



Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 07:34 am BST

PDB ID : 2VZ7
Title : Crystal structure of the YC-17-bound PikC D50N mutant
Authors : Li, S.; Sherman, D.H.; Podust, L.M.
Deposited on : 2008-07-30
Resolution : 3.20 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

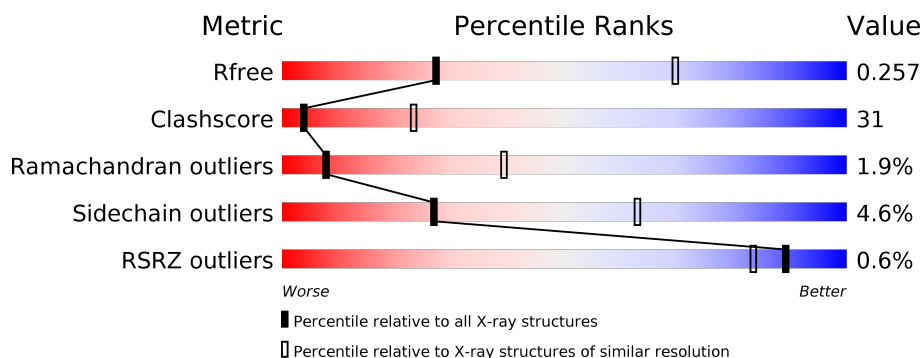
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1133 (3.20-3.20)
Clashscore	141614	1253 (3.20-3.20)
Ramachandran outliers	138981	1234 (3.20-3.20)
Sidechain outliers	138945	1233 (3.20-3.20)
RSRZ outliers	127900	1095 (3.20-3.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	436	<div> <div>%</div> <div> <div></div> <div>46%</div> <div>41%</div> <div>9%</div> </div> </div>
1	B	436	<div> <div></div> <div>48%</div> <div>39%</div> <div>10%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	PXI	A	1408	X	-	-	-
3	PXI	B	1408	X	-	-	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6441 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.

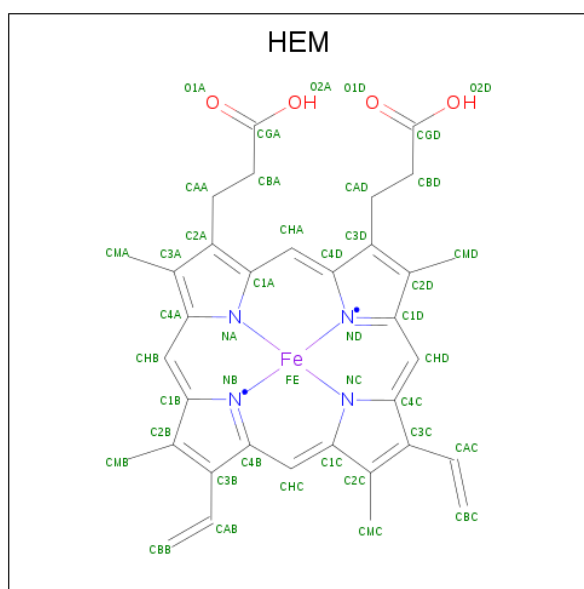
- Molecule 1 is a protein called CYTOCHROME P450 MONOOXYGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	396	Total	C	N	O	S	0	0	0
			3068	1937	552	566	13			
1	B	393	Total	C	N	O	S	0	0	0
			3048	1924	546	565	13			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	50	ASN	ASP	engineered mutation	UNP O87605
B	50	ASN	ASP	engineered mutation	UNP O87605

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



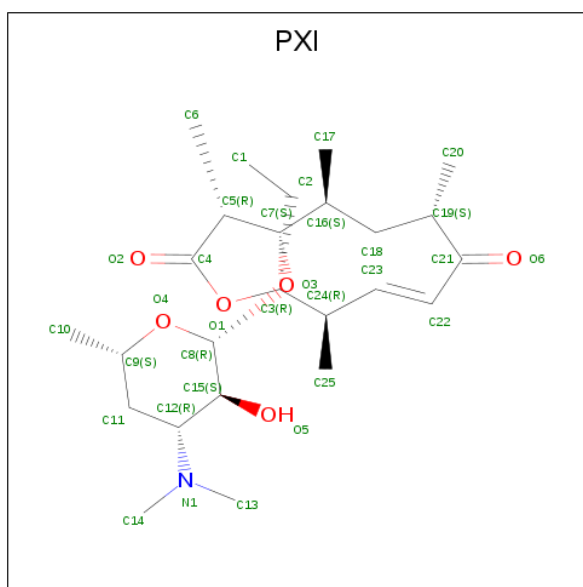
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

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

- Molecule 3 is 4-{{4-(DIMETHYLAMINO)-3-HYDROXY-6-METHYLTETRAHYDRO-2H-PYRAN-2-YL}OXY}-12-ETHYL-3,5,7,11-TETRAMETHYLOXACYCLODODEC-9-ENE-2,8-DIONE (three-letter code: PXI) (formula: C₂₅H₄₃NO₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			32	25	1	6		
3	B	1	Total	C	N	O	0	0
			32	25	1	6		

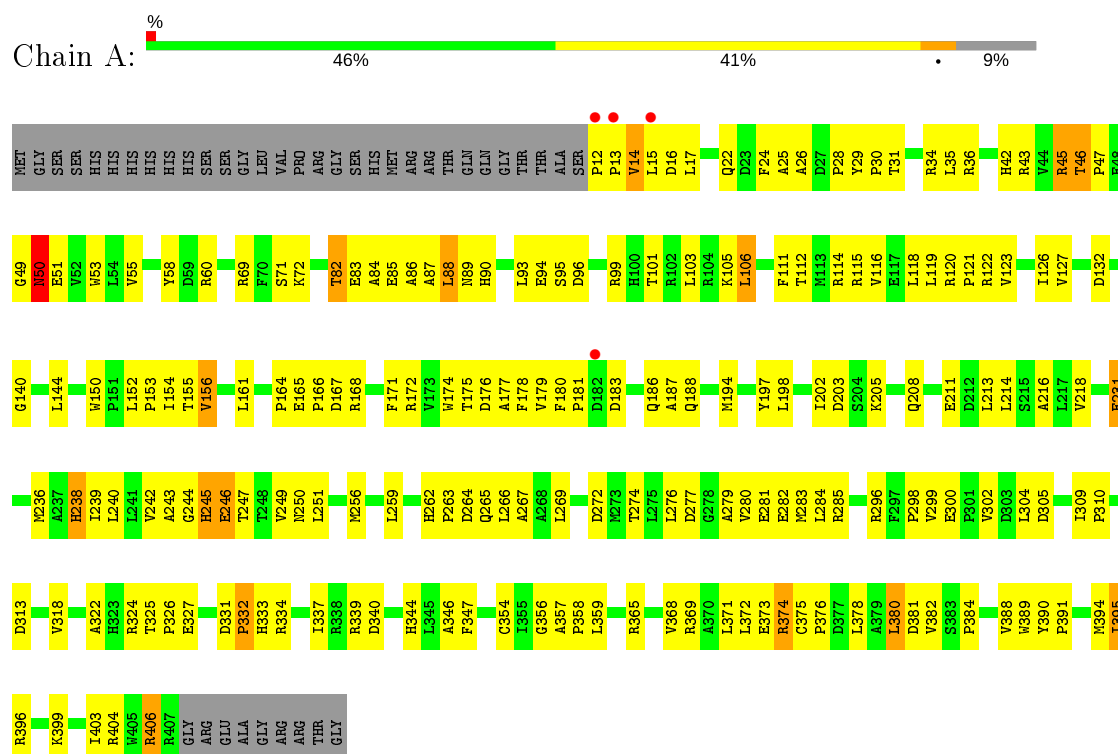
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	81	Total O	0	0
			81 81		
4	B	94	Total O	0	0
			94 94		

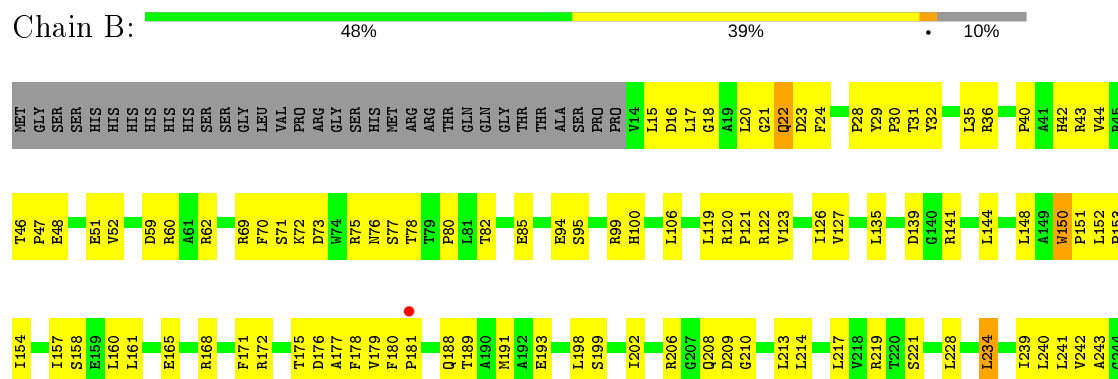
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: CYTOCHROME P450 MONOOXYGENASE



• Molecule 1: CYTOCHROME P450 MONOOXYGENASE



ALA GLY ARG ARG THR THR GLY	D321	R245
	R324	E246
	T325	T247
	P326	T248
	F329	L251
	P330	L252
	D331	A253
	P332	N254
	H333	G255
	R334	N256
F335	L259	
D336	H262	
I337	D263	
R338	D264	
R339	Q265	
H344	L266	
L345	A267	
A346	A268	
H352	L269	
R351	R270	
R365	L276	
I366	D277	
R369	V280	
A370	E281	
L371	E282	
L372	N283	
E373	R285	
R374	V286	
C375	E287	
P376	V290	
A379	T294	
L380	V295	
D381	R296	
V382	F297	
S383	F298	
P384	V302	
L387	D303	
N392	L304	
P393	D305	
M394	T309	
I395	P310	
R396	A311	
I403	G312	
R406	D313	
ARG	L316	
GLY	V317	
ARG	V318	
ARG	L319	
GLY	A320	

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	59.86Å 109.32Å 153.04Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.67 – 3.20 49.66 – 3.19	Depositor EDS
% Data completeness (in resolution range)	94.5 (49.67-3.20) 94.2 (49.66-3.19)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.77 (at 3.19Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.179 , 0.260 0.175 , 0.257	Depositor DCC
R_{free} test set	1718 reflections (9.96%)	wwPDB-VP
Wilson B-factor (Å ²)	48.6	Xtriage
Anisotropy	0.414	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 56.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6441	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 13.64% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0/3140	0.62	1/4285 (0.0%)
1	B	0.35	0/3119	0.59	0/4256
All	All	0.35	0/6259	0.61	1/8541 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	12	PRO	N-CA-CB	5.63	110.06	103.30

There are no chirality outliers.

There are no planarity outliers.

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	3068	0	3036	189	0
1	B	3048	0	3016	184	0
2	A	43	0	30	8	0
2	B	43	0	30	1	0
3	A	32	0	43	15	0
3	B	32	0	43	11	0
4	A	81	0	0	9	0
4	B	94	0	0	11	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	6441	0	6198	384	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 31.

All (384) 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)
3:A:1408:PXI:H16	3:A:1408:PXI:H22	1.38	1.01
1:A:43:ARG:HH11	1:A:43:ARG:HB3	1.23	1.01
1:B:43:ARG:HH12	1:B:51:GLU:HB3	1.21	1.00
1:A:43:ARG:NH1	1:A:43:ARG:HB3	1.80	0.97
1:B:384:PRO:HA	1:B:387:LEU:HD12	1.49	0.92
1:B:265:GLN:HE22	1:B:337:ILE:H	1.15	0.90
1:A:88:LEU:H	1:A:88:LEU:HD12	1.40	0.85
1:B:36:ARG:HH12	1:B:326:PRO:HD2	1.41	0.84
1:A:43:ARG:HH22	1:A:51:GLU:HB3	1.43	0.83
1:A:165:GLU:HG3	1:A:166:PRO:HD3	1.61	0.81
1:A:179:VAL:HG12	1:A:180:PHE:H	1.46	0.80
1:B:72:LYS:NZ	1:B:72:LYS:HB3	1.96	0.80
1:A:88:LEU:N	1:A:88:LEU:HD12	1.97	0.80
1:A:177:ALA:HA	1:A:181:PRO:HB3	1.64	0.79
1:A:36:ARG:HH12	1:A:326:PRO:HD2	1.46	0.78
1:B:150:TRP:CH2	1:B:172:ARG:HB2	2.18	0.78
1:A:175:THR:HG21	1:A:245:HIS:HD1	1.47	0.78
1:B:82:THR:HG23	1:B:85:GLU:OE1	1.85	0.77
1:A:177:ALA:HB1	1:A:187:ALA:HA	1.68	0.75
3:A:1408:PXI:H22	3:A:1408:PXI:C16	2.17	0.74
1:A:179:VAL:HG12	1:A:180:PHE:N	2.02	0.74
1:B:296:ARG:HG3	1:B:296:ARG:HH11	1.52	0.74
1:A:259:LEU:HD21	1:A:269:LEU:HD23	1.70	0.74
1:B:251:LEU:HD13	1:B:284:LEU:HD23	1.69	0.73
1:A:35:LEU:HB3	1:A:42:HIS:CD2	2.23	0.72
1:A:50:ASN:HD22	1:A:50:ASN:N	1.86	0.72
1:B:43:ARG:NH1	1:B:51:GLU:HB3	2.01	0.71
1:A:45:ARG:HB2	1:A:45:ARG:NH1	2.06	0.71
1:B:265:GLN:NE2	1:B:337:ILE:H	1.88	0.71
1:A:380:LEU:HD22	1:A:382:VAL:H	1.55	0.71
1:A:36:ARG:NH1	1:A:326:PRO:HD2	2.06	0.70
1:B:175:THR:HG21	1:B:245:HIS:HD2	1.55	0.69
1:B:152:LEU:HB3	1:B:153:PRO:HD3	1.73	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:318:VAL:HG12	1:B:321:ASP:H	1.57	0.69
1:A:24:PHE:O	1:A:28:PRO:HG3	1.92	0.69
1:B:269:LEU:HD11	1:B:276:LEU:HA	1.74	0.69
1:A:152:LEU:HB3	1:A:153:PRO:HD3	1.74	0.69
1:A:171:PHE:O	1:A:175:THR:HG23	1.93	0.68
1:A:45:ARG:HB2	1:A:45:ARG:HH11	1.58	0.68
1:A:285:ARG:NH2	1:A:332:PRO:O	2.28	0.67
1:B:178:PHE:CE2	1:B:191:MET:HG3	2.30	0.67
1:B:265:GLN:HE22	1:B:337:ILE:N	1.92	0.66
1:B:384:PRO:HA	1:B:387:LEU:CD1	2.25	0.66
1:A:123:VAL:O	1:A:127:VAL:HG23	1.95	0.66
1:B:72:LYS:HZ3	1:B:72:LYS:HB3	1.60	0.66
1:B:94:GLU:OE1	3:B:1408:PXI:H143	1.95	0.65
1:A:331:ASP:OD1	1:A:334:ARG:HD2	1.96	0.65
1:A:205:LYS:HE3	1:A:214:LEU:HD23	1.78	0.65
1:A:103:LEU:O	1:A:106:LEU:HB2	1.96	0.65
1:B:248:THR:O	1:B:252:ILE:HG13	1.96	0.65
1:B:36:ARG:NH1	1:B:326:PRO:HD2	2.12	0.65
3:B:1408:PXI:C22	3:B:1408:PXI:H16	2.27	0.65
1:B:144:LEU:HB2	1:B:403:ILE:HD13	1.79	0.65
1:B:344:HIS:HD2	1:B:346:ALA:CB	2.10	0.64
1:B:119:LEU:HD11	1:B:160:LEU:HD21	1.78	0.64
1:B:206:ARG:HH11	1:B:206:ARG:HG3	1.61	0.64
1:A:161:LEU:O	1:A:214:LEU:HB2	1.97	0.64
1:A:279:ALA:O	1:A:283:MET:HG3	1.98	0.64
1:B:296:ARG:HG3	1:B:296:ARG:NH1	2.12	0.64
1:B:32:TYR:HB3	1:B:321:ASP:HB2	1.80	0.64
1:A:16:ASP:OD2	1:A:45:ARG:HD2	1.98	0.63
1:A:72:LYS:HB3	1:A:72:LYS:NZ	2.13	0.63
1:B:71:SER:CB	1:B:76:ASN:HD22	2.10	0.63
1:A:238:HIS:O	1:A:242:VAL:HG22	1.98	0.63
1:B:36:ARG:HH12	1:B:325:THR:HA	1.63	0.63
1:B:139:ASP:OD2	1:B:141:ARG:HB2	1.98	0.62
1:B:165:GLU:HA	1:B:168:ARG:NE	2.14	0.62
1:A:214:LEU:O	1:A:218:VAL:HG23	1.99	0.62
1:B:189:THR:O	1:B:193:GLU:HG3	1.99	0.62
1:A:272:ASP:OD1	1:A:274:THR:HG23	1.99	0.62
1:A:31:THR:HA	1:A:34:ARG:NH1	2.15	0.62
1:B:344:HIS:CD2	1:B:346:ALA:HB3	2.35	0.61
1:A:122:ARG:O	1:A:126:ILE:HG13	2.01	0.61
2:A:1407:HEM:HBB2	2:A:1407:HEM:HMB2	1.81	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:144:LEU:HD21	1:A:256:MET:HG3	1.82	0.61
3:A:1408:PXI:C4	3:A:1408:PXI:H173	2.30	0.61
1:A:242:VAL:HG23	3:A:1408:PXI:H6C3	1.83	0.61
1:A:34:ARG:HH11	1:A:34:ARG:HB3	1.66	0.61
1:A:84:ALA:HB3	1:A:188:GLN:HE22	1.65	0.61
1:A:277:ASP:O	1:A:280:VAL:HG12	2.01	0.60
1:B:29:TYR:N	1:B:30:PRO:HD2	2.17	0.60
1:A:179:VAL:O	1:A:181:PRO:HD3	2.02	0.60
1:A:390:TYR:HE1	1:A:399:LYS:HA	1.67	0.60
1:B:82:THR:HG21	1:B:188:GLN:OE1	2.01	0.60
1:A:82:THR:OG1	1:A:85:GLU:HG3	2.02	0.60
1:B:36:ARG:NH1	1:B:325:THR:HA	2.16	0.59
1:A:116:VAL:HG11	1:A:359:LEU:HA	1.85	0.59
1:B:324:ARG:HA	1:B:332:PRO:HB2	1.85	0.59
1:B:29:TYR:HD1	1:B:29:TYR:H	1.51	0.59
1:A:304:LEU:HD12	1:A:309:ILE:HD11	1.85	0.59
1:A:388:VAL:HG13	1:A:399:LYS:HB2	1.85	0.58
1:A:13:PRO:O	1:A:14:VAL:HB	2.03	0.58
1:B:141:ARG:HH11	1:B:141:ARG:HB2	1.68	0.58
1:A:101:THR:O	1:A:105:LYS:HG3	2.03	0.58
1:A:150:TRP:O	1:A:153:PRO:HD2	2.04	0.58
1:B:396:ARG:HG2	1:B:396:ARG:HH11	1.69	0.58
3:B:1408:PXI:C4	3:B:1408:PXI:H173	2.33	0.58
1:B:338:ARG:HD3	4:B:2076:HOH:O	2.03	0.58
1:A:132:ASP:OD1	1:A:374:ARG:NH2	2.31	0.58
1:A:395:ILE:HG23	1:A:396:ARG:N	2.19	0.58
1:A:88:LEU:CD1	1:A:88:LEU:H	2.15	0.58
1:B:178:PHE:HE2	1:B:191:MET:HG3	1.66	0.57
1:A:339:ARG:HG2	1:A:340:ASP:N	2.19	0.57
1:B:251:LEU:HD22	1:B:284:LEU:HD21	1.86	0.57
1:B:265:GLN:HE21	1:B:337:ILE:HG23	1.69	0.57
1:B:95:SER:CB	1:B:99:ARG:HD2	2.34	0.57
1:B:304:LEU:HD12	1:B:309:ILE:HD11	1.85	0.57
1:B:29:TYR:HD2	1:B:324:ARG:HG3	1.69	0.57
1:A:395:ILE:HG12	3:A:1408:PXI:H23	1.87	0.57
1:A:269:LEU:HD23	1:A:372:LEU:HD11	1.86	0.57
1:A:165:GLU:CG	1:A:166:PRO:HD3	2.33	0.57
1:A:34:ARG:CB	1:A:34:ARG:NH1	2.68	0.56
1:B:35:LEU:O	1:B:42:HIS:NE2	2.38	0.56
1:A:211:GLU:HA	1:A:211:GLU:OE1	2.04	0.56
1:B:285:ARG:NH2	1:B:332:PRO:O	2.38	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:265:GLN:NE2	1:B:337:ILE:N	2.51	0.56
1:A:150:TRP:CZ2	1:A:172:ARG:HB2	2.41	0.56
1:B:344:HIS:HD2	1:B:346:ALA:HB3	1.71	0.56
1:B:406:ARG:HA	4:B:2091:HOH:O	2.05	0.56
1:B:122:ARG:O	1:B:126:ILE:HG13	2.06	0.56
1:B:382:VAL:HG22	1:B:383:SER:N	2.21	0.56
1:B:395:ILE:HG23	1:B:396:ARG:N	2.21	0.56
1:B:265:GLN:NE2	1:B:337:ILE:HG23	2.20	0.56
1:A:277:ASP:O	1:A:281:GLU:HG2	2.06	0.55
1:B:264:ASP:O	1:B:267:ALA:HB3	2.06	0.55
1:A:356:GLY:HA2	4:A:2071:HOH:O	2.06	0.55
1:B:269:LEU:HD23	1:B:372:LEU:HD11	1.89	0.55
1:B:177:ALA:O	1:B:181:PRO:HB3	2.06	0.55
1:A:236:MET:O	1:A:239:ILE:HG22	2.07	0.55
1:A:87:ALA:HB3	1:A:88:LEU:HD12	1.89	0.55
1:B:298:PRO:HB3	1:B:302:VAL:HG23	1.88	0.55
1:B:150:TRP:HZ3	1:B:245:HIS:NE2	2.05	0.54
1:A:179:VAL:CG1	1:A:180:PHE:H	2.18	0.54
1:A:29:TYR:HB2	1:A:30:PRO:HD3	1.88	0.54
1:B:127:VAL:HG21	1:B:366:ILE:HG22	1.89	0.54
1:B:175:THR:HG21	1:B:245:HIS:CD2	2.41	0.54
1:B:381:ASP:OD1	1:B:381:ASP:O	2.26	0.54
1:A:357:ALA:HB3	1:A:358:PRO:HD3	1.89	0.54
1:A:93:LEU:HB2	2:A:1407:HEM:HAD2	1.89	0.54
1:B:175:THR:CG2	1:B:246:GLU:HG2	2.38	0.54
1:B:334:ARG:HD3	4:B:2072:HOH:O	2.08	0.54
1:A:179:VAL:CG1	1:A:180:PHE:N	2.72	0.53
1:B:210:GLY:O	1:B:219:ARG:NH2	2.41	0.53
1:B:254:ASN:HB3	1:B:287:GLU:HB3	1.90	0.53
1:B:70:PHE:HB3	1:B:296:ARG:HB3	1.91	0.53
1:B:280:VAL:O	1:B:283:MET:HB2	2.08	0.53
1:B:60:ARG:HD2	1:B:304:LEU:HD22	1.91	0.53
1:A:58:TYR:HA	1:A:322:ALA:CB	2.38	0.53
1:A:120:ARG:HB3	1:A:121:PRO:HD3	1.90	0.53
1:B:95:SER:HB3	1:B:99:ARG:HH11	1.73	0.53
1:A:282:GLU:CD	1:A:339:ARG:HD2	2.29	0.53
1:B:259:LEU:HD21	1:B:372:LEU:HD11	1.91	0.53
1:B:379:ALA:HB1	4:B:2088:HOH:O	2.09	0.53
1:B:161:LEU:HD22	1:B:241:LEU:CD2	2.39	0.52
3:A:1408:PXI:H16	3:A:1408:PXI:C22	2.27	0.52
1:A:395:ILE:CG2	1:A:396:ARG:N	2.72	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:46:THR:HG23	1:A:50:ASN:O	2.08	0.52
1:A:58:TYR:HA	1:A:322:ALA:HB1	1.92	0.52
1:A:279:ALA:CB	1:A:337:ILE:HB	2.40	0.52
1:A:384:PRO:HD3	4:B:2062:HOH:O	2.08	0.52
1:A:244:GLY:HA3	4:A:2055:HOH:O	2.10	0.51
1:A:50:ASN:N	1:A:50:ASN:ND2	2.56	0.51
1:B:243:ALA:HA	3:B:1408:PXI:O2	2.09	0.51
1:B:165:GLU:HA	1:B:168:ARG:HE	1.75	0.51
1:B:361:ARG:O	1:B:365:ARG:HG3	2.11	0.51
1:B:78:THR:O	1:B:80:PRO:HD3	2.11	0.51
1:A:55:VAL:O	1:A:318:VAL:HG23	2.10	0.51
1:B:171:PHE:CZ	1:B:241:LEU:HD13	2.46	0.51
1:A:198:LEU:O	1:A:202:ILE:HG13	2.10	0.51
1:A:266:LEU:HD11	1:A:372:LEU:HD22	1.92	0.51
1:B:72:LYS:HZ2	1:B:72:LYS:HB3	1.75	0.51
1:B:310:PRO:HA	4:B:2068:HOH:O	2.10	0.51
1:A:112:THR:O	1:A:116:VAL:HG23	2.10	0.51
1:A:15:LEU:CD1	1:A:17:LEU:HD23	2.41	0.51
1:A:231:GLU:OE1	1:A:231:GLU:N	2.44	0.50
1:B:144:LEU:HD21	1:B:256:MET:HG3	1.92	0.50
1:A:118:LEU:HA	4:A:2026:HOH:O	2.11	0.50
1:A:281:GLU:O	1:A:284:LEU:HB2	2.11	0.50
1:A:34:ARG:HB2	1:A:34:ARG:NH1	2.26	0.50
1:A:394:MET:HG2	3:A:1408:PXI:H251	1.93	0.50
1:A:153:PRO:HG2	1:A:249:VAL:HG22	1.93	0.50
1:B:150:TRP:CZ3	1:B:245:HIS:NE2	2.80	0.50
1:B:395:ILE:CG2	1:B:396:ARG:N	2.74	0.50
1:B:29:TYR:N	1:B:29:TYR:CD1	2.79	0.50
1:B:282:GLU:HG2	1:B:336:ASP:O	2.11	0.50
1:B:135:LEU:HD12	1:B:374:ARG:NH2	2.26	0.50
1:B:221:SER:HB2	1:B:228:LEU:HD23	1.94	0.50
1:A:34:ARG:HH11	1:A:34:ARG:CB	2.24	0.50
1:B:154:ILE:HB	1:B:245:HIS:HE1	1.77	0.49
1:A:175:THR:HG21	1:A:245:HIS:ND1	2.24	0.49
1:A:45:ARG:HG3	1:A:45:ARG:O	2.12	0.49
1:B:120:ARG:HB3	1:B:121:PRO:HD3	1.94	0.49
1:B:46:THR:HG23	1:B:52:VAL:HG22	1.94	0.49
1:A:327:GLU:OE1	1:A:327:GLU:N	2.43	0.49
1:A:53:TRP:CE2	1:A:309:ILE:HG23	2.47	0.49
1:A:69:ARG:O	1:A:298:PRO:HA	2.13	0.49
1:A:378:LEU:HD12	1:A:404:ARG:O	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:198:LEU:O	1:B:199:SER:C	2.51	0.49
1:A:164:PRO:HA	4:A:2035:HOH:O	2.12	0.49
1:B:82:THR:OG1	1:B:85:GLU:HG3	2.13	0.49
1:A:13:PRO:O	1:A:14:VAL:CB	2.60	0.49
1:B:172:ARG:HG2	1:B:176:ASP:OD2	2.12	0.49
1:A:167:ASP:OD1	1:A:197:TYR:OH	2.31	0.49
1:A:154:ILE:HD13	1:A:245:HIS:NE2	2.28	0.49
1:B:179:VAL:HA	3:B:1408:PXI:H201	1.94	0.48
1:A:381:ASP:HB2	1:A:403:ILE:CA	2.43	0.48
1:A:90:HIS:CD2	1:A:99:ARG:HH12	2.31	0.48
1:B:148:LEU:C	1:B:151:PRO:HD2	2.33	0.48
1:A:247:THR:HB	2:A:1407:HEM:C2B	2.48	0.48
1:A:246:GLU:O	1:A:250:ASN:ND2	2.47	0.48
1:A:265:GLN:HE21	1:A:337:ILE:HG23	1.79	0.48
1:A:95:SER:HB3	1:A:99:ARG:HH21	1.78	0.48
1:B:374:ARG:O	1:B:376:PRO:HD3	2.14	0.48
1:A:43:ARG:CB	1:A:43:ARG:HH11	2.10	0.48
1:B:282:GLU:HA	1:B:282:GLU:OE1	2.14	0.48
1:A:282:GLU:OE1	1:A:285:ARG:HD3	2.13	0.48
1:B:285:ARG:NH1	1:B:339:ARG:NE	2.61	0.48
1:A:380:LEU:HD22	1:A:382:VAL:O	2.14	0.48
1:A:49:GLY:O	1:A:50:ASN:O	2.32	0.48
1:B:59:ASP:HB2	4:B:2008:HOH:O	2.14	0.48
1:B:77:SER:HB2	1:B:297:PHE:CE1	2.49	0.48
1:A:279:ALA:HB2	1:A:337:ILE:HB	1.96	0.48
1:B:43:ARG:HH11	1:B:43:ARG:HG2	1.79	0.48
1:A:310:PRO:HB2	1:A:313:ASP:OD2	2.12	0.47
1:A:94:GLU:OE2	3:A:1408:PXI:C14	2.62	0.47
1:A:144:LEU:O	1:A:144:LEU:HG	2.13	0.47
1:A:239:ILE:HD12	3:A:1408:PXI:H6C2	1.96	0.47
1:B:40:PRO:HB2	1:B:60:ARG:HG2	1.94	0.47
1:B:21:GLY:C	1:B:23:ASP:H	2.17	0.47
1:A:357:ALA:N	1:A:358:PRO:CD	2.77	0.47
1:B:262:HIS:CD2	1:B:335:PHE:HB3	2.49	0.47
1:B:15:LEU:HD12	1:B:16:ASP:N	2.29	0.47
1:B:28:PRO:HD2	1:B:29:TYR:CD1	2.50	0.47
1:B:392:ASN:OD1	1:B:393:PRO:HD2	2.13	0.47
1:B:29:TYR:OH	1:B:396:ARG:NH1	2.47	0.47
1:B:161:LEU:CD2	1:B:241:LEU:HG	2.45	0.47
1:A:94:GLU:OE2	3:A:1408:PXI:H142	2.15	0.47
1:A:34:ARG:CZ	1:A:34:ARG:HB2	2.45	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:83:GLU:CB	4:A:2015:HOH:O	2.64	0.46
1:B:263:PRO:HD2	4:B:2062:HOH:O	2.15	0.46
1:B:310:PRO:HB2	1:B:313:ASP:OD1	2.15	0.46
1:A:49:GLY:C	1:A:50:ASN:HD22	2.16	0.46
1:B:17:LEU:HD11	1:B:44:VAL:HG21	1.97	0.46
1:A:344:HIS:HD2	1:A:346:ALA:CB	2.29	0.46
1:B:161:LEU:O	1:B:214:LEU:HB2	2.15	0.46
1:B:78:THR:HG23	1:B:312:GLY:HA3	1.98	0.46
1:A:115:ARG:O	1:A:118:LEU:HB2	2.16	0.46
1:A:17:LEU:HB2	1:A:47:PRO:HD3	1.97	0.46
1:A:339:ARG:CG	1:A:340:ASP:N	2.79	0.46
1:A:208:GLN:NE2	1:A:208:GLN:HA	2.31	0.46
1:A:26:ALA:HA	1:A:389:TRP:CD1	2.51	0.46
1:A:69:ARG:NH1	1:A:302:VAL:HG22	2.31	0.46
1:B:28:PRO:C	1:B:30:PRO:HD2	2.35	0.46
1:B:94:GLU:OE2	3:B:1408:PXI:C14	2.64	0.46
1:B:285:ARG:HD2	1:B:329:PHE:CD2	2.51	0.46
1:B:30:PRO:HG2	1:B:31:THR:H	1.81	0.46
1:B:150:TRP:CZ3	1:B:245:HIS:CE1	3.03	0.45
1:B:69:ARG:HB3	4:B:2011:HOH:O	2.16	0.45
3:A:1408:PXI:C22	3:A:1408:PXI:C16	2.92	0.45
1:A:243:ALA:HA	3:A:1408:PXI:O2	2.15	0.45
1:A:231:GLU:CD	1:A:231:GLU:H	2.20	0.45
1:A:309:ILE:HA	1:A:310:PRO:HD3	1.82	0.45
1:A:376:PRO:HB2	1:A:406:ARG:CB	2.47	0.45
1:B:286:TYR:CD2	1:B:287:GLU:HG2	2.52	0.45
1:A:155:THR:O	1:A:156:VAL:C	2.55	0.45
1:A:36:ARG:HH12	1:A:326:PRO:CD	2.23	0.45
1:B:242:VAL:HG12	3:B:1408:PXI:H6C3	1.98	0.45
1:A:25:ALA:HB2	1:A:391:PRO:HA	1.99	0.45
1:A:264:ASP:O	1:A:267:ALA:HB3	2.16	0.45
1:A:296:ARG:HG3	1:A:296:ARG:HH11	1.82	0.45
1:A:194:MET:CE	1:A:238:HIS:HD2	2.30	0.45
1:B:290:VAL:HG22	3:B:1408:PXI:H252	1.99	0.45
1:B:46:THR:HG23	1:B:52:VAL:CG2	2.47	0.45
1:B:296:ARG:HH22	2:B:1407:HEM:CGA	2.30	0.45
1:B:344:HIS:HD2	1:B:346:ALA:HB2	1.78	0.45
1:A:281:GLU:HA	1:A:284:LEU:HD12	1.99	0.45
1:A:245:HIS:HB3	1:A:246:GLU:H	1.50	0.45
1:B:135:LEU:HD12	1:B:374:ARG:HH21	1.81	0.45
1:B:319:LEU:HD13	1:B:345:LEU:O	2.16	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:299:VAL:HG23	1:A:300:GLU:HG2	1.98	0.44
1:B:382:VAL:HG22	1:B:383:SER:H	1.82	0.44
1:B:202:ILE:HD11	1:B:234:LEU:HA	2.00	0.44
1:B:330:PRO:HB2	4:B:2069:HOH:O	2.17	0.44
1:B:251:LEU:HD13	1:B:284:LEU:CD2	2.44	0.44
1:A:247:THR:HB	2:A:1407:HEM:C3B	2.52	0.44
1:A:150:TRP:CE2	1:A:172:ARG:HD2	2.53	0.44
1:A:304:LEU:HD23	1:A:304:LEU:HA	1.87	0.44
1:A:251:LEU:HD22	2:A:1407:HEM:HBB1	1.99	0.44
1:B:148:LEU:O	1:B:151:PRO:HD2	2.18	0.44
1:B:295:TYR:HB2	4:B:2014:HOH:O	2.17	0.44
1:B:324:ARG:HA	1:B:324:ARG:HD3	1.77	0.44
1:A:251:LEU:HB2	2:A:1407:HEM:HBB1	1.99	0.44
1:A:280:VAL:HG11	1:A:365:ARG:HE	1.83	0.44
1:B:276:LEU:O	1:B:277:ASP:C	2.55	0.44
1:B:18:GLY:HA3	1:B:47:PRO:HA	1.99	0.44
1:A:71:SER:OG	1:A:96:ASP:OD1	2.29	0.44
1:B:239:ILE:HG23	1:B:240:LEU:N	2.33	0.44
1:B:242:VAL:CG1	3:B:1408:PXI:H172	2.48	0.43
1:B:259:LEU:HD11	1:B:266:LEU:HA	2.00	0.43
1:A:371:LEU:O	1:A:375:CYS:HB2	2.17	0.43
1:B:206:ARG:HH11	1:B:206:ARG:CG	2.29	0.43
1:B:381:ASP:O	1:B:382:VAL:HB	2.18	0.43
1:B:62:ARG:HB2	1:B:345:LEU:HD21	1.99	0.43
1:B:73:ASP:OD1	1:B:75:ARG:HG3	2.17	0.43
1:A:194:MET:HE3	1:A:238:HIS:HD2	1.82	0.43
1:A:85:GLU:O	1:A:87:ALA:N	2.51	0.43
1:B:277:ASP:OD1	1:B:365:ARG:HD3	2.17	0.43
1:A:94:GLU:OE1	3:A:1408:PXI:H143	2.18	0.43
1:A:174:TRP:O	1:A:175:THR:C	2.56	0.43
1:A:239:ILE:HG23	1:A:240:LEU:N	2.33	0.43
1:B:141:ARG:CB	1:B:141:ARG:NH1	2.81	0.43
1:B:161:LEU:HD12	1:B:161:LEU:N	2.34	0.43
1:B:304:LEU:O	1:B:305:ASP:C	2.57	0.43
1:B:144:LEU:HD11	1:B:371:LEU:HD11	2.00	0.43
1:A:119:LEU:O	1:A:123:VAL:HG23	2.19	0.43
1:A:266:LEU:CD1	1:A:372:LEU:HD22	2.49	0.43
1:B:29:TYR:N	1:B:30:PRO:CD	2.82	0.43
1:B:281:GLU:OE1	1:B:344:HIS:CE1	2.72	0.43
1:A:111:PHE:HB3	4:A:2071:HOH:O	2.19	0.43
1:A:140:GLY:O	1:A:404:ARG:HA	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:17:LEU:HD22	1:A:24:PHE:CZ	2.54	0.43
1:A:390:TYR:HA	1:A:391:PRO:HD3	1.87	0.43
1:A:31:THR:HA	1:A:34:ARG:HH11	1.83	0.42
1:B:214:LEU:O	1:B:217:LEU:HB2	2.19	0.42
1:B:95:SER:HB2	1:B:99:ARG:HD2	2.01	0.42
1:A:150:TRP:CH2	1:A:245:HIS:CE1	3.07	0.42
1:A:276:LEU:HD22	1:A:369:ARG:HH21	1.84	0.42
1:A:388:VAL:HG23	4:A:2001:HOH:O	2.18	0.42
1:B:85:GLU:HG2	1:B:191:MET:SD	2.59	0.42
1:B:396:ARG:CG	1:B:396:ARG:HH11	2.30	0.42
2:A:1407:HEM:HBB2	2:A:1407:HEM:CMB	2.48	0.42
1:A:15:LEU:HD12	1:A:17:LEU:HD23	2.02	0.42
1:B:21:GLY:C	1:B:23:ASP:N	2.72	0.42
1:A:181:PRO:HG3	1:A:187:ALA:HB2	2.01	0.42
1:B:119:LEU:O	1:B:123:VAL:HG23	2.20	0.42
1:B:213:LEU:O	1:B:214:LEU:C	2.57	0.42
1:B:298:PRO:HB3	1:B:302:VAL:CG2	2.48	0.42
1:B:304:LEU:HD12	1:B:309:ILE:CD1	2.50	0.42
1:A:358:PRO:HD2	4:A:2071:HOH:O	2.19	0.42
1:A:72:LYS:HB3	1:A:72:LYS:HZ2	1.82	0.42
1:A:213:LEU:O	1:A:216:ALA:HB3	2.19	0.42
1:B:150:TRP:CZ3	1:B:172:ARG:HB2	2.52	0.42
1:B:295:TYR:CE1	1:B:316:LEU:HD21	2.54	0.42
2:A:1407:HEM:CBB	2:A:1407:HEM:HMB2	2.47	0.42
1:A:90:HIS:HD2	1:A:99:ARG:NH1	2.17	0.42
1:B:48:GLU:H	1:B:48:GLU:CD	2.22	0.42
1:A:165:GLU:HA	1:A:168:ARG:NE	2.35	0.42
1:A:282:GLU:OE2	1:A:339:ARG:HD2	2.19	0.42
1:A:354:CYS:SG	1:A:357:ALA:N	2.93	0.42
1:B:157:ILE:HG23	1:B:158:SER:N	2.35	0.42
1:B:330:PRO:O	1:B:331:ASP:C	2.58	0.42
1:A:183:ASP:O	1:A:186:GLN:HB3	2.20	0.42
1:A:239:ILE:CD1	3:A:1408:PXI:H6C2	2.50	0.41
1:B:22:GLN:H	1:B:22:GLN:NE2	2.19	0.41
1:A:324:ARG:O	1:A:326:PRO:HD3	2.21	0.41
1:B:208:GLN:NE2	1:B:208:GLN:HA	2.36	0.41
1:B:254:ASN:ND2	1:B:396:ARG:O	2.47	0.41
1:A:282:GLU:OE1	1:A:339:ARG:HD2	2.21	0.41
1:B:94:GLU:CD	3:B:1408:PXI:H143	2.41	0.41
1:B:76:ASN:O	1:B:312:GLY:HA2	2.21	0.41
1:A:282:GLU:HA	1:A:282:GLU:OE1	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:24:PHE:O	1:B:28:PRO:HB3	2.20	0.41
1:B:72:LYS:HE3	1:B:294:THR:OG1	2.20	0.41
1:B:239:ILE:CG2	1:B:240:LEU:N	2.83	0.41
1:A:101:THR:HG22	1:A:105:LYS:HE3	2.03	0.41
1:B:154:ILE:HB	1:B:245:HIS:CE1	2.56	0.41
1:A:178:PHE:HB3	3:A:1408:PXI:H203	2.03	0.41
1:B:141:ARG:NH1	1:B:141:ARG:HB2	2.35	0.41
1:B:304:LEU:HA	1:B:304:LEU:HD23	1.89	0.41
1:A:84:ALA:HB3	1:A:188:GLN:NE2	2.34	0.41
1:A:368:VAL:O	1:A:369:ARG:C	2.60	0.40
1:A:90:HIS:CD2	1:A:99:ARG:NH1	2.89	0.40
1:A:299:VAL:HA	4:A:2063:HOH:O	2.20	0.40
1:B:371:LEU:O	1:B:375:CYS:HB2	2.21	0.40
1:A:325:THR:HA	1:A:326:PRO:HD2	1.92	0.40
1:A:347:PHE:HB3	1:A:354:CYS:HB3	2.03	0.40
1:B:206:ARG:NH1	1:B:206:ARG:CG	2.85	0.40
3:B:1408:PXI:H142	3:B:1408:PXI:O5	2.21	0.40
1:B:176:ASP:HB3	1:B:180:PHE:CE1	2.57	0.40
1:B:100:HIS:CD2	1:B:352:HIS:CE1	3.10	0.40
1:A:262:HIS:N	1:A:263:PRO:HD3	2.37	0.40
1:A:395:ILE:HG23	1:A:396:ARG:H	1.87	0.40
1:A:82:THR:HG23	1:A:85:GLU:CD	2.42	0.40
1:B:331:ASP:O	1:B:339:ARG:NH2	2.54	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	394/436 (90%)	330 (84%)	54 (14%)	10 (2%)	5 32
1	B	391/436 (90%)	337 (86%)	49 (12%)	5 (1%)	12 47

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	785/872 (90%)	667 (85%)	103 (13%)	15 (2%)	8	39

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	50	ASN
1	A	14	VAL
1	A	374	ARG
1	B	245	HIS
1	B	305	ASP
1	A	305	ASP
1	A	406	ARG
1	B	294	THR
1	B	382	VAL
1	A	86	ALA
1	A	246	GLU
1	A	333	HIS
1	B	369	ARG
1	A	156	VAL
1	A	332	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	319/355 (90%)	301 (94%)	18 (6%)	21	57
1	B	318/355 (90%)	307 (96%)	11 (4%)	36	69
All	All	637/710 (90%)	608 (95%)	29 (5%)	27	63

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	GLN
1	A	45	ARG
1	A	46	THR

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Mol	Chain	Res	Type
1	A	50	ASN
1	A	60	ARG
1	A	82	THR
1	A	88	LEU
1	A	89	ASN
1	A	106	LEU
1	A	114	ARG
1	A	176	ASP
1	A	203	ASP
1	A	231	GLU
1	A	238	HIS
1	A	245	HIS
1	A	373	GLU
1	A	380	LEU
1	A	395	ILE
1	B	20	LEU
1	B	22	GLN
1	B	106	LEU
1	B	150	TRP
1	B	209	ASP
1	B	234	LEU
1	B	270	ARG
1	B	277	ASP
1	B	285	ARG
1	B	394	MET
1	B	395	ILE

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

Mol	Chain	Res	Type
1	A	50	ASN
1	A	90	HIS
1	A	188	GLN
1	A	208	GLN
1	A	238	HIS
1	A	265	GLN
1	A	349	HIS
1	B	22	GLN
1	B	76	ASN
1	B	188	GLN
1	B	208	GLN
1	B	245	HIS

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Mol	Chain	Res	Type
1	B	265	GLN
1	B	333	HIS
1	B	344	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 carbohydrates 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	PXI	B	1408	-	33,33,33	1.59	9 (27%)	40,47,47	2.41	13 (32%)
2	HEM	A	1407	1	27,50,50	1.79	5 (18%)	17,82,82	1.47	2 (11%)
3	PXI	A	1408	-	33,33,33	1.82	8 (24%)	40,47,47	2.36	15 (37%)
2	HEM	B	1407	-	27,50,50	1.84	9 (33%)	17,82,82	1.25	2 (11%)

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	PXI	B	1408	-	5/5/13/15	10/43/59/59	0/1/2/2
2	HEM	A	1407	1	-	2/6/54/54	-
3	PXI	A	1408	-	5/5/13/15	7/43/59/59	0/1/2/2
2	HEM	B	1407	-	-	0/6/54/54	-

All (31) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1408	PXI	C19-C21	5.05	1.59	1.51
3	A	1408	PXI	C5-C4	4.58	1.61	1.51
2	A	1407	HEM	C3B-CAB	-4.49	1.38	1.47
2	A	1407	HEM	C3C-CAC	-4.13	1.39	1.47
2	B	1407	HEM	C3B-CAB	-3.85	1.40	1.47
2	A	1407	HEM	C3B-C2B	-3.77	1.35	1.40
3	B	1408	PXI	C5-C4	3.68	1.59	1.51
2	B	1407	HEM	C3C-CAC	-3.58	1.40	1.47
3	A	1408	PXI	O1-C3	3.56	1.52	1.46
3	B	1408	PXI	O1-C3	3.36	1.51	1.46
3	B	1408	PXI	C24-C3	3.26	1.61	1.53
3	A	1408	PXI	C5-C7	3.00	1.62	1.55
2	B	1407	HEM	C3C-C2C	-2.96	1.36	1.40
3	A	1408	PXI	O3-C8	2.86	1.49	1.41
3	A	1408	PXI	C15-C12	2.81	1.59	1.53
2	B	1407	HEM	CBB-CAB	2.70	1.47	1.29
3	B	1408	PXI	C5-C7	2.67	1.61	1.55
2	B	1407	HEM	CBC-CAC	2.63	1.46	1.29
3	B	1408	PXI	C15-C12	2.54	1.59	1.53
2	B	1407	HEM	C4A-NA	2.53	1.41	1.36
2	B	1407	HEM	C3B-C2B	-2.51	1.36	1.40
3	B	1408	PXI	O3-C8	2.51	1.48	1.41
2	A	1407	HEM	CBB-CAB	2.37	1.45	1.29
2	A	1407	HEM	CBC-CAC	2.35	1.44	1.29
3	A	1408	PXI	O4-C8	2.26	1.47	1.41
3	B	1408	PXI	C19-C21	2.16	1.54	1.51
3	B	1408	PXI	O4-C8	2.15	1.47	1.41
2	B	1407	HEM	C1C-C2C	2.10	1.47	1.42
3	A	1408	PXI	O1-C4	2.10	1.39	1.34
2	B	1407	HEM	C4D-C3D	2.03	1.47	1.42
3	B	1408	PXI	O1-C4	2.01	1.39	1.34

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1408	PXI	C8-O3-C7	-7.91	98.40	117.96
3	A	1408	PXI	C8-O3-C7	-7.43	99.57	117.96
3	B	1408	PXI	O1-C4-C5	5.66	123.99	111.56
3	A	1408	PXI	O1-C4-C5	5.08	122.71	111.56
3	B	1408	PXI	O1-C3-C2	4.92	114.77	106.92
3	A	1408	PXI	O1-C3-C2	4.83	114.64	106.92
3	A	1408	PXI	C3-O1-C4	3.91	123.76	117.89
3	B	1408	PXI	C6-C5-C7	-3.78	104.34	112.92
3	A	1408	PXI	C10-C9-C11	-3.47	107.95	113.40
3	B	1408	PXI	C11-C12-N1	-3.35	106.20	115.67
2	A	1407	HEM	C1D-C2D-C3D	3.06	109.12	107.00
3	A	1408	PXI	C11-C12-N1	-3.03	107.12	115.67
3	A	1408	PXI	C6-C5-C7	-3.02	106.08	112.92
3	B	1408	PXI	O1-C4-O2	-2.98	118.38	123.94
2	A	1407	HEM	CMC-C2C-C3C	2.97	130.24	124.68
3	B	1408	PXI	C10-C9-C11	-2.87	108.90	113.40
3	A	1408	PXI	O1-C4-O2	-2.86	118.60	123.94
3	B	1408	PXI	C17-C16-C18	-2.84	106.43	110.69
3	B	1408	PXI	C7-C5-C4	2.75	115.64	110.01
3	A	1408	PXI	C2-C3-C24	-2.52	106.33	114.96
3	B	1408	PXI	C18-C16-C7	2.45	117.67	111.64
3	B	1408	PXI	C11-C12-C15	-2.43	106.46	109.97
3	A	1408	PXI	C18-C16-C7	2.42	117.59	111.64
3	B	1408	PXI	O2-C4-C5	-2.41	117.62	124.08
2	B	1407	HEM	CBD-CAD-C3D	2.37	116.85	112.48
3	A	1408	PXI	C7-C5-C4	2.34	114.79	110.01
2	B	1407	HEM	CMA-C3A-C4A	-2.23	125.03	128.46
3	B	1408	PXI	C3-O1-C4	2.20	121.19	117.89
3	A	1408	PXI	C11-C12-C15	-2.19	106.81	109.97
3	A	1408	PXI	C14-N1-C13	-2.03	104.37	110.38
3	A	1408	PXI	C13-N1-C12	-2.02	107.04	113.11
3	A	1408	PXI	O2-C4-C5	-2.02	118.69	124.08

All (10) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	B	1408	PXI	C12
3	B	1408	PXI	C15
3	B	1408	PXI	C8
3	B	1408	PXI	C9
3	B	1408	PXI	C19
3	A	1408	PXI	C12
3	A	1408	PXI	C15

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Mol	Chain	Res	Type	Atom
3	A	1408	PXI	C8
3	A	1408	PXI	C9
3	A	1408	PXI	C19

All (19) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	1408	PXI	C23-C24-C3-C2
3	B	1408	PXI	C25-C24-C3-C2
3	B	1408	PXI	C23-C24-C3-O1
3	B	1408	PXI	C25-C24-C3-O1
3	B	1408	PXI	C4-C5-C7-C16
3	B	1408	PXI	C6-C5-C7-O3
2	A	1407	HEM	C2D-C3D-CAD-CBD
2	A	1407	HEM	C4D-C3D-CAD-CBD
3	A	1408	PXI	C4-C5-C7-C16
3	A	1408	PXI	C6-C5-C7-O3
3	A	1408	PXI	C18-C19-C21-O6
3	A	1408	PXI	C18-C19-C21-C22
3	B	1408	PXI	C22-C23-C24-C25
3	B	1408	PXI	C11-C12-N1-C14
3	A	1408	PXI	C11-C12-N1-C14
3	A	1408	PXI	O4-C8-O3-C7
3	A	1408	PXI	C15-C8-O3-C7
3	B	1408	PXI	O4-C8-O3-C7
3	B	1408	PXI	C15-C8-O3-C7

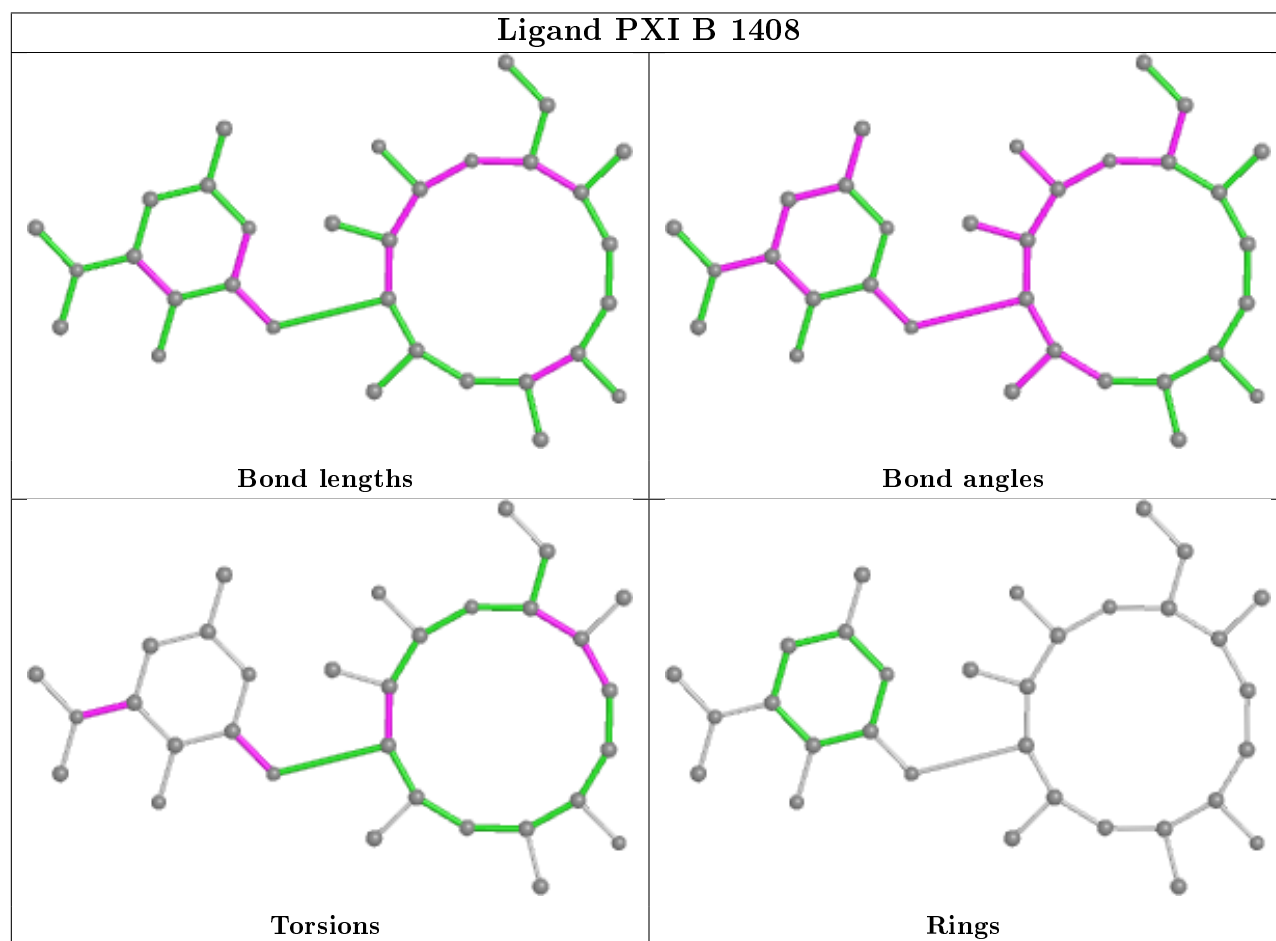
There are no ring outliers.

4 monomers are involved in 35 short contacts:

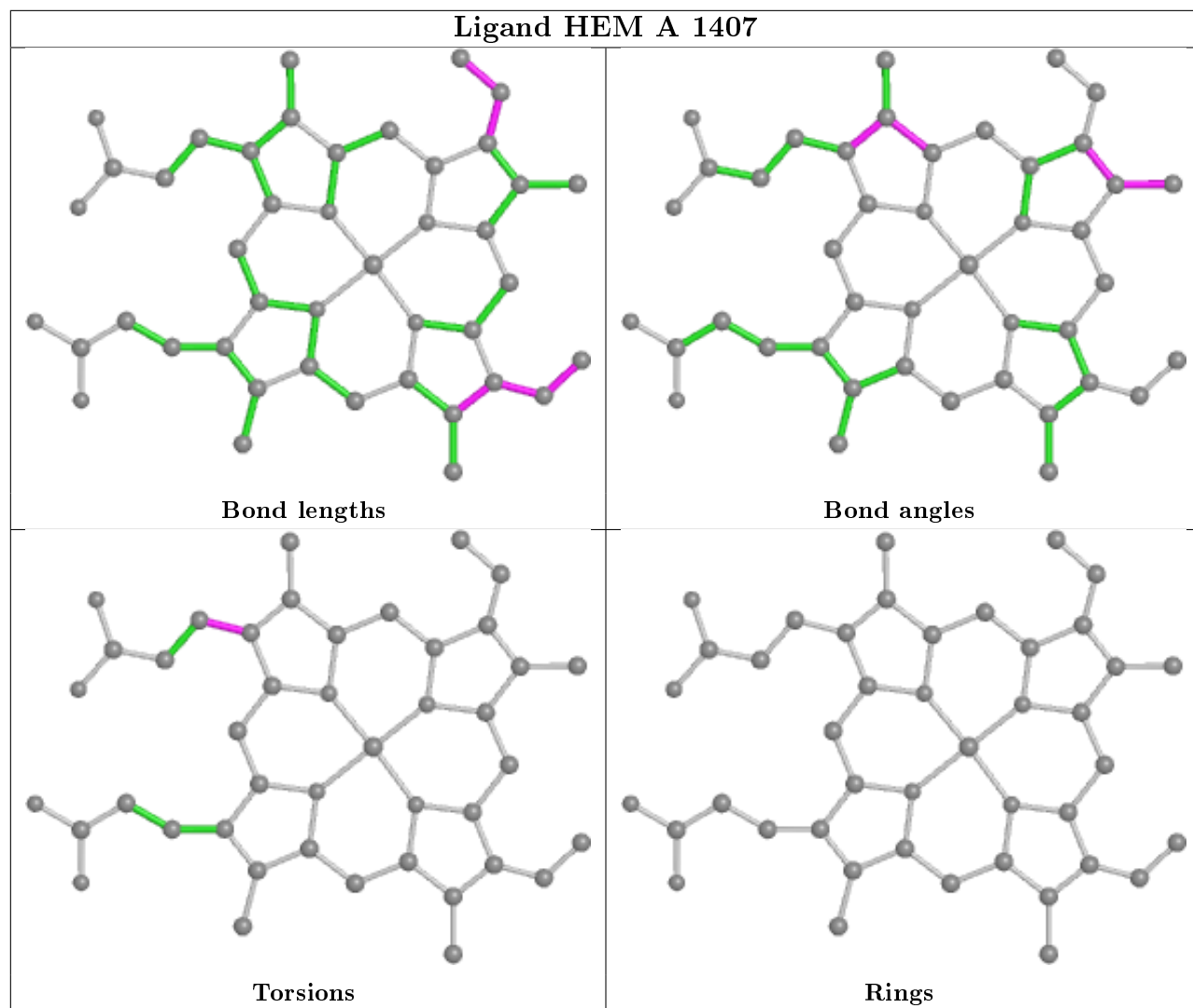
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1408	PXI	11	0
2	A	1407	HEM	8	0
3	A	1408	PXI	15	0
2	B	1407	HEM	1	0

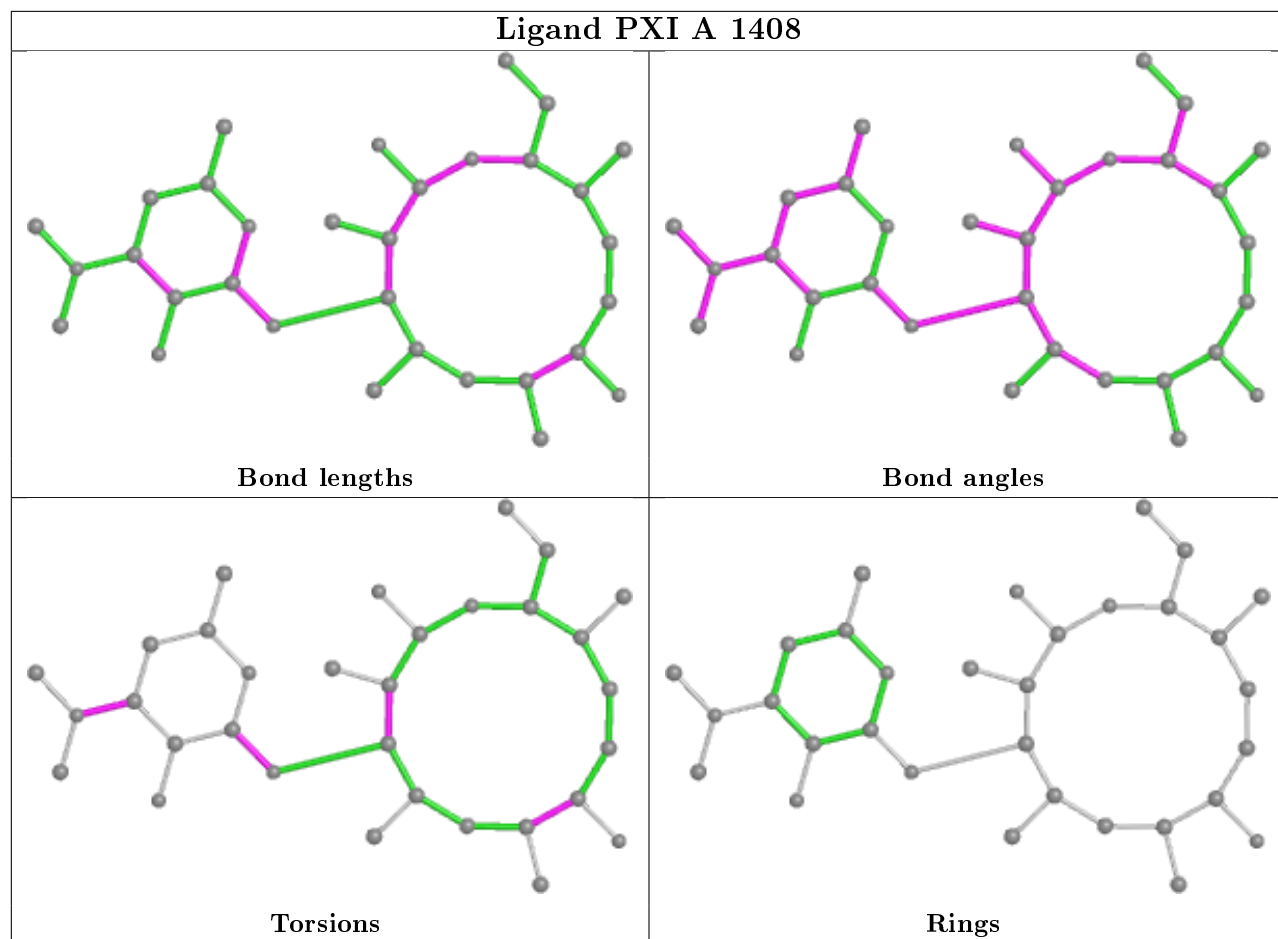
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be

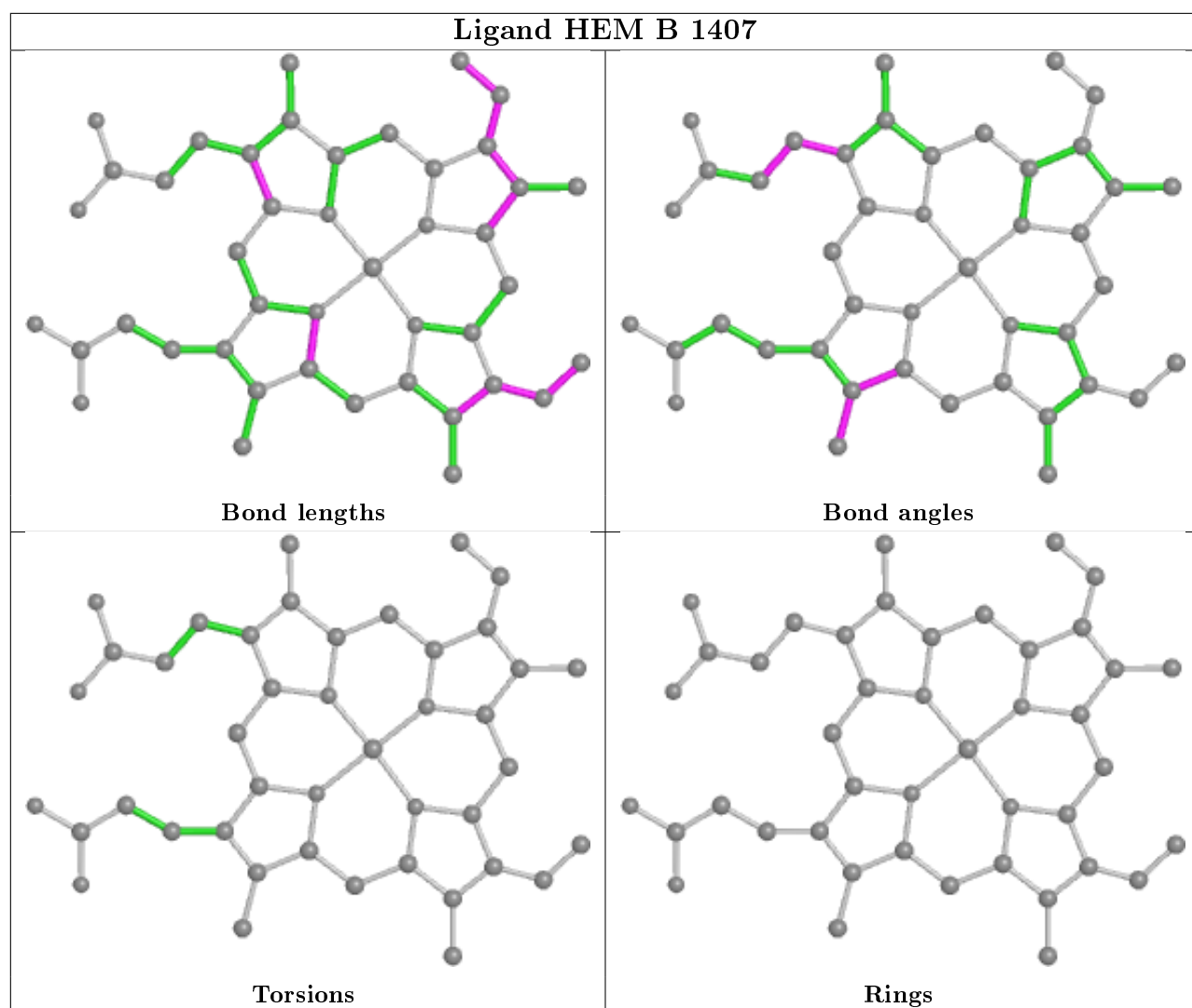
highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



Ligand HEM A 1407







5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

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

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	396/436 (90%)	-0.28	4 (1%) 82 72	26, 44, 71, 93	0
1	B	393/436 (90%)	-0.41	1 (0%) 94 92	19, 44, 73, 92	0
All	All	789/872 (90%)	-0.35	5 (0%) 89 83	19, 44, 72, 93	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	12	PRO	2.6
1	A	13	PRO	2.6
1	A	182	ASP	2.4
1	B	181	PRO	2.2
1	A	15	LEU	2.0

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

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

6.3 Carbohydrates [i](#)

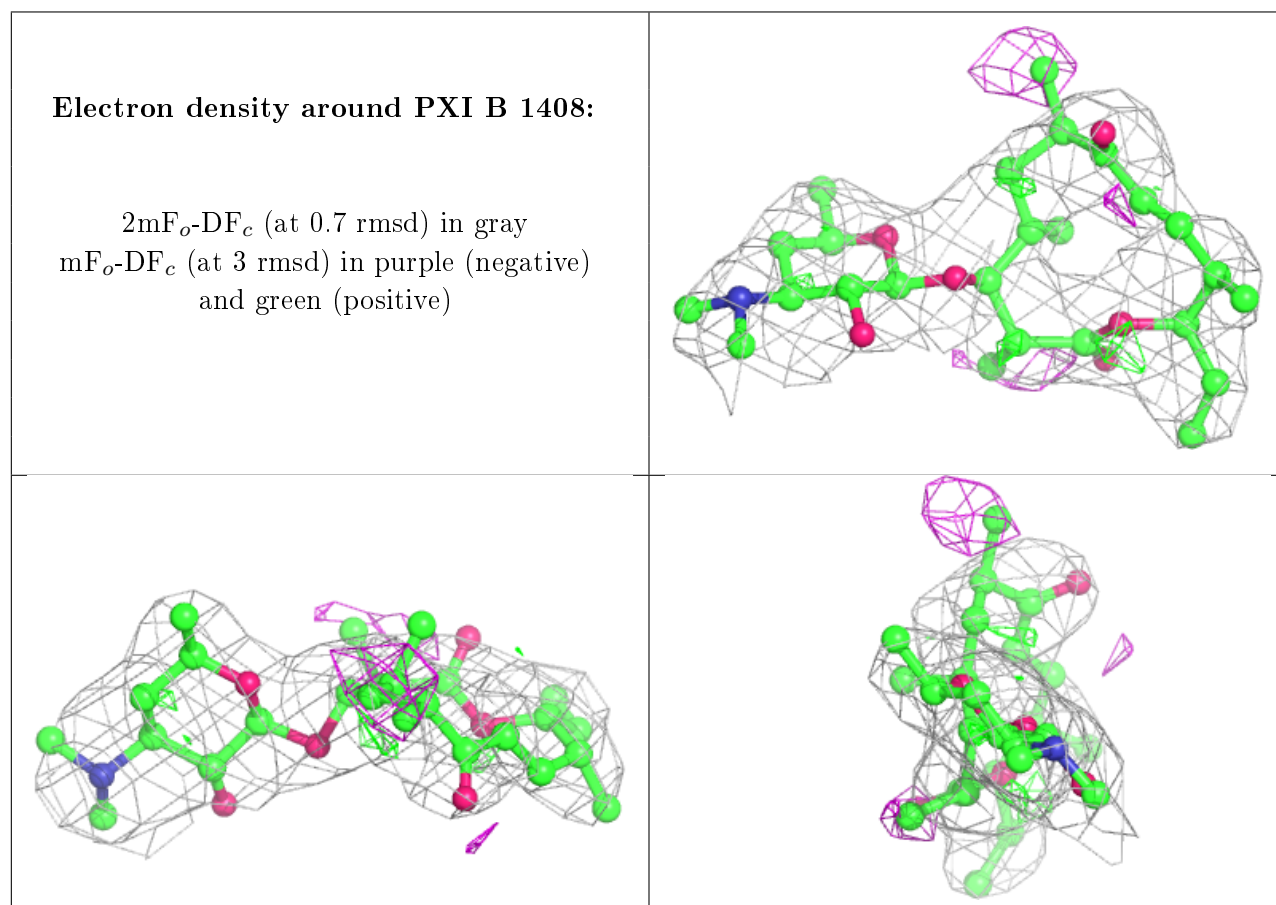
There are no carbohydrates in this entry.

6.4 Ligands [i](#)

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

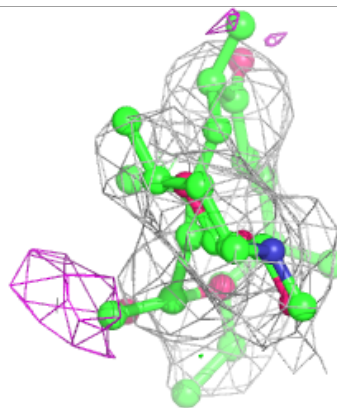
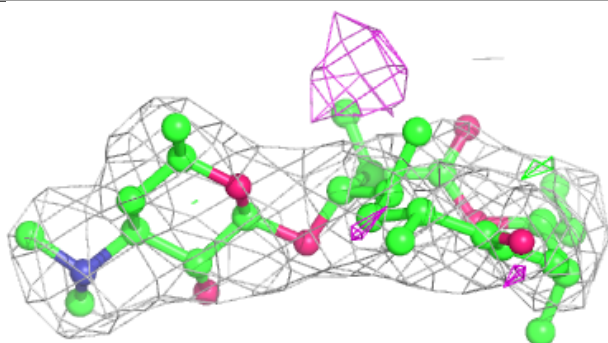
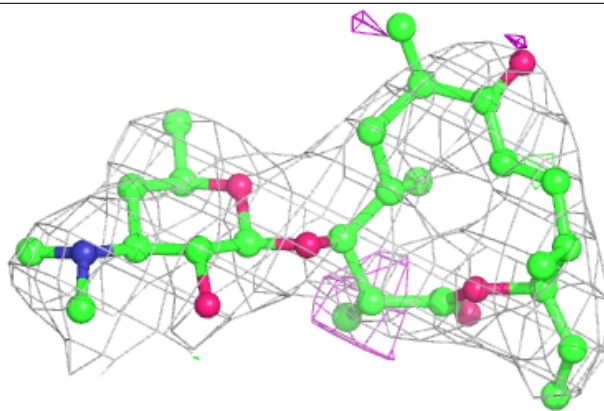
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	PXI	B	1408	32/32	0.91	0.31	36,42,49,52	0
3	PXI	A	1408	32/32	0.93	0.32	38,48,55,56	0
2	HEM	A	1407	43/43	0.98	0.16	25,30,32,33	0
2	HEM	B	1407	43/43	0.98	0.14	24,29,31,33	0

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.



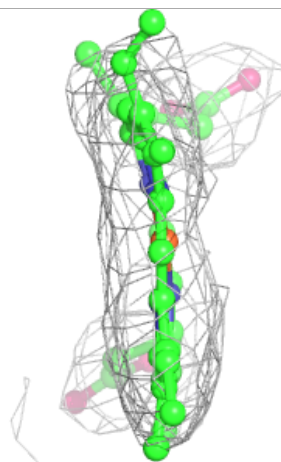
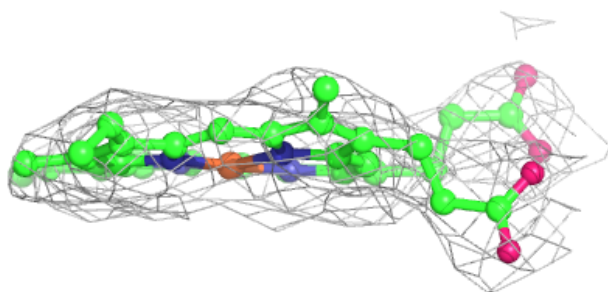
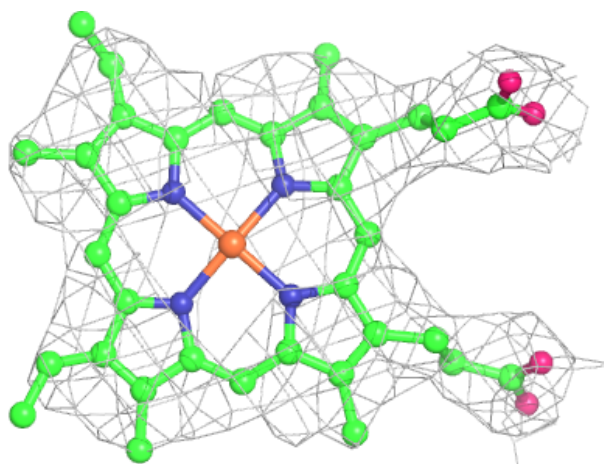
Electron density around PXI A 1408:

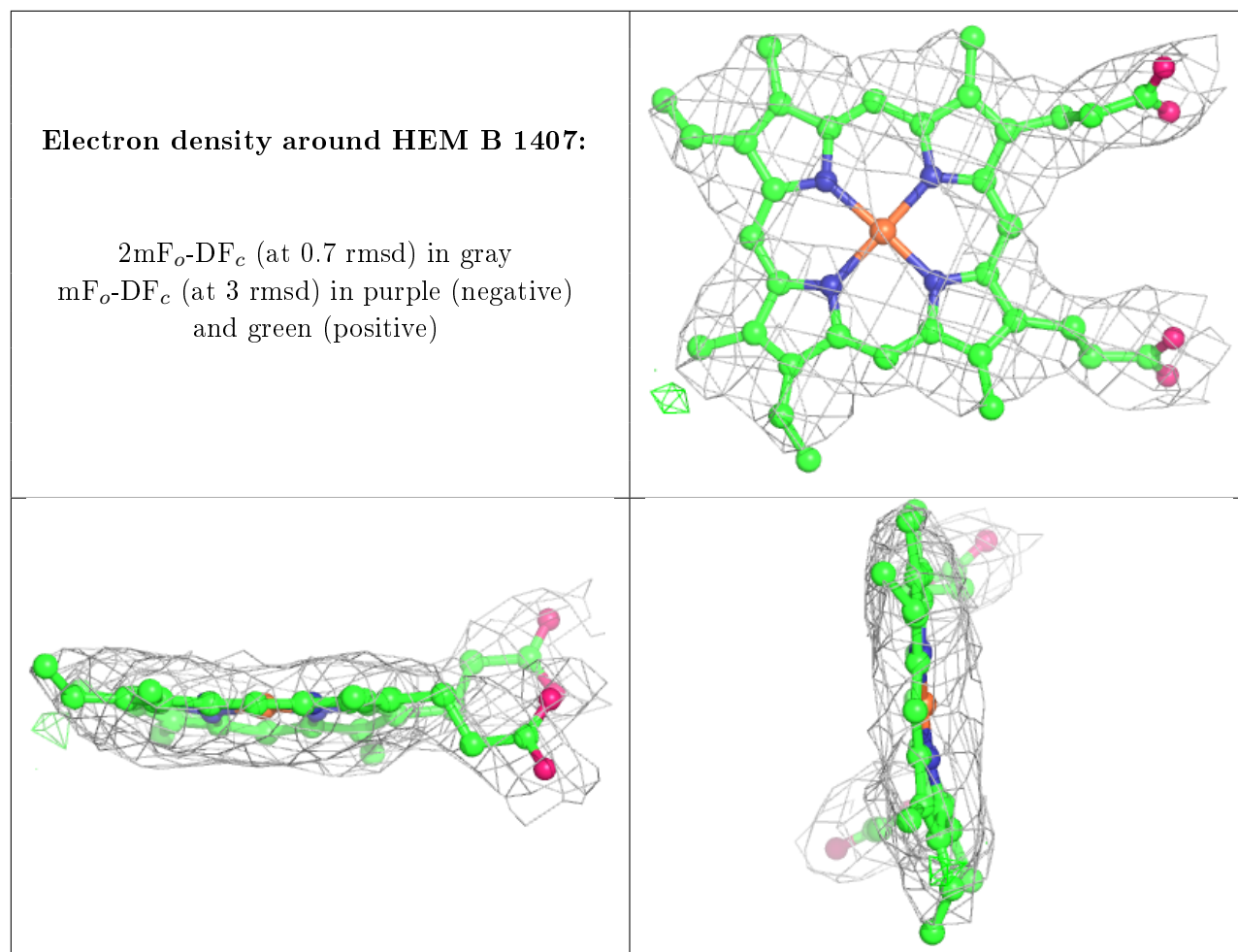
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around HEM A 1407:

$2mF_o-DF_c$ (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.