



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 02:31 PM UTC

PDB ID : 5FSA / pdb_00005fsa
Title : Crystal structure of sterol 14-alpha demethylase (CYP51) from a pathogenic yeast *Candida albicans* in complex with the antifungal drug posaconazole
Authors : Hargrove, T.Y.; Wawrzak, Z.; Friggeri, L.; Lepesheva, G.I.
Deposited on : 2016-01-02
Resolution : 2.86 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

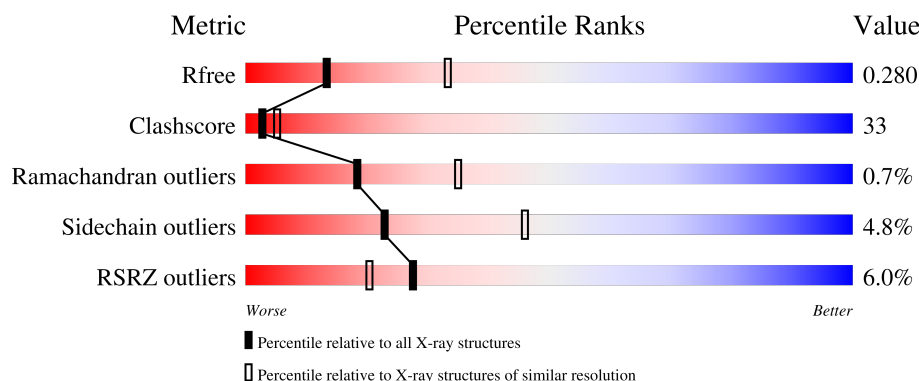
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.86 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	490	6% 56% 38% ..
1	B	490	6% 55% 37% 6% ..

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8163 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 CYP51 VARIANT1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	484	Total	C	N	O	S	0	0	0
			3929	2539	652	722	16			
1	B	483	Total	C	N	O	S	0	0	0
			3919	2532	650	721	16			

There are 22 discrepancies between the modelled and reference sequences:

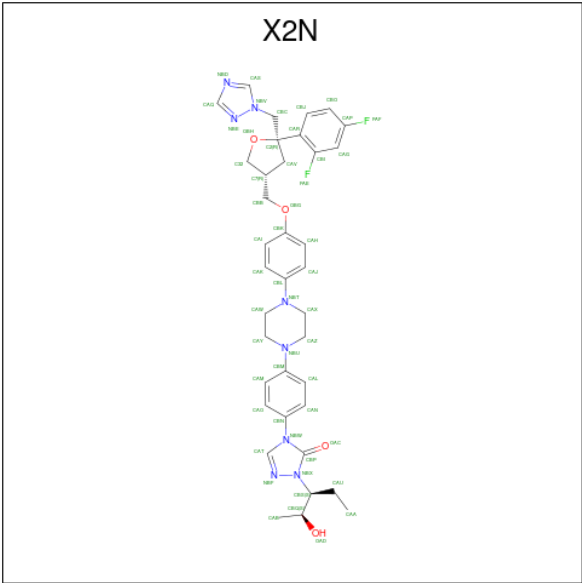
Chain	Residue	Modelled	Actual	Comment	Reference
A	43	MET	-	expression tag	UNP Q9P4W0
A	44	ALA	-	expression tag	UNP Q9P4W0
A	45	LYS	-	expression tag	UNP Q9P4W0
A	46	LYS	-	expression tag	UNP Q9P4W0
A	47	THR	-	expression tag	UNP Q9P4W0
A	48	PRO	-	expression tag	UNP Q9P4W0
A	529	HIS	-	expression tag	UNP Q9P4W0
A	530	HIS	-	expression tag	UNP Q9P4W0
A	531	HIS	-	expression tag	UNP Q9P4W0
A	532	HIS	-	expression tag	UNP Q9P4W0
A	263	LEU	SER	engineered mutation	UNP Q9P4W0
B	43	MET	-	expression tag	UNP Q9P4W0
B	44	ALA	-	expression tag	UNP Q9P4W0
B	45	LYS	-	expression tag	UNP Q9P4W0
B	46	LYS	-	expression tag	UNP Q9P4W0
B	47	THR	-	expression tag	UNP Q9P4W0
B	48	PRO	-	expression tag	UNP Q9P4W0
B	529	HIS	-	expression tag	UNP Q9P4W0
B	530	HIS	-	expression tag	UNP Q9P4W0
B	531	HIS	-	expression tag	UNP Q9P4W0
B	532	HIS	-	expression tag	UNP Q9P4W0
B	263	LEU	SER	engineered mutation	UNP Q9P4W0

- 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 POSACONAZOLE (CCD ID: X2N) (formula: $C_{37}H_{42}F_2N_8O_4$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	F	N	O	0	0
			51	37	2	8	4		
3	B	1	Total	C	F	N	O	0	0
			51	37	2	8	4		

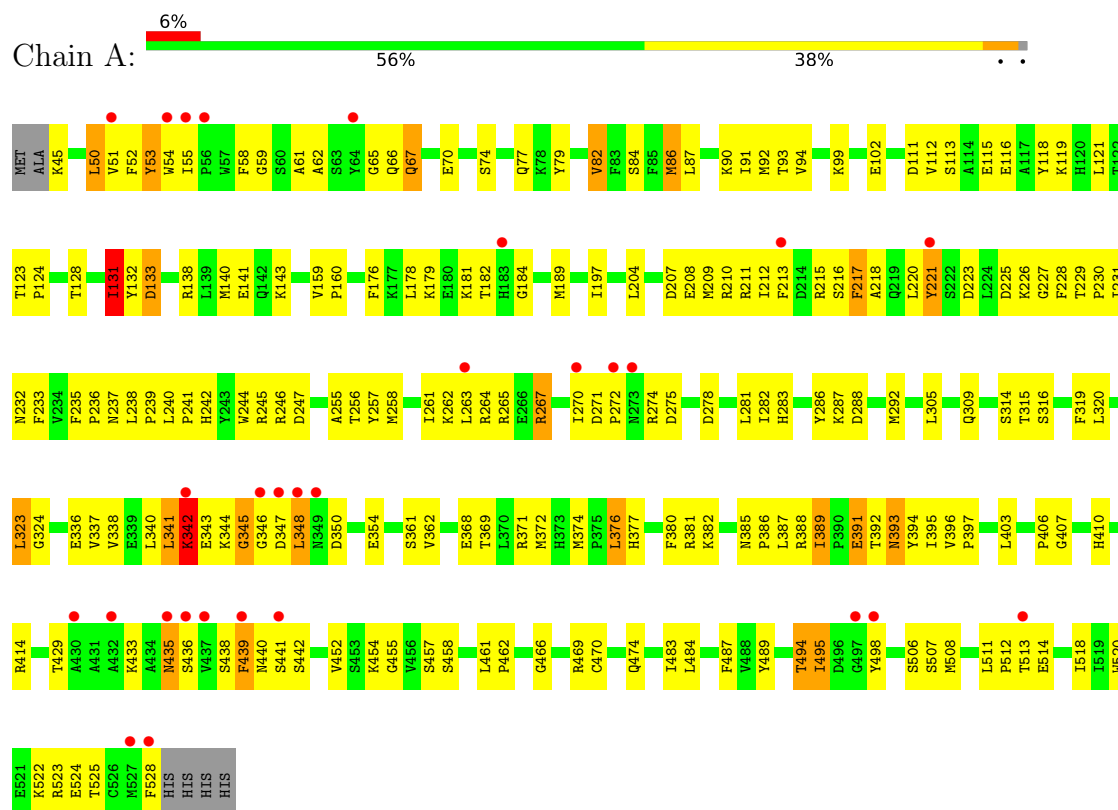
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	61	Total 61	O 61	0	0
4	B	66	Total 66	O 66	0	0

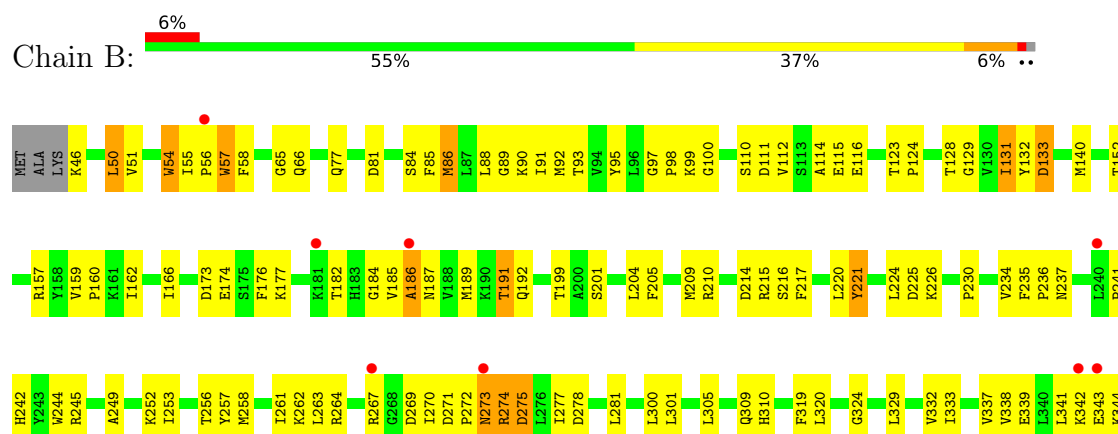
3 Residue-property plots

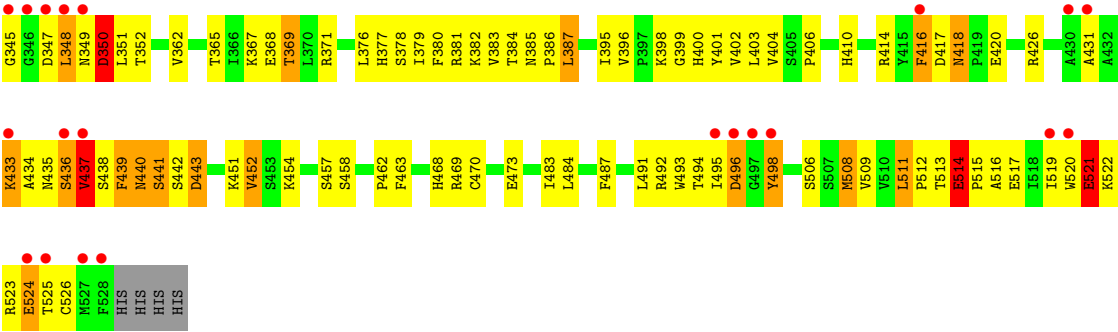
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: CYP51 VARIANT1



• Molecule 1: CYP51 VARIANT1





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	179.59Å 72.91Å 79.71Å 90.00° 96.13° 90.00°	Depositor
Resolution (Å)	50.39 – 2.86 50.39 – 2.86	Depositor EDS
% Data completeness (in resolution range)	99.3 (50.39-2.86) 99.3 (50.39-2.86)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.18 (at 2.86Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.241 , 0.256 0.243 , 0.280	Depositor DCC
R_{free} test set	1235 reflections (5.20%)	wwPDB-VP
Wilson B-factor (Å ²)	66.1	Xtriage
Anisotropy	0.142	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 60.8	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	0.92	EDS
Total number of atoms	8163	wwPDB-VP
Average B, all atoms (Å ²)	73.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.07% 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, X2N

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	0.34	0/4040	0.88	13/5475 (0.2%)
1	B	0.38	0/4030	0.94	18/5462 (0.3%)
All	All	0.36	0/8070	0.91	31/10937 (0.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

There are no bond length outliers.

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	343	GLU	N-CA-C	9.13	124.62	111.56
1	A	271	ASP	CA-C-N	-8.13	111.38	119.76
1	A	271	ASP	C-N-CA	-8.13	111.38	119.76
1	A	435	ASN	N-CA-C	-8.01	97.67	108.24
1	B	514	GLU	N-CA-C	7.92	124.23	113.77
1	B	441	SER	N-CA-C	7.21	119.04	111.03
1	B	496	ASP	N-CA-C	-7.12	104.55	113.23
1	A	343	GLU	N-CA-C	7.01	118.92	111.28
1	B	437	VAL	CB-CA-C	-6.88	102.91	111.92
1	B	186	ALA	N-CA-C	6.85	120.62	109.59
1	B	56	PRO	N-CA-C	6.47	122.72	113.47
1	B	521	GLU	N-CA-C	-6.40	100.78	110.70
1	B	437	VAL	N-CA-C	-6.30	105.59	111.45
1	A	133	ASP	N-CA-C	-6.22	104.05	112.26

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	131	ILE	N-CA-C	6.12	122.06	109.34
1	A	345	GLY	N-CA-C	-5.98	99.01	113.18
1	A	391	GLU	N-CA-C	-5.87	104.72	112.24
1	B	97	GLY	CA-C-N	5.68	125.15	119.24
1	B	97	GLY	C-N-CA	5.68	125.15	119.24
1	B	350	ASP	N-CA-C	5.66	122.86	110.80
1	A	389	ILE	CA-C-N	-5.62	114.95	120.52
1	A	389	ILE	C-N-CA	-5.62	114.95	120.52
1	A	380	PHE	N-CA-C	5.51	118.54	109.72
1	B	57	TRP	N-CA-CB	-5.26	102.30	110.98
1	A	348	LEU	N-CA-C	-5.24	105.57	111.28
1	B	498	TYR	N-CA-C	-5.23	102.45	110.14
1	B	133	ASP	N-CA-C	-5.17	105.43	112.26
1	A	131	ILE	N-CA-C	5.17	120.09	109.34
1	B	275	ASP	N-CA-C	5.15	115.84	108.74
1	B	416	PHE	N-CA-C	5.12	116.08	108.60
1	A	495	ILE	N-CA-C	5.08	119.89	109.34

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	342	LYS	Peptide

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3929	0	3865	226	1
1	B	3919	0	3847	298	0
2	A	43	0	30	5	0
2	B	43	0	30	10	0
3	A	51	0	42	5	0
3	B	51	0	42	6	0
4	A	61	0	0	2	0
4	B	66	0	0	3	0
All	All	8163	0	7856	526	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 33.

All (526) 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:272:PRO:HB2	1:A:274:ARG:HG3	1.32	1.09
1:B:523:ARG:HG3	1:B:524:GLU:H	1.25	1.00
1:A:121:LEU:HD11	1:A:508:MET:HE1	1.45	0.99
1:B:506:SER:HB2	3:B:590:X2N:HAT	1.47	0.97
1:B:263:LEU:O	1:B:267:ARG:HG2	1.64	0.96
1:B:184:GLY:HA2	1:B:521:GLU:HB3	1.45	0.96
1:B:226:LYS:HG3	1:B:242:HIS:NE2	1.81	0.94
1:B:271:ASP:N	1:B:272:PRO:CD	2.30	0.93
1:B:491:LEU:HA	1:B:523:ARG:HB3	1.51	0.93
1:B:221:TYR:O	1:B:225:ASP:HB2	1.69	0.92
1:B:270:ILE:C	1:B:272:PRO:HD2	1.95	0.92
1:A:213:PHE:HA	1:A:217:PHE:CE2	2.04	0.91
1:B:111:ASP:C	1:B:384:THR:HG22	1.96	0.91
1:B:272:PRO:HB2	1:B:278:ASP:OD2	1.69	0.91
1:B:271:ASP:N	1:B:272:PRO:HD2	1.87	0.90
1:B:272:PRO:HB3	1:B:274:ARG:NE	1.86	0.90
1:A:212:ILE:C	1:A:217:PHE:HE2	1.80	0.90
1:B:272:PRO:HB3	1:B:274:ARG:CG	2.02	0.89
1:B:523:ARG:HG3	1:B:524:GLU:N	1.88	0.88
1:B:514:GLU:CG	1:B:515:PRO:HD3	2.05	0.87
1:A:265:ARG:HA	1:A:270:ILE:HD11	1.56	0.87
1:B:439:PHE:O	1:B:440:ASN:HB2	1.72	0.86
1:A:438:SER:HB3	1:A:458:SER:HB3	1.56	0.86
1:B:173:ASP:O	1:B:177:LYS:HG2	1.76	0.85
1:B:493:TRP:HA	1:B:520:TRP:O	1.75	0.85
1:B:90:LYS:HE3	1:B:401:TYR:CE1	2.10	0.85
1:B:329:LEU:O	1:B:332:VAL:HG12	1.79	0.82
1:A:86:MET:HA	1:A:91:ILE:HA	1.60	0.81
1:B:514:GLU:HG2	1:B:515:PRO:HD3	1.62	0.81
1:B:204:LEU:O	1:B:275:ASP:HB2	1.80	0.81
1:B:319:PHE:HB3	1:B:369:THR:HG21	1.62	0.80
1:B:93:THR:OG1	1:B:400:HIS:CE1	2.35	0.79
1:B:508:MET:HA	1:B:508:MET:HE2	1.64	0.79
1:B:174:GLU:HA	1:B:177:LYS:HE3	1.65	0.79
1:B:173:ASP:C	1:B:177:LYS:HE2	2.08	0.79
1:B:437:VAL:HG12	1:B:438:SER:N	1.96	0.78
1:A:52:PHE:HB3	1:B:54:TRP:CH2	2.18	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:220:LEU:HD22	1:B:252:LYS:HD3	1.66	0.77
1:B:46:LYS:O	1:B:395:ILE:HG22	1.86	0.76
1:B:494:THR:HG22	1:B:495:ILE:N	2.00	0.76
1:B:496:ASP:HB2	1:B:519:ILE:HG13	1.68	0.76
1:A:115:GLU:O	1:A:119:LYS:HG3	1.86	0.76
1:A:354:GLU:N	1:A:354:GLU:OE1	2.17	0.76
1:A:212:ILE:C	1:A:217:PHE:CE2	2.65	0.74
1:B:205:PHE:O	1:B:264:ARG:NH2	2.20	0.74
2:A:580:HEM:CMB	2:A:580:HEM:HBB2	2.18	0.74
1:A:113:SER:HB2	1:A:115:GLU:OE1	1.86	0.73
1:B:420:GLU:OE1	1:B:420:GLU:N	2.20	0.73
1:A:319:PHE:O	1:A:323:LEU:HB2	1.89	0.73
1:B:112:VAL:HG13	1:B:381:ARG:HB3	1.70	0.73
1:B:435:ASN:OD1	1:B:436:SER:N	2.22	0.73
1:A:257:TYR:O	1:A:261:ILE:HG13	1.89	0.72
1:B:128:THR:HG23	1:B:133:ASP:OD2	1.88	0.72
1:B:523:ARG:CG	1:B:524:GLU:H	2.02	0.72
1:B:523:ARG:HG2	1:B:523:ARG:HH11	1.54	0.72
1:B:398:LYS:HG3	1:B:399:GLY:N	2.03	0.72
1:B:416:PHE:HD1	1:B:417:ASP:H	1.36	0.72
1:A:438:SER:CB	1:A:458:SER:HB3	2.20	0.72
1:A:223:ASP:HB3	1:A:245:ARG:HG3	1.70	0.72
2:A:580:HEM:HBB2	2:A:580:HEM:HMB2	1.72	0.72
1:B:50:LEU:HD13	1:B:50:LEU:C	2.16	0.71
1:B:494:THR:HG22	1:B:495:ILE:H	1.53	0.71
1:A:337:VAL:O	1:A:341:LEU:HD22	1.90	0.71
1:A:213:PHE:N	1:A:217:PHE:HE2	1.88	0.71
1:B:272:PRO:HB3	1:B:274:ARG:HG2	1.73	0.71
1:A:128:THR:HG23	1:A:133:ASP:OD2	1.91	0.70
1:B:348:LEU:O	1:B:349:ASN:HB2	1.91	0.70
1:A:215:ARG:HG2	1:A:216:SER:H	1.56	0.70
1:A:53:TYR:HB3	1:A:59:GLY:HA2	1.74	0.70
1:A:121:LEU:CD1	1:A:508:MET:HE1	2.20	0.70
1:A:264:ARG:NH1	1:A:278:ASP:OD1	2.25	0.70
1:A:121:LEU:HD11	1:A:508:MET:CE	2.21	0.70
1:B:438:SER:HB3	1:B:458:SER:HB2	1.74	0.70
1:B:270:ILE:HD13	1:B:281:LEU:HD13	1.75	0.69
1:B:525:THR:OG1	1:B:526:CYS:N	2.21	0.69
1:B:523:ARG:HG2	1:B:523:ARG:NH1	2.07	0.69
1:A:495:ILE:HG22	1:A:495:ILE:O	1.91	0.69
1:B:495:ILE:O	1:B:496:ASP:HB3	1.93	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:324:GLY:C	1:A:495:ILE:HG12	2.18	0.69
1:B:514:GLU:HG3	1:B:515:PRO:HD3	1.73	0.69
1:A:270:ILE:HD13	1:A:281:LEU:HD13	1.75	0.69
1:A:272:PRO:HB2	1:A:274:ARG:CG	2.17	0.69
1:B:264:ARG:HG2	1:B:267:ARG:HH21	1.58	0.69
1:B:348:LEU:O	1:B:348:LEU:HD13	1.92	0.69
1:A:348:LEU:C	1:A:348:LEU:HD13	2.17	0.69
1:A:52:PHE:HB3	1:B:54:TRP:CZ3	2.28	0.68
1:A:62:ALA:O	1:A:66:GLN:HG3	1.93	0.68
1:A:228:PHE:CD1	1:A:508:MET:HE3	2.28	0.68
1:B:348:LEU:O	1:B:348:LEU:HD22	1.93	0.68
1:B:111:ASP:C	1:B:384:THR:CG2	2.66	0.68
1:B:438:SER:HB3	1:B:458:SER:CB	2.24	0.68
1:B:493:TRP:CA	1:B:520:TRP:O	2.41	0.68
1:B:55:ILE:CG2	1:B:58:PHE:HD2	2.05	0.68
1:A:138:ARG:NH2	1:A:288:ASP:OD2	2.26	0.67
1:A:438:SER:CB	1:A:458:SER:CB	2.71	0.67
2:B:580:HEM:HBB2	2:B:580:HEM:CMB	2.23	0.67
1:A:54:TRP:HA	1:B:54:TRP:HZ2	1.59	0.67
1:B:309:GLN:HE21	1:B:310:HIS:CE1	2.12	0.67
2:B:580:HEM:HBB2	2:B:580:HEM:HMB2	1.76	0.67
1:B:377:HIS:CE1	1:B:378:SER:HG	2.12	0.67
1:B:90:LYS:CE	1:B:401:TYR:CE1	2.78	0.67
1:B:438:SER:CB	1:B:458:SER:HB2	2.26	0.66
1:B:496:ASP:HB2	1:B:519:ILE:CG1	2.25	0.66
1:B:157:ARG:NH2	1:B:273:ASN:O	2.29	0.66
1:A:176:PHE:O	1:A:181:LYS:HE3	1.95	0.66
1:B:221:TYR:OH	1:B:305:LEU:HD23	1.96	0.66
1:B:513:THR:HG22	1:B:514:GLU:HG2	1.77	0.66
1:A:392:THR:HG22	1:A:393:ASN:N	2.11	0.66
1:A:267:ARG:HH22	1:A:274:ARG:HH12	1.43	0.65
1:A:348:LEU:HD13	1:A:348:LEU:O	1.96	0.65
1:B:517:GLU:OE2	1:B:519:ILE:HD11	1.96	0.65
1:B:383:VAL:CG2	1:B:400:HIS:O	2.45	0.65
1:B:468:HIS:HD2	2:B:580:HEM:O2D	1.80	0.65
1:B:494:THR:CG2	1:B:495:ILE:H	2.10	0.65
1:A:389:ILE:HD12	1:A:389:ILE:H	1.62	0.64
1:B:513:THR:HG22	1:B:514:GLU:N	2.12	0.64
1:A:116:GLU:N	1:A:116:GLU:OE1	2.30	0.64
1:A:246:ARG:NH1	1:A:247:ASP:OD1	2.30	0.64
1:B:493:TRP:CB	1:B:520:TRP:O	2.45	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:213:PHE:CA	1:A:217:PHE:CE2	2.78	0.64
1:B:272:PRO:HA	4:B:2035:HOH:O	1.98	0.64
1:B:93:THR:HG1	1:B:400:HIS:CE1	2.14	0.64
1:A:391:GLU:O	1:A:391:GLU:HG3	1.98	0.64
1:B:333:ILE:O	1:B:337:VAL:HG23	1.98	0.64
3:B:590:X2N:HAA2	3:B:590:X2N:OAD	1.98	0.63
1:B:379:ILE:HB	1:B:404:VAL:CG1	2.29	0.63
1:A:513:THR:HG22	1:A:514:GLU:N	2.11	0.63
1:B:57:TRP:O	1:B:88:LEU:HA	1.98	0.63
1:B:386:PRO:O	1:B:387:LEU:HD12	1.99	0.63
1:A:51:VAL:O	1:A:53:TYR:CE1	2.52	0.63
1:A:102:GLU:HG2	1:A:457:SER:OG	1.99	0.62
1:B:191:THR:HG22	1:B:192:GLN:N	2.14	0.62
1:A:439:PHE:O	1:A:440:ASN:HB2	1.98	0.62
1:B:50:LEU:HD13	1:B:51:VAL:O	1.99	0.62
1:B:418:ASN:N	1:B:418:ASN:HD22	1.96	0.62
1:A:212:ILE:O	1:A:217:PHE:CE2	2.52	0.62
1:A:362:VAL:HG13	1:A:484:LEU:HD23	1.81	0.62
1:A:212:ILE:HG22	1:A:217:PHE:CE2	2.35	0.62
1:B:272:PRO:HB3	1:B:274:ARG:CD	2.29	0.62
1:B:441:SER:O	1:B:442:SER:OG	2.16	0.62
1:A:229:THR:N	1:A:508:MET:HE2	2.15	0.62
1:B:523:ARG:O	1:B:524:GLU:HB2	1.98	0.61
1:B:383:VAL:HG23	1:B:400:HIS:O	2.00	0.61
1:A:438:SER:HB2	1:A:458:SER:CB	2.30	0.61
1:B:380:PHE:HE1	1:B:403:LEU:HD12	1.65	0.61
1:B:496:ASP:O	1:B:498:TYR:CD1	2.53	0.61
1:B:192:GLN:NE2	1:B:320:LEU:HD11	2.14	0.61
1:B:241:PRO:HA	1:B:244:TRP:CE3	2.35	0.61
1:A:441:SER:O	1:A:455:GLY:HA3	1.99	0.61
3:B:590:X2N:NBF	3:B:590:X2N:HAA3	2.14	0.61
1:B:184:GLY:HA3	1:B:519:ILE:O	2.01	0.60
1:B:513:THR:HG22	1:B:514:GLU:H	1.66	0.60
1:A:213:PHE:C	1:A:215:ARG:H	2.07	0.60
1:B:495:ILE:O	1:B:495:ILE:HG22	2.00	0.60
1:A:212:ILE:O	1:A:217:PHE:CD2	2.54	0.60
1:B:523:ARG:O	1:B:524:GLU:CB	2.49	0.60
1:B:174:GLU:HA	1:B:177:LYS:CE	2.30	0.60
1:B:494:THR:H	1:B:520:TRP:HB2	1.67	0.60
1:A:392:THR:CG2	1:A:394:TYR:CD1	2.84	0.60
1:B:272:PRO:CB	1:B:274:ARG:CG	2.76	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:270:ILE:HG22	1:A:270:ILE:O	2.01	0.60
1:A:54:TRP:HA	1:B:54:TRP:CZ2	2.37	0.60
1:B:185:VAL:HG12	1:B:185:VAL:O	2.00	0.60
1:A:212:ILE:HG22	1:A:217:PHE:CD2	2.37	0.59
1:A:228:PHE:C	1:A:508:MET:HE2	2.28	0.59
1:A:433:LYS:HE3	1:A:474:GLN:HG2	1.83	0.59
1:B:220:LEU:HD13	1:B:252:LYS:NZ	2.17	0.59
1:A:315:THR:HG23	1:A:369:THR:CG2	2.31	0.59
1:B:272:PRO:CB	1:B:274:ARG:HG2	2.32	0.59
2:A:580:HEM:HMB2	2:A:580:HEM:CBB	2.32	0.59
1:A:221:TYR:OH	1:A:305:LEU:HD23	2.03	0.58
1:A:347:ASP:OD2	1:A:350:ASP:OD1	2.21	0.58
1:B:509:VAL:O	1:B:511:LEU:HD22	2.02	0.58
1:A:55:ILE:O	1:A:59:GLY:N	2.33	0.58
1:A:112:VAL:HB	1:A:381:ARG:HB3	1.85	0.58
1:B:443:ASP:O	1:B:454:LYS:N	2.35	0.58
1:A:116:GLU:HB3	1:A:382:LYS:HD3	1.85	0.58
1:B:66:GLN:OE1	1:B:506:SER:HB3	2.03	0.58
1:B:184:GLY:HA2	1:B:521:GLU:CB	2.26	0.58
1:B:90:LYS:HE3	1:B:401:TYR:CZ	2.37	0.58
1:B:162:ILE:HD13	1:B:199:THR:HB	1.86	0.58
1:A:226:LYS:HG2	1:A:242:HIS:NE2	2.18	0.58
1:B:131:ILE:HG22	1:B:131:ILE:O	2.03	0.58
1:B:191:THR:CG2	1:B:192:GLN:NE2	2.67	0.58
2:A:580:HEM:HBC2	2:A:580:HEM:CMC	2.32	0.57
1:A:227:GLY:HA3	1:A:246:ARG:HG3	1.85	0.57
1:B:157:ARG:NH2	1:B:274:ARG:HA	2.18	0.57
1:A:111:ASP:HB3	1:A:387:LEU:CD2	2.34	0.57
1:B:494:THR:N	1:B:520:TRP:HB2	2.20	0.57
1:B:128:THR:HA	1:B:133:ASP:OD2	2.05	0.57
1:B:439:PHE:O	1:B:440:ASN:CB	2.49	0.57
1:B:338:VAL:O	1:B:341:LEU:O	2.21	0.57
1:A:159:VAL:HB	1:A:160:PRO:HD3	1.86	0.57
1:B:224:LEU:HD22	1:B:253:ILE:HD12	1.86	0.56
1:A:45:LYS:O	1:A:393:ASN:O	2.23	0.56
1:A:102:GLU:CG	1:A:457:SER:OG	2.53	0.56
1:B:367:LYS:HZ3	1:B:431:ALA:HA	1.69	0.56
1:B:55:ILE:HG22	1:B:58:PHE:HD2	1.69	0.56
1:A:115:GLU:CD	1:A:115:GLU:H	2.12	0.56
1:B:381:ARG:NH2	2:B:580:HEM:O2A	2.38	0.56
1:B:513:THR:CG2	1:B:514:GLU:H	2.18	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:204:LEU:O	1:A:275:ASP:HB2	2.04	0.56
1:B:110:SER:O	1:B:384:THR:HG23	2.06	0.56
1:A:176:PHE:HB3	1:A:178:LEU:CD1	2.36	0.55
1:A:213:PHE:CA	1:A:217:PHE:HE2	2.19	0.55
1:B:191:THR:CG2	1:B:192:GLN:N	2.70	0.55
1:B:379:ILE:HB	1:B:404:VAL:HG13	1.87	0.55
1:A:438:SER:HB3	1:A:458:SER:CB	2.27	0.55
1:B:92:MET:HE1	1:B:380:PHE:CD1	2.42	0.55
1:B:242:HIS:HA	1:B:245:ARG:NH2	2.21	0.55
1:B:342:LYS:NZ	1:B:348:LEU:HG	2.22	0.55
1:A:221:TYR:O	1:A:225:ASP:HB2	2.07	0.55
1:A:207:ASP:O	1:A:210:ARG:HG3	2.07	0.55
1:A:494:THR:O	1:A:495:ILE:HD13	2.07	0.55
1:B:267:ARG:NH1	1:B:269:ASP:HB3	2.22	0.55
1:A:392:THR:CG2	1:A:393:ASN:N	2.70	0.54
1:A:392:THR:HG22	1:A:394:TYR:CD1	2.42	0.54
1:A:228:PHE:CG	1:A:508:MET:HE3	2.43	0.54
1:B:324:GLY:C	1:B:495:ILE:HG12	2.33	0.54
1:A:316:SER:O	1:A:320:LEU:HG	2.08	0.54
1:B:494:THR:OG1	1:B:520:TRP:CE3	2.61	0.54
1:B:376:LEU:HD22	1:B:379:ILE:HD11	1.89	0.54
3:B:590:X2N:NBF	3:B:590:X2N:CAA	2.71	0.54
1:A:346:GLY:HA3	1:A:489:TYR:OH	2.08	0.54
1:B:95:TYR:OH	1:B:99:LYS:HE3	2.08	0.54
1:B:443:ASP:OD1	1:B:454:LYS:HD3	2.08	0.54
1:A:182:THR:HA	1:A:522:LYS:HE2	1.90	0.53
1:B:57:TRP:O	1:B:57:TRP:CD2	2.61	0.53
1:B:186:ALA:O	1:B:517:GLU:HB2	2.08	0.53
1:B:404:VAL:HG22	1:B:406:PRO:HD3	1.89	0.53
1:B:92:MET:CE	1:B:403:LEU:HD12	2.39	0.53
1:B:110:SER:O	1:B:384:THR:CG2	2.56	0.53
1:A:255:ALA:HA	1:A:258:MET:HE3	1.90	0.53
1:A:52:PHE:HE2	1:B:55:ILE:HD11	1.73	0.53
1:B:342:LYS:HZ3	1:B:348:LEU:HG	1.73	0.53
1:A:393:ASN:C	1:A:394:TYR:HD1	2.16	0.53
1:A:513:THR:CG2	1:A:514:GLU:N	2.72	0.53
1:B:115:GLU:HG2	1:B:116:GLU:OE1	2.09	0.53
1:B:362:VAL:HG13	1:B:484:LEU:HD23	1.90	0.53
1:B:159:VAL:HB	1:B:160:PRO:HD3	1.90	0.53
1:A:208:GLU:O	1:A:211:ARG:HG2	2.08	0.53
1:B:51:VAL:HG22	1:B:85:PHE:HA	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:494:THR:HB	1:B:520:TRP:CE3	2.43	0.53
1:B:270:ILE:O	1:B:270:ILE:HG22	2.09	0.52
1:B:157:ARG:CZ	1:B:274:ARG:HA	2.39	0.52
1:B:439:PHE:N	4:B:2051:HOH:O	2.38	0.52
1:B:470:CYS:HB2	2:B:580:HEM:C1A	2.44	0.52
1:A:184:GLY:HA3	1:A:520:TRP:CE2	2.45	0.52
1:B:192:GLN:NE2	1:B:320:LEU:CD1	2.72	0.52
1:A:111:ASP:HB3	1:A:387:LEU:HD21	1.91	0.52
1:B:230:PRO:HD3	1:B:508:MET:HE3	1.92	0.52
1:A:348:LEU:C	1:A:348:LEU:CD1	2.82	0.52
1:B:50:LEU:HD13	1:B:51:VAL:N	2.24	0.52
1:B:380:PHE:HE1	1:B:403:LEU:CD1	2.23	0.52
1:B:77:GLN:HA	1:B:77:GLN:OE1	2.10	0.51
1:B:272:PRO:CB	1:B:278:ASP:OD2	2.51	0.51
1:B:377:HIS:CE1	1:B:378:SER:OG	2.63	0.51
1:A:62:ALA:O	1:A:66:GLN:CG	2.58	0.51
1:A:82:VAL:HA	1:A:94:VAL:O	2.10	0.51
1:B:513:THR:CG2	1:B:514:GLU:N	2.73	0.51
1:A:341:LEU:N	1:A:341:LEU:HD13	2.24	0.51
1:B:347:ASP:O	1:B:348:LEU:HB3	2.10	0.51
1:B:418:ASN:HD22	1:B:418:ASN:H	1.57	0.51
1:A:213:PHE:CE1	1:A:218:ALA:HB1	2.46	0.51
1:A:438:SER:CB	1:A:458:SER:HA	2.40	0.51
1:B:377:HIS:CE1	3:B:590:X2N:HAY1	2.46	0.51
1:B:174:GLU:N	1:B:177:LYS:HE2	2.26	0.51
1:A:324:GLY:C	1:A:495:ILE:CG1	2.82	0.51
1:B:189:MET:HE1	1:B:512:PRO:HD2	1.93	0.51
1:A:315:THR:HG23	1:A:369:THR:HG23	1.92	0.51
1:A:495:ILE:HD11	1:A:518:ILE:HG22	1.92	0.51
1:B:162:ILE:O	1:B:166:ILE:HG13	2.10	0.51
1:B:514:GLU:CG	1:B:515:PRO:CD	2.85	0.51
1:A:513:THR:HG22	1:A:514:GLU:H	1.76	0.50
1:B:417:ASP:O	1:B:426:ARG:NH2	2.44	0.50
1:B:342:LYS:HZ3	1:B:342:LYS:HB2	1.75	0.50
1:A:336:GLU:OE1	1:A:361:SER:OG	2.29	0.50
1:B:258:MET:O	1:B:262:LYS:HG2	2.12	0.50
1:B:344:LYS:HG3	1:B:345:GLY:N	2.27	0.50
1:A:208:GLU:O	1:A:212:ILE:HG12	2.10	0.50
1:A:315:THR:HG23	1:A:369:THR:HG21	1.92	0.50
1:A:393:ASN:N	1:A:393:ASN:OD1	2.34	0.50
1:B:277:ILE:O	1:B:281:LEU:HG	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:386:PRO:O	1:A:388:ARG:NH1	2.45	0.50
1:B:112:VAL:N	1:B:384:THR:HG22	2.27	0.50
1:B:201:SER:O	1:B:205:PHE:HB2	2.11	0.50
1:B:365:THR:O	1:B:369:THR:HG22	2.12	0.50
1:A:278:ASP:HA	1:A:281:LEU:HD12	1.94	0.49
1:B:92:MET:HE2	1:B:403:LEU:HD12	1.94	0.49
1:B:189:MET:HE1	1:B:511:LEU:HB3	1.94	0.49
1:A:128:THR:CG2	1:A:133:ASP:OD2	2.58	0.49
1:B:50:LEU:C	1:B:50:LEU:CD1	2.83	0.49
1:B:483:ILE:O	1:B:487:PHE:HD1	1.94	0.49
1:B:494:THR:CB	1:B:520:TRP:CE3	2.95	0.49
1:B:376:LEU:HD22	1:B:379:ILE:CD1	2.43	0.49
1:B:187:ASN:OD1	1:B:516:ALA:O	2.31	0.49
1:B:272:PRO:O	1:B:273:ASN:OD1	2.30	0.49
1:A:239:PRO:O	1:A:244:TRP:CZ2	2.66	0.49
2:A:580:HEM:HBC2	2:A:580:HEM:HMC2	1.92	0.49
1:A:54:TRP:CA	1:B:54:TRP:HZ2	2.26	0.49
1:B:494:THR:HB	1:B:520:TRP:CD2	2.48	0.49
1:A:53:TYR:CB	1:A:59:GLY:HA2	2.41	0.49
2:B:580:HEM:HMB2	2:B:580:HEM:CBB	2.42	0.49
1:A:213:PHE:CE1	1:A:218:ALA:CB	2.96	0.49
1:A:513:THR:CG2	1:A:514:GLU:H	2.26	0.49
1:A:189:MET:HE1	1:A:511:LEU:HB3	1.94	0.49
1:B:50:LEU:CD1	1:B:51:VAL:O	2.61	0.49
1:A:51:VAL:HG22	1:A:84:SER:O	2.13	0.48
1:A:61:ALA:O	1:A:65:GLY:N	2.37	0.48
1:A:226:LYS:HG2	1:A:242:HIS:CD2	2.47	0.48
1:A:176:PHE:HB3	1:A:178:LEU:HD13	1.95	0.48
1:A:403:LEU:C	1:A:403:LEU:HD23	2.38	0.48
1:B:123:THR:HB	1:B:124:PRO:HD3	1.95	0.48
1:A:522:LYS:HE3	1:A:525:THR:HA	1.95	0.48
1:B:272:PRO:CA	1:B:274:ARG:HG2	2.42	0.48
1:B:329:LEU:O	1:B:332:VAL:CG1	2.55	0.48
1:B:368:GLU:HA	1:B:368:GLU:OE1	2.12	0.48
1:B:234:VAL:HG23	1:B:235:PHE:N	2.28	0.48
1:B:270:ILE:CA	1:B:272:PRO:HD2	2.42	0.48
1:B:438:SER:CB	1:B:458:SER:CB	2.88	0.48
1:A:414:ARG:O	1:A:414:ARG:HG3	2.13	0.48
1:B:418:ASN:N	1:B:418:ASN:ND2	2.62	0.48
1:A:209:MET:HE1	1:A:256:THR:HG22	1.96	0.48
1:B:95:TYR:CG	1:B:100:GLY:HA2	2.48	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:483:ILE:O	1:A:487:PHE:HD1	1.96	0.48
1:A:93:THR:O	1:A:93:THR:HG23	2.14	0.48
1:B:225:ASP:OD1	1:B:310:HIS:HE1	1.97	0.48
1:B:380:PHE:CE1	1:B:403:LEU:HD12	2.47	0.48
1:B:494:THR:CG2	1:B:495:ILE:N	2.64	0.47
1:A:506:SER:HB2	3:A:590:X2N:HAT	1.95	0.47
1:A:189:MET:CE	1:A:512:PRO:HG2	2.44	0.47
1:A:229:THR:O	1:A:232:ASN:ND2	2.33	0.47
1:B:309:GLN:NE2	1:B:310:HIS:ND1	2.57	0.47
1:A:319:PHE:CD1	1:A:319:PHE:C	2.93	0.47
1:B:140:MET:HE2	1:B:140:MET:HA	1.96	0.47
1:A:216:SER:HA	4:A:2027:HOH:O	2.15	0.47
1:B:263:LEU:O	1:B:267:ARG:CG	2.51	0.47
1:A:338:VAL:HA	1:A:341:LEU:HD22	1.97	0.47
1:B:57:TRP:O	1:B:57:TRP:CE3	2.68	0.47
1:B:65:GLY:HA2	3:B:590:X2N:HAO	1.97	0.47
1:A:376:LEU:HD11	3:A:590:X2N:HAH	1.97	0.47
1:A:141:GLU:OE2	1:A:287:LYS:HG3	2.15	0.46
1:A:344:LYS:C	1:A:345:GLY:O	2.56	0.46
1:A:396:VAL:O	1:A:397:PRO:C	2.56	0.46
1:B:514:GLU:HG3	1:B:515:PRO:CD	2.42	0.46
1:A:286:TYR:OH	1:A:292:MET:HG2	2.15	0.46
1:B:98:PRO:HB3	1:B:457:SER:O	2.15	0.46
1:A:116:GLU:CB	1:A:382:LYS:HD3	2.46	0.46
1:A:436:SER:OG	1:A:442:SER:HB2	2.16	0.46
1:B:220:LEU:HD13	1:B:252:LYS:CE	2.44	0.46
1:B:386:PRO:HA	1:B:396:VAL:O	2.15	0.46
1:B:496:ASP:HB2	1:B:519:ILE:CB	2.46	0.46
1:A:237:ASN:C	1:A:238:LEU:HD12	2.40	0.46
1:A:240:LEU:HB3	1:A:241:PRO:HD2	1.97	0.46
1:A:374:MET:HE3	1:A:462:PRO:O	2.15	0.46
1:B:182:THR:O	1:B:182:THR:HG22	2.15	0.46
1:B:272:PRO:O	1:B:273:ASN:C	2.59	0.46
1:B:309:GLN:NE2	1:B:310:HIS:CE1	2.82	0.46
1:B:403:LEU:HD23	1:B:403:LEU:C	2.41	0.46
1:B:438:SER:CB	1:B:458:SER:CA	2.94	0.46
1:A:258:MET:O	1:A:262:LYS:HG2	2.16	0.46
1:B:438:SER:HB2	1:B:458:SER:HA	1.97	0.46
1:A:341:LEU:HG	1:A:347:ASP:HA	1.98	0.46
1:A:385:ASN:O	1:A:387:LEU:CD2	2.64	0.46
1:B:217:PHE:HA	1:B:220:LEU:HD12	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:74:SER:O	1:A:77:GLN:HG2	2.15	0.46
1:A:113:SER:OG	1:A:116:GLU:OE1	2.32	0.46
1:A:429:THR:O	1:A:429:THR:HG22	2.16	0.46
1:B:214:ASP:O	1:B:215:ARG:HD3	2.16	0.46
1:B:270:ILE:CD1	1:B:281:LEU:HD13	2.44	0.46
1:B:508:MET:HA	1:B:508:MET:CE	2.39	0.46
1:A:386:PRO:C	1:A:387:LEU:HD23	2.41	0.45
1:B:111:ASP:CA	1:B:384:THR:CG2	2.94	0.45
1:B:176:PHE:O	1:B:177:LYS:C	2.59	0.45
1:B:451:LYS:HD3	1:B:451:LYS:N	2.31	0.45
1:A:58:PHE:HZ	1:A:62:ALA:HB2	1.82	0.45
1:B:496:ASP:O	1:B:498:TYR:HD1	1.98	0.45
1:A:236:PRO:O	1:A:246:ARG:NH2	2.49	0.45
1:A:315:THR:CG2	1:A:369:THR:HG21	2.47	0.45
1:A:406:PRO:O	1:A:410:HIS:CD2	2.70	0.45
1:B:272:PRO:C	1:B:274:ARG:HG2	2.41	0.45
1:A:438:SER:HB2	1:A:458:SER:HB2	1.98	0.45
1:B:349:ASN:O	1:B:350:ASP:HB2	2.17	0.45
1:B:91:ILE:HG23	1:B:400:HIS:CD2	2.52	0.45
1:B:271:ASP:N	1:B:272:PRO:HD3	2.27	0.45
1:B:469:ARG:NH2	1:B:473:GLU:OE1	2.31	0.45
1:B:111:ASP:HA	1:B:384:THR:HG23	1.99	0.45
1:B:221:TYR:O	1:B:225:ASP:N	2.44	0.45
1:B:261:ILE:HG23	1:B:281:LEU:HD21	1.99	0.45
1:B:492:ARG:NH1	4:B:2060:HOH:O	2.49	0.45
1:A:118:TYR:HD2	1:A:132:TYR:CZ	2.35	0.45
1:A:121:LEU:HD22	1:A:233:PHE:CZ	2.51	0.45
1:A:469:ARG:NH1	1:A:470:CYS:O	2.47	0.45
1:B:90:LYS:CE	1:B:401:TYR:CZ	2.99	0.45
1:B:376:LEU:CD2	1:B:379:ILE:CD1	2.95	0.45
1:A:87:LEU:N	1:A:90:LYS:O	2.47	0.45
1:A:371:ARG:O	1:A:410:HIS:HB3	2.16	0.45
1:A:439:PHE:CD1	1:A:440:ASN:N	2.80	0.45
1:A:239:PRO:O	1:A:244:TRP:CH2	2.71	0.44
1:A:389:ILE:HD12	1:A:389:ILE:N	2.29	0.44
1:B:344:LYS:HG3	1:B:345:GLY:H	1.82	0.44
1:B:348:LEU:O	1:B:349:ASN:CB	2.61	0.44
1:B:384:THR:HG23	1:B:385:ASN:N	2.32	0.44
1:B:436:SER:HB3	1:B:442:SER:HB2	1.98	0.44
1:A:217:PHE:HB2	1:A:220:LEU:HD12	1.99	0.44
1:A:438:SER:CB	1:A:458:SER:CA	2.95	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:215:ARG:C	1:B:217:PHE:N	2.74	0.44
1:B:470:CYS:HA	2:B:580:HEM:C4D	2.52	0.44
1:A:52:PHE:CE2	1:B:55:ILE:HD11	2.52	0.44
1:B:496:ASP:CB	1:B:519:ILE:HG13	2.42	0.44
1:A:230:PRO:O	1:A:233:PHE:HD2	2.00	0.44
1:B:133:ASP:O	1:B:133:ASP:OD1	2.35	0.44
1:A:51:VAL:HG11	1:A:79:TYR:CE2	2.53	0.44
1:A:189:MET:HE2	1:A:512:PRO:HG2	1.98	0.44
1:B:272:PRO:HB3	1:B:274:ARG:HG3	1.92	0.44
1:B:210:ARG:HD3	1:B:210:ARG:C	2.43	0.44
1:B:371:ARG:O	1:B:410:HIS:HB3	2.18	0.44
1:A:213:PHE:C	1:A:215:ARG:N	2.72	0.44
1:B:436:SER:CB	1:B:442:SER:HB2	2.48	0.44
1:B:439:PHE:HD1	1:B:439:PHE:HA	1.69	0.44
1:A:53:TYR:C	1:B:54:TRP:CZ2	2.96	0.43
1:A:115:GLU:OE1	1:A:116:GLU:OE1	2.35	0.43
1:A:392:THR:CG2	1:A:394:TYR:CG	3.01	0.43
3:A:590:X2N:HAO	3:A:590:X2N:OAC	2.17	0.43
1:A:84:SER:HA	1:A:92:MET:O	2.18	0.43
1:A:407:GLY:O	1:A:410:HIS:HB2	2.18	0.43
1:B:309:GLN:HG3	1:B:310:HIS:N	2.32	0.43
1:A:229:THR:OG1	1:A:231:ILE:HD12	2.19	0.43
1:A:53:TYR:C	1:B:54:TRP:HZ2	2.26	0.43
1:B:215:ARG:O	1:B:217:PHE:N	2.52	0.43
1:A:132:TYR:HA	4:A:2013:HOH:O	2.19	0.43
1:B:383:VAL:CG2	1:B:402:VAL:HG23	2.49	0.43
1:A:111:ASP:HB3	1:A:387:LEU:HD22	1.99	0.43
1:B:272:PRO:CB	1:B:274:ARG:NE	2.72	0.43
1:A:324:GLY:CA	1:A:495:ILE:HG12	2.49	0.43
1:A:388:ARG:HG2	1:A:395:ILE:CD1	2.49	0.43
1:A:394:TYR:CD1	1:A:394:TYR:N	2.87	0.43
1:B:93:THR:OG1	1:B:400:HIS:NE2	2.37	0.43
1:B:264:ARG:CG	1:B:267:ARG:HH21	2.28	0.43
1:B:272:PRO:CB	1:B:274:ARG:HG3	2.48	0.43
1:A:99:LYS:HA	1:A:99:LYS:HE2	2.00	0.43
1:B:264:ARG:HA	1:B:267:ARG:NE	2.34	0.43
1:B:116:GLU:HB3	1:B:382:LYS:HD3	2.01	0.43
1:A:229:THR:C	1:A:231:ILE:H	2.26	0.43
1:A:272:PRO:C	1:A:274:ARG:H	2.26	0.43
1:B:470:CYS:HB2	2:B:580:HEM:NA	2.34	0.43
1:A:131:ILE:O	1:A:132:TYR:CB	2.67	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:466:GLY:O	1:A:469:ARG:HB2	2.18	0.42
1:A:520:TRP:HZ3	1:A:522:LYS:HG2	1.84	0.42
1:B:416:PHE:CD1	1:B:417:ASP:N	2.70	0.42
1:A:123:THR:N	1:A:124:PRO:HD2	2.34	0.42
1:A:347:ASP:CG	1:A:350:ASP:OD1	2.61	0.42
1:B:257:TYR:O	1:B:261:ILE:HG13	2.19	0.42
1:B:438:SER:OG	1:B:458:SER:HB2	2.19	0.42
1:B:462:PRO:HG2	1:B:463:PHE:CD2	2.54	0.42
1:A:131:ILE:O	1:A:132:TYR:HB2	2.18	0.42
1:A:230:PRO:HG2	1:A:507:SER:HA	2.01	0.42
1:A:377:HIS:NE2	3:A:590:X2N:HAZ2	2.34	0.42
1:B:131:ILE:O	1:B:132:TYR:CB	2.68	0.42
1:A:263:LEU:O	1:A:267:ARG:HB3	2.19	0.42
1:B:236:PRO:HB2	1:B:237:ASN:OD1	2.20	0.42
1:A:272:PRO:C	1:A:274:ARG:N	2.77	0.42
1:A:324:GLY:O	1:A:495:ILE:HG12	2.19	0.42
1:B:128:THR:HG22	1:B:129:GLY:N	2.34	0.42
1:B:351:LEU:HD23	1:B:352:THR:N	2.35	0.42
1:A:197:ILE:HD13	1:A:309:GLN:HG3	2.01	0.42
1:A:371:ARG:NH2	1:A:372:MET:HE2	2.34	0.42
1:A:495:ILE:O	1:A:495:ILE:CG2	2.60	0.42
1:B:381:ARG:HH21	2:B:580:HEM:CGA	2.31	0.42
1:A:385:ASN:O	1:A:387:LEU:HD23	2.20	0.42
1:B:114:ALA:HB3	1:B:468:HIS:NE2	2.34	0.42
1:B:379:ILE:N	1:B:379:ILE:HD12	2.35	0.42
1:A:179:LYS:HE2	1:A:528:PHE:HB3	2.01	0.42
1:A:340:LEU:C	1:A:340:LEU:HD23	2.45	0.42
1:B:92:MET:HB3	1:B:92:MET:HE3	1.69	0.42
1:B:270:ILE:HD13	1:B:281:LEU:CD1	2.47	0.42
1:B:496:ASP:HB2	1:B:519:ILE:HB	2.01	0.42
1:B:184:GLY:CA	1:B:521:GLU:H	2.33	0.42
1:B:344:LYS:O	1:B:345:GLY:C	2.62	0.42
1:A:50:LEU:HA	1:A:84:SER:HB2	2.02	0.41
1:A:55:ILE:O	1:A:58:PHE:HB3	2.20	0.41
1:A:67:GLN:HG3	1:A:70:GLU:CD	2.45	0.41
1:B:131:ILE:HD11	1:B:300:LEU:HD12	2.01	0.41
1:A:282:ILE:HG13	1:A:283:HIS:CD2	2.55	0.41
1:B:438:SER:CB	1:B:458:SER:HA	2.50	0.41
1:A:215:ARG:HG2	1:A:216:SER:N	2.29	0.41
1:A:231:ILE:HG23	1:A:235:PHE:HD2	1.85	0.41
1:A:233:PHE:HZ	3:A:590:X2N:CAK	2.34	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:523:ARG:C	1:A:524:GLU:HG3	2.45	0.41
1:B:272:PRO:HB3	1:B:274:ARG:CZ	2.48	0.41
1:A:209:MET:HE2	1:A:257:TYR:CD1	2.56	0.41
1:A:229:THR:C	1:A:231:ILE:N	2.79	0.41
1:A:261:ILE:HG23	1:A:281:LEU:HD21	2.03	0.41
1:B:81:ASP:OD1	1:B:81:ASP:N	2.53	0.41
1:B:452:VAL:HG21	1:B:469:ARG:HD2	2.02	0.41
1:A:53:TYR:O	1:B:54:TRP:CZ2	2.74	0.41
1:A:435:ASN:OD1	1:A:436:SER:N	2.47	0.41
1:A:51:VAL:CG1	1:A:79:TYR:CE2	3.02	0.41
1:A:140:MET:O	1:A:143:LYS:HB2	2.21	0.41
1:B:277:ILE:HG12	1:B:301:LEU:HD21	2.02	0.41
1:B:492:ARG:HB3	1:B:522:LYS:HE3	2.03	0.41
1:B:521:GLU:OE1	1:B:521:GLU:O	2.39	0.41
1:A:121:LEU:HD11	1:A:508:MET:SD	2.61	0.41
1:B:214:ASP:O	1:B:215:ARG:CG	2.69	0.41
1:A:392:THR:CG2	1:A:393:ASN:H	2.33	0.41
1:B:344:LYS:C	1:B:347:ASP:H	2.29	0.41
1:B:348:LEU:O	1:B:348:LEU:CD2	2.67	0.41
1:B:469:ARG:NH1	1:B:470:CYS:O	2.50	0.41
1:A:454:LYS:HD3	1:A:455:GLY:N	2.36	0.41
1:B:214:ASP:O	1:B:215:ARG:HG2	2.21	0.41
1:B:272:PRO:HG3	1:B:274:ARG:NH2	2.36	0.41
1:A:319:PHE:C	1:A:319:PHE:HD1	2.29	0.40
1:A:324:GLY:HA3	1:A:495:ILE:CG1	2.51	0.40
1:A:368:GLU:HA	1:A:368:GLU:OE1	2.21	0.40
1:B:51:VAL:HG13	1:B:84:SER:O	2.20	0.40
1:B:55:ILE:HG21	1:B:58:PHE:HD2	1.80	0.40
1:B:272:PRO:HG3	1:B:274:ARG:CZ	2.51	0.40
1:A:316:SER:HA	1:A:319:PHE:CZ	2.56	0.40
1:B:86:MET:HE2	1:B:89:GLY:H	1.87	0.40
1:B:184:GLY:HA3	1:B:521:GLU:H	1.87	0.40
1:B:468:HIS:CD2	2:B:580:HEM:O2D	2.68	0.40
1:B:187:ASN:C	1:B:189:MET:N	2.79	0.40
1:B:209:MET:HE1	1:B:256:THR:HG22	2.03	0.40
1:B:249:ALA:O	1:B:253:ILE:HG13	2.21	0.40
1:A:341:LEU:C	1:A:342:LYS:HG3	2.46	0.40
1:B:433:LYS:O	1:B:434:ALA:HB2	2.22	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:A:53:TYR:OH	1:A:237:ASN:O[2_756]	2.03	0.17

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	482/490 (98%)	455 (94%)	25 (5%)	2 (0%)	30	48
1	B	481/490 (98%)	448 (93%)	28 (6%)	5 (1%)	12	25
All	All	963/980 (98%)	903 (94%)	53 (6%)	7 (1%)	18	35

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	524	GLU
1	A	342	LYS
1	B	440	ASN
1	B	216	SER
1	B	350	ASP
1	B	273	ASN
1	A	131	ILE

5.3.2 Protein sidechains [i](#)

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	431/436 (99%)	413 (96%)	18 (4%)	26	51

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	430/436 (99%)	407 (95%)	23 (5%)	20	42
All	All	861/872 (99%)	820 (95%)	41 (5%)	23	46

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	50	LEU
1	A	53	TYR
1	A	67	GLN
1	A	82	VAL
1	A	86	MET
1	A	217	PHE
1	A	221	TYR
1	A	267	ARG
1	A	314	SER
1	A	323	LEU
1	A	341	LEU
1	A	376	LEU
1	A	393	ASN
1	A	439	PHE
1	A	452	VAL
1	A	461	LEU
1	A	494	THR
1	A	498	TYR
1	B	50	LEU
1	B	54	TRP
1	B	86	MET
1	B	152	THR
1	B	191	THR
1	B	221	TYR
1	B	274	ARG
1	B	339	GLU
1	B	348	LEU
1	B	369	THR
1	B	387	LEU
1	B	414	ARG
1	B	418	ASN
1	B	433	LYS
1	B	436	SER
1	B	437	VAL
1	B	439	PHE
1	B	443	ASP

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Mol	Chain	Res	Type
1	B	452	VAL
1	B	508	MET
1	B	511	LEU
1	B	514	GLU
1	B	521	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	120	HIS
1	A	187	ASN
1	A	490	ASN
1	B	187	ASN
1	B	192	GLN
1	B	273	ASN
1	B	295	GLN
1	B	309	GLN
1	B	310	HIS
1	B	357	GLN
1	B	363	ASN
1	B	418	ASN

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	$\# Z > 2$
3	X2N	A	590	2	55,57,57	1.79	9 (16%)	68,82,82	2.27	24 (35%)
2	HEM	A	580	1,3	50,50,50	2.13	20 (40%)	67,82,82	1.78	17 (25%)
3	X2N	B	590	2	55,57,57	1.75	8 (14%)	68,82,82	2.32	28 (41%)
2	HEM	B	580	1,3	50,50,50	1.91	17 (34%)	67,82,82	2.04	20 (29%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	X2N	A	590	2	-	10/38/59/59	0/7/7/7
2	HEM	A	580	1,3	-	0/14/54/54	-
3	X2N	B	590	2	-	8/38/59/59	0/7/7/7
2	HEM	B	580	1,3	-	4/14/54/54	-

All (54) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	590	X2N	CAR-CBI	7.14	1.48	1.39
3	A	590	X2N	CAR-CBI	7.07	1.48	1.39
3	A	590	X2N	CBN-NBW	-5.91	1.34	1.44
3	B	590	X2N	CBN-NBW	-5.62	1.34	1.44
2	A	580	HEM	C4D-ND	-5.23	1.31	1.40
3	A	590	X2N	CBP-NBW	-4.95	1.33	1.39
2	A	580	HEM	FE-ND	-4.66	1.80	1.94
3	B	590	X2N	CBP-NBW	-4.50	1.33	1.39
2	B	580	HEM	FE-ND	-4.28	1.81	1.94
2	A	580	HEM	C1B-NB	-4.10	1.33	1.40
2	A	580	HEM	C1D-ND	-3.99	1.31	1.38
2	B	580	HEM	C4D-ND	-3.94	1.33	1.40
3	A	590	X2N	CAT-NBF	3.85	1.36	1.30
2	B	580	HEM	C1B-NB	-3.83	1.33	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	580	HEM	C1C-C2C	-3.76	1.37	1.45
3	B	590	X2N	CAT-NBF	3.74	1.36	1.30
2	B	580	HEM	C1C-C2C	-3.73	1.38	1.45
2	A	580	HEM	FE-NB	3.67	2.06	1.94
2	A	580	HEM	C3C-C4C	-3.60	1.39	1.46
2	B	580	HEM	FE-NB	3.46	2.05	1.94
3	A	590	X2N	OAC-CBP	3.45	1.29	1.22
3	B	590	X2N	OAC-CBP	3.39	1.28	1.22
2	B	580	HEM	C1D-ND	-3.20	1.32	1.38
2	B	580	HEM	C3C-C4C	-3.19	1.40	1.46
2	B	580	HEM	C4B-NB	-3.16	1.32	1.38
2	A	580	HEM	C4B-NB	-3.07	1.32	1.38
2	A	580	HEM	C1C-NC	-2.91	1.34	1.39
2	B	580	HEM	C1C-NC	-2.81	1.34	1.39
2	A	580	HEM	C1A-C2A	-2.79	1.39	1.44
2	A	580	HEM	C4C-NC	-2.69	1.34	1.39
2	A	580	HEM	C1A-NA	-2.60	1.34	1.39
2	B	580	HEM	C4C-NC	-2.59	1.34	1.39
2	B	580	HEM	C3D-C2D	-2.57	1.31	1.36
2	A	580	HEM	C4A-NA	-2.45	1.35	1.39
2	B	580	HEM	FE-NC	2.44	2.03	1.95
2	A	580	HEM	O2D-CGD	-2.43	1.22	1.30
2	A	580	HEM	O2A-CGA	-2.40	1.22	1.30
2	A	580	HEM	FE-NC	2.37	2.03	1.95
3	B	590	X2N	CAT-NBW	-2.35	1.34	1.37
2	A	580	HEM	C2A-C3A	-2.30	1.32	1.38
2	B	580	HEM	O2A-CGA	-2.30	1.23	1.30
2	A	580	HEM	CHA-C1A	-2.27	1.34	1.39
3	A	590	X2N	CAT-NBW	-2.27	1.34	1.37
2	A	580	HEM	C3D-C2D	-2.24	1.31	1.36
3	A	590	X2N	NBV-NBE	2.24	1.40	1.36
3	B	590	X2N	NBV-NBE	2.23	1.40	1.36
3	A	590	X2N	NBX-NBF	2.22	1.41	1.37
2	B	580	HEM	CHA-C1A	-2.17	1.34	1.39
2	A	580	HEM	C1B-C2B	-2.16	1.40	1.44
2	B	580	HEM	C1A-C2A	-2.16	1.40	1.44
3	B	590	X2N	NBX-NBF	2.16	1.41	1.37
2	B	580	HEM	O2D-CGD	-2.13	1.23	1.30
2	B	580	HEM	C2A-C3A	-2.09	1.33	1.38
3	A	590	X2N	CAQ-NBE	2.00	1.35	1.32

All (89) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	590	X2N	CBP-NBX-NBF	-7.42	107.04	112.87
3	B	590	X2N	CBP-NBX-NBF	-7.39	107.07	112.87
2	B	580	HEM	CHC-C4B-NB	6.64	131.57	124.42
3	B	590	X2N	CAO-CBN-NBW	5.51	126.81	119.62
3	A	590	X2N	CAT-NBW-CBP	5.35	111.23	107.27
2	B	580	HEM	CHA-C4D-ND	5.31	130.94	124.37
2	A	580	HEM	CHD-C1D-ND	5.25	130.07	124.42
2	B	580	HEM	CHD-C1D-ND	5.00	129.81	124.42
3	B	590	X2N	NBW-CBP-NBX	4.87	108.37	102.88
3	A	590	X2N	NBW-CBP-NBX	4.66	108.14	102.88
2	A	580	HEM	CHC-C4B-NB	4.58	129.35	124.42
3	B	590	X2N	CAT-NBW-CBP	4.43	110.54	107.27
2	A	580	HEM	CHA-C4D-ND	4.04	129.37	124.37
3	B	590	X2N	CBS-NBX-CBP	3.96	135.79	125.26
3	B	590	X2N	CAG-CBI-CAR	-3.93	119.60	123.88
2	A	580	HEM	CHB-C1B-NB	3.93	129.22	124.37
3	A	590	X2N	C2-CAR-CBI	-3.91	119.07	122.74
3	A	590	X2N	CBS-NBX-CBP	3.83	135.45	125.26
3	A	590	X2N	CAG-CBI-CAR	-3.74	119.81	123.88
3	B	590	X2N	CAL-CBM-NBU	-3.66	116.30	121.39
2	B	580	HEM	CHA-C4D-C3D	-3.61	118.56	125.23
3	A	590	X2N	CBC-NBV-NBE	-3.61	116.00	121.25
3	B	590	X2N	CBS-NBX-NBF	-3.58	117.14	121.28
3	A	590	X2N	CAT-NBF-NBX	3.56	107.20	103.68
3	A	590	X2N	CAX-CAZ-NBU	-3.38	103.67	110.78
3	A	590	X2N	CAL-CBM-NBU	-3.38	116.70	121.39
2	A	580	HEM	CHD-C1D-C2D	-3.36	119.72	125.03
2	B	580	HEM	C1B-NB-C4B	3.32	109.14	105.21
3	B	590	X2N	CAT-NBF-NBX	3.30	106.94	103.68
3	A	590	X2N	CBS-NBX-NBF	-3.28	117.50	121.28
3	B	590	X2N	CBC-NBV-NBE	-3.25	116.52	121.25
3	A	590	X2N	CBJ-CBO-CAP	3.23	121.69	118.38
3	B	590	X2N	CAX-NBT-CBL	-3.21	109.36	118.11
2	B	580	HEM	CAD-C3D-C4D	3.14	130.18	124.70
3	A	590	X2N	CAX-NBT-CBL	-3.14	109.56	118.11
2	A	580	HEM	CHA-C4D-C3D	-3.04	119.62	125.23
3	B	590	X2N	CBJ-CBO-CAP	3.04	121.50	118.38
2	B	580	HEM	CHB-C1B-NB	2.97	128.04	124.37
3	B	590	X2N	C2-CAR-CBI	-2.97	119.95	122.74
2	B	580	HEM	CHD-C4C-NC	2.96	127.67	124.45
3	A	590	X2N	CBI-CAG-CAP	2.94	119.80	116.67
2	B	580	HEM	C1A-CHA-C4D	-2.89	119.44	126.25
3	A	590	X2N	CAO-CBN-NBW	2.89	123.39	119.62

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	590	X2N	CAM-CBM-NBU	2.87	125.38	121.39
3	B	590	X2N	CBI-CAG-CAP	2.87	119.72	116.67
2	A	580	HEM	O2A-CGA-O1A	-2.80	116.14	123.33
3	A	590	X2N	CAW-CAY-NBU	-2.79	104.91	110.78
3	A	590	X2N	NBW-CAT-NBF	-2.79	106.16	111.42
3	B	590	X2N	CBN-NBW-CBP	2.75	128.61	125.87
2	B	580	HEM	CAA-C2A-C1A	2.73	130.28	124.94
2	B	580	HEM	O2A-CGA-O1A	-2.72	116.33	123.33
2	B	580	HEM	CHD-C1D-C2D	-2.65	120.84	125.03
3	B	590	X2N	CBN-NBW-CAT	-2.63	120.83	125.28
2	B	580	HEM	O2D-CGD-O1D	-2.61	116.63	123.33
3	B	590	X2N	CAJ-CBL-NBT	-2.60	117.78	121.39
3	B	590	X2N	CAZ-NBU-CBM	-2.60	111.03	118.11
3	B	590	X2N	CAX-CAZ-NBU	-2.56	105.40	110.78
2	B	580	HEM	C4A-C3A-C2A	2.56	109.75	106.82
2	A	580	HEM	C1B-NB-C4B	2.55	108.23	105.21
2	B	580	HEM	O2D-CGD-CBD	2.53	122.01	114.00
2	A	580	HEM	C4C-CHD-C1D	-2.51	120.69	126.02
3	A	590	X2N	CAZ-NBU-CBM	-2.50	111.28	118.11
3	B	590	X2N	CAS-NBD-CAQ	2.50	109.61	102.85
2	B	580	HEM	C4C-CHD-C1D	-2.49	120.72	126.02
3	A	590	X2N	CAS-NBD-CAQ	2.47	109.53	102.85
2	B	580	HEM	CHC-C4B-C3B	-2.43	120.22	125.07
3	A	590	X2N	CBC-NBV-CAS	2.39	133.51	128.82
2	A	580	HEM	O2A-CGA-CBA	2.39	121.56	114.00
2	A	580	HEM	CHD-C4C-NC	2.39	127.05	124.45
2	B	580	HEM	C1C-CHC-C4B	-2.35	121.03	126.02
2	A	580	HEM	CHB-C1B-C2B	-2.31	120.39	126.95
3	B	590	X2N	NBW-CAT-NBF	-2.30	107.08	111.42
2	A	580	HEM	C1A-CHA-C4D	-2.25	120.95	126.25
2	A	580	HEM	C4A-CHB-C1B	-2.24	120.97	126.25
3	B	590	X2N	CAL-CAN-CBN	2.19	123.07	120.30
3	B	590	X2N	CAO-CBN-CAN	-2.18	115.02	119.18
2	A	580	HEM	CBD-CAD-C3D	2.17	118.52	112.53
3	B	590	X2N	CAM-CAO-CBN	2.16	123.04	120.30
3	A	590	X2N	NBE-CAQ-NBD	-2.14	108.80	114.90
3	A	590	X2N	CAJ-CBL-NBT	-2.12	118.45	121.39
3	B	590	X2N	NBE-CAQ-NBD	-2.11	108.88	114.90
3	A	590	X2N	CAM-CBM-NBU	2.09	124.29	121.39
2	B	580	HEM	CBD-CAD-C3D	2.09	118.30	112.53
3	B	590	X2N	CBC-NBV-CAS	2.07	132.88	128.82
2	A	580	HEM	O2D-CGD-O1D	-2.06	118.04	123.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	580	HEM	C1C-CHC-C4B	-2.05	121.66	126.02
3	B	590	X2N	NBV-CAS-NBD	-2.04	106.21	110.42
3	A	590	X2N	CBN-NBW-CBP	-2.01	123.86	125.87
2	B	580	HEM	CMD-C2D-C1D	2.00	128.16	125.03

There are no chirality outliers.

All (22) torsion outliers are listed below:

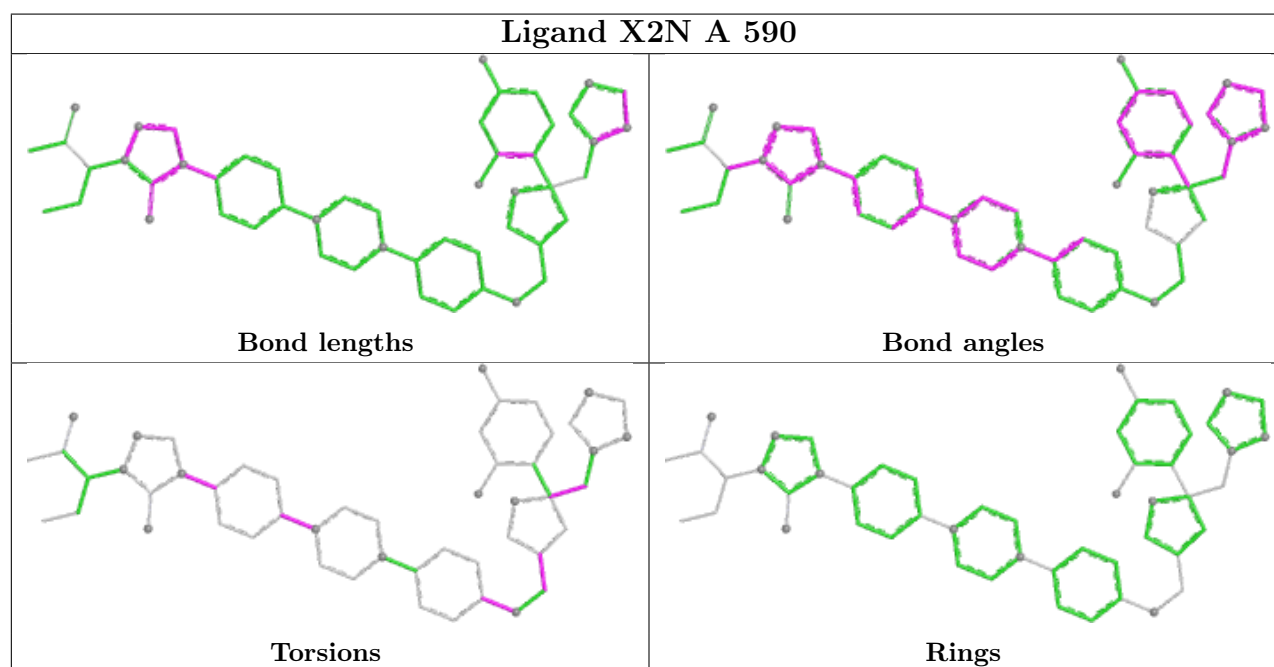
Mol	Chain	Res	Type	Atoms
3	A	590	X2N	CAV-C7-CBB-OBG
3	B	590	X2N	CAV-C7-CBB-OBG
3	B	590	X2N	CAH-CBK-OBG-CBB
3	B	590	X2N	CAI-CBK-OBG-CBB
3	A	590	X2N	CAH-CBK-OBG-CBB
3	A	590	X2N	CAI-CBK-OBG-CBB
3	A	590	X2N	OBH-C2-CBC-NBV
3	A	590	X2N	CAR-C2-CBC-NBV
3	B	590	X2N	CAR-C2-CBC-NBV
3	A	590	X2N	C32-C7-CBB-OBG
3	B	590	X2N	C32-C7-CBB-OBG
3	A	590	X2N	CAO-CBN-NBW-CAT
3	B	590	X2N	CAA-CAU-CBS-CBQ
2	B	580	HEM	C2D-C3D-CAD-CBD
3	A	590	X2N	CAM-CBM-NBU-CAZ
3	A	590	X2N	CAN-CBN-NBW-CAT
3	A	590	X2N	CAL-CBM-NBU-CAZ
2	B	580	HEM	C4D-C3D-CAD-CBD
2	B	580	HEM	C4C-C3C-CAC-CBC
3	B	590	X2N	CAK-CBL-NBT-CAX
2	B	580	HEM	C2C-C3C-CAC-CBC
3	B	590	X2N	CBC-C2-CAR-CBI

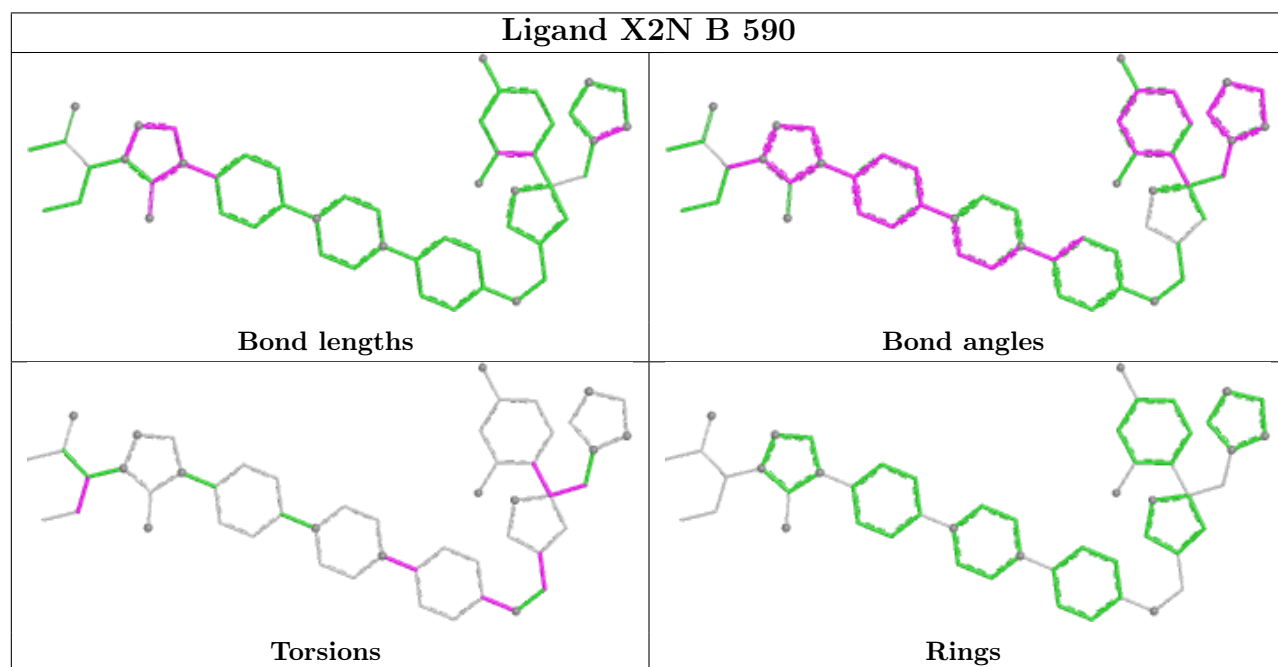
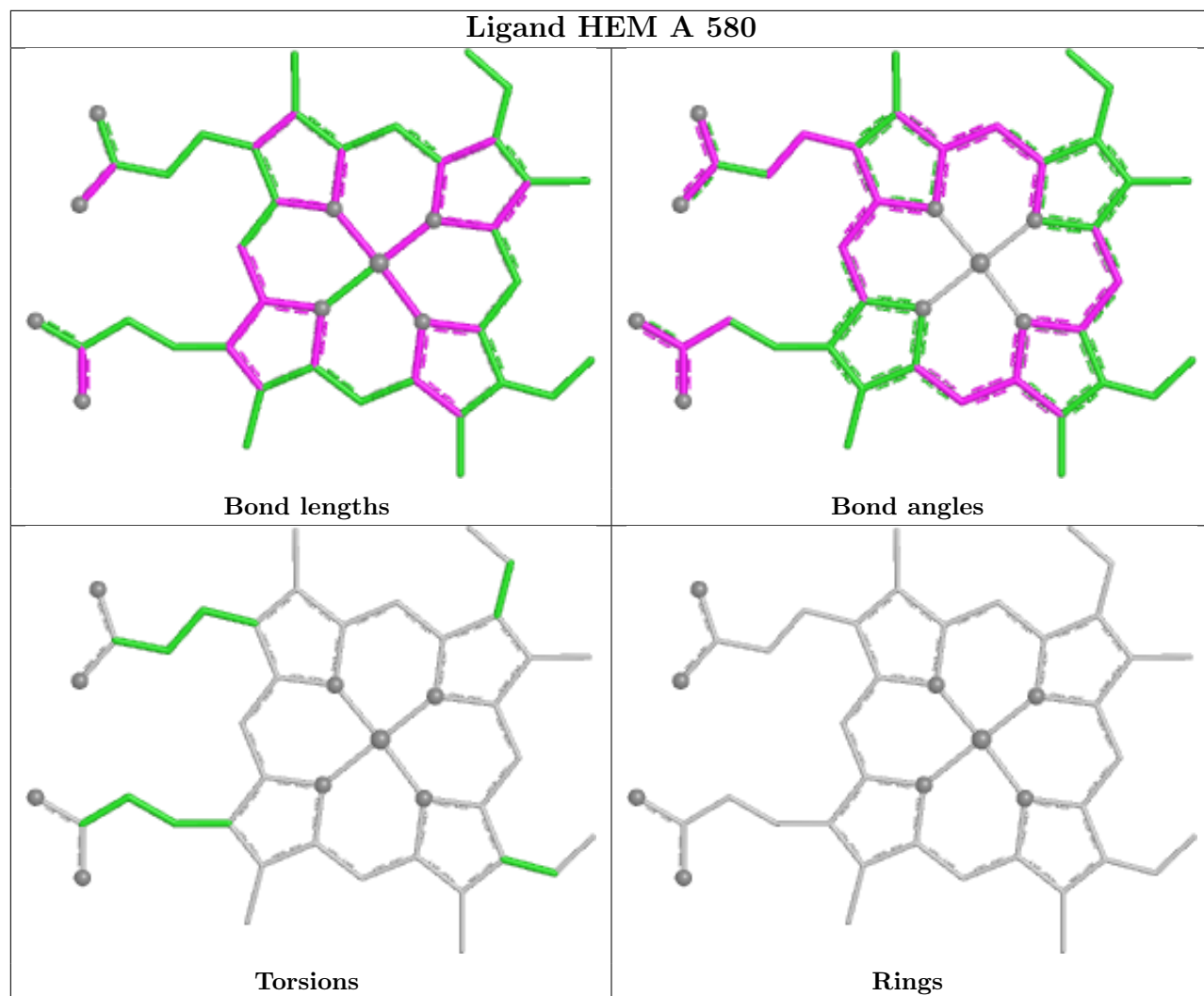
There are no ring outliers.

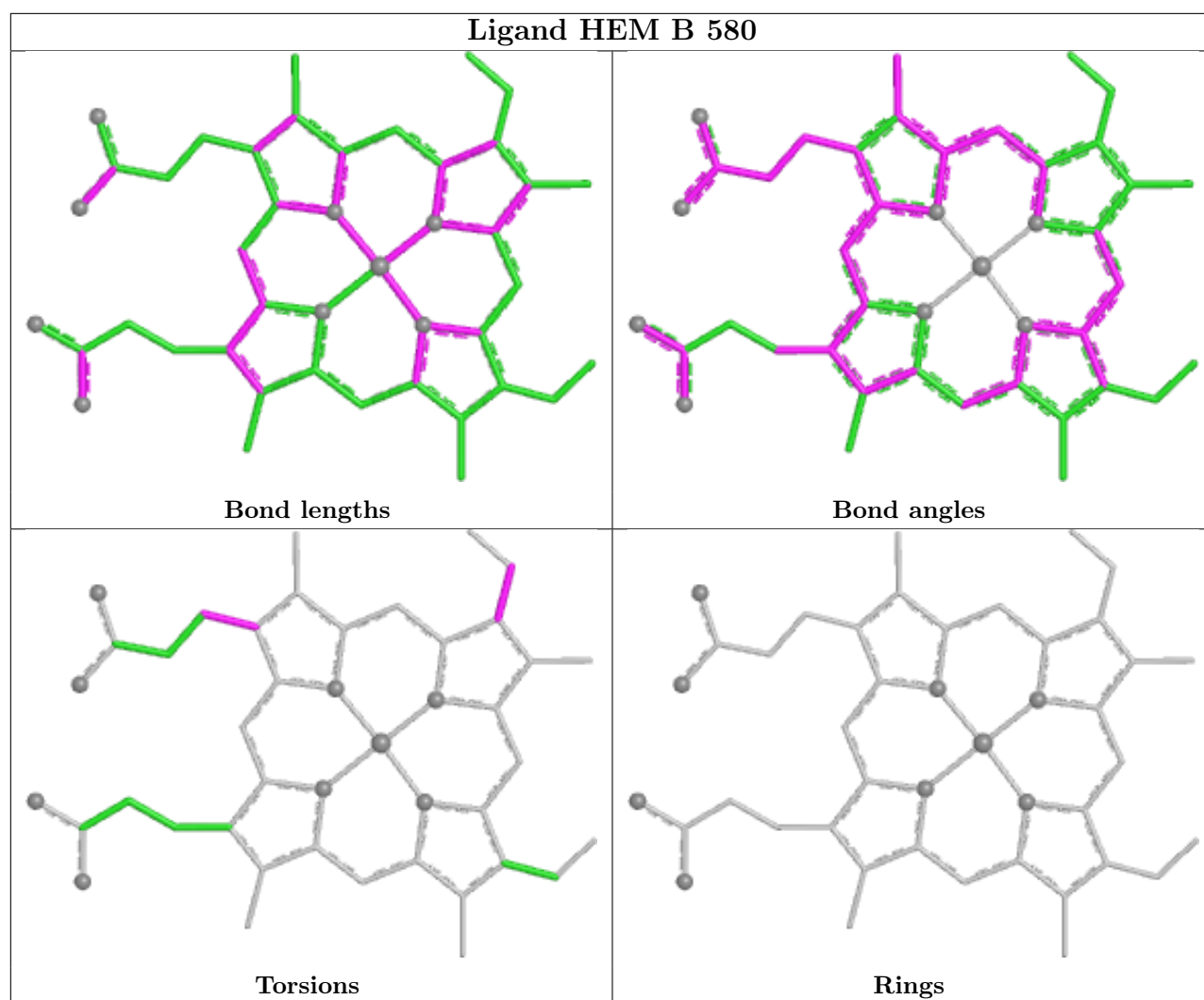
4 monomers are involved in 26 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	590	X2N	5	0
2	A	580	HEM	5	0
3	B	590	X2N	6	0
2	B	580	HEM	10	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	484/490 (98%)	0.31	29 (5%)	27 20	36, 63, 130, 189	0
1	B	483/490 (98%)	0.39	29 (6%)	27 20	38, 65, 136, 187	0
All	All	967/980 (98%)	0.35	58 (5%)	27 20	36, 64, 134, 189	0

All (58) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	520	TRP	6.5
1	A	349	ASN	5.0
1	A	272	PRO	4.1
1	B	527	MET	3.9
1	B	498	TYR	3.6
1	A	437	VAL	3.6
1	A	498	TYR	3.4
1	B	430	ALA	3.3
1	B	348	LEU	3.3
1	A	213	PHE	3.2
1	B	519	ILE	3.2
1	A	430	ALA	3.2
1	A	346	GLY	3.2
1	A	56	PRO	3.1
1	A	441	SER	3.1
1	A	527	MET	3.0
1	A	497	GLY	3.0
1	B	186	ALA	2.9
1	B	497	GLY	2.9
1	B	56	PRO	2.8
1	A	347	ASP	2.8
1	B	342	LYS	2.8
1	B	431	ALA	2.8
1	B	347	ASP	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	433	LYS	2.8
1	B	436	SER	2.8
1	B	496	ASP	2.7
1	A	51	VAL	2.7
1	B	437	VAL	2.6
1	B	524	GLU	2.6
1	B	528	PHE	2.5
1	A	183	HIS	2.5
1	A	348	LEU	2.5
1	A	270	ILE	2.4
1	A	436	SER	2.4
1	B	240	LEU	2.3
1	B	181	LYS	2.3
1	A	54	TRP	2.3
1	A	513	THR	2.3
1	A	64	TYR	2.3
1	A	435	ASN	2.2
1	B	346	GLY	2.2
1	A	528	PHE	2.2
1	A	273	ASN	2.2
1	B	349	ASN	2.2
1	B	416	PHE	2.2
1	A	342	LYS	2.1
1	B	525	THR	2.1
1	A	221	TYR	2.1
1	B	273	ASN	2.1
1	B	345	GLY	2.1
1	A	263	LEU	2.1
1	A	432	ALA	2.1
1	B	267	ARG	2.1
1	A	439	PHE	2.1
1	A	55	ILE	2.1
1	B	343	GLU	2.1
1	B	495	ILE	2.0

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

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](#)

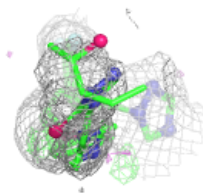
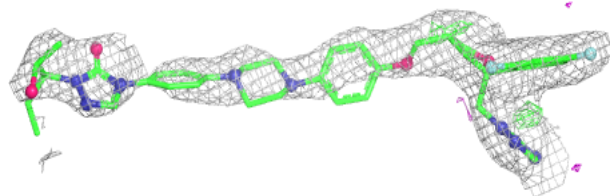
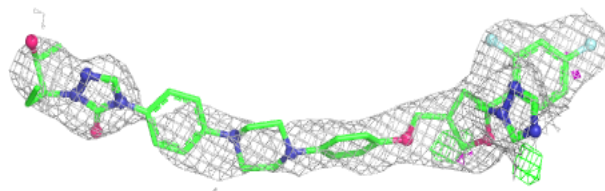
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	X2N	B	590	51/51	0.85	0.14	45,99,118,122	0
3	X2N	A	590	51/51	0.91	0.11	46,74,103,111	0
2	HEM	A	580	43/43	0.96	0.08	34,45,51,53	0
2	HEM	B	580	43/43	0.97	0.08	44,50,55,60	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.

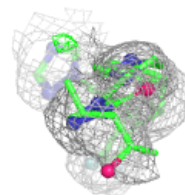
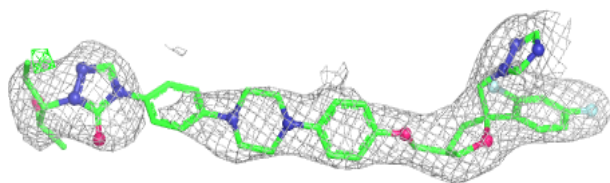
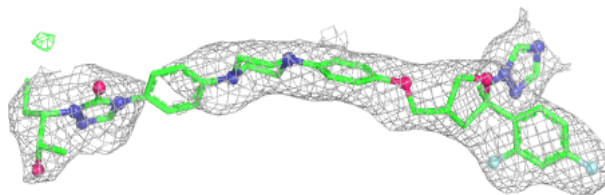
Electron density around X2N B 590:

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



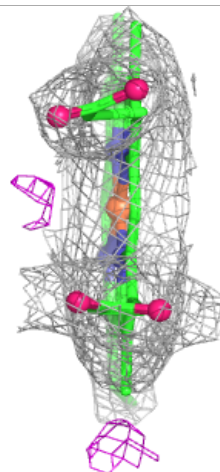
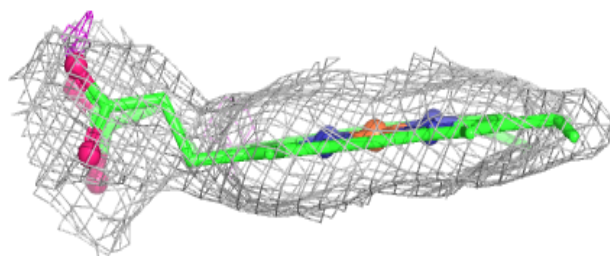
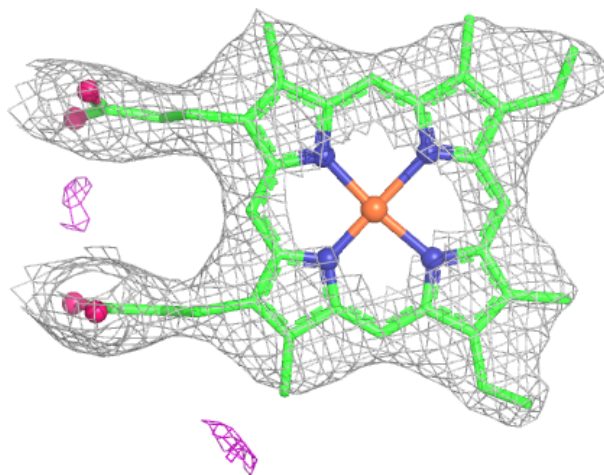
Electron density around X2N A 590:

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



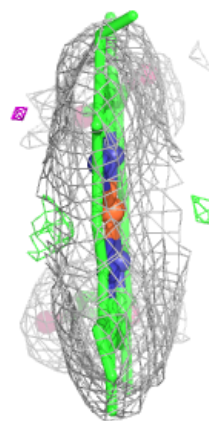
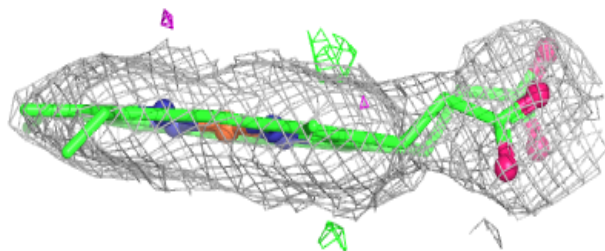
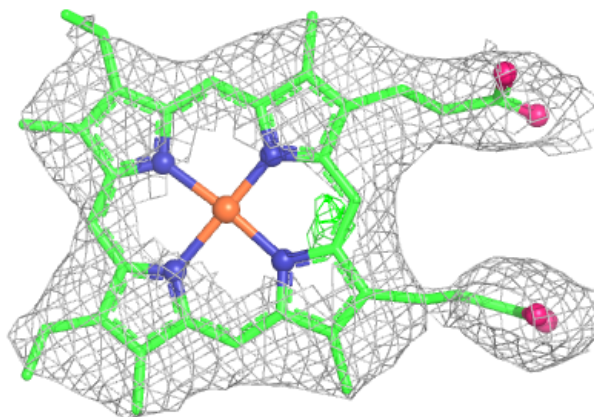
Electron density around HEM A 580:

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



Electron density around HEM B 580:

$2mF_o-DF_c$ (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.