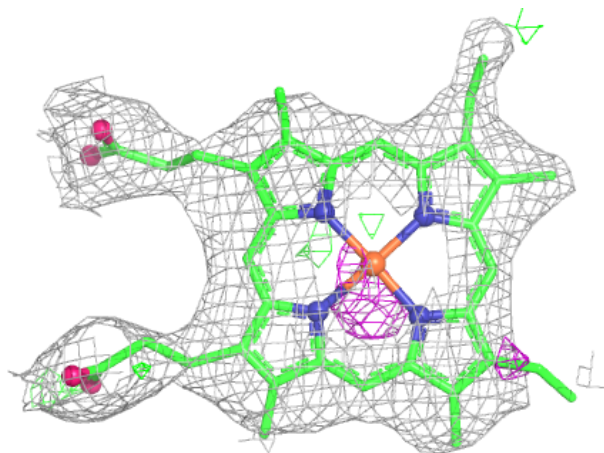
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

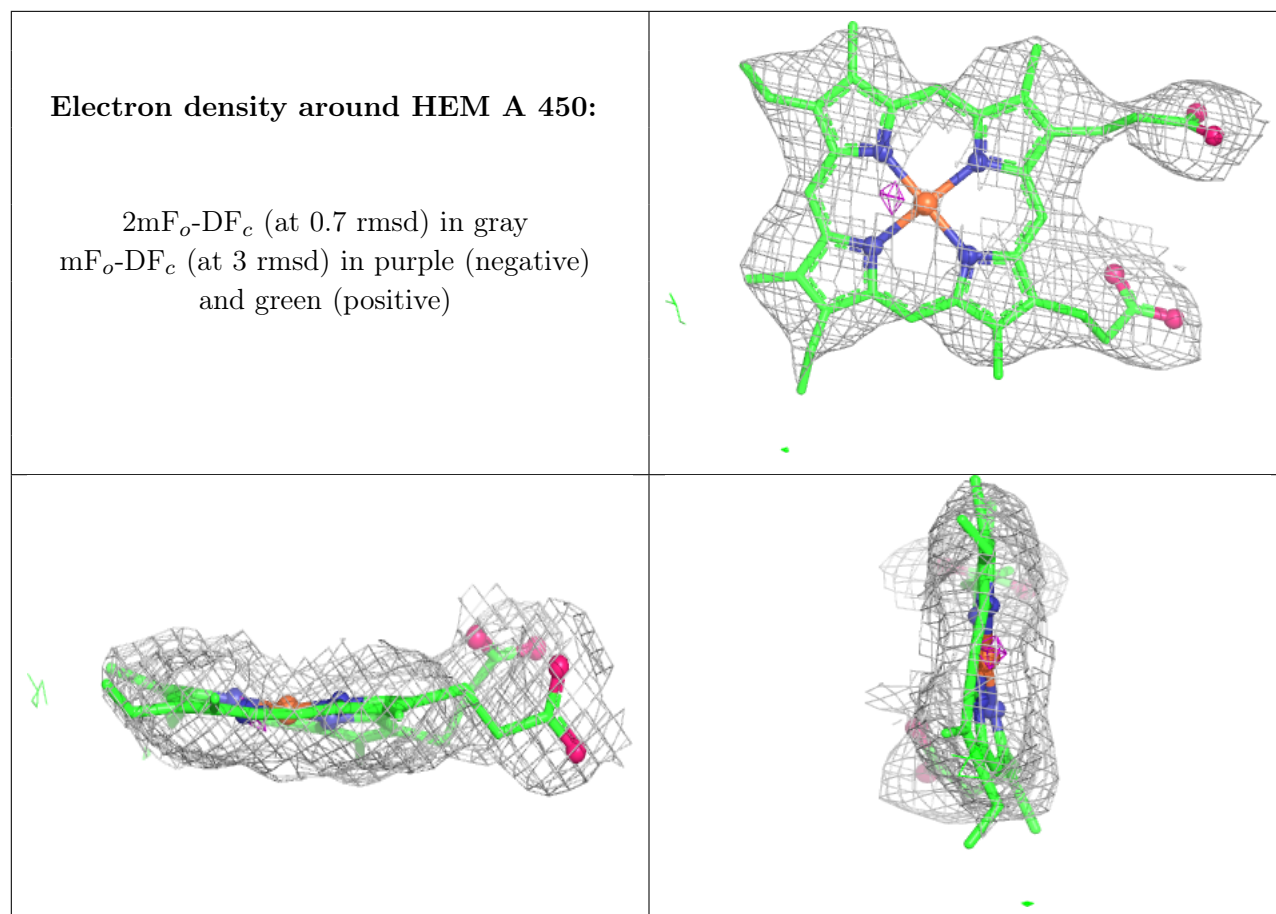
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	CNL	B	500	11/11	0.89	0.22	53,55,56,56	0
3	CNL	A	500	11/11	0.90	0.19	58,59,60,61	0
2	HEM	B	450	43/43	0.93	0.12	37,49,61,63	0
2	HEM	A	450	43/43	0.94	0.12	46,50,56,60	0

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 HEM B 450:

$2mF_o-DF_c$ (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.