



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 10:24 AM UTC

PDB ID : 2NZA / pdb_00002nza
Title : Structure and Function Studies of Cytochrome P450 158A1 from *Streptomyces coelicolor* A3(2)
Authors : Zhao, B.; Waterman, M.R.
Deposited on : 2006-11-22
Resolution : 2.90 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

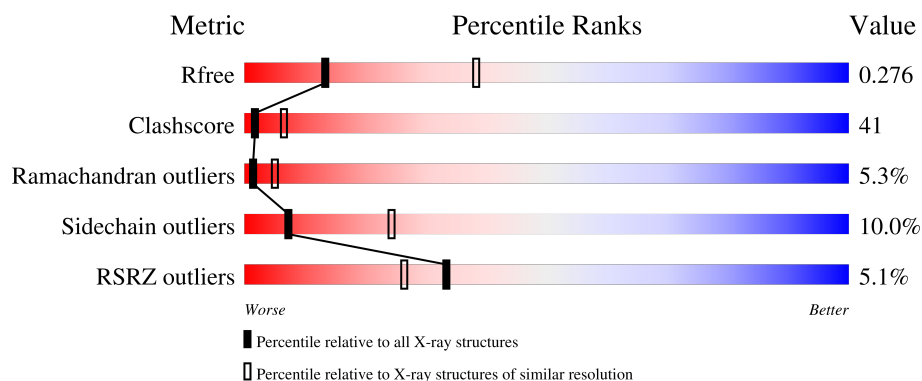
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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	413	
1	B	413	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6371 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 CYP158A1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	397	Total	C	N	O	S	0	0	0
			3065	1920	581	554	10			
1	B	388	Total	C	N	O	S	0	0	0
			2992	1872	566	544	10			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	408	HIS	-	expression tag	UNP Q9KZF5
A	409	HIS	-	expression tag	UNP Q9KZF5
A	410	HIS	-	expression tag	UNP Q9KZF5
A	411	HIS	-	expression tag	UNP Q9KZF5
A	412	HIS	-	expression tag	UNP Q9KZF5
A	413	HIS	-	expression tag	UNP Q9KZF5
B	408	HIS	-	expression tag	UNP Q9KZF5
B	409	HIS	-	expression tag	UNP Q9KZF5
B	410	HIS	-	expression tag	UNP Q9KZF5
B	411	HIS	-	expression tag	UNP Q9KZF5
B	412	HIS	-	expression tag	UNP Q9KZF5
B	413	HIS	-	expression tag	UNP Q9KZF5

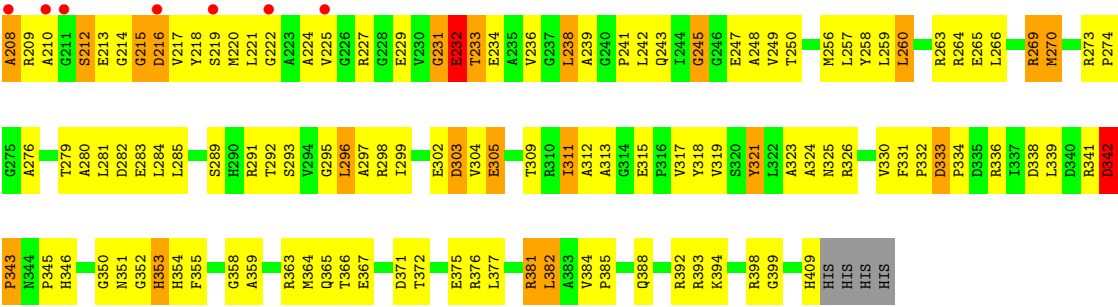
- Molecule 2 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



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

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	117	Total	O	0	0
			117	117		
3	B	111	Total	O	0	0
			111	111		



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	186.61Å 44.35Å 130.62Å 90.00° 98.23° 90.00°	Depositor
Resolution (Å)	10.00 – 2.90 10.00 – 2.90	Depositor EDS
% Data completeness (in resolution range)	85.7 (10.00-2.90) 84.5 (10.00-2.90)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.37 (at 2.81Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.197 , 0.267 0.220 , 0.276	Depositor DCC
R_{free} test set	1144 reflections (4.84%)	wwPDB-VP
Wilson B-factor (Å ²)	20.3	Xtriage
Anisotropy	0.296	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.43 , 66.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	6371	wwPDB-VP
Average B, all atoms (Å ²)	14.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.37% 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

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	0/3136	1.10	18/4271 (0.4%)
1	B	0.48	0/3060	1.10	26/4165 (0.6%)
All	All	0.50	0/6196	1.10	44/8436 (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	142	GLY	CA-C-N	12.77	133.54	120.38
1	B	142	GLY	C-N-CA	12.77	133.54	120.38
1	A	142	GLY	CA-C-N	10.29	130.98	120.38
1	A	142	GLY	C-N-CA	10.29	130.98	120.38
1	A	180	GLN	N-CA-C	-9.86	99.99	112.90
1	B	143	PRO	N-CA-C	8.78	121.41	110.70
1	A	234	GLU	N-CA-C	-8.55	103.29	114.31
1	A	143	PRO	N-CA-C	8.40	120.95	110.70
1	A	183	SER	N-CA-C	-8.35	102.12	111.14
1	B	208	ALA	N-CA-C	-7.84	100.42	110.53
1	B	162	GLU	N-CA-C	-7.44	103.11	111.14
1	B	23	ASP	N-CA-C	-7.27	104.35	112.57
1	A	240	GLY	CA-C-N	6.92	126.73	119.05
1	A	240	GLY	C-N-CA	6.92	126.73	119.05
1	A	287	TRP	N-CA-C	6.79	118.68	111.28
1	B	181	ILE	N-CA-C	-6.70	103.90	113.07
1	B	342	ASP	CA-C-N	-6.69	111.47	119.84
1	B	342	ASP	C-N-CA	-6.69	111.47	119.84
1	B	143	PRO	CA-C-N	-6.67	111.50	119.84
1	B	143	PRO	C-N-CA	-6.67	111.50	119.84
1	A	342	ASP	CA-C-N	-6.56	112.94	119.56
1	A	342	ASP	C-N-CA	-6.56	112.94	119.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	13	PRO	N-CA-CB	6.19	109.81	103.00
1	A	208	ALA	N-CA-C	-6.19	97.62	110.80
1	A	13	PRO	N-CA-CB	6.14	109.76	103.00
1	B	112	ALA	N-CA-C	6.12	119.84	112.38
1	B	382	LEU	N-CA-C	-6.08	100.59	110.20
1	B	200	GLY	N-CA-C	-5.98	105.56	112.50
1	B	321	TYR	N-CA-C	-5.65	105.27	111.82
1	A	333	ASP	CA-C-N	5.64	125.75	119.32
1	A	333	ASP	C-N-CA	5.64	125.75	119.32
1	A	33	LEU	N-CA-C	-5.61	105.08	111.14
1	B	189	GLU	N-CA-C	-5.58	106.51	113.38
1	B	69	ASP	CA-C-N	5.57	126.81	119.84
1	B	69	ASP	C-N-CA	5.57	126.81	119.84
1	B	180	GLN	N-CA-C	-5.41	103.48	111.81
1	B	333	ASP	CA-C-N	5.39	126.57	119.84
1	B	333	ASP	C-N-CA	5.39	126.57	119.84
1	B	215	GLY	N-CA-C	-5.31	106.35	112.73
1	B	30	ASP	CA-C-N	5.19	124.81	119.05
1	B	30	ASP	C-N-CA	5.19	124.81	119.05
1	A	215	GLY	N-CA-C	-5.19	108.64	114.40
1	B	207	ARG	N-CA-C	-5.13	105.58	111.07
1	A	189	GLU	N-CA-C	-5.03	106.83	113.17

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	3065	0	3047	256	0
1	B	2992	0	2969	244	0
2	A	43	0	30	2	0
2	B	43	0	30	2	0
3	A	117	0	0	21	0
3	B	111	0	0	19	0
All	All	6371	0	6076	498	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 41.

All (498) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:289:SER:HB2	1:B:398:ARG:HH21	1.20	1.06
1:A:283:GLU:HG2	1:A:339:LEU:H	1.27	0.99
1:A:182:ILE:HD13	1:A:186:GLY:H	1.28	0.98
1:B:304:VAL:HG12	1:B:305:GLU:H	1.28	0.95
1:B:64:LYS:HD2	1:B:351:ASN:OD1	1.70	0.91
1:A:165:GLY:HA3	1:A:216:ASP:OD1	1.70	0.91
1:B:58:THR:HG23	1:B:323:ALA:HB1	1.53	0.91
1:A:124:ARG:HB3	1:A:125:PRO:HD3	1.52	0.90
1:B:209:ARG:HB3	1:B:219:SER:HA	1.54	0.89
1:A:33:LEU:HD21	1:A:323:ALA:HB2	1.55	0.88
1:A:85:LEU:HD22	1:A:293:SER:HB3	1.56	0.88
1:A:208:ALA:O	1:A:210:ALA:N	2.06	0.87
1:B:33:LEU:HD21	1:B:323:ALA:HB2	1.59	0.84
1:A:392:ARG:HG3	1:A:399:GLY:O	1.78	0.83
1:B:289:SER:CB	1:B:398:ARG:HH21	1.91	0.83
1:A:85:LEU:HD23	1:A:396:MET:HE2	1.61	0.82
1:B:37:MET:HG2	1:B:58:THR:HG21	1.60	0.81
1:A:381:ARG:HD3	3:A:435:HOH:O	1.80	0.81
1:B:289:SER:HB2	1:B:398:ARG:NH2	1.95	0.81
1:B:351:ASN:ND2	1:B:352:GLY:H	1.79	0.80
1:A:193:ARG:HG2	1:A:196:ARG:NH2	1.95	0.80
1:B:143:PRO:HB2	1:B:144:PRO:HD3	1.60	0.80
1:B:175:HIS:O	1:B:179:ARG:HG3	1.82	0.79
1:A:93:PRO:HB2	1:A:233:THR:O	1.82	0.79
1:B:63:VAL:HG13	1:B:319:VAL:HB	1.64	0.79
1:B:22:LEU:HG	1:B:24:LEU:HD12	1.65	0.79
1:B:202:ILE:O	1:B:206:VAL:HG23	1.83	0.78
1:A:80:ARG:NH1	1:A:314:GLY:O	2.17	0.78
1:B:14:PRO:HA	3:B:452:HOH:O	1.83	0.77
1:A:181:ILE:HD12	1:A:181:ILE:H	1.48	0.76
1:A:41:PRO:HB3	1:A:59:ARG:NH1	2.00	0.76
1:A:291:ARG:NH2	1:A:396:MET:HG3	2.00	0.76
1:A:304:VAL:O	1:A:311:ILE:HG22	1.86	0.76
1:B:342:ASP:HB2	1:B:343:PRO:HD3	1.66	0.75
1:B:342:ASP:CB	1:B:343:PRO:HD3	2.17	0.75
2:B:430:HEM:HHA	2:B:430:HEM:HBA2	1.70	0.74
1:A:338:ASP:HB3	3:A:445:HOH:O	1.86	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:41:PRO:HA	1:B:59:ARG:HD3	1.70	0.74
1:B:363:ARG:O	1:B:367:GLU:HG3	1.87	0.73
1:B:304:VAL:HG12	1:B:305:GLU:N	2.02	0.73
1:A:342:ASP:OD2	1:A:343:PRO:HD2	1.88	0.73
1:B:182:ILE:HG23	1:B:186:GLY:O	1.89	0.72
1:A:193:ARG:HG2	1:A:196:ARG:HH21	1.51	0.72
1:A:62:ASP:O	1:A:66:ILE:HG13	1.88	0.72
1:B:63:VAL:O	1:B:67:THR:HG22	1.88	0.72
1:A:210:ALA:HA	1:A:222:GLY:O	1.89	0.71
1:A:182:ILE:HD13	1:A:186:GLY:N	2.04	0.71
1:B:351:ASN:HD22	1:B:352:GLY:H	1.37	0.71
1:A:255:GLN:NE2	1:A:288:ILE:HG23	2.06	0.71
1:B:108:ARG:HG2	1:B:108:ARG:HH11	1.55	0.71
1:B:342:ASP:CG	1:B:343:PRO:HD3	2.16	0.70
1:B:37:MET:HG2	1:B:58:THR:CG2	2.21	0.70
1:A:179:ARG:N	1:A:179:ARG:HD2	2.06	0.70
1:A:291:ARG:HH22	1:A:396:MET:HG3	1.57	0.70
1:B:303:ASP:HB2	1:B:313:ALA:HB2	1.73	0.69
1:B:342:ASP:OD2	1:B:343:PRO:HD3	1.91	0.69
1:A:67:THR:HG22	1:A:298:ARG:NH2	2.08	0.69
1:A:179:ARG:HH11	1:A:179:ARG:HA	1.58	0.69
1:B:205:THR:HG23	1:B:209:ARG:NH1	2.08	0.69
1:B:205:THR:HG23	1:B:209:ARG:CZ	2.23	0.68
1:B:209:ARG:HA	1:B:212:SER:HB2	1.75	0.68
1:A:67:THR:HG22	1:A:298:ARG:CZ	2.24	0.68
1:A:98:PHE:CE2	1:A:296:LEU:HD23	2.29	0.67
1:A:47:LEU:HD12	1:A:83:THR:O	1.95	0.67
1:A:176:SER:O	1:A:179:ARG:HD3	1.94	0.67
1:A:82:ILE:HD12	1:A:82:ILE:H	1.60	0.67
1:A:182:ILE:HG22	1:A:183:SER:H	1.59	0.67
1:A:207:ARG:HG3	1:A:207:ARG:HH11	1.58	0.67
1:A:181:ILE:HD12	1:A:181:ILE:N	2.10	0.66
1:B:91:PRO:HB3	1:B:98:PHE:CE1	2.29	0.66
1:A:37:MET:HG2	1:A:58:THR:CG2	2.25	0.66
1:A:207:ARG:HG3	1:A:207:ARG:NH1	2.10	0.66
1:A:286:ARG:HH21	1:A:341:ARG:HH21	1.42	0.66
1:B:165:GLY:HA3	1:B:216:ASP:OD2	1.96	0.66
1:B:138:ILE:HG13	1:B:139:LEU:N	2.10	0.66
1:A:230:VAL:O	1:A:231:GLY:O	2.14	0.66
1:A:385:PRO:HD2	1:A:388:GLN:HG3	1.78	0.66
1:B:33:LEU:HD23	1:B:33:LEU:O	1.96	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:269:ARG:CG	1:B:269:ARG:HH11	2.09	0.65
1:B:23:ASP:HA	3:B:432:HOH:O	1.97	0.65
1:B:155:PHE:HB3	1:B:156:PRO:HD3	1.77	0.65
1:A:37:MET:HG2	1:A:58:THR:HG21	1.77	0.65
1:A:263:ARG:HB3	1:A:265:GLU:OE2	1.96	0.65
1:A:363:ARG:O	1:A:367:GLU:HG3	1.97	0.65
1:A:225:VAL:HG13	1:A:231:GLY:HA2	1.78	0.65
1:B:351:ASN:ND2	1:B:352:GLY:N	2.45	0.65
1:A:41:PRO:HB3	1:A:59:ARG:HH12	1.62	0.64
1:B:91:PRO:HB3	1:B:98:PHE:CD1	2.32	0.64
1:A:342:ASP:HB2	1:A:343:PRO:CD	2.27	0.64
1:B:67:THR:HB	3:B:457:HOH:O	1.96	0.64
1:A:135:VAL:HB	1:A:376:ARG:NH1	2.12	0.64
1:B:207:ARG:O	1:B:210:ALA:HB2	1.97	0.64
1:A:221:LEU:O	1:A:225:VAL:HG23	1.98	0.64
1:B:58:THR:HG23	1:B:323:ALA:CB	2.26	0.64
1:B:249:VAL:HG11	1:B:365:GLN:HE21	1.62	0.64
1:B:321:TYR:CZ	1:B:350:GLY:HA2	2.33	0.63
1:B:33:LEU:O	1:B:37:MET:HG3	1.97	0.63
1:B:76:GLU:HB3	1:B:299:ILE:HG13	1.80	0.63
1:A:144:PRO:O	1:B:196:ARG:NH1	2.32	0.63
1:A:249:VAL:HG21	1:A:365:GLN:NE2	2.13	0.63
1:A:182:ILE:CD1	1:A:185:SER:HA	2.29	0.62
1:A:342:ASP:HB2	1:A:343:PRO:HD3	1.80	0.62
1:B:47:LEU:HB3	3:B:509:HOH:O	1.99	0.62
1:A:133:GLY:O	1:A:136:ASP:HB2	2.00	0.62
1:A:286:ARG:NH2	1:A:341:ARG:NH2	2.48	0.62
1:B:297:ALA:HB2	1:B:318:TYR:CE2	2.35	0.62
1:B:342:ASP:CG	1:B:343:PRO:CD	2.73	0.62
1:A:115:PHE:HB3	3:A:432:HOH:O	2.00	0.62
1:B:124:ARG:HB3	1:B:125:PRO:HD3	1.81	0.62
1:B:375:GLU:HG3	1:B:376:ARG:HG2	1.82	0.62
1:B:44:ARG:HD2	1:B:53:TRP:CZ3	2.36	0.61
1:B:138:ILE:HD11	1:B:377:LEU:HD21	1.82	0.61
1:A:387:GLU:CD	1:A:387:GLU:H	2.08	0.61
1:B:74:ARG:HG2	1:B:97:ALA:O	1.99	0.61
1:B:394:LYS:HE3	3:B:494:HOH:O	2.00	0.61
1:A:231:GLY:O	1:A:232:GLU:O	2.18	0.61
1:A:303:ASP:HA	1:A:311:ILE:O	2.00	0.61
1:A:180:GLN:CA	1:A:180:GLN:HE21	2.14	0.61
1:B:33:LEU:HD23	1:B:37:MET:HG3	1.83	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:159:VAL:O	1:B:163:VAL:HG23	2.01	0.60
1:A:342:ASP:CB	1:A:343:PRO:CD	2.79	0.60
1:B:105:ASN:O	1:B:106:ARG:C	2.43	0.60
1:A:270:MET:HE1	1:A:280:ALA:HB1	1.83	0.60
1:B:375:GLU:HG3	1:B:376:ARG:N	2.16	0.60
1:A:211:GLY:H	1:A:223:ALA:HA	1.67	0.60
1:A:178:THR:C	1:A:179:ARG:HD2	2.26	0.59
1:A:286:ARG:HH21	1:A:341:ARG:NH2	2.00	0.59
1:A:182:ILE:HD13	1:A:185:SER:HA	1.83	0.59
1:A:208:ALA:C	1:A:210:ALA:H	2.11	0.59
1:A:207:ARG:HH12	1:A:232:GLU:CD	2.10	0.59
1:B:22:LEU:CG	1:B:24:LEU:HD12	2.32	0.59
1:B:153:GLU:HA	1:B:250:THR:CG2	2.32	0.59
1:B:260:LEU:HD21	1:B:270:MET:HE2	1.83	0.59
1:A:224:ALA:CA	1:A:227:ARG:HH21	2.16	0.59
1:B:79:GLN:O	1:B:80:ARG:HG2	2.02	0.58
1:A:180:GLN:HE21	1:A:180:GLN:HA	1.67	0.58
1:A:59:ARG:O	1:A:63:VAL:HG23	2.02	0.58
1:A:177:TRP:C	1:A:179:ARG:H	2.10	0.58
1:B:259:LEU:O	1:B:263:ARG:HB2	2.04	0.58
1:A:271:ARG:HD3	3:A:440:HOH:O	2.03	0.58
1:A:300:ALA:HB1	3:A:538:HOH:O	2.04	0.58
1:B:342:ASP:OD2	1:B:343:PRO:CD	2.52	0.58
1:B:62:ASP:O	1:B:66:ILE:HG13	2.03	0.58
1:A:249:VAL:O	1:A:253:VAL:HG23	2.03	0.58
1:B:64:LYS:O	1:B:64:LYS:HG3	2.04	0.57
1:A:131:LEU:O	1:A:135:VAL:HG23	2.04	0.57
1:B:153:GLU:N	1:B:154:PRO:HD2	2.20	0.57
1:A:63:VAL:O	1:A:67:THR:OG1	2.22	0.57
1:B:304:VAL:CG1	1:B:305:GLU:H	2.12	0.57
1:B:208:ALA:O	1:B:209:ARG:HB2	2.02	0.57
1:B:182:ILE:HG22	1:B:185:SER:C	2.30	0.57
1:B:203:THR:O	1:B:207:ARG:HD3	2.05	0.57
1:A:181:ILE:C	1:A:182:ILE:HG13	2.29	0.56
1:A:232:GLU:C	1:A:234:GLU:H	2.12	0.56
1:A:74:ARG:HG3	1:A:74:ARG:O	2.05	0.56
1:A:155:PHE:CD2	1:A:369:LEU:HD21	2.40	0.56
1:A:228:GLY:HA2	3:A:461:HOH:O	2.05	0.56
1:B:241:PRO:HG2	3:B:522:HOH:O	2.06	0.56
1:B:311:ILE:HG22	1:B:311:ILE:O	2.06	0.56
1:A:44:ARG:HG2	1:A:53:TRP:CZ3	2.41	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:165:GLY:HA3	1:B:216:ASP:CG	2.30	0.56
1:B:22:LEU:HD21	1:B:24:LEU:HB2	1.88	0.56
1:B:149:GLU:O	1:B:149:GLU:HG2	2.06	0.56
1:B:183:SER:C	1:B:185:SER:H	2.12	0.56
1:B:295:GLY:HA3	3:B:513:HOH:O	2.06	0.55
1:B:338:ASP:O	1:B:339:LEU:HB2	2.06	0.55
1:A:138:ILE:HD11	1:A:377:LEU:HD11	1.88	0.55
1:A:167:PRO:O	1:A:171:ARG:HG2	2.05	0.55
1:A:181:ILE:H	1:A:181:ILE:CD1	2.15	0.55
1:A:385:PRO:HD2	1:A:388:GLN:CG	2.36	0.55
1:B:168:ALA:O	1:B:171:ARG:HG2	2.06	0.55
1:B:392:ARG:HG3	1:B:399:GLY:O	2.07	0.55
1:A:28:GLU:N	1:A:28:GLU:OE1	2.40	0.55
1:A:171:ARG:HG3	1:A:171:ARG:HH11	1.72	0.55
1:A:60:TYR:CD2	1:A:330:VAL:HG21	2.42	0.55
1:B:269:ARG:HH11	1:B:269:ARG:HG2	1.70	0.55
1:B:58:THR:CG2	1:B:323:ALA:HB1	2.34	0.55
1:A:76:GLU:OE1	1:A:80:ARG:HD2	2.07	0.54
1:A:249:VAL:HG21	1:A:365:GLN:HE21	1.71	0.54
1:A:283:GLU:HG2	1:A:339:LEU:N	2.10	0.54
1:A:252:ASN:HD21	1:A:256:MET:HE3	1.73	0.54
1:A:286:ARG:NH2	1:A:341:ARG:HH21	2.02	0.54
1:A:108:ARG:HH11	1:A:108:ARG:HG2	1.73	0.54
1:A:124:ARG:HB3	1:A:125:PRO:CD	2.33	0.54
1:B:342:ASP:CB	1:B:343:PRO:CD	2.85	0.54
1:A:60:TYR:CG	1:A:330:VAL:HG21	2.43	0.53
1:A:209:ARG:HD3	1:A:218:TYR:C	2.33	0.53
1:B:351:ASN:HD22	1:B:352:GLY:N	2.06	0.53
1:B:232:GLU:O	1:B:234:GLU:N	2.41	0.53
1:A:37:MET:HE3	1:A:323:ALA:HB1	1.89	0.53
1:A:99:ALA:HB1	1:A:104:HIS:HA	1.90	0.53
1:B:153:GLU:OE2	1:B:157:ILE:HD11	2.07	0.53
1:B:152:LEU:HD11	1:B:257:LEU:HD12	1.89	0.53
1:A:99:ALA:HB3	1:A:104:HIS:CD2	2.44	0.53
1:B:269:ARG:HH21	1:B:273:ARG:CZ	2.21	0.53
1:B:59:ARG:O	1:B:63:VAL:HG23	2.08	0.53
1:A:143:PRO:HB2	1:A:144:PRO:HD3	1.90	0.52
1:B:257:LEU:O	1:B:258:TYR:C	2.51	0.52
1:A:126:ARG:O	1:A:130:ILE:HG13	2.10	0.52
1:A:205:THR:HG22	1:A:209:ARG:HD2	1.91	0.52
1:B:213:GLU:CD	1:B:214:GLY:N	2.68	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:372:THR:O	1:B:375:GLU:HG2	2.09	0.52
1:A:77:VAL:HG22	1:A:299:ILE:HG12	1.91	0.52
1:A:91:PRO:HB3	1:A:98:PHE:CE1	2.44	0.52
1:A:100:ASP:O	1:A:103:ASP:HB2	2.09	0.52
1:B:205:THR:HG22	1:B:218:TYR:HB3	1.90	0.52
1:B:242:LEU:O	1:B:243:GLN:C	2.53	0.52
1:B:376:ARG:HD2	3:B:526:HOH:O	2.08	0.52
1:B:393:ARG:NH2	3:B:448:HOH:O	2.42	0.52
1:A:83:THR:O	1:A:83:THR:HG23	2.08	0.52
1:B:175:HIS:O	1:B:178:THR:HG22	2.09	0.52
1:B:311:ILE:O	1:B:311:ILE:CG2	2.58	0.52
1:A:176:SER:C	1:A:179:ARG:HD3	2.35	0.52
1:A:192:GLU:OE2	1:A:195:LYS:HE2	2.10	0.52
1:A:126:ARG:O	1:A:126:ARG:HD2	2.09	0.52
1:B:33:LEU:HD23	1:B:33:LEU:C	2.35	0.52
1:B:42:LEU:HD23	1:B:309:THR:HG21	1.91	0.52
1:A:286:ARG:HG3	1:A:325:ASN:HB3	1.91	0.52
1:A:144:PRO:HB2	1:B:196:ARG:NH1	2.25	0.52
1:A:192:GLU:O	1:A:196:ARG:HG3	2.10	0.51
1:A:77:VAL:CG2	1:A:299:ILE:HG12	2.40	0.51
1:A:149:GLU:O	1:A:149:GLU:HG2	2.11	0.51
1:B:117:VAL:O	1:B:121:LYS:HG3	2.10	0.51
1:B:172:GLU:O	1:B:176:SER:HB2	2.10	0.51
1:B:274:PRO:C	1:B:276:ALA:H	2.17	0.51
1:A:105:ASN:ND2	3:A:460:HOH:O	2.42	0.51
1:A:152:LEU:O	1:A:250:THR:HG23	2.10	0.51
1:A:182:ILE:HG21	1:A:186:GLY:N	2.25	0.51
1:A:232:GLU:C	1:A:233:THR:HG23	2.35	0.51
1:A:207:ARG:C	1:A:208:ALA:O	2.53	0.51
1:B:78:THR:HG23	1:B:91:PRO:HG2	1.92	0.51
1:B:231:GLY:O	1:B:232:GLU:O	2.29	0.51
1:A:69:ASP:HB3	1:A:72:PHE:HD1	1.76	0.51
1:B:201:TRP:HZ3	1:B:218:TYR:CE1	2.29	0.51
1:A:16:VAL:HG11	1:A:46:ARG:NH1	2.26	0.51
1:A:187:GLY:N	1:A:190:ALA:HB3	2.26	0.51
1:B:77:VAL:HG22	1:B:299:ILE:HD11	1.93	0.51
1:B:215:GLY:O	1:B:220:MET:HE2	2.10	0.51
1:A:206:VAL:HG13	1:A:222:GLY:HA2	1.92	0.50
1:B:192:GLU:O	1:B:196:ARG:HG3	2.12	0.50
1:B:256:MET:HE1	1:B:281:LEU:HD22	1.92	0.50
1:A:187:GLY:O	1:A:190:ALA:HB3	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:291:ARG:HG3	1:B:293:SER:O	2.11	0.50
1:B:302:GLU:O	1:B:303:ASP:HB3	2.11	0.50
1:A:178:THR:HG22	1:A:178:THR:O	2.11	0.50
1:A:85:LEU:HD22	1:A:293:SER:CB	2.36	0.50
1:A:37:MET:HE1	1:A:326:ARG:HB2	1.92	0.50
1:B:303:ASP:CB	1:B:313:ALA:HB2	2.42	0.50
1:B:331:PHE:HE2	1:B:345:PRO:HD2	1.77	0.50
1:A:283:GLU:CG	1:A:339:LEU:H	2.10	0.50
1:A:363:ARG:HD3	3:A:450:HOH:O	2.12	0.50
1:A:177:TRP:O	1:A:180:GLN:OE1	2.30	0.50
1:A:195:LYS:HE3	3:A:491:HOH:O	2.12	0.50
1:B:330:VAL:O	1:B:332:PRO:HD3	2.11	0.50
1:A:28:GLU:O	1:A:398:ARG:NH2	2.42	0.49
1:A:224:ALA:N	1:A:227:ARG:HH21	2.10	0.49
1:B:224:ALA:HB1	1:B:229:GLU:HB2	1.94	0.49
1:A:63:VAL:HG13	1:A:319:VAL:HB	1.95	0.49
1:A:305:GLU:O	1:A:305:GLU:HG3	2.12	0.49
1:A:328:PRO:HG3	1:A:334:PRO:HG3	1.94	0.49
1:B:203:THR:OG1	1:B:236:VAL:HG21	2.12	0.49
1:B:354:HIS:O	2:B:430:HEM:HBA1	2.11	0.49
1:B:21:ALA:O	1:B:22:LEU:HB2	2.12	0.49
1:B:120:THR:HG23	1:B:364:MET:HE3	1.95	0.49
1:B:168:ALA:HA	1:B:171:ARG:HE	1.78	0.49
1:A:224:ALA:HA	1:A:227:ARG:HH21	1.78	0.49
1:B:22:LEU:CD2	1:B:24:LEU:HB2	2.42	0.49
1:B:225:VAL:HB	3:B:466:HOH:O	2.12	0.49
1:A:14:PRO:HB3	3:A:525:HOH:O	2.12	0.49
1:A:67:THR:HG21	1:A:321:TYR:OH	2.12	0.49
1:A:199:TYR:HB3	1:A:236:VAL:HG11	1.94	0.49
1:A:252:ASN:ND2	1:A:256:MET:HE3	2.27	0.49
1:A:123:LEU:O	1:A:124:ARG:C	2.55	0.49
1:A:225:VAL:HG13	1:A:232:GLU:N	2.28	0.49
1:B:63:VAL:O	1:B:67:THR:CG2	2.60	0.49
1:B:105:ASN:O	1:B:107:LEU:N	2.46	0.49
1:B:108:ARG:HG2	1:B:108:ARG:NH1	2.23	0.49
1:A:77:VAL:HG12	1:A:77:VAL:O	2.12	0.49
1:A:349:TYR:CE2	2:A:430:HEM:HBB2	2.48	0.49
1:A:273:ARG:NH1	3:A:446:HOH:O	2.27	0.48
1:A:283:GLU:HG2	1:A:338:ASP:H	1.78	0.48
1:B:283:GLU:HG2	1:B:339:LEU:H	1.77	0.48
1:A:192:GLU:HG3	1:A:196:ARG:NE	2.28	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:109:ARG:HD2	1:B:109:ARG:O	2.13	0.48
1:B:215:GLY:O	1:B:216:ASP:C	2.55	0.48
1:B:333:ASP:N	1:B:334:PRO:CD	2.76	0.48
1:A:302:GLU:HG3	1:A:302:GLU:O	2.12	0.48
1:A:341:ARG:O	1:A:342:ASP:O	2.32	0.48
1:B:29:PHE:CE2	1:B:33:LEU:HD22	2.47	0.48
1:B:67:THR:HG21	1:B:351:ASN:OD1	2.13	0.48
1:A:193:ARG:CG	1:A:196:ARG:HH21	2.24	0.48
1:A:214:GLY:C	1:A:216:ASP:H	2.19	0.48
1:A:225:VAL:HG13	1:A:232:GLU:H	1.77	0.48
1:A:292:THR:HG22	1:A:398:ARG:HG2	1.95	0.48
1:A:198:LEU:HD21	1:A:243:GLN:HG2	1.96	0.48
1:B:109:ARG:HD2	1:B:109:ARG:C	2.39	0.48
1:A:20:PRO:HB2	1:A:22:LEU:HD23	1.96	0.48
1:A:180:GLN:CA	1:A:180:GLN:NE2	2.76	0.48
1:B:199:TYR:HB3	1:B:236:VAL:HG11	1.95	0.48
1:A:256:MET:HB3	1:A:284:LEU:HD13	1.95	0.48
1:B:229:GLU:HA	3:B:502:HOH:O	2.13	0.48
1:B:170:ASP:CG	1:B:173:ARG:HH11	2.23	0.47
1:B:203:THR:HG23	1:B:232:GLU:OE2	2.15	0.47
1:B:312:ALA:HB3	1:B:315:GLU:HG3	1.97	0.47
1:A:92:ARG:HG2	1:A:93:PRO:HD2	1.95	0.47
1:A:203:THR:OG1	1:A:236:VAL:HG21	2.15	0.47
1:B:198:LEU:O	1:B:202:ILE:HG13	2.13	0.47
1:A:99:ALA:HB1	1:A:104:HIS:CA	2.44	0.47
1:A:33:LEU:HD23	1:A:37:MET:HG3	1.95	0.47
1:A:74:ARG:HG2	1:A:97:ALA:O	2.15	0.47
1:A:155:PHE:HD2	1:A:369:LEU:HD21	1.79	0.47
1:B:241:PRO:O	1:B:245:GLY:N	2.44	0.47
1:A:177:TRP:C	1:A:179:ARG:N	2.73	0.47
1:A:202:ILE:O	1:A:206:VAL:HG23	2.15	0.47
1:B:108:ARG:HD2	3:B:482:HOH:O	2.15	0.47
1:B:220:MET:N	3:B:498:HOH:O	2.47	0.47
1:A:124:ARG:CB	1:A:125:PRO:HD3	2.35	0.47
1:A:182:ILE:HD12	1:A:185:SER:HA	1.97	0.47
1:A:236:VAL:HG12	1:A:236:VAL:O	2.14	0.47
1:B:63:VAL:C	1:B:65:ALA:H	2.22	0.47
1:B:99:ALA:HB3	1:B:104:HIS:HD2	1.79	0.47
1:B:155:PHE:CZ	1:B:159:VAL:HG21	2.49	0.47
1:B:184:THR:O	1:B:185:SER:HB2	2.14	0.47
1:B:358:GLY:O	1:B:359:ALA:C	2.57	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:137:GLY:O	1:A:141:GLU:HG3	2.15	0.47
1:B:367:GLU:O	1:B:371:ASP:HB2	2.15	0.47
1:A:154:PRO:HD3	3:A:540:HOH:O	2.14	0.47
1:B:203:THR:HG22	1:B:207:ARG:HE	1.80	0.47
1:B:269:ARG:CG	1:B:269:ARG:NH1	2.70	0.47
1:B:93:PRO:O	1:B:234:GLU:HA	2.15	0.47
1:A:274:PRO:O	1:A:275:GLY:C	2.56	0.46
1:A:124:ARG:HD2	1:A:364:MET:HG2	1.96	0.46
1:A:296:LEU:N	1:A:296:LEU:HD12	2.29	0.46
1:B:375:GLU:CG	1:B:376:ARG:N	2.77	0.46
1:B:59:ARG:N	1:B:59:ARG:HD2	2.30	0.46
1:B:201:TRP:CZ3	1:B:218:TYR:CE1	3.04	0.46
1:A:155:PHE:HB3	1:A:156:PRO:HD3	1.98	0.46
1:A:256:MET:HB3	1:A:284:LEU:CD1	2.45	0.46
1:B:212:SER:OG	1:B:213:GLU:N	2.44	0.46
1:A:190:ALA:C	1:A:192:GLU:H	2.24	0.46
1:A:153:GLU:H	1:A:154:PRO:HD2	1.79	0.46
1:A:280:ALA:HA	1:A:339:LEU:HA	1.98	0.46
1:B:280:ALA:HA	1:B:339:LEU:HA	1.98	0.46
1:A:230:VAL:O	1:A:231:GLY:C	2.59	0.46
1:A:338:ASP:C	1:A:340:ASP:H	2.24	0.46
1:B:97:ALA:HB1	1:B:296:LEU:CD2	2.46	0.46
1:A:99:ALA:HB3	1:A:104:HIS:HD2	1.81	0.46
1:B:14:PRO:HG2	1:B:44:ARG:NE	2.31	0.46
1:B:167:PRO:O	1:B:170:ASP:N	2.43	0.46
1:A:123:LEU:HD23	1:A:364:MET:CE	2.46	0.46
1:A:209:ARG:N	3:A:464:HOH:O	2.49	0.46
1:A:352:GLY:C	1:A:354:HIS:H	2.24	0.46
1:B:136:ASP:OD1	1:B:376:ARG:NH2	2.45	0.46
1:B:303:ASP:HB2	1:B:313:ALA:CB	2.43	0.46
1:A:178:THR:O	1:A:179:ARG:NH1	2.49	0.45
1:A:207:ARG:HH11	1:A:207:ARG:CG	2.24	0.45
1:A:297:ALA:HB2	1:A:318:TYR:CE2	2.50	0.45
1:A:321:TYR:CZ	1:A:350:GLY:HA2	2.51	0.45
1:B:15:PRO:HD2	3:B:452:HOH:O	2.15	0.45
1:B:247:GLU:HG3	1:B:248:ALA:N	2.31	0.45
1:A:270:MET:O	1:A:277:ARG:NH1	2.49	0.45
1:B:97:ALA:O	1:B:354:HIS:CE1	2.68	0.45
1:B:208:ALA:N	3:B:495:HOH:O	2.50	0.45
1:B:212:SER:OG	1:B:219:SER:HB3	2.17	0.45
1:A:153:GLU:N	1:A:154:PRO:HD2	2.32	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:208:ALA:HB2	3:B:495:HOH:O	2.16	0.45
1:B:177:TRP:C	1:B:179:ARG:H	2.25	0.45
1:B:353:HIS:C	1:B:355:PHE:H	2.24	0.45
1:A:226:GLY:HA2	3:A:545:HOH:O	2.17	0.45
1:B:14:PRO:O	1:B:15:PRO:C	2.60	0.45
1:B:259:LEU:O	1:B:266:LEU:HD12	2.16	0.45
1:A:28:GLU:OE1	1:A:28:GLU:CA	2.64	0.45
1:B:221:LEU:O	1:B:222:GLY:C	2.59	0.45
1:B:225:VAL:C	1:B:227:ARG:H	2.25	0.45
1:B:283:GLU:CD	1:B:341:ARG:HE	2.25	0.45
1:A:373:LEU:HA	1:A:373:LEU:HD23	1.79	0.45
1:B:71:ARG:NH1	1:B:304:VAL:HG13	2.32	0.45
1:B:167:PRO:O	1:B:169:ALA:N	2.49	0.45
1:A:138:ILE:HG13	1:A:139:LEU:N	2.31	0.44
1:A:214:GLY:C	1:A:216:ASP:N	2.72	0.44
1:B:78:THR:HG23	1:B:91:PRO:CG	2.46	0.44
1:B:177:TRP:C	1:B:179:ARG:N	2.75	0.44
1:A:295:GLY:O	1:A:318:TYR:HE2	2.00	0.44
1:B:107:LEU:O	1:B:238:LEU:HD11	2.17	0.44
1:A:342:ASP:CG	1:A:343:PRO:N	2.75	0.44
1:A:143:PRO:O	1:A:145:ALA:N	2.49	0.44
1:A:186:GLY:C	1:A:190:ALA:CB	2.91	0.44
1:A:265:GLU:H	1:A:265:GLU:HG3	1.43	0.44
1:A:341:ARG:HB3	1:A:344:ASN:HB2	1.99	0.44
1:B:22:LEU:CD1	1:B:24:LEU:HD12	2.47	0.44
1:B:291:ARG:HD3	1:B:293:SER:O	2.17	0.44
1:A:82:ILE:HD12	1:A:82:ILE:N	2.30	0.44
1:B:20:PRO:O	1:B:21:ALA:O	2.36	0.44
1:B:385:PRO:HD2	1:B:388:GLN:HG3	1.99	0.44
1:A:152:LEU:HD11	1:A:257:LEU:HD12	1.99	0.44
1:A:181:ILE:N	1:A:181:ILE:CD1	2.79	0.44
1:A:342:ASP:CG	1:A:343:PRO:CD	2.91	0.44
1:B:72:PHE:HB3	1:B:298:ARG:HB3	2.00	0.44
1:A:146:ASP:OD1	1:A:146:ASP:C	2.59	0.44
1:B:105:ASN:O	1:B:108:ARG:N	2.50	0.44
1:B:153:GLU:HA	1:B:250:THR:HG23	1.99	0.44
1:A:69:ASP:HB3	1:A:72:PHE:CD1	2.53	0.44
1:A:189:GLU:HA	3:A:488:HOH:O	2.16	0.44
1:A:203:THR:CG2	1:A:207:ARG:NH1	2.81	0.44
1:A:69:ASP:OD2	1:A:71:ARG:NH2	2.51	0.44
1:A:198:LEU:O	1:A:202:ILE:HG13	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:114:ALA:HB2	1:B:220:MET:HE1	2.00	0.43
1:B:215:GLY:O	1:B:217:VAL:N	2.51	0.43
1:B:239:ALA:HA	1:B:242:LEU:HD12	2.00	0.43
1:A:146:ASP:HA	1:A:403:LEU:O	2.17	0.43
1:A:156:PRO:O	1:A:160:VAL:HG23	2.18	0.43
1:A:209:ARG:HG3	1:A:209:ARG:HH11	1.83	0.43
1:B:103:ASP:CG	1:B:106:ARG:HH21	2.26	0.43
1:B:283:GLU:OE1	1:B:283:GLU:HA	2.18	0.43
1:A:60:TYR:HB3	3:A:437:HOH:O	2.17	0.43
1:B:37:MET:CG	1:B:58:THR:HG21	2.41	0.43
1:B:198:LEU:HD23	1:B:198:LEU:C	2.43	0.43
1:B:232:GLU:O	1:B:233:THR:HB	2.18	0.43
1:B:296:LEU:N	1:B:296:LEU:CD1	2.81	0.43
1:A:115:PHE:CD1	1:A:115:PHE:N	2.86	0.43
1:A:192:GLU:HG3	1:A:196:ARG:HE	1.83	0.43
1:A:203:THR:HG23	1:A:207:ARG:NH1	2.33	0.43
1:A:302:GLU:O	1:A:303:ASP:O	2.37	0.43
1:A:15:PRO:HG2	1:A:17:ARG:CZ	2.48	0.43
1:A:37:MET:HG2	1:A:58:THR:HG23	1.97	0.43
1:A:108:ARG:HH11	1:A:112:ALA:HB2	1.82	0.43
1:A:171:ARG:HG3	1:A:171:ARG:NH1	2.33	0.43
1:B:36:LEU:HD12	1:B:36:LEU:HA	1.90	0.43
1:B:67:THR:HG21	1:B:351:ASN:HB3	2.00	0.43
1:B:209:ARG:HD3	1:B:219:SER:OG	2.18	0.43
1:B:321:TYR:O	1:B:325:ASN:N	2.51	0.43
1:A:81:GLN:OE1	1:A:81:GLN:HA	2.18	0.43
1:A:83:THR:O	1:A:83:THR:CG2	2.65	0.43
1:A:239:ALA:O	1:A:240:GLY:C	2.61	0.43
1:B:153:GLU:OE1	1:B:179:ARG:NH2	2.52	0.43
1:A:103:ASP:O	1:A:104:HIS:C	2.61	0.43
1:A:270:MET:CA	1:A:270:MET:HE2	2.48	0.43
1:B:168:ALA:HA	1:B:171:ARG:NE	2.33	0.43
1:B:249:VAL:HG11	1:B:365:GLN:NE2	2.32	0.43
1:A:58:THR:HG23	1:A:323:ALA:HB1	2.01	0.43
1:A:164:MET:HA	1:A:217:VAL:HB	2.00	0.43
1:B:321:TYR:HA	1:B:324:ALA:HB3	1.99	0.43
1:B:326:ARG:HB3	3:B:489:HOH:O	2.18	0.43
1:B:333:ASP:OD1	1:B:336:ARG:HD3	2.19	0.43
1:A:403:LEU:HD23	1:A:403:LEU:HA	1.90	0.42
1:B:109:ARG:O	1:B:110:ALA:C	2.62	0.42
1:B:212:SER:CB	1:B:219:SER:HB3	2.50	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:349:TYR:HE2	2:A:430:HEM:HBB2	1.82	0.42
1:A:291:ARG:HD2	3:A:473:HOH:O	2.19	0.42
1:B:131:LEU:HD21	1:B:372:THR:HB	2.01	0.42
1:B:143:PRO:CG	1:B:409:HIS:HB2	2.49	0.42
1:B:221:LEU:O	1:B:225:VAL:HG23	2.18	0.42
1:B:216:ASP:O	1:B:217:VAL:C	2.62	0.42
1:A:244:ILE:O	1:A:244:ILE:HG22	2.20	0.42
1:B:47:LEU:O	1:B:48:PRO:O	2.38	0.42
1:B:212:SER:HB3	1:B:219:SER:HB3	2.02	0.42
1:B:22:LEU:HD13	1:B:292:THR:HA	2.01	0.42
1:B:79:GLN:C	1:B:80:ARG:HG2	2.44	0.42
1:B:173:ARG:HH12	1:B:201:TRP:HE1	1.67	0.41
1:B:312:ALA:HB3	1:B:315:GLU:CD	2.45	0.41
1:A:47:LEU:HD12	1:A:84:ARG:HA	2.02	0.41
1:A:105:ASN:HA	3:A:460:HOH:O	2.20	0.41
1:A:106:ARG:HA	1:A:109:ARG:HH21	1.84	0.41
1:A:355:PHE:CD2	1:A:356:CYS:N	2.88	0.41
1:A:131:LEU:HD21	1:A:372:THR:HB	2.03	0.41
1:B:135:VAL:O	1:B:136:ASP:C	2.63	0.41
1:B:259:LEU:HD22	1:B:263:ARG:HD2	2.02	0.41
1:A:180:GLN:C	1:A:182:ILE:HG13	2.45	0.41
1:B:167:PRO:CG	1:B:209:ARG:NH2	2.83	0.41
1:B:381:ARG:HD3	3:B:503:HOH:O	2.21	0.41
1:B:164:MET:CE	1:B:239:ALA:HB1	2.51	0.41
1:B:342:ASP:N	1:B:342:ASP:OD1	2.54	0.41
1:A:297:ALA:HB2	1:A:318:TYR:CZ	2.55	0.41
1:A:336:ARG:NH2	1:A:338:ASP:OD2	2.53	0.41
1:B:101:GLN:OE1	1:B:104:HIS:HB3	2.19	0.41
1:B:183:SER:C	1:B:185:SER:N	2.78	0.41
1:A:123:LEU:HD23	1:A:364:MET:HE1	2.03	0.41
1:B:164:MET:O	1:B:216:ASP:OD2	2.38	0.41
1:A:92:ARG:HG2	1:A:93:PRO:CD	2.50	0.41
1:A:333:ASP:N	1:A:334:PRO:CD	2.84	0.41
1:A:353:HIS:HB3	3:A:436:HOH:O	2.20	0.41
1:B:198:LEU:HD23	1:B:202:ILE:HG13	2.02	0.41
1:B:384:VAL:HB	1:B:385:PRO:CD	2.51	0.41
1:A:283:GLU:HA	1:A:283:GLU:OE1	2.20	0.41
1:A:33:LEU:HD23	1:A:33:LEU:C	2.46	0.41
1:A:73:GLY:HA3	3:A:526:HOH:O	2.21	0.40
1:A:186:GLY:O	1:B:79:GLN:OE1	2.39	0.40
1:A:401:ARG:NH2	3:A:531:HOH:O	2.52	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:22:LEU:HD12	1:B:30:ASP:OD1	2.21	0.40
1:B:126:ARG:O	1:B:130:ILE:HG13	2.21	0.40
1:B:353:HIS:CD2	1:B:353:HIS:N	2.88	0.40
1:B:208:ALA:C	1:B:210:ALA:H	2.28	0.40
1:B:279:THR:O	1:B:282:ASP:HB2	2.22	0.40
1:B:319:VAL:HG21	3:B:457:HOH:O	2.22	0.40
1:A:101:GLN:OE1	1:A:105:ASN:HB2	2.22	0.40
1:B:285:LEU:HA	1:B:285:LEU:HD23	1.90	0.40
1:B:264:ARG:O	1:B:265:GLU:C	2.64	0.40
1:B:231:GLY:C	1:B:232:GLU:O	2.62	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	395/413 (96%)	335 (85%)	41 (10%)	19 (5%)	2	7
1	B	384/413 (93%)	305 (79%)	57 (15%)	22 (6%)	1	4
All	All	779/826 (94%)	640 (82%)	98 (13%)	41 (5%)	1	5

All (41) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	182	ILE
1	A	208	ALA
1	A	209	ARG
1	A	213	GLU
1	A	231	GLY
1	A	232	GLU
1	A	276	ALA
1	A	303	ASP

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Mol	Chain	Res	Type
1	A	342	ASP
1	B	21	ALA
1	B	48	PRO
1	B	168	ALA
1	B	232	GLU
1	B	342	ASP
1	A	211	GLY
1	A	214	GLY
1	B	22	LEU
1	B	71	ARG
1	B	143	PRO
1	B	216	ASP
1	B	231	GLY
1	B	245	GLY
1	A	190	ALA
1	B	106	ARG
1	B	176	SER
1	A	103	ASP
1	A	302	GLU
1	B	105	ASN
1	B	183	SER
1	B	212	SER
1	B	233	THR
1	B	343	PRO
1	A	167	PRO
1	A	230	VAL
1	A	293	SER
1	B	346	HIS
1	B	93	PRO
1	A	144	PRO
1	A	306	VAL
1	B	20	PRO
1	B	70	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	310/329 (94%)	274 (88%)	36 (12%)	5	18
1	B	303/329 (92%)	278 (92%)	25 (8%)	10	32
All	All	613/658 (93%)	552 (90%)	61 (10%)	7	24

All (61) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	LEU
1	A	24	LEU
1	A	28	GLU
1	A	36	LEU
1	A	51	GLU
1	A	59	ARG
1	A	67	THR
1	A	74	ARG
1	A	76	GLU
1	A	82	ILE
1	A	102	PRO
1	A	103	ASP
1	A	105	ASN
1	A	143	PRO
1	A	151	VAL
1	A	172	GLU
1	A	179	ARG
1	A	180	GLN
1	A	181	ILE
1	A	182	ILE
1	A	204	GLU
1	A	207	ARG
1	A	213	GLU
1	A	232	GLU
1	A	233	THR
1	A	238	LEU
1	A	265	GLU
1	A	269	ARG
1	A	277	ARG
1	A	285	LEU
1	A	294	VAL
1	A	343	PRO
1	A	366	THR
1	A	387	GLU
1	A	395	THR

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Mol	Chain	Res	Type
1	A	396	MET
1	B	36	LEU
1	B	48	PRO
1	B	76	GLU
1	B	109	ARG
1	B	117	VAL
1	B	143	PRO
1	B	149	GLU
1	B	160	VAL
1	B	198	LEU
1	B	207	ARG
1	B	232	GLU
1	B	238	LEU
1	B	260	LEU
1	B	269	ARG
1	B	270	MET
1	B	284	LEU
1	B	296	LEU
1	B	303	ASP
1	B	305	GLU
1	B	311	ILE
1	B	317	VAL
1	B	353	HIS
1	B	366	THR
1	B	381	ARG
1	B	382	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	68	ASN
1	A	105	ASN
1	A	180	GLN
1	A	252	ASN
1	A	255	GLN
1	A	351	ASN
1	A	353	HIS
1	A	365	GLN
1	B	68	ASN
1	B	79	GLN
1	B	105	ASN
1	B	290	HIS

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Mol	Chain	Res	Type
1	B	307	HIS
1	B	365	GLN
1	B	408	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](#)

2 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	B	430	1,3	50,50,50	1.71	11 (22%)	67,82,82	0.98	5 (7%)
2	HEM	A	430	1	50,50,50	1.63	11 (22%)	67,82,82	0.88	2 (2%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	HEM	B	430	1,3	-	8/14/54/54	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	HEM	A	430	1	-	3/14/54/54	-

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	430	HEM	CBB-CAB	5.04	1.54	1.30
2	A	430	HEM	CBC-CAC	4.98	1.54	1.30
2	B	430	HEM	CBC-CAC	4.92	1.54	1.30
2	A	430	HEM	CBB-CAB	4.80	1.53	1.30
2	A	430	HEM	CAC-C3C	3.36	1.56	1.47
2	B	430	HEM	CAC-C3C	3.17	1.55	1.47
2	B	430	HEM	FE-NA	2.97	2.05	1.95
2	B	430	HEM	C1A-C2A	2.61	1.49	1.44
2	B	430	HEM	CMC-C2C	2.47	1.55	1.50
2	B	430	HEM	FE-NB	2.47	2.02	1.94
2	B	430	HEM	CAB-C3B	2.43	1.53	1.47
2	A	430	HEM	FE-NA	2.35	2.02	1.95
2	A	430	HEM	FE-NB	2.32	2.02	1.94
2	A	430	HEM	CMC-C2C	2.30	1.55	1.50
2	B	430	HEM	CAA-C2A	2.26	1.57	1.51
2	A	430	HEM	FE-NC	2.21	2.02	1.95
2	A	430	HEM	O2A-CGA	-2.13	1.23	1.30
2	A	430	HEM	CAB-C3B	2.11	1.53	1.47
2	B	430	HEM	O1A-CGA	2.11	1.29	1.22
2	A	430	HEM	CMD-C2D	2.10	1.55	1.50
2	A	430	HEM	O1A-CGA	2.06	1.28	1.22
2	B	430	HEM	C2A-C3A	-2.04	1.33	1.38

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	430	HEM	CBB-CAB-C3B	-2.70	114.02	127.53
2	B	430	HEM	O1A-CGA-CBA	-2.66	114.66	123.09
2	A	430	HEM	CBB-CAB-C3B	-2.66	114.26	127.53
2	A	430	HEM	O1A-CGA-CBA	-2.44	115.36	123.09
2	B	430	HEM	CBC-CAC-C3C	-2.30	116.01	127.53
2	B	430	HEM	CAA-C2A-C1A	2.10	129.05	124.94
2	B	430	HEM	CBA-CAA-C2A	2.08	118.29	112.53

There are no chirality outliers.

All (11) torsion outliers are listed below:

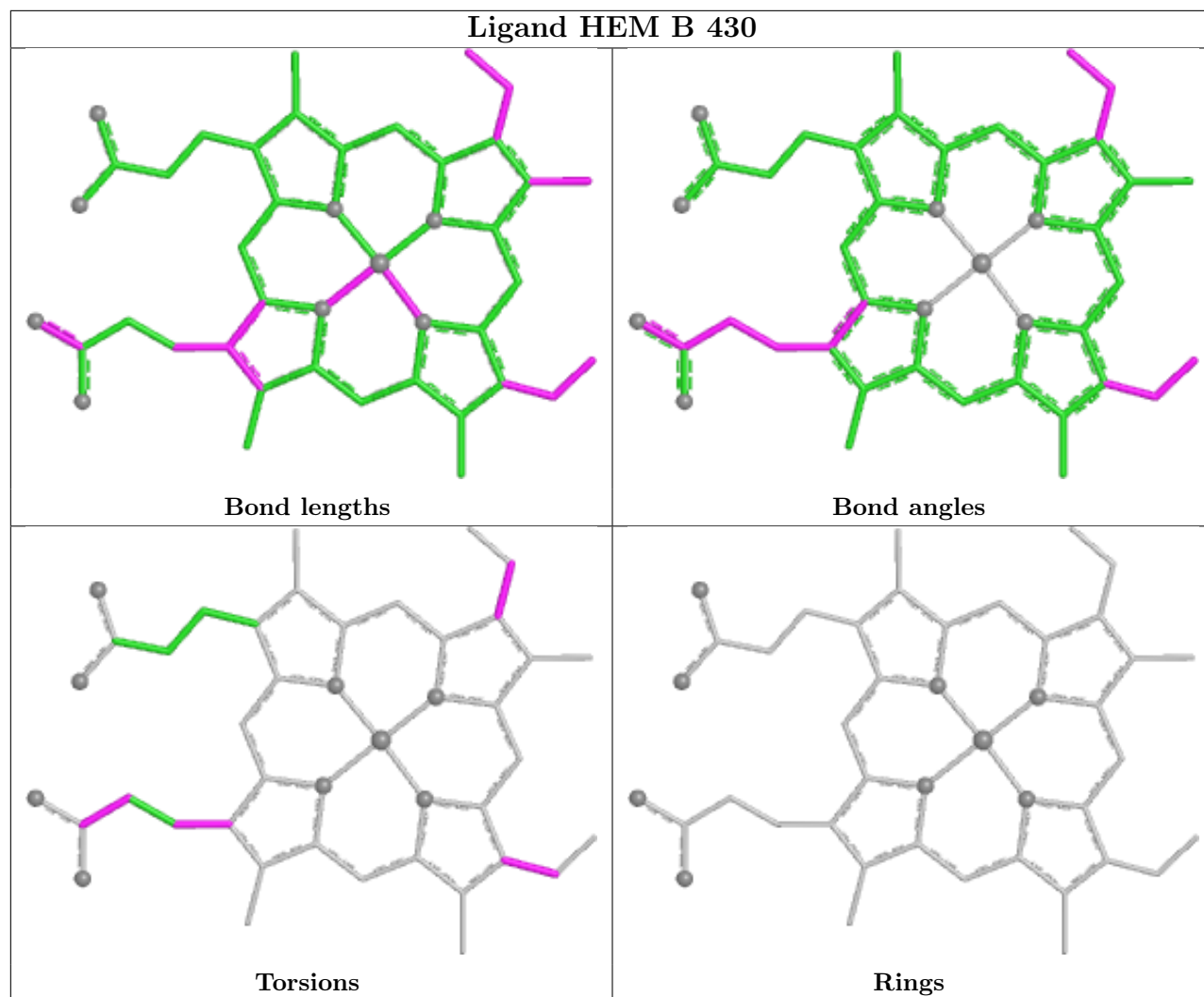
Mol	Chain	Res	Type	Atoms
2	A	430	HEM	C2B-C3B-CAB-CBB
2	A	430	HEM	C4B-C3B-CAB-CBB
2	B	430	HEM	C2B-C3B-CAB-CBB
2	B	430	HEM	C1A-C2A-CAA-CBA
2	B	430	HEM	C3A-C2A-CAA-CBA
2	B	430	HEM	C4B-C3B-CAB-CBB
2	B	430	HEM	C2C-C3C-CAC-CBC
2	B	430	HEM	CAA-CBA-CGA-O2A
2	B	430	HEM	CAA-CBA-CGA-O1A
2	B	430	HEM	C4C-C3C-CAC-CBC
2	A	430	HEM	CAA-CBA-CGA-O1A

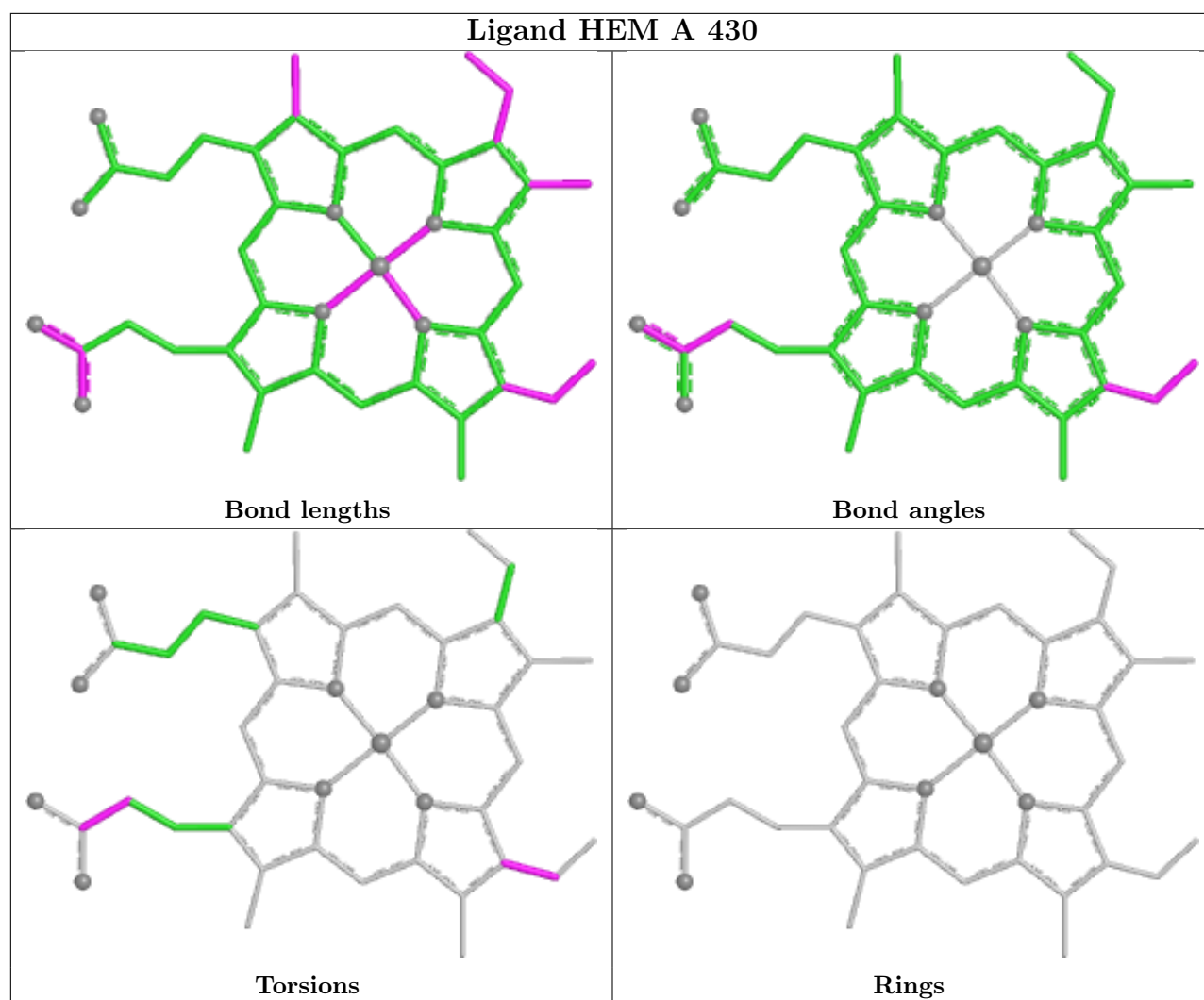
There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	430	HEM	2	0
2	A	430	HEM	2	0

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





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	397/413 (96%)	0.05	14 (3%) 47 38	2, 8, 34, 73	0
1	B	388/413 (93%)	0.27	26 (6%) 24 19	2, 14, 48, 66	0
All	All	785/826 (95%)	0.16	40 (5%) 33 26	2, 10, 44, 73	0

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	184	THR	6.8
1	A	188	ALA	5.0
1	B	183	SER	4.8
1	B	184	THR	4.8
1	A	182	ILE	4.3
1	B	185	SER	4.2
1	B	14	PRO	3.9
1	B	50	GLY	3.9
1	B	181	ILE	3.7
1	B	22	LEU	3.6
1	A	143	PRO	3.6
1	B	182	ILE	3.5
1	A	14	PRO	3.3
1	A	185	SER	3.2
1	A	231	GLY	3.1
1	B	210	ALA	3.0
1	A	189	GLU	3.0
1	A	343	PRO	2.9
1	B	189	GLU	2.8
1	B	28	GLU	2.8
1	A	228	GLY	2.7
1	B	211	GLY	2.5
1	B	187	GLY	2.4
1	B	225	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	187	GLY	2.3
1	A	186	GLY	2.3
1	A	216	ASP	2.3
1	B	191	ALA	2.3
1	B	186	GLY	2.3
1	B	172	GLU	2.3
1	B	194	ALA	2.3
1	B	180	GLN	2.3
1	B	216	ASP	2.2
1	B	18	ASP	2.2
1	A	183	SER	2.2
1	B	208	ALA	2.2
1	B	143	PRO	2.1
1	B	222	GLY	2.1
1	B	219	SER	2.1
1	B	65	ALA	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](#)

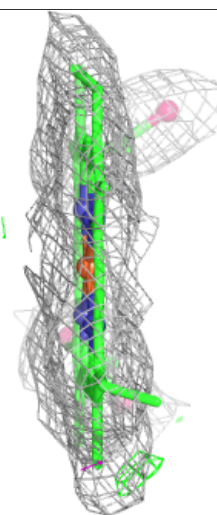
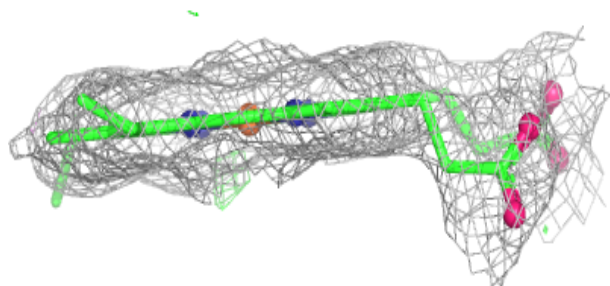
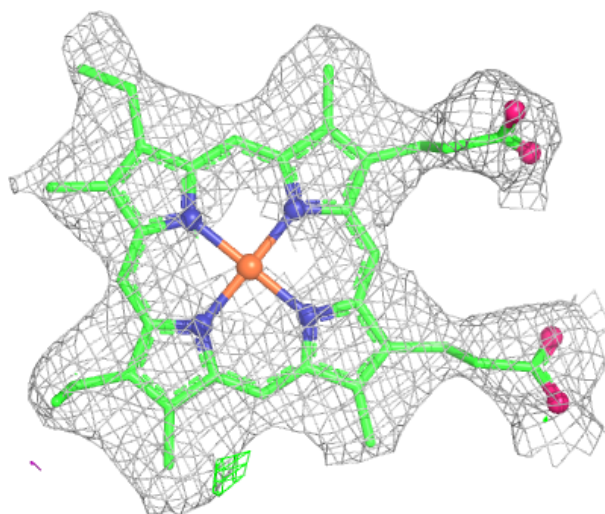
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

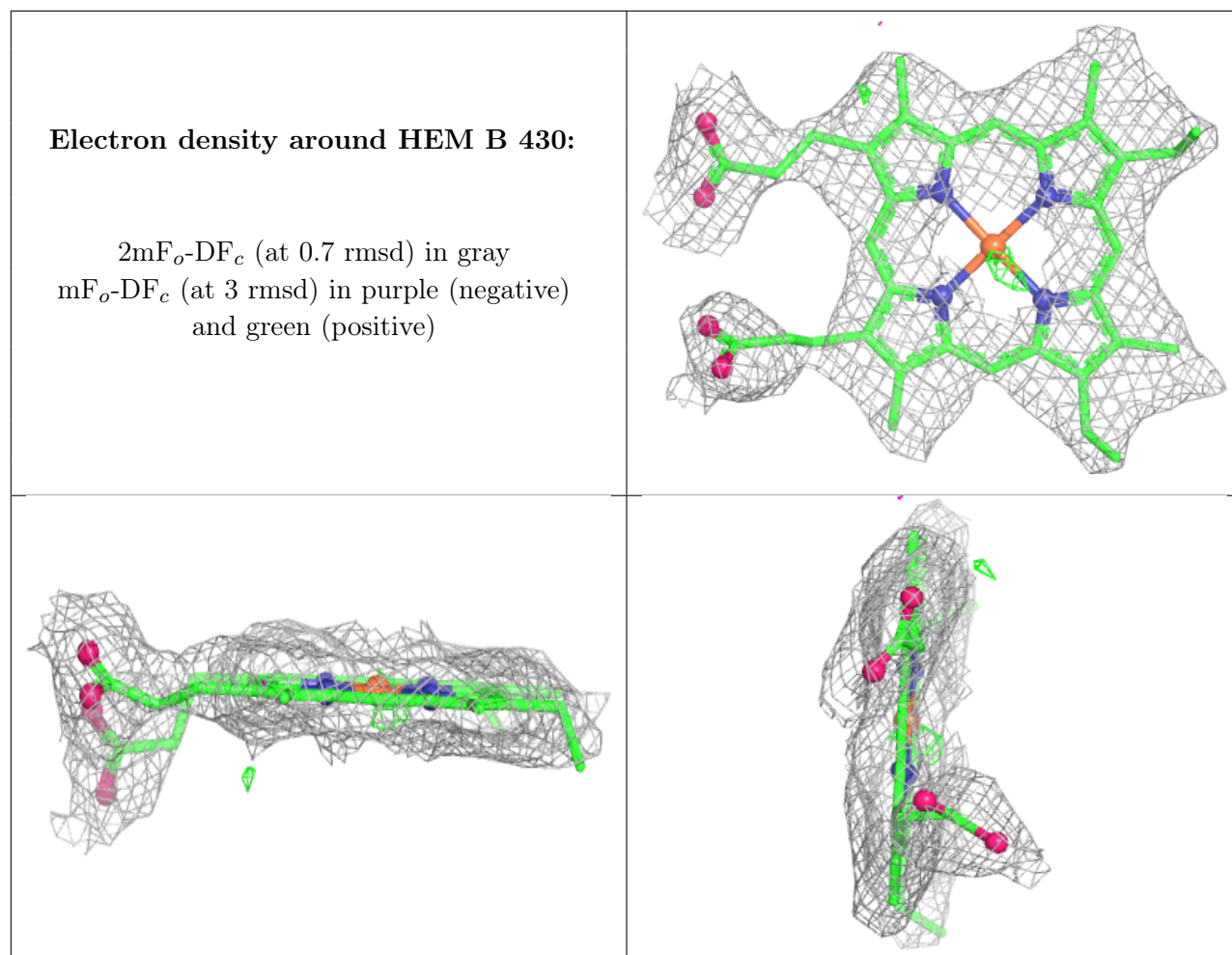
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	HEM	A	430	43/43	0.94	0.09	1,6,7,8	0
2	HEM	B	430	43/43	0.94	0.10	10,13,16,21	0

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 A 430:

$2mF_o-DF_c$ (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.