



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 5, 2026 – 05:56 PM EDT

PDB ID : 3E7T / pdb_00003e7t
Title : Structure of murine iNOS oxygenase domain with inhibitor AR-C102222
Authors : Garcin, E.D.; Arvai, A.S.; Rosenfeld, R.J.; Kroeger, M.D.; Crane, B.R.; Andersson, G.; Andrews, G.; Hamley, P.J.; Mallinder, P.R.; Nicholls, D.J.; St-Gallay, S.A.; Tinker, A.C.; Gensmantel, N.P.; Mete, A.; Cheshire, D.R.; Connolly, S.; Stuehr, D.J.; Aberg, A.; Wallace, A.V.; Tainer, J.A.; Getzoff, E.D.
Deposited on : 2008-08-18
Resolution : 2.60 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](#) ①) were used in the production of this report:

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

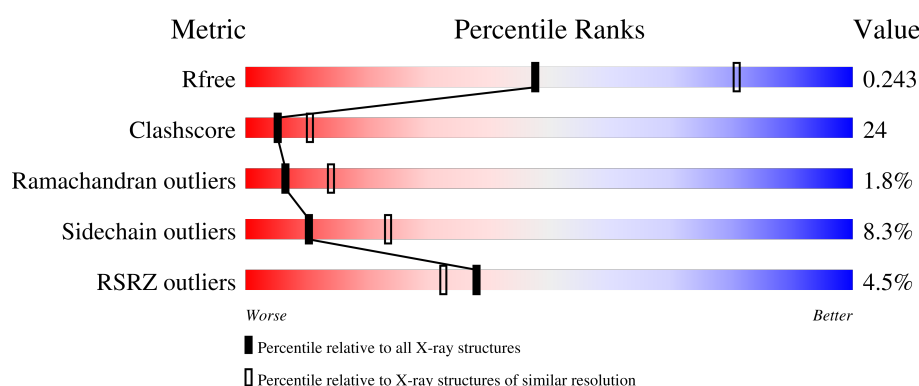
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	433	4% 56% 34% 5% 5%
1	B	433	5% 55% 35% 5% 5%

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 7241 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 Nitric oxide synthase, inducible.

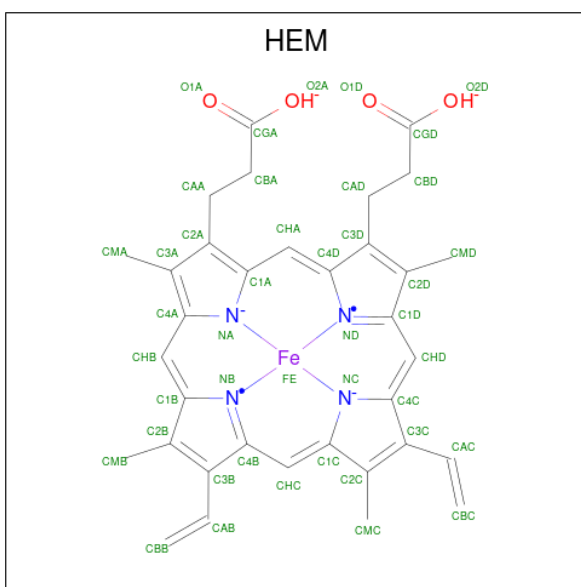
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	413	Total	C	N	O	S	0	0	0
			3368	2159	581	608	20			
1	B	413	Total	C	N	O	S	0	0	0
			3368	2159	581	608	20			

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



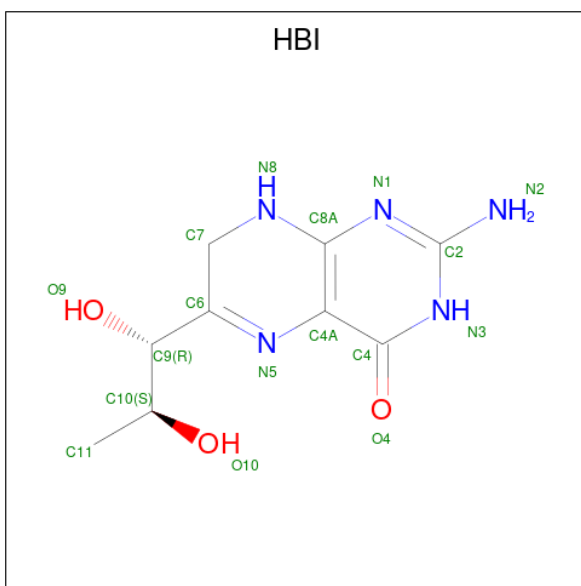
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: C₃₄H₃₂FeN₄O₄).



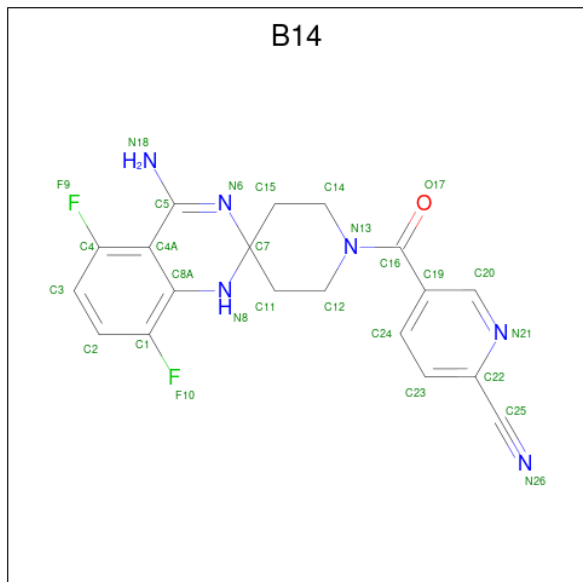
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
3	B	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 4 is 7,8-DIHYDROBIOPTERIN (CCD ID: HBI) (formula: $\text{C}_9\text{H}_{13}\text{N}_5\text{O}_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total 17	C 9	N 5	O 3	0	0
4	B	1	Total 17	C 9	N 5	O 3	0	0

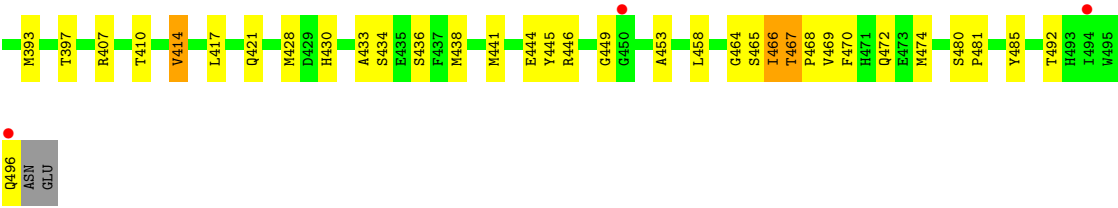
- Molecule 5 is 1-(6-CYANO-3-PYRIDYLCARBONYL)-5',8'-DIFLUOROSPIRO[PIPERIDINE-4,2'(1'H)-QUINAZOLINE]-4'-AMINE (CCD ID: B14) (formula: C₁₉H₁₆F₂N₆O).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	F	N	O	0	0
			28	19	2	6	1		
5	B	1	Total	C	F	N	O	0	0
			28	19	2	6	1		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	154	Total	O	0	0
			154	154		
6	B	170	Total	O	0	0
			170	170		



4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	214.16Å 214.16Å 116.95Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	19.99 – 2.60 19.99 – 2.60	Depositor EDS
% Data completeness (in resolution range)	90.6 (19.99-2.60) 90.7 (19.99-2.60)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.95 (at 2.61Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.232 , 0.263 0.220 , 0.243	Depositor DCC
R_{free} test set	2359 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	43.7	Xtriage
Anisotropy	0.624	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 47.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7241	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.17% 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: HBI, SO4, B14, HEM

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.45	0/3466	0.94	8/4712 (0.2%)
1	B	0.46	0/3466	0.93	12/4712 (0.3%)
All	All	0.45	0/6932	0.94	20/9424 (0.2%)

There are no bond length outliers.

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	348	ASN	N-CA-C	8.75	122.45	112.57
1	B	348	ASN	N-CA-C	8.57	123.04	112.59
1	A	485	TYR	N-CA-C	-7.71	100.14	110.55
1	A	125	ASP	N-CA-C	-7.22	106.39	114.62
1	A	122	GLY	CA-C-N	6.99	128.58	119.84
1	A	122	GLY	C-N-CA	6.99	128.58	119.84
1	B	485	TYR	N-CA-C	-6.33	101.42	110.59
1	A	367	TYR	N-CA-C	6.04	118.50	110.53
1	B	188	TRP	N-CA-C	-5.99	104.67	111.14
1	B	367	TYR	N-CA-C	5.77	118.64	110.50
1	A	365	GLY	N-CA-C	-5.73	101.66	110.77
1	B	234	ILE	N-CA-C	5.61	116.64	109.30
1	B	253	LEU	N-CA-C	-5.43	99.88	108.73
1	A	234	ILE	N-CA-C	5.35	117.16	109.51
1	B	87	GLY	N-CA-C	-5.30	108.32	115.21
1	B	386	LEU	N-CA-C	5.30	116.74	111.07
1	B	251	PHE	N-CA-C	-5.23	100.72	109.24
1	B	365	GLY	N-CA-C	-5.16	100.95	113.18
1	B	393	MET	N-CA-C	-5.15	107.04	113.38
1	B	466	ILE	N-CA-C	-5.06	106.65	112.98

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3368	0	3265	158	0
1	B	3368	0	3265	169	0
2	A	5	0	0	0	0
3	A	43	0	30	1	0
3	B	43	0	30	0	0
4	A	17	0	13	2	0
4	B	17	0	13	1	0
5	A	28	0	16	0	0
5	B	28	0	16	0	0
6	A	154	0	0	14	0
6	B	170	0	0	19	2
All	All	7241	0	6648	330	2

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 24.

All (330) 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:134:LEU:HD11	1:A:164:THR:HG22	1.33	1.04
1:B:438:MET:HE3	1:B:469:VAL:HG12	1.41	1.02
1:B:149:LYS:HD3	1:B:149:LYS:H	1.23	1.01
1:B:195:ILE:HG13	6:B:2143:HOH:O	1.71	0.90
1:B:110:LEU:HD23	1:B:110:LEU:H	1.36	0.88
1:A:407:ARG:HD3	6:A:1155:HOH:O	1.73	0.88
1:A:217:MET:HE2	1:A:305:ALA:HB2	1.57	0.87
1:A:134:LEU:O	1:A:138:ILE:HG12	1.76	0.86
1:B:99:THR:HG22	1:B:100:SER:H	1.41	0.85
1:A:368:MET:HE2	1:A:433:ALA:CB	2.10	0.82
1:B:149:LYS:HD3	1:B:149:LYS:N	1.95	0.82
1:B:99:THR:HG22	1:B:100:SER:N	1.94	0.82
1:B:301:LEU:HD13	1:B:315:ILE:HD11	1.62	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:266:MET:HE1	1:A:293:TYR:HD1	1.46	0.78
1:B:428:MET:HE3	6:B:2107:HOH:O	1.82	0.78
1:A:132:GLU:O	1:A:135:PRO:HD2	1.83	0.78
1:B:161:GLU:HG2	1:B:165:LYS:HE3	1.64	0.77
1:B:215:GLN:O	1:B:219:GLN:HG3	1.83	0.77
1:A:134:LEU:CD1	1:A:164:THR:HG22	2.14	0.77
1:A:438:MET:HE3	1:A:469:VAL:HG12	1.68	0.76
1:B:304:GLN:O	1:B:304:GLN:HG3	1.86	0.74
1:A:141:ILE:CD1	1:A:163:VAL:HG21	2.18	0.74
1:A:221:ILE:HG21	1:A:301:LEU:HD21	1.69	0.73
1:A:215:GLN:HE21	1:A:219:GLN:HE21	1.37	0.71
1:B:271:ILE:HD13	1:B:278:LEU:HD11	1.71	0.71
1:B:99:THR:CG2	1:B:100:SER:H	2.02	0.71
1:B:441:MET:HE1	1:B:472:GLN:CD	2.16	0.70
1:A:445:TYR:HA	1:A:450:GLY:H	1.56	0.70
1:B:177:LEU:O	1:B:181:ILE:HG12	1.91	0.70
1:B:249:HIS:C	1:B:306:ASP:O	2.35	0.69
1:B:149:LYS:H	1:B:149:LYS:CD	2.04	0.68
1:B:272:ARG:HH12	1:B:295:ARG:HG3	1.58	0.68
1:A:180:LEU:HB2	6:A:1086:HOH:O	1.93	0.67
1:B:428:MET:HE2	1:B:433:ALA:CB	2.25	0.67
1:A:123:PRO:HD3	1:A:487:ILE:HD12	1.77	0.66
1:A:224:HIS:HD2	1:A:239:THR:HG22	1.60	0.66
1:B:182:PHE:CZ	1:B:186:MET:HE3	2.30	0.66
1:B:441:MET:HE1	1:B:472:GLN:HG2	1.78	0.65
1:A:141:ILE:HD13	1:A:163:VAL:HG21	1.78	0.65
1:A:249:HIS:C	1:A:306:ASP:O	2.40	0.65
1:B:272:ARG:NH1	1:B:295:ARG:HG3	2.12	0.65
1:A:215:GLN:HG3	1:A:219:GLN:NE2	2.11	0.65
1:A:124:ARG:HG2	1:A:124:ARG:HH11	1.61	0.64
1:B:99:THR:CG2	1:B:100:SER:N	2.61	0.64
1:B:163:VAL:O	1:B:167:ILE:HG13	1.97	0.64
1:B:213:THR:OG1	1:B:216:GLU:HG3	1.97	0.64
1:B:285:ILE:HD11	1:B:291:PRO:HB3	1.78	0.64
1:A:411:GLU:O	1:A:414:VAL:HG22	1.96	0.64
1:B:127:PRO:HG3	1:B:246:ASP:HA	1.80	0.64
1:B:330:TYR:HD2	1:B:332:TRP:CZ2	2.16	0.64
1:A:494:ILE:O	1:A:495:TRP:HB2	1.99	0.63
1:A:134:LEU:HD11	1:A:164:THR:CG2	2.20	0.62
1:A:271:ILE:HD13	1:A:278:LEU:HD11	1.81	0.62
1:B:285:ILE:CD1	1:B:291:PRO:HB3	2.30	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:281:THR:HB	6:B:2134:HOH:O	2.00	0.61
1:A:241:PHE:HB3	1:A:242:PRO:CD	2.29	0.61
1:B:258:LEU:HB3	1:B:259:ILE:HD12	1.82	0.61
1:A:159:ARG:O	1:A:163:VAL:HG23	2.01	0.61
1:B:221:ILE:HG21	1:B:301:LEU:HD21	1.83	0.61
1:A:224:HIS:CD2	1:A:239:THR:HG22	2.36	0.61
1:A:266:MET:HE1	1:A:293:TYR:CD1	2.33	0.60
1:A:281:THR:O	1:A:285:ILE:HG12	2.02	0.60
1:A:332:TRP:CE3	1:A:392:ARG:HD2	2.37	0.60
1:B:244:ARG:HA	6:B:2018:HOH:O	2.02	0.59
1:A:438:MET:HE3	1:A:469:VAL:HA	1.84	0.59
1:B:441:MET:HE1	1:B:472:GLN:CG	2.31	0.59
1:A:163:VAL:O	1:A:167:ILE:HG13	2.02	0.59
1:A:223:ARG:HG3	1:A:223:ARG:HH11	1.68	0.58
1:A:266:MET:SD	1:A:272:ARG:HD3	2.43	0.58
1:A:217:MET:HE3	1:A:303:LEU:HB3	1.86	0.58
1:B:224:HIS:HE1	1:B:364:ASN:OD1	1.87	0.58
1:B:188:TRP:CE3	1:B:200:TRP:HA	2.39	0.58
1:A:80:ARG:NH2	6:A:1051:HOH:O	2.36	0.57
1:B:243:GLN:HB3	1:B:358:PHE:CE2	2.38	0.57
1:B:78:TYR:CD1	1:B:78:TYR:C	2.82	0.57
1:A:445:TYR:O	1:A:449:GLY:HA2	2.05	0.57
1:A:480:SER:HA	1:A:481:PRO:C	2.30	0.57
1:A:134:LEU:HB3	1:A:135:PRO:HD3	1.86	0.57
1:A:193:ARG:HD2	6:A:1008:HOH:O	2.05	0.57
1:B:259:ILE:HD12	1:B:259:ILE:N	2.19	0.57
1:A:141:ILE:HD11	1:A:163:VAL:HG21	1.86	0.57
1:B:467:THR:CG2	1:B:469:VAL:HG22	2.35	0.56
1:A:272:ARG:HH21	1:A:294:GLY:HA2	1.70	0.56
1:B:378:CYS:SG	1:B:386:LEU:HD23	2.44	0.56
1:A:368:MET:HE2	1:A:433:ALA:HB1	1.87	0.56
1:A:467:THR:CG2	1:A:469:VAL:HG22	2.35	0.56
1:A:303:LEU:N	1:A:303:LEU:CD1	2.68	0.56
1:A:124:ARG:HG2	1:A:124:ARG:NH1	2.18	0.56
1:B:428:MET:HE2	1:B:433:ALA:HA	1.88	0.56
1:A:124:ARG:HB3	1:A:247:GLY:HA3	1.88	0.55
1:B:266:MET:HB3	1:B:267:PRO:HD2	1.88	0.55
1:B:375:ARG:NH1	1:B:379:ASP:OD2	2.40	0.55
1:B:149:LYS:C	1:B:150:GLU:HG2	2.31	0.55
1:B:428:MET:HE2	1:B:433:ALA:HB2	1.87	0.55
1:B:141:ILE:HD11	1:B:163:VAL:HG11	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:318:ASP:HA	6:B:2164:HOH:O	2.07	0.55
1:A:195:ILE:HG21	1:A:368:MET:HE3	1.89	0.55
1:A:239:THR:CG2	1:A:362:PRO:HG2	2.36	0.54
1:A:375:ARG:O	1:A:379:ASP:HB2	2.07	0.54
1:A:304:GLN:O	1:A:304:GLN:HG3	2.07	0.54
1:B:385:ILE:HA	6:B:2127:HOH:O	2.06	0.54
1:B:182:PHE:CE2	1:B:186:MET:HE3	2.43	0.54
1:A:217:MET:CE	1:A:303:LEU:HB3	2.37	0.54
1:B:129:PRO:HG2	1:B:132:GLU:HG2	1.88	0.54
1:B:239:THR:O	1:B:361:CYS:HA	2.08	0.54
1:A:295:ARG:HD2	6:A:1144:HOH:O	2.06	0.53
1:A:241:PHE:HB3	1:A:242:PRO:HD2	1.90	0.53
1:A:271:ILE:HG23	1:A:278:LEU:HD21	1.90	0.53
1:B:96:HIS:C	1:B:98:ALA:H	2.14	0.53
1:B:272:ARG:HH11	1:B:272:ARG:HG2	1.73	0.53
1:A:217:MET:HE1	1:A:304:GLN:C	2.34	0.53
1:B:130:LEU:HD21	1:B:167:ILE:HG22	1.90	0.53
1:A:496:GLN:H	1:A:496:GLN:NE2	2.07	0.53
1:B:134:LEU:HB3	1:B:135:PRO:HD3	1.89	0.53
1:A:246:ASP:CG	1:A:248:LYS:H	2.16	0.53
1:A:338:LEU:HD21	1:A:385:ILE:HD13	1.90	0.53
1:B:141:ILE:CD1	1:B:163:VAL:HG21	2.40	0.52
1:A:258:LEU:HB2	1:A:345:ALA:HB3	1.92	0.52
1:B:128:THR:HG22	6:B:2131:HOH:O	2.09	0.52
1:B:464:GLY:O	1:B:467:THR:HB	2.09	0.52
1:B:246:ASP:CG	1:B:247:GLY:H	2.18	0.52
1:B:382:ARG:HG3	1:B:382:ARG:HH11	1.75	0.52
1:B:407:ARG:HG3	6:B:2042:HOH:O	2.08	0.52
1:A:145:TYR:HE2	1:A:159:ARG:HG2	1.75	0.52
1:A:185:LYS:HE2	1:A:201:SER:O	2.09	0.52
1:A:194:CYS:O	1:A:197:ARG:HG3	2.10	0.52
1:B:374:VAL:O	1:B:378:CYS:HB2	2.09	0.52
1:A:133:LEU:C	1:A:133:LEU:HD23	2.35	0.51
1:B:176:THR:OG1	1:B:179:GLU:HG3	2.09	0.51
1:A:116:PRO:HG2	1:A:119:LEU:HB2	1.91	0.51
1:A:350:LEU:HD21	1:A:357:GLU:HB2	1.91	0.51
1:B:480:SER:HA	1:B:481:PRO:C	2.35	0.51
1:A:153:ILE:O	1:A:157:LEU:HD13	2.11	0.51
1:A:252:ARG:HD3	1:A:359:PRO:HB2	1.92	0.51
1:A:410:THR:O	1:A:414:VAL:HG13	2.10	0.51
1:B:114:MET:HE2	1:B:114:MET:HA	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:152:LYS:HD2	1:B:155:GLU:OE2	2.11	0.50
1:B:210:ASN:HD22	1:B:211:CYS:N	2.08	0.50
1:A:144:TYR:O	1:A:147:SER:HB3	2.11	0.50
1:B:428:MET:HE2	1:B:433:ALA:CA	2.41	0.50
1:B:466:ILE:O	1:B:466:ILE:HG22	2.11	0.50
1:A:303:LEU:O	1:A:310:PRO:HA	2.11	0.50
1:B:252:ARG:HD3	1:B:359:PRO:HB2	1.93	0.50
1:B:330:TYR:N	1:B:330:TYR:CD1	2.79	0.50
1:A:238:ILE:HG13	1:A:363:PHE:HB3	1.92	0.50
1:B:161:GLU:O	1:B:165:LYS:HG3	2.12	0.50
1:B:445:TYR:O	1:B:449:GLY:HA2	2.11	0.50
3:A:901:HEM:HBC2	3:A:901:HEM:HMC2	1.94	0.50
1:A:285:ILE:CD1	1:A:291:PRO:HB3	2.42	0.49
1:B:267:PRO:C	1:B:269:GLY:H	2.19	0.49
1:A:190:ASN:O	1:A:192:PRO:HD3	2.12	0.49
1:B:264:TYR:HB3	6:B:2161:HOH:O	2.11	0.49
1:B:175:LEU:H	1:B:175:LEU:HD22	1.77	0.49
1:B:257:GLN:OE1	1:B:260:ARG:NH1	2.44	0.49
1:A:298:VAL:HG21	1:A:320:VAL:HG11	1.93	0.49
1:A:405:LYS:O	1:A:409:VAL:HG23	2.13	0.49
1:A:301:LEU:HB3	1:A:303:LEU:CD1	2.43	0.49
1:A:384:ASN:HA	6:A:1048:HOH:O	2.12	0.49
1:B:110:LEU:H	1:B:110:LEU:CD2	2.16	0.49
1:B:368:MET:HE1	6:B:2140:HOH:O	2.12	0.49
1:A:386:LEU:HD22	1:A:404:TRP:CZ3	2.48	0.49
1:B:145:TYR:HB3	6:B:2152:HOH:O	2.12	0.49
1:A:460:PRO:HG3	6:A:1152:HOH:O	2.11	0.49
1:A:195:ILE:HG21	1:A:368:MET:CE	2.43	0.49
1:A:272:ARG:HG2	1:A:272:ARG:HH11	1.78	0.48
1:A:441:MET:HE1	1:A:474:MET:HB3	1.95	0.48
4:A:902:HBI:O10	4:A:902:HBI:H71	2.12	0.48
1:B:130:LEU:HD11	1:B:134:LEU:HD23	1.95	0.48
1:A:224:HIS:HE1	1:A:364:ASN:OD1	1.96	0.48
1:A:494:ILE:HG23	6:A:1099:HOH:O	2.14	0.48
1:B:113:ILE:C	1:B:114:MET:HE2	2.39	0.48
1:B:331:GLU:OE1	1:B:331:GLU:HA	2.14	0.48
1:B:332:TRP:O	1:B:335:GLU:HB2	2.12	0.48
1:B:241:PHE:HB3	1:B:242:PRO:HD2	1.95	0.48
1:A:332:TRP:O	1:A:335:GLU:HB2	2.13	0.48
1:A:350:LEU:C	1:A:350:LEU:HD23	2.39	0.48
1:B:304:GLN:NE2	6:B:2135:HOH:O	2.47	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:350:LEU:C	1:B:350:LEU:HD23	2.39	0.48
1:A:197:ARG:NH2	1:A:452:PRO:O	2.47	0.48
1:B:246:ASP:CG	1:B:247:GLY:N	2.72	0.48
1:B:110:LEU:HD23	1:B:110:LEU:N	2.17	0.48
1:B:141:ILE:HD11	1:B:163:VAL:HG21	1.96	0.47
1:A:132:GLU:C	1:A:135:PRO:HD2	2.39	0.47
1:A:301:LEU:HB3	1:A:303:LEU:HD11	1.96	0.47
1:A:438:MET:HE3	1:A:469:VAL:CA	2.44	0.47
1:A:494:ILE:O	1:A:495:TRP:CB	2.60	0.47
1:B:166:GLU:OE2	1:B:173:TYR:HA	2.14	0.47
1:B:330:TYR:HD2	1:B:332:TRP:HZ2	1.60	0.47
1:B:130:LEU:HD13	1:B:134:LEU:HB2	1.97	0.47
1:B:252:ARG:CD	1:B:359:PRO:HB2	2.44	0.47
1:A:89:ILE:O	1:A:90:LEU:HD23	2.15	0.47
1:A:164:THR:O	1:A:168:GLU:HG2	2.15	0.47
1:A:368:MET:HE3	1:A:428:MET:HE3	1.97	0.47
1:A:438:MET:HE3	1:A:469:VAL:CG1	2.41	0.47
1:A:467:THR:HG21	1:A:469:VAL:HG22	1.97	0.47
1:B:113:ILE:O	1:B:114:MET:HE2	2.15	0.47
1:B:241:PHE:HB3	1:B:242:PRO:CD	2.45	0.47
1:B:492:THR:HG23	1:B:492:THR:O	2.15	0.47
1:A:239:THR:HG23	1:A:362:PRO:HG2	1.97	0.46
1:B:204:GLN:HB3	1:B:237:ALA:HB2	1.97	0.46
1:B:348:ASN:C	1:B:348:ASN:HD22	2.23	0.46
1:B:438:MET:HE2	1:B:438:MET:HA	1.98	0.46
1:A:153:ILE:H	1:A:153:ILE:HG13	1.43	0.46
1:B:258:LEU:HB2	1:B:345:ALA:HB3	1.96	0.46
1:B:434:SER:HB3	1:B:467:THR:HG23	1.98	0.46
1:A:264:TYR:HB2	1:A:266:MET:CE	2.44	0.46
1:B:410:THR:O	1:B:414:VAL:HG13	2.16	0.46
1:A:438:MET:HG3	1:A:468:PRO:HB2	1.97	0.46
1:B:130:LEU:CD2	1:B:167:ILE:HG22	2.46	0.46
1:A:138:ILE:HG23	1:A:160:LEU:HD22	1.98	0.46
1:A:251:PHE:O	1:A:360:ALA:HB2	2.16	0.46
1:B:333:PHE:O	1:B:336:LEU:HB2	2.16	0.46
1:A:385:ILE:HA	6:A:1068:HOH:O	2.15	0.46
1:B:116:PRO:HG2	1:B:119:LEU:HB2	1.97	0.46
1:B:133:LEU:HD13	1:B:133:LEU:O	2.15	0.46
1:B:188:TRP:CD2	1:B:200:TRP:HA	2.50	0.46
1:B:438:MET:HE3	1:B:469:VAL:CG1	2.30	0.46
1:B:446:ARG:HD3	1:B:446:ARG:C	2.41	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:124:ARG:HH21	1:B:128:THR:HB	1.80	0.46
1:A:145:TYR:C	1:A:147:SER:H	2.25	0.45
1:B:133:LEU:HD13	1:B:133:LEU:C	2.41	0.45
1:B:330:TYR:N	1:B:330:TYR:HD1	2.15	0.45
1:B:154:GLU:HB3	6:B:2094:HOH:O	2.16	0.45
1:B:256:SER:HA	1:B:348:ASN:OD1	2.15	0.45
1:A:99:THR:HG22	1:A:100:SER:N	2.32	0.45
1:B:163:VAL:HG13	1:B:173:TYR:CD2	2.52	0.45
1:A:78:TYR:CD1	1:A:78:TYR:C	2.95	0.45
1:B:295:ARG:HG2	6:B:2162:HOH:O	2.16	0.45
1:B:304:GLN:HA	6:B:2075:HOH:O	2.15	0.45
1:B:327:HIS:CG	1:B:328:PRO:HD2	2.52	0.45
1:A:133:LEU:HD23	1:A:133:LEU:O	2.17	0.45
1:A:244:ARG:HD2	1:A:357:GLU:OE2	2.16	0.45
1:A:360:ALA:HA	6:A:1033:HOH:O	2.17	0.45
1:B:175:LEU:HD21	1:B:356:LEU:HD11	1.97	0.45
1:B:210:ASN:ND2	1:B:210:ASN:H	2.14	0.45
1:A:89:ILE:HG22	1:A:90:LEU:N	2.32	0.45
1:A:188:TRP:CE3	1:A:200:TRP:HA	2.51	0.45
1:A:305:ALA:O	1:A:307:GLY:N	2.50	0.44
1:A:444:GLU:HA	1:A:444:GLU:OE1	2.17	0.44
1:B:249:HIS:HA	1:B:306:ASP:O	2.17	0.44
1:B:428:MET:HA	6:B:2119:HOH:O	2.17	0.44
1:A:154:GLU:HG2	1:A:155:GLU:H	1.82	0.44
1:A:344:PRO:O	1:A:364:ASN:HB2	2.18	0.44
1:A:371:GLU:O	1:A:375:ARG:HB2	2.17	0.44
1:A:375:ARG:NH1	1:A:379:ASP:OD2	2.51	0.44
1:B:164:THR:O	1:B:168:GLU:HG3	2.16	0.44
1:B:177:LEU:C	1:B:177:LEU:HD13	2.42	0.44
1:B:453:ALA:HB1	1:B:458:LEU:CD1	2.48	0.44
1:A:123:PRO:CD	1:A:487:ILE:HD12	2.47	0.44
1:B:121:ARG:HA	1:B:121:ARG:HD3	1.75	0.44
1:A:191:ALA:O	1:A:197:ARG:NH1	2.51	0.44
1:A:494:ILE:CD1	1:A:496:GLN:HE22	2.30	0.44
1:B:387:GLU:HG3	1:B:397:THR:HG21	2.00	0.44
1:A:221:ILE:CG2	1:A:301:LEU:HD21	2.43	0.44
1:A:264:TYR:CE1	1:A:293:TYR:HA	2.52	0.44
1:A:301:LEU:HD13	1:A:315:ILE:HD11	2.00	0.44
1:B:136:HIS:O	1:B:139:GLU:HB3	2.18	0.44
1:B:372:ILE:HA	1:B:376:ASP:OD2	2.17	0.44
1:A:309:ASP:OD1	1:A:495:TRP:HA	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:266:MET:HE2	1:B:272:ARG:HD2	1.99	0.44
1:A:444:GLU:HB3	1:A:451:CYS:HB2	2.00	0.44
1:B:250:ASP:O	1:B:304:GLN:NE2	2.51	0.43
1:B:175:LEU:HD22	1:B:175:LEU:N	2.33	0.43
1:B:258:LEU:C	1:B:259:ILE:HD12	2.42	0.43
1:B:397:THR:O	1:B:397:THR:CG2	2.66	0.43
1:A:148:PHE:CD1	1:A:149:LYS:N	2.87	0.43
1:A:249:HIS:HA	1:A:306:ASP:O	2.18	0.43
1:A:470:PHE:CD1	1:A:470:PHE:C	2.97	0.43
1:A:154:GLU:H	1:A:154:GLU:CD	2.24	0.43
1:A:434:SER:HB3	1:A:468:PRO:HD2	1.99	0.43
1:B:144:TYR:CE2	1:B:179:GLU:HA	2.54	0.43
1:B:465:SER:HA	1:B:470:PHE:CG	2.53	0.43
1:A:215:GLN:HG3	1:A:219:GLN:HE21	1.82	0.43
1:B:95:HIS:O	1:B:98:ALA:HB2	2.19	0.43
1:B:166:GLU:O	1:B:170:THR:OG1	2.31	0.43
1:A:134:LEU:CD1	1:A:138:ILE:HD11	2.48	0.43
1:B:210:ASN:HD22	1:B:210:ASN:C	2.27	0.43
1:B:375:ARG:O	1:B:379:ASP:HB2	2.19	0.43
1:B:438:MET:HG3	1:B:468:PRO:HB2	1.99	0.43
1:A:272:ARG:NH1	6:A:1119:HOH:O	2.51	0.43
1:B:249:HIS:CA	1:B:306:ASP:O	2.67	0.43
1:B:254:TRP:HB2	1:B:302:VAL:HB	2.00	0.43
1:A:456:ILE:HG22	4:A:902:HBI:H8	1.84	0.42
1:B:397:THR:O	1:B:397:THR:HG22	2.19	0.42
1:A:430:HIS:ND1	1:A:430:HIS:N	2.66	0.42
1:A:494:ILE:CG2	6:A:1099:HOH:O	2.67	0.42
1:A:176:THR:OG1	1:A:179:GLU:HG3	2.19	0.42
1:A:333:PHE:HB3	6:A:1060:HOH:O	2.20	0.42
1:B:191:ALA:O	1:B:197:ARG:HD3	2.20	0.42
1:A:308:GLN:HA	1:A:308:GLN:NE2	2.35	0.42
1:A:494:ILE:HG13	1:A:495:TRP:N	2.34	0.42
1:B:78:TYR:C	1:B:78:TYR:HD1	2.25	0.42
1:B:436:SER:HB3	6:B:2080:HOH:O	2.20	0.42
1:B:467:THR:HG21	1:B:469:VAL:HG22	2.00	0.42
1:A:224:HIS:HD2	1:A:239:THR:CG2	2.32	0.42
1:A:349:MET:HE2	1:A:483:TYR:HB3	2.02	0.42
1:B:149:LYS:O	1:B:150:GLU:C	2.62	0.42
1:B:161:GLU:CG	1:B:165:LYS:HE3	2.42	0.42
1:B:353:VAL:HG13	1:B:356:LEU:HD11	2.02	0.42
1:B:434:SER:CB	1:B:467:THR:HG23	2.50	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:320:VAL:HB	6:A:1032:HOH:O	2.19	0.41
4:B:902:HBI:H71	4:B:902:HBI:O10	2.19	0.41
1:A:148:PHE:HD1	1:A:149:LYS:N	2.19	0.41
1:A:137:ALA:O	1:A:141:ILE:HG13	2.20	0.41
1:B:428:MET:CE	1:B:433:ALA:HA	2.50	0.41
1:B:430:HIS:ND1	1:B:430:HIS:N	2.69	0.41
1:B:272:ARG:NH1	1:B:272:ARG:HG2	2.36	0.41
1:A:250:ASP:N	1:A:306:ASP:O	2.53	0.41
1:B:294:GLY:N	1:B:297:ASP:OD2	2.53	0.41
1:A:411:GLU:HA	1:A:414:VAL:HG22	2.02	0.41
1:B:467:THR:HG22	1:B:469:VAL:HG22	2.02	0.41
1:A:110:LEU:N	1:A:110:LEU:HD22	2.35	0.41
1:A:249:HIS:CA	1:A:306:ASP:O	2.68	0.41
1:B:217:MET:HB3	1:B:303:LEU:HD13	2.03	0.41
1:A:152:LYS:HD2	1:A:155:GLU:OE2	2.21	0.41
1:A:368:MET:HE2	1:A:433:ALA:HB3	2.00	0.41
1:B:116:PRO:CG	1:B:119:LEU:HD22	2.50	0.41
1:B:224:HIS:HD2	1:B:239:THR:OG1	2.03	0.41
1:B:376:ASP:HA	1:B:382:ARG:NH1	2.36	0.41
1:A:94:LEU:O	1:A:96:HIS:N	2.54	0.40
1:A:246:ASP:CG	1:A:247:GLY:N	2.79	0.40
1:B:194:CYS:O	1:B:197:ARG:HG3	2.21	0.40
1:B:257:GLN:HE21	1:B:257:GLN:HB3	1.58	0.40
1:B:348:ASN:HB2	6:B:2136:HOH:O	2.20	0.40
1:A:346:VAL:HB	1:A:363:PHE:CE1	2.56	0.40
1:B:198:ILE:HG13	6:B:2003:HOH:O	2.21	0.40
1:A:148:PHE:HD1	1:A:149:LYS:O	2.02	0.40
1:A:252:ARG:HB2	1:A:304:GLN:HG2	2.04	0.40
1:A:285:ILE:HD11	1:A:291:PRO:HB3	2.03	0.40
1:B:188:TRP:CZ3	1:B:200:TRP:HA	2.56	0.40

All (2) 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)
6:B:2118:HOH:O	6:B:2118:HOH:O[9_766]	0.50	1.70
6:B:2102:HOH:O	6:B:2102:HOH:O[9_766]	0.51	1.69

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	409/433 (94%)	363 (89%)	38 (9%)	8 (2%)	6	12
1	B	409/433 (94%)	372 (91%)	30 (7%)	7 (2%)	7	15
All	All	818/866 (94%)	735 (90%)	68 (8%)	15 (2%)	6	14

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	151	ALA
1	A	494	ILE
1	A	495	TRP
1	A	200	TRP
1	A	95	HIS
1	A	306	ASP
1	B	150	GLU
1	B	200	TRP
1	B	306	ASP
1	A	150	GLU
1	B	308	GLN
1	B	307	GLY
1	B	268	ASP
1	B	384	ASN
1	A	307	GLY

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	361/381 (95%)	330 (91%)	31 (9%)	10	22
1	B	361/381 (95%)	332 (92%)	29 (8%)	11	25
All	All	722/762 (95%)	662 (92%)	60 (8%)	10	23

All (60) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	110	LEU
1	A	119	LEU
1	A	124	ARG
1	A	125	ASP
1	A	129	PRO
1	A	153	ILE
1	A	161	GLU
1	A	164	THR
1	A	170	THR
1	A	178	ASP
1	A	180	LEU
1	A	257	GLN
1	A	258	LEU
1	A	301	LEU
1	A	303	LEU
1	A	334	GLN
1	A	336	LEU
1	A	338	LEU
1	A	348	ASN
1	A	386	LEU
1	A	387	GLU
1	A	395	LEU
1	A	397	THR
1	A	400	LEU
1	A	417	LEU
1	A	444	GLU
1	A	446	ARG
1	A	467	THR
1	A	469	VAL
1	A	495	TRP
1	A	496	GLN
1	B	119	LEU
1	B	125	ASP
1	B	126	LYS
1	B	128	THR

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Mol	Chain	Res	Type
1	B	130	LEU
1	B	134	LEU
1	B	149	LYS
1	B	175	LEU
1	B	178	ASP
1	B	210	ASN
1	B	226	LEU
1	B	257	GLN
1	B	258	LEU
1	B	259	ILE
1	B	301	LEU
1	B	304	GLN
1	B	330	TYR
1	B	336	LEU
1	B	348	ASN
1	B	356	LEU
1	B	375	ARG
1	B	386	LEU
1	B	414	VAL
1	B	417	LEU
1	B	421	GLN
1	B	444	GLU
1	B	467	THR
1	B	474	MET
1	B	496	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	96	HIS
1	A	115	ASN
1	A	202	ASN
1	A	219	GLN
1	A	224	HIS
1	A	231	ASN
1	A	233	ASN
1	A	243	GLN
1	A	308	GLN
1	A	334	GLN
1	A	348	ASN
1	A	421	GLN
1	A	486	GLN

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Mol	Chain	Res	Type
1	A	496	GLN
1	B	77	GLN
1	B	142	ASN
1	B	199	GLN
1	B	210	ASN
1	B	215	GLN
1	B	224	HIS
1	B	231	ASN
1	B	233	ASN
1	B	348	ASN
1	B	418	HIS
1	B	476	ASN
1	B	486	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

7 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	$\# Z > 2$
4	HBI	B	902	-	15,18,18	3.20	4 (26%)	14,26,26	1.92	4 (28%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	SO4	A	903	-	4,4,4	0.39	0	6,6,6	0.14	0
4	HBI	A	902	-	15,18,18	3.29	4 (26%)	14,26,26	1.87	4 (28%)
3	HEM	A	901	1	50,50,50	1.77	10 (20%)	67,82,82	1.21	6 (8%)
5	B14	A	906	-	27,31,31	2.57	12 (44%)	37,46,46	1.91	11 (29%)
5	B14	B	906	-	27,31,31	2.57	13 (48%)	37,46,46	1.92	11 (29%)
3	HEM	B	901	1	50,50,50	1.66	7 (14%)	67,82,82	1.20	8 (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
4	HBI	B	902	-	-	0/4/17/17	0/2/2/2
4	HBI	A	902	-	-	0/4/17/17	0/2/2/2
3	HEM	A	901	1	-	2/14/54/54	-
5	B14	A	906	-	-	2/8/36/36	0/3/4/4
5	B14	B	906	-	-	2/8/36/36	0/3/4/4
3	HEM	B	901	1	-	0/14/54/54	-

All (50) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	902	HBI	C6-N5	10.62	1.45	1.28
4	B	902	HBI	C6-N5	10.13	1.44	1.28
5	B	906	B14	C25-N26	8.56	1.33	1.14
5	A	906	B14	C25-N26	8.39	1.33	1.14
4	A	902	HBI	C7-N8	-4.78	1.38	1.45
4	B	902	HBI	C7-N8	-4.75	1.38	1.45
3	A	901	HEM	FE-NB	4.67	2.09	1.94
3	B	901	HEM	FE-NB	4.63	2.09	1.94
3	A	901	HEM	FE-NC	4.32	2.09	1.95
3	B	901	HEM	FE-ND	4.07	2.07	1.94
3	A	901	HEM	FE-ND	3.94	2.07	1.94
5	B	906	B14	C20-N21	3.66	1.41	1.34
3	B	901	HEM	FE-NC	3.65	2.07	1.95
3	B	901	HEM	FE-NA	3.49	2.06	1.95
3	A	901	HEM	C1A-C2A	3.43	1.51	1.44
4	B	902	HBI	C7-C6	-3.31	1.46	1.49
5	B	906	B14	C19-C16	-3.29	1.44	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	906	B14	C19-C16	-3.27	1.44	1.50
5	B	906	B14	C20-C19	3.23	1.44	1.39
5	A	906	B14	C22-C25	-3.22	1.36	1.45
5	A	906	B14	C20-C19	3.19	1.44	1.39
5	A	906	B14	C8A-C1	3.18	1.43	1.38
5	B	906	B14	C22-C25	-3.17	1.36	1.45
5	A	906	B14	C20-N21	3.14	1.40	1.34
3	A	901	HEM	CAB-C3B	-2.99	1.39	1.47
4	A	902	HBI	C7-C6	-2.92	1.46	1.49
3	A	901	HEM	C4D-C3D	2.80	1.49	1.45
4	A	902	HBI	C4-N3	-2.80	1.33	1.38
3	A	901	HEM	FE-NA	2.70	2.04	1.95
4	B	902	HBI	C4-N3	-2.69	1.33	1.38
5	B	906	B14	C4A-C4	2.69	1.43	1.39
3	B	901	HEM	CAC-C3C	-2.67	1.40	1.47
5	A	906	B14	C4A-C4	2.65	1.43	1.39
3	B	901	HEM	C3B-C4B	2.62	1.50	1.44
5	B	906	B14	C16-N13	2.56	1.40	1.34
5	A	906	B14	C5-N6	2.49	1.35	1.29
3	A	901	HEM	CHA-C4D	2.47	1.43	1.38
5	A	906	B14	C16-N13	2.40	1.39	1.34
5	B	906	B14	C8A-C1	2.39	1.42	1.38
5	B	906	B14	C5-N6	2.39	1.35	1.29
5	A	906	B14	C22-N21	2.28	1.40	1.35
5	B	906	B14	F9-C4	-2.26	1.29	1.35
5	A	906	B14	C24-C19	2.21	1.42	1.39
5	B	906	B14	F10-C1	-2.20	1.29	1.35
5	B	906	B14	C5-N18	-2.20	1.30	1.34
5	A	906	B14	F10-C1	-2.13	1.29	1.35
3	A	901	HEM	CBC-CAC	2.13	1.40	1.30
3	B	901	HEM	CAB-C3B	-2.11	1.41	1.47
3	A	901	HEM	CAC-C3C	-2.07	1.41	1.47
5	B	906	B14	C22-N21	2.04	1.39	1.35

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	906	B14	C14-N13-C12	-4.94	102.60	112.68
5	A	906	B14	C14-N13-C12	-4.94	102.62	112.68
5	B	906	B14	C14-C15-C7	-4.53	106.50	112.55
5	A	906	B14	C14-C15-C7	-4.31	106.80	112.55
4	B	902	HBI	C2-N1-C8A	3.96	120.34	113.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	906	B14	C19-C20-N21	-3.95	117.78	123.59
4	A	902	HBI	C2-N1-C8A	3.90	120.24	113.36
5	B	906	B14	C19-C20-N21	-3.79	118.01	123.59
3	B	901	HEM	CMD-C2D-C1D	3.66	130.75	125.03
3	A	901	HEM	CMD-C2D-C1D	3.43	130.39	125.03
4	B	902	HBI	O4-C4-C4A	-3.40	117.57	126.53
4	A	902	HBI	O4-C4-C4A	-3.28	117.87	126.53
5	A	906	B14	C20-N21-C22	3.24	121.84	116.73
3	A	901	HEM	CAA-C2A-C1A	3.05	130.90	124.94
5	B	906	B14	C20-N21-C22	2.91	121.33	116.73
5	B	906	B14	C15-C14-N13	-2.76	105.66	110.82
3	A	901	HEM	CHC-C4B-NB	-2.76	121.45	124.42
4	B	902	HBI	C2-N3-C4	-2.72	120.18	125.11
4	A	902	HBI	C2-N3-C4	-2.69	120.23	125.11
3	B	901	HEM	C4B-C3B-C2B	-2.62	104.87	107.28
5	A	906	B14	C15-C14-N13	-2.62	105.93	110.82
5	B	906	B14	C15-C7-N8	2.57	114.49	110.47
3	B	901	HEM	CHC-C4B-NB	-2.55	121.68	124.42
5	A	906	B14	C3-C4-C4A	-2.51	118.88	123.48
5	B	906	B14	C8A-C4A-C5	-2.44	114.87	117.41
4	B	902	HBI	C4A-C4-N3	2.43	119.45	113.25
5	A	906	B14	C12-C11-C7	-2.40	109.34	112.55
4	A	902	HBI	C4A-C4-N3	2.40	119.37	113.25
5	B	906	B14	C3-C4-C4A	-2.35	119.17	123.48
5	B	906	B14	C12-C11-C7	-2.34	109.42	112.55
5	A	906	B14	C8A-C4A-C5	-2.29	115.03	117.41
5	A	906	B14	C11-C12-N13	-2.27	106.59	110.82
3	A	901	HEM	CAB-C3B-C2B	-2.26	121.08	128.43
3	B	901	HEM	CAD-C3D-C2D	-2.24	123.68	127.87
3	A	901	HEM	C4C-C3C-C2C	-2.23	104.88	106.81
3	A	901	HEM	CAD-C3D-C4D	2.16	128.47	124.70
3	B	901	HEM	C3B-C4B-NB	2.16	111.02	109.47
5	A	906	B14	C15-C7-N8	2.14	113.81	110.47
3	B	901	HEM	CAB-C3B-C2B	-2.12	121.53	128.43
3	B	901	HEM	CAA-C2A-C1A	2.12	129.08	124.94
5	B	906	B14	C11-C12-N13	-2.08	106.93	110.82
3	B	901	HEM	CAB-C3B-C4B	2.07	133.55	124.39
5	B	906	B14	F9-C4-C4A	2.03	121.71	118.15
5	A	906	B14	C12-N13-C16	-2.01	116.56	122.79

There are no chirality outliers.

All (6) torsion outliers are listed below:

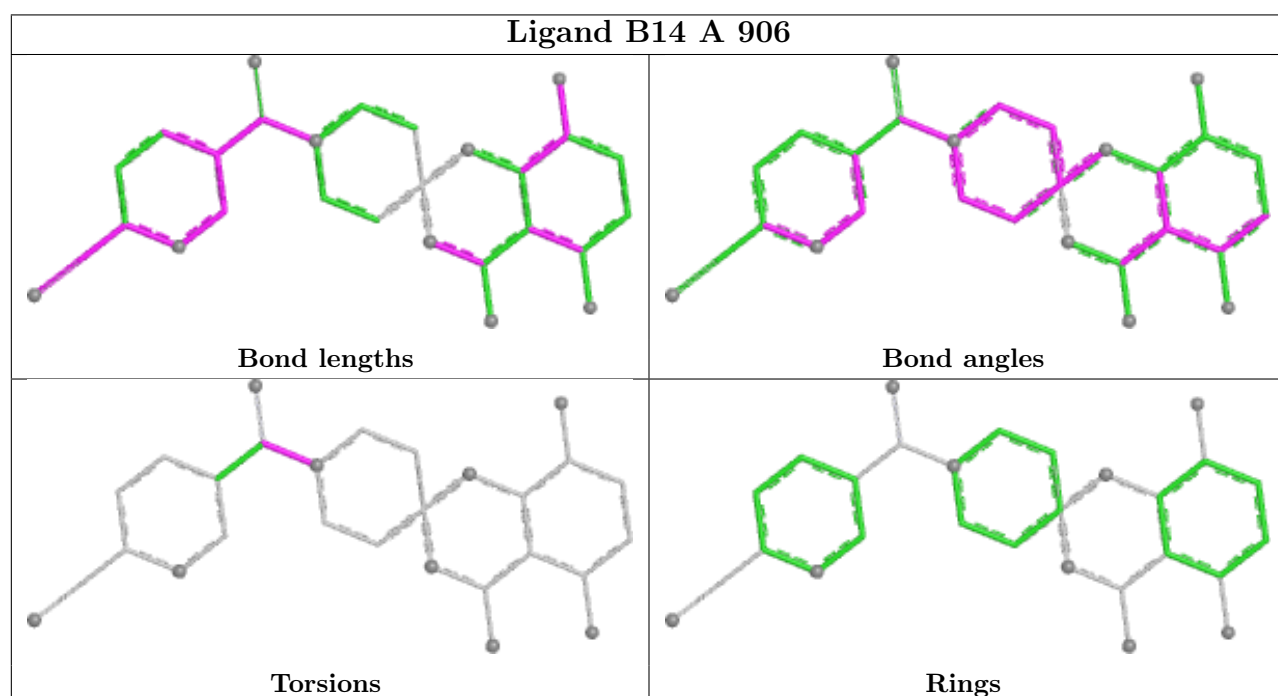
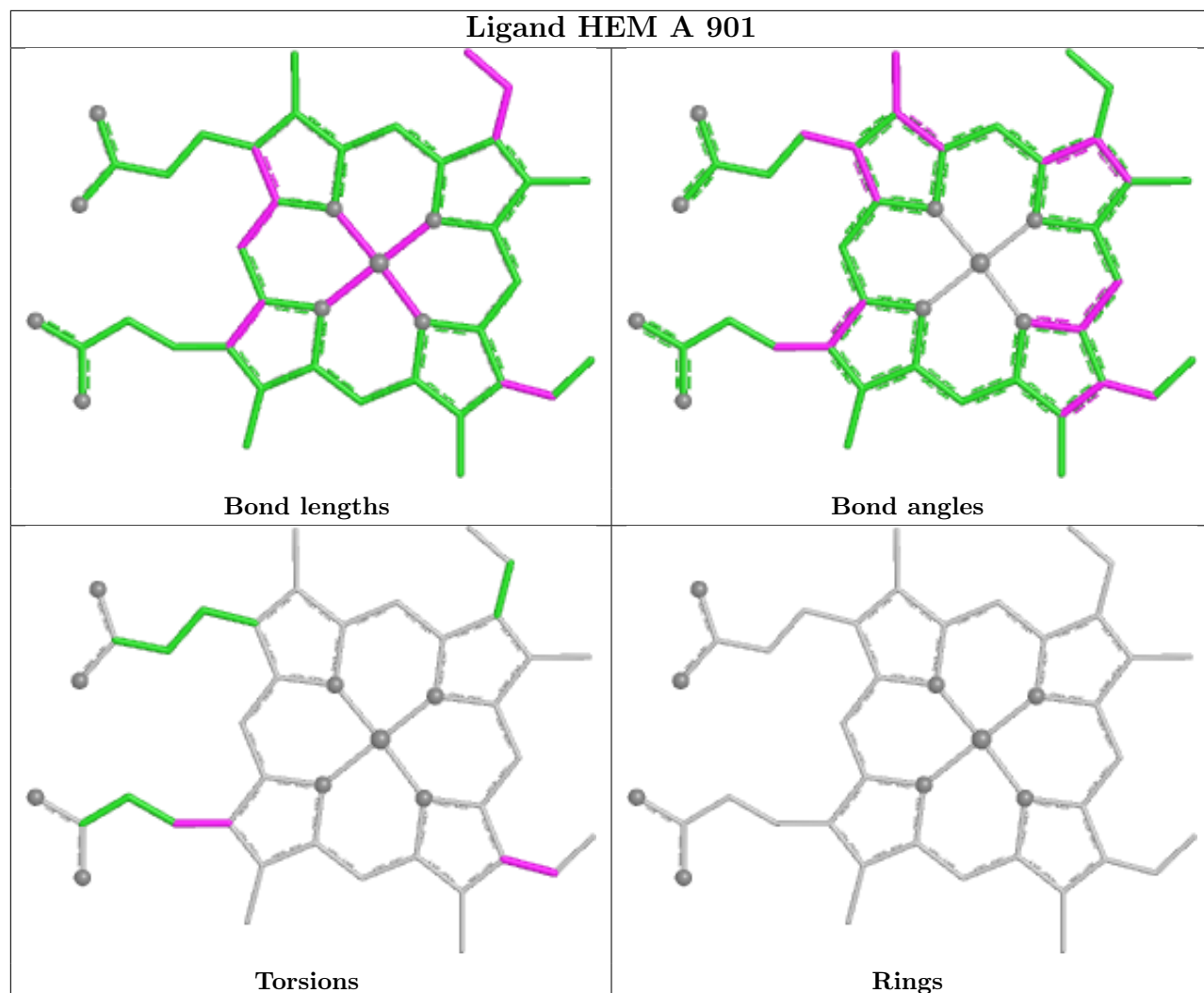
Mol	Chain	Res	Type	Atoms
5	A	906	B14	O17-C16-N13-C14
5	B	906	B14	O17-C16-N13-C14
5	A	906	B14	C19-C16-N13-C14
5	B	906	B14	C19-C16-N13-C14
3	A	901	HEM	C1A-C2A-CAA-CBA
3	A	901	HEM	C4B-C3B-CAB-CBB

There are no ring outliers.

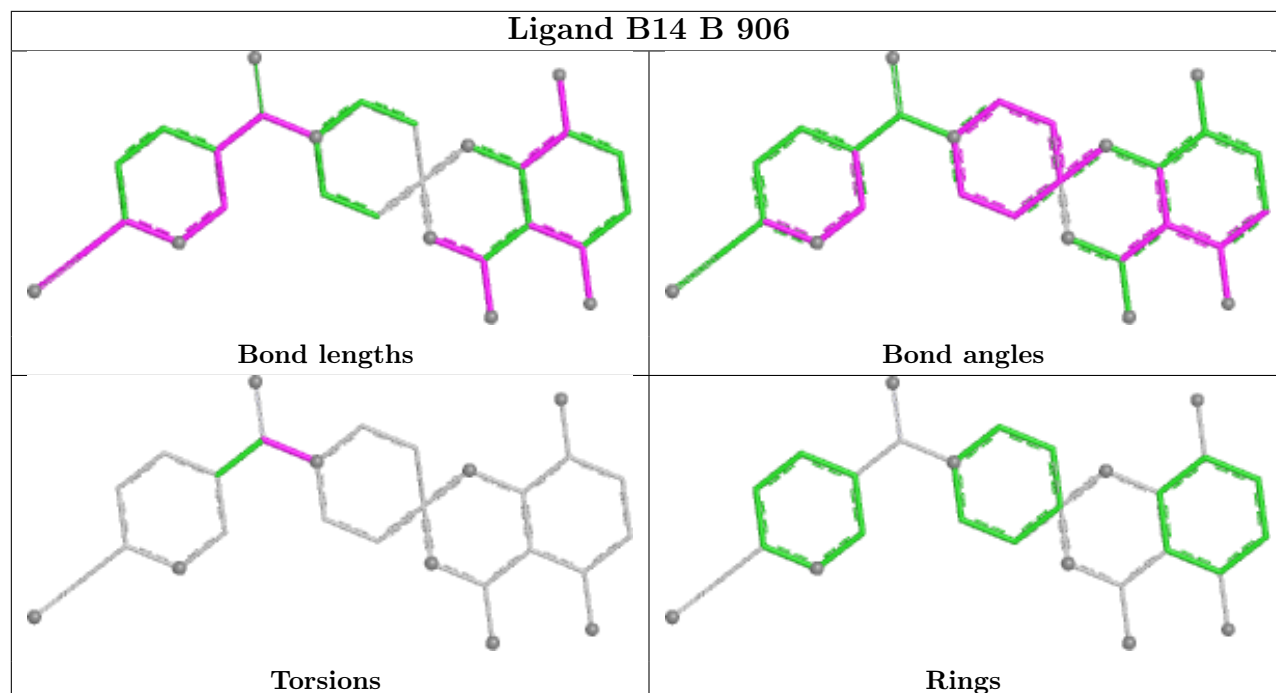
3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	902	HBI	1	0
4	A	902	HBI	2	0
3	A	901	HEM	1	0

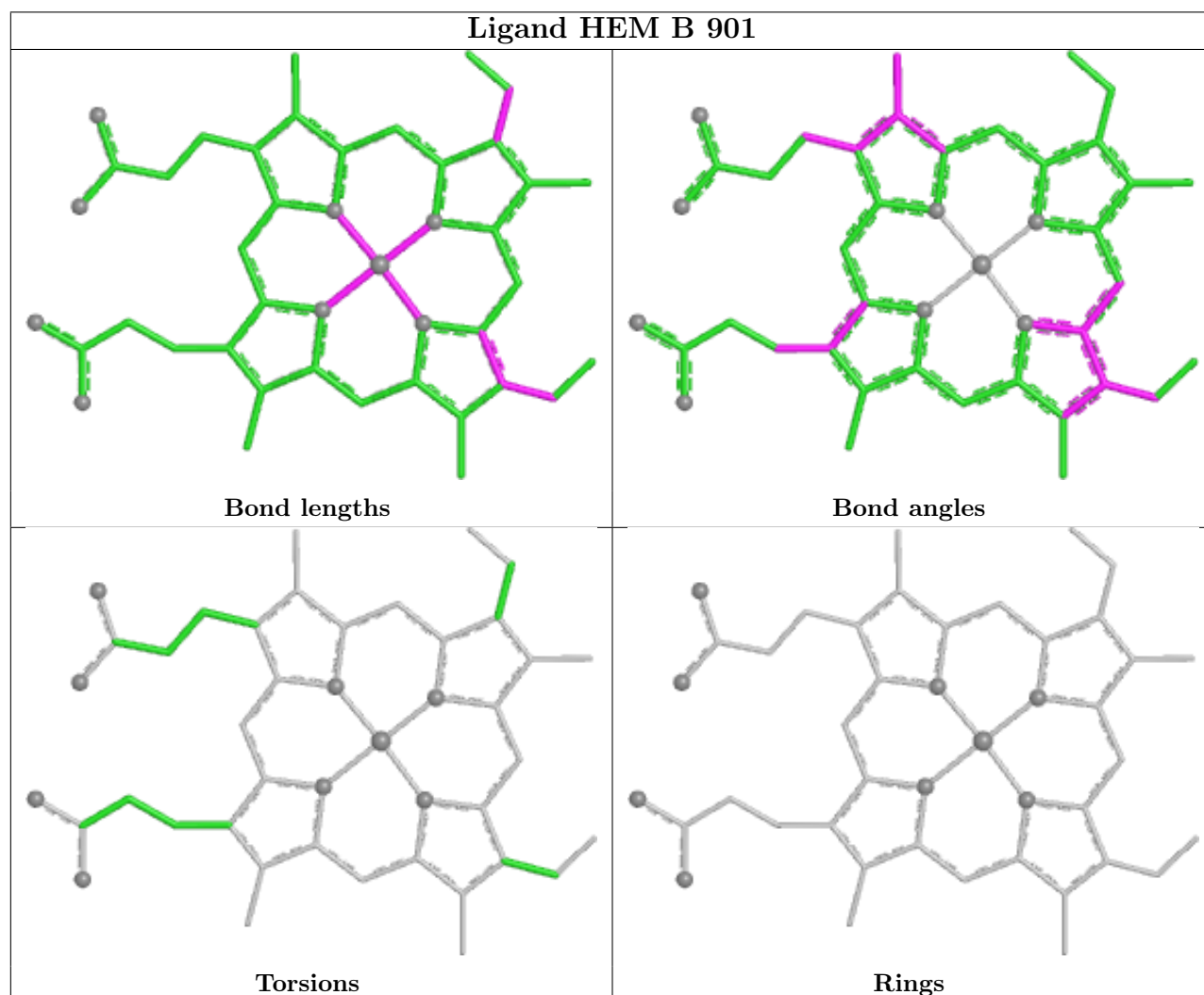
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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 B14 B 906



Ligand HEM B 901



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	413/433 (95%)	0.20	16 (3%) 43 38	25, 48, 89, 111	0
1	B	413/433 (95%)	0.13	21 (5%) 33 28	27, 49, 83, 108	0
All	All	826/866 (95%)	0.16	37 (4%) 38 32	25, 49, 86, 111	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	110	LEU	4.3
1	B	108	SER	4.2
1	B	110	LEU	4.0
1	B	114	MET	4.0
1	A	109	CYS	3.9
1	A	114	MET	3.8
1	A	171	GLY	3.7
1	B	148	PHE	3.4
1	A	494	ILE	3.3
1	B	109	CYS	3.2
1	B	494	ILE	3.2
1	A	108	SER	3.2
1	A	148	PHE	3.2
1	B	496	GLN	3.1
1	A	495	TRP	3.1
1	A	111	GLY	2.8
1	B	111	GLY	2.7
1	A	146	GLY	2.6
1	B	113	ILE	2.6
1	B	130	LEU	2.5
1	B	247	GLY	2.5
1	B	125	ASP	2.4
1	A	147	SER	2.4
1	A	152	LYS	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	450	GLY	2.4
1	B	78	TYR	2.3
1	B	100	SER	2.3
1	A	112	SER	2.3
1	B	152	LYS	2.2
1	A	307	GLY	2.2
1	B	131	GLU	2.1
1	B	272	ARG	2.1
1	A	308	GLN	2.1
1	B	171	GLY	2.1
1	B	290	LYS	2.0
1	B	99	THR	2.0
1	A	91	HIS	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

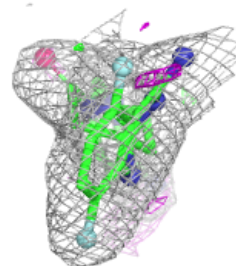
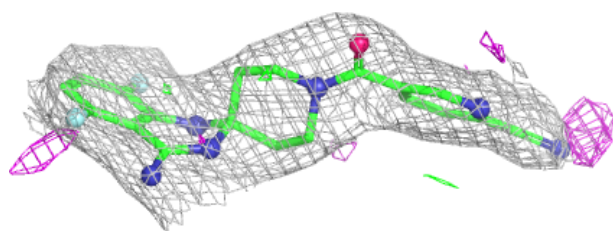
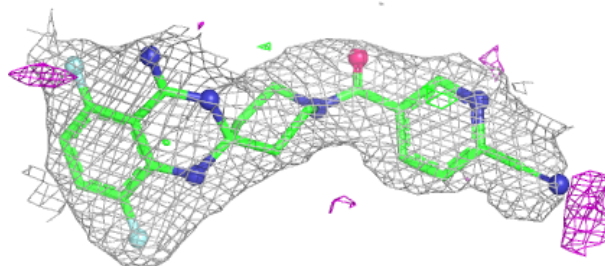
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	HBI	B	902	17/17	0.83	0.20	78,80,86,87	0
4	HBI	A	902	17/17	0.86	0.17	78,81,84,84	0
2	SO4	A	903	5/5	0.92	0.13	93,93,93,94	0
5	B14	A	906	28/28	0.95	0.08	25,29,34,38	0
5	B14	B	906	28/28	0.95	0.08	30,32,37,43	0
3	HEM	A	901	43/43	0.98	0.06	24,27,34,39	0
3	HEM	B	901	43/43	0.98	0.06	25,29,34,37	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.

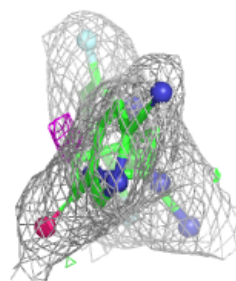
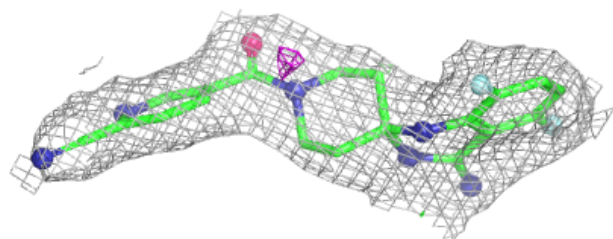
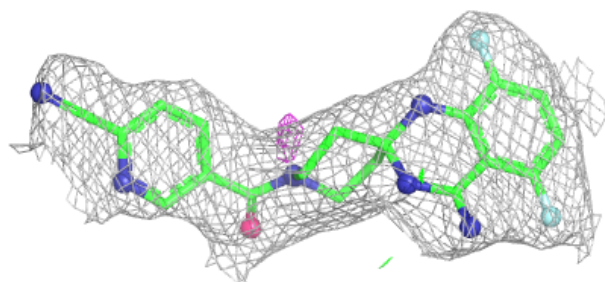
Electron density around B14 A 906:

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



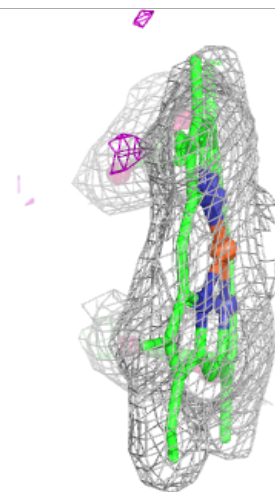
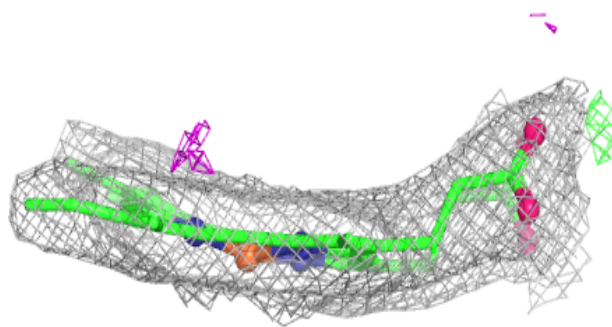
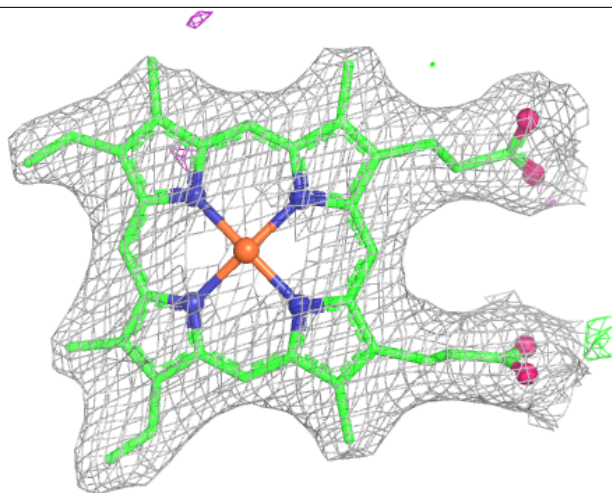
Electron density around B14 B 906:

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



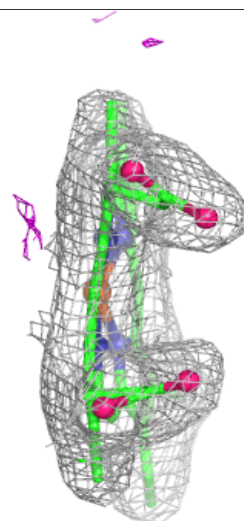
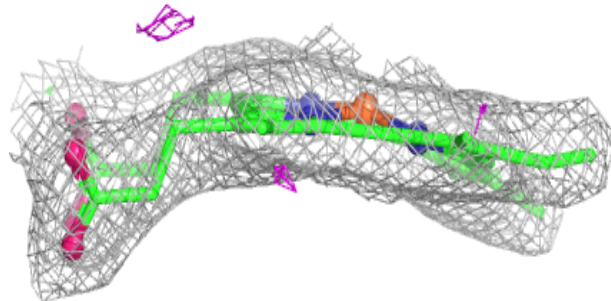
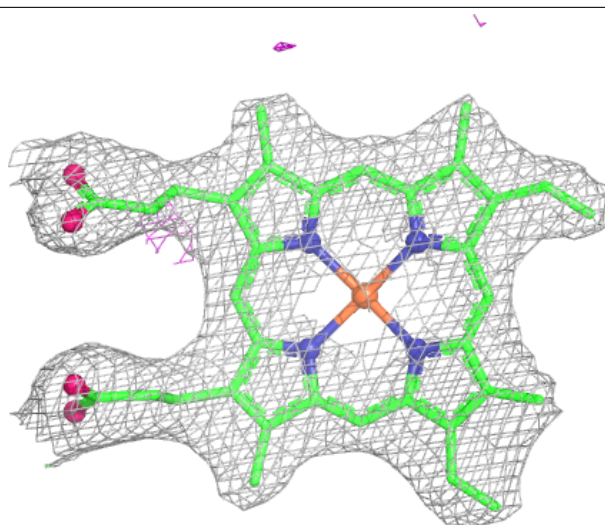
Electron density around HEM A 901:

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



Electron density around HEM B 901:

$2mF_o-DF_c$ (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.