Full wwPDB X-ray Structure Validation Report (i)

Jun 23, 2024 – 08:01 AM EDT

PDB ID	:	4UX6
Title	:	The discovery of novel, potent and highly selective inhibitors of inducible nitric oxide synthase (iNOS)
Authors	:	Cheshire, D.R.; Andrews, G.; Beaton, H.G.; Birkinshaw, T.; Boughton-Smith, N.; Connolly, S.; Cook, T.R.; Cooper, A.; Cooper, S.L.; Cox, D.; Dixon, J.; Gensmantel, N.; Hamley, P.J.; Harrison, R.; Hartopp, P.; Kack, H.; Luker, T.; Mete, A.; Millichip, I.; Nicholls, D.J.; Pimm, A.D.; St-Gallay, S.A.; Wallace, A.V.
Deposited on	:	2014-08-19
Resolution	:	3.00 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 (i) 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 (i)) were used in the production of this report:

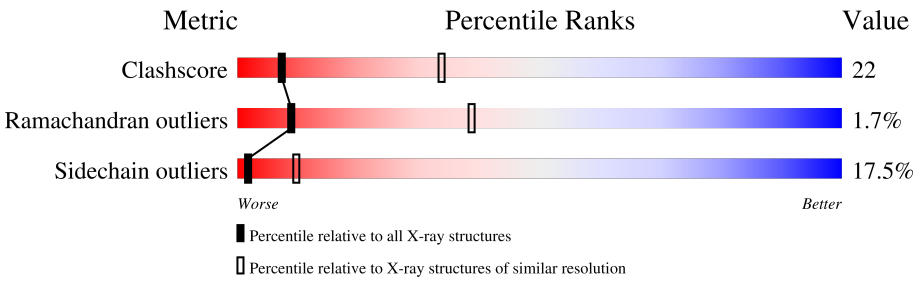
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	141614	2416 (3.00-3.00)
Ramachandran outliers	138981	2333 (3.00-3.00)
Sidechain outliers	138945	2336 (3.00-3.00)

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 <=5%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	24	17% 71% 12%
2	B	389	54% 38% 8%

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
4	H4B	B	1498	X	-	-	-

Validation Pipeline (wwPDB-VP) : 2.37.1

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 3451 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.

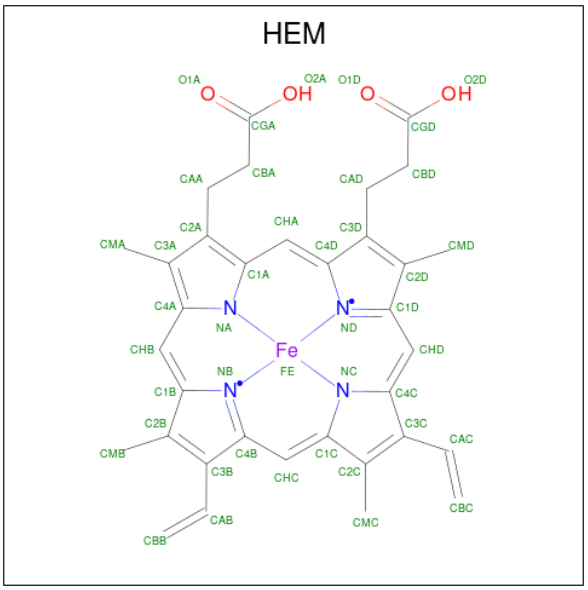
- Molecule 1 is a protein called NITRIC OXIDE SYNTHASE, INDUCIBLE.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
1	A	24	197	124	38	35	0	0	0

- Molecule 2 is a protein called NITRIC OXIDE SYNTHASE, INDUCIBLE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	389	3171	2035	543	573	20	0	0	0

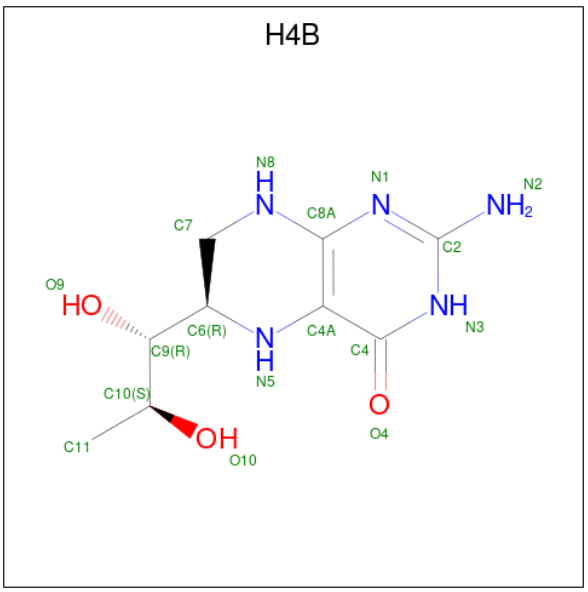
- Molecule 3 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



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

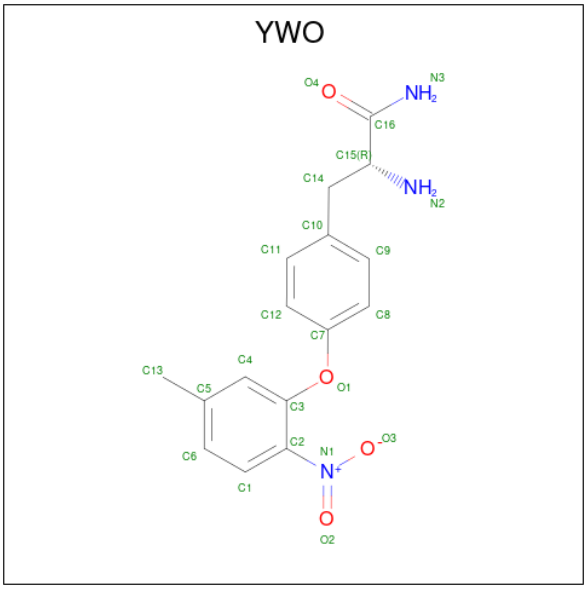
- Molecule 4 is 5,6,7,8-TETRAHYDROBIOPTERIN (three-letter code: H4B) (formula:

$C_9H_{15}N_5O_3$).



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

- Molecule 5 is O-(5-methyl-2-nitrophenyl)-D-tyrosinamide (three-letter code: YWO) (formula: $C_{16}H_{17}N_3O_4$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
5	B	1	23	16	3	4	0	0

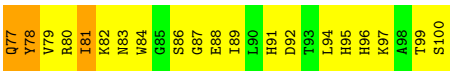
3 Residue-property plots (i)

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. 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. 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.

Note EDS was not executed.

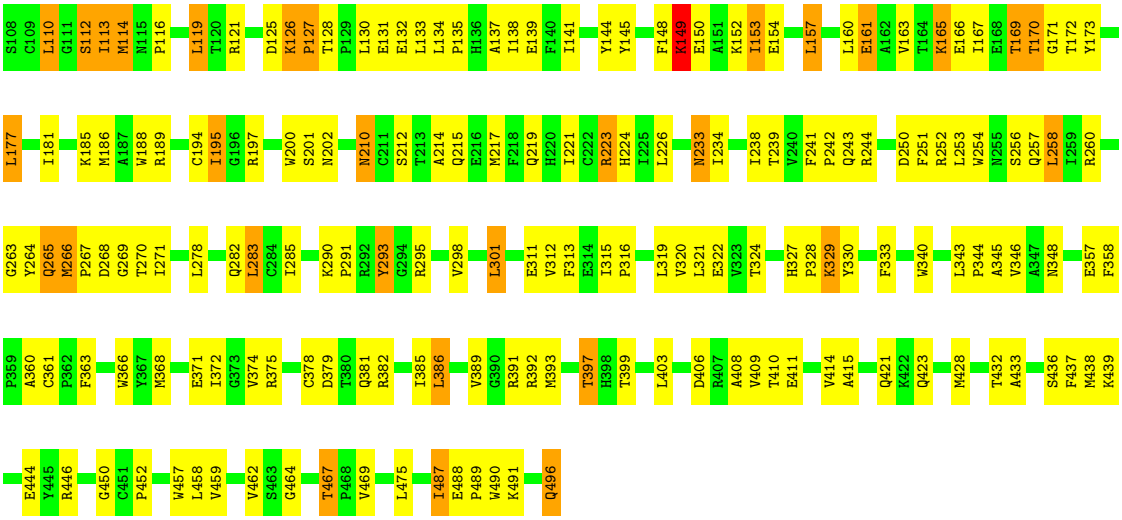
• Molecule 1: NITRIC OXIDE SYNTHASE, INDUCIBLE

Chain A: 17% 71% 12%



• Molecule 2: NITRIC OXIDE SYNTHASE, INDUCIBLE

Chain B: 54% 38% 8%



4 Data and refinement statistics (i)

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α, β, γ	213.55Å 213.55Å 115.99Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 3.00	Depositor
% Data completeness (in resolution range)	95.0 (20.00-3.00)	Depositor
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.230 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3451	wwPDB-VP
Average B, all atoms (Å^2)	32.0	wwPDB-VP

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, YWO, H4B

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	0.53	0/202	0.69	0/272
2	B	0.57	0/3264	0.82	5/4440 (0.1%)
All	All	0.57	0/3466	0.81	5/4712 (0.1%)

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
2	B	0	1

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	223	ARG	NE-CZ-NH1	14.73	127.67	120.30
2	B	223	ARG	NE-CZ-NH2	-12.10	114.25	120.30
2	B	223	ARG	CD-NE-CZ	6.72	133.01	123.60
2	B	149	LYS	N-CA-C	-6.08	94.59	111.00
2	B	368	MET	N-CA-C	-5.21	96.94	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	145	TYR	Sidechain

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	197	0	190	19	0
2	B	3171	0	3075	129	0
3	B	43	0	30	4	0
4	B	17	0	13	4	0
5	B	23	0	17	8	0
All	All	3451	0	3325	152	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 22.

All (152) 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)
2:B:114:MET:HG2	4:B:1498:H4B:H112	1.22	1.15
2:B:438:MET:HE3	2:B:469:VAL:HG12	1.17	1.13
2:B:271:ILE:HD13	2:B:278:LEU:HD11	1.44	0.99
2:B:210:ASN:H	2:B:210:ASN:HD22	1.18	0.88
2:B:327:HIS:ND1	2:B:328:PRO:HD2	1.89	0.88
2:B:265:GLN:O	2:B:265:GLN:HG3	1.76	0.85
1:A:77:GLN:HE21	1:A:77:GLN:HA	1.45	0.82
2:B:188:TRP:CE3	2:B:200:TRP:HA	2.15	0.81
2:B:411:GLU:O	2:B:414:VAL:HG22	1.81	0.79
1:A:77:GLN:HA	1:A:77:GLN:NE2	1.98	0.79
2:B:195:ILE:N	2:B:195:ILE:HD13	1.99	0.76
2:B:328:PRO:O	2:B:329:LYS:HD3	1.85	0.75
2:B:210:ASN:H	2:B:210:ASN:ND2	1.83	0.75
2:B:210:ASN:HD22	2:B:210:ASN:N	1.86	0.71
1:A:77:GLN:O	1:A:96:HIS:HE1	1.73	0.71
5:B:1499:YWO:HN2	5:B:1499:YWO:H9	1.55	0.71
2:B:266:MET:HB2	2:B:270:THR:O	1.92	0.70
2:B:110:LEU:HB3	2:B:113:ILE:HD13	1.75	0.69
2:B:438:MET:CE	2:B:469:VAL:HG12	2.10	0.69
5:B:1499:YWO:H9	5:B:1499:YWO:N2	2.08	0.68
2:B:410:THR:O	2:B:414:VAL:HG13	1.94	0.67
2:B:165:LYS:O	2:B:169:THR:HG23	1.95	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:114:MET:HG2	4:B:1498:H4B:C11	2.14	0.66
2:B:285:ILE:HD11	2:B:291:PRO:HB3	1.77	0.66
2:B:437:PHE:CE2	2:B:458:LEU:HD13	2.31	0.66
2:B:257:GLN:HB2	2:B:345:ALA:O	1.96	0.65
2:B:137:ALA:O	2:B:141:ILE:HD12	1.97	0.64
2:B:233:ASN:C	2:B:233:ASN:HD22	2.00	0.64
2:B:135:PRO:HA	2:B:138:ILE:HD12	1.79	0.63
2:B:141:ILE:HD11	2:B:163:VAL:HG21	1.79	0.63
2:B:238:ILE:HD12	2:B:363:PHE:HB3	1.81	0.62
2:B:241:PHE:HB3	2:B:242:PRO:CD	2.29	0.61
2:B:327:HIS:ND1	2:B:328:PRO:CD	2.64	0.61
2:B:149:LYS:HD3	2:B:149:LYS:H	1.66	0.60
1:A:86:SER:OG	1:A:88:GLU:HB2	2.02	0.59
2:B:215:GLN:HG3	2:B:219:GLN:NE2	2.17	0.59
2:B:285:ILE:CD1	2:B:291:PRO:HB3	2.32	0.59
2:B:149:LYS:H	2:B:149:LYS:CD	2.16	0.59
2:B:188:TRP:CD2	2:B:200:TRP:HA	2.38	0.59
2:B:301:LEU:HD22	2:B:315:ILE:HG12	1.84	0.59
2:B:371:GLU:HG3	5:B:1499:YWO:O3	2.01	0.58
1:A:78:TYR:CD1	1:A:78:TYR:C	2.77	0.58
2:B:385:ILE:O	2:B:389:VAL:HG23	2.03	0.58
2:B:375:ARG:NH1	2:B:379:ASP:OD2	2.37	0.57
2:B:197:ARG:NH2	2:B:452:PRO:O	2.32	0.57
2:B:221:ILE:HG21	2:B:301:LEU:HD21	1.87	0.56
2:B:257:GLN:OE1	5:B:1499:YWO:N2	2.38	0.56
2:B:186:MET:HE1	2:B:189:ARG:HH11	1.71	0.56
2:B:438:MET:HE3	2:B:469:VAL:CG1	2.12	0.56
2:B:327:HIS:CE1	2:B:328:PRO:HD2	2.40	0.56
2:B:243:GLN:HB3	2:B:358:PHE:CE1	2.41	0.55
2:B:194:CYS:HB3	2:B:197:ARG:HD2	1.90	0.54
2:B:374:VAL:O	2:B:378:CYS:HB2	2.07	0.54
2:B:252:ARG:NH2	2:B:489:PRO:HD3	2.22	0.54
1:A:82:LYS:HG2	1:A:83:ASN:N	2.23	0.53
2:B:264:TYR:CE2	2:B:293:TYR:HA	2.43	0.53
2:B:112:SER:O	4:B:1498:H4B:H111	2.06	0.53
2:B:433:ALA:O	2:B:436:SER:HB3	2.09	0.53
2:B:340:TRP:CD1	2:B:372:ILE:HG12	2.45	0.52
2:B:496:GLN:CD	2:B:496:GLN:C	2.67	0.52
2:B:298:VAL:HG21	2:B:320:VAL:HG11	1.91	0.52
2:B:133:LEU:HD13	2:B:133:LEU:O	2.10	0.52
2:B:167:ILE:O	2:B:171:GLY:N	2.42	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:188:TRP:CZ3	2:B:200:TRP:HA	2.46	0.51
1:A:87:GLY:O	1:A:89:ILE:HD12	2.11	0.50
1:A:78:TYR:CD1	1:A:78:TYR:O	2.64	0.50
2:B:311:GLU:HB3	2:B:313:PHE:CE1	2.45	0.50
2:B:321:LEU:HD23	2:B:322:GLU:N	2.27	0.50
2:B:126:LYS:O	2:B:127:PRO:O	2.30	0.50
2:B:366:TRP:H	3:B:1497:HEM:HAB	1.77	0.50
2:B:130:LEU:N	2:B:130:LEU:HD23	2.27	0.49
2:B:233:ASN:C	2:B:233:ASN:ND2	2.65	0.49
2:B:241:PHE:HB3	2:B:242:PRO:HD2	1.95	0.49
2:B:254:TRP:CE2	2:B:489:PRO:HB2	2.48	0.49
2:B:258:LEU:HD22	2:B:345:ALA:HB1	1.95	0.49
2:B:239:THR:O	2:B:361:CYS:HA	2.14	0.48
2:B:185:LYS:O	2:B:188:TRP:HB3	2.14	0.48
2:B:321:LEU:HD23	2:B:321:LEU:C	2.34	0.48
2:B:467:THR:CG2	2:B:469:VAL:HG22	2.44	0.48
2:B:134:LEU:CD1	2:B:138:ILE:HD11	2.44	0.48
2:B:316:PRO:HD2	2:B:319:LEU:HD12	1.96	0.48
2:B:141:ILE:CD1	2:B:163:VAL:HG21	2.44	0.48
2:B:327:HIS:CD2	2:B:330:TYR:HD2	2.32	0.47
2:B:242:PRO:HG2	2:B:251:PHE:CZ	2.48	0.47
2:B:163:VAL:HG12	2:B:167:ILE:HD12	1.95	0.47
2:B:116:PRO:HG2	2:B:119:LEU:HB2	1.97	0.47
2:B:157:LEU:N	2:B:157:LEU:HD23	2.29	0.47
2:B:161:GLU:OE2	2:B:165:LYS:NZ	2.48	0.47
2:B:263:GLY:O	2:B:278:LEU:HD23	2.14	0.47
1:A:78:TYR:CE1	1:A:91:HIS:CD2	3.03	0.47
2:B:195:ILE:HD13	2:B:195:ILE:H	1.75	0.47
2:B:177:LEU:HD13	2:B:181:ILE:HD12	1.96	0.46
1:A:79:VAL:HG23	1:A:95:HIS:CE1	2.51	0.46
2:B:133:LEU:HD13	2:B:133:LEU:C	2.36	0.46
2:B:224:HIS:HD2	2:B:239:THR:OG1	1.98	0.46
1:A:78:TYR:C	1:A:78:TYR:HD1	2.18	0.46
1:A:94:LEU:HB3	2:B:450:GLY:HA3	1.98	0.46
1:A:92:ASP:OD1	1:A:95:HIS:CD2	2.68	0.46
3:B:1497:HEM:O1A	4:B:1498:H4B:N3	2.45	0.46
1:A:92:ASP:OD1	1:A:95:HIS:HD2	1.99	0.46
2:B:153:ILE:H	2:B:153:ILE:HD13	1.80	0.46
2:B:149:LYS:O	2:B:150:GLU:HB2	2.16	0.45
1:A:82:LYS:NZ	1:A:84:TRP:CZ3	2.82	0.45
2:B:126:LYS:O	2:B:127:PRO:C	2.55	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:381:GLN:O	2:B:382:ARG:HD2	2.17	0.45
2:B:125:ASP:C	2:B:126:LYS:HG2	2.37	0.45
2:B:163:VAL:HG13	2:B:173:TYR:CD2	2.51	0.45
2:B:489:PRO:O	2:B:491:LYS:N	2.50	0.45
2:B:114:MET:CE	2:B:457:TRP:HZ2	2.29	0.44
2:B:244:ARG:HD2	2:B:357:GLU:OE1	2.18	0.44
2:B:346:VAL:HG23	5:B:1499:YWO:H4	1.99	0.44
2:B:214:ALA:O	2:B:217:MET:HB2	2.17	0.44
2:B:487:ILE:CD1	2:B:487:ILE:N	2.80	0.44
2:B:153:ILE:H	2:B:153:ILE:CD1	2.31	0.44
2:B:408:ALA:O	2:B:411:GLU:N	2.47	0.43
2:B:488:GLU:OE2	2:B:491:LYS:HE3	2.18	0.43
2:B:234:ILE:H	2:B:234:ILE:HG12	1.61	0.43
2:B:258:LEU:HB2	2:B:345:ALA:HB3	1.99	0.43
2:B:265:GLN:NE2	2:B:266:MET:O	2.46	0.43
2:B:397:THR:O	2:B:397:THR:CG2	2.67	0.43
1:A:95:HIS:CD2	1:A:95:HIS:H	2.36	0.43
2:B:386:LEU:HD23	2:B:386:LEU:HA	1.89	0.43
2:B:144:TYR:CE2	2:B:148:PHE:HZ	2.36	0.43
2:B:167:ILE:HA	2:B:172:THR:O	2.18	0.42
2:B:282:GLN:O	2:B:283:LEU:C	2.57	0.42
2:B:343:LEU:HA	2:B:344:PRO:HD3	1.89	0.42
2:B:188:TRP:CZ2	3:B:1497:HEM:HMC1	2.54	0.42
2:B:195:ILE:HG22	2:B:437:PHE:HB2	2.00	0.42
2:B:459:VAL:HG22	2:B:469:VAL:HG23	2.01	0.42
2:B:126:LYS:C	2:B:127:PRO:O	2.58	0.42
2:B:165:LYS:HD2	2:B:165:LYS:HA	1.82	0.42
5:B:1499:YWO:N2	5:B:1499:YWO:C9	2.82	0.42
2:B:251:PHE:O	2:B:360:ALA:HB2	2.19	0.42
2:B:195:ILE:N	2:B:195:ILE:CD1	2.69	0.41
2:B:393:MET:HE1	2:B:411:GLU:HG2	2.02	0.41
2:B:489:PRO:C	2:B:491:LYS:H	2.24	0.41
3:B:1497:HEM:O2D	5:B:1499:YWO:H11	2.20	0.41
1:A:81:ILE:HD13	1:A:81:ILE:HA	1.74	0.41
2:B:312:VAL:C	2:B:313:PHE:CD1	2.93	0.41
2:B:333:PHE:CE2	2:B:415:ALA:HB1	2.55	0.41
2:B:391:ARG:C	2:B:393:MET:H	2.24	0.41
2:B:244:ARG:NH2	2:B:250:ASP:OD2	2.43	0.41
2:B:166:GLU:O	2:B:170:THR:N	2.51	0.41
2:B:363:PHE:HE1	5:B:1499:YWO:H13A	1.86	0.41
2:B:138:ILE:HG23	2:B:160:LEU:CD2	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:167:ILE:O	2:B:171:GLY:CA	2.69	0.41
2:B:428:MET:CE	2:B:432:THR:HG22	2.51	0.41
2:B:121:ARG:HA	2:B:121:ARG:HD3	1.91	0.40
1:A:77:GLN:NE2	1:A:77:GLN:CA	2.78	0.40
1:A:86:SER:C	1:A:88:GLU:H	2.25	0.40
2:B:267:PRO:C	2:B:269:GLY:H	2.25	0.40
2:B:327:HIS:HB2	2:B:333:PHE:CG	2.57	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	22/24 (92%)	21 (96%)	1 (4%)	0	100	100
2	B	387/389 (100%)	339 (88%)	41 (11%)	7 (2%)	8	37
All	All	409/413 (99%)	360 (88%)	42 (10%)	7 (2%)	9	39

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	293	TYR
2	B	464	GLY
2	B	268	ASP
2	B	392	ARG
2	B	397	THR
2	B	127	PRO
2	B	490	TRP

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	21/21 (100%)	14 (67%)	7 (33%)	0 1
2	B	340/340 (100%)	284 (84%)	56 (16%)	2 11
All	All	361/361 (100%)	298 (82%)	63 (18%)	2 10

All (63) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	77	GLN
1	A	78	TYR
1	A	80	ARG
1	A	81	ILE
1	A	97	LYS
1	A	99	THR
1	A	100	SER
2	B	110	LEU
2	B	112	SER
2	B	113	ILE
2	B	114	MET
2	B	119	LEU
2	B	126	LYS
2	B	128	THR
2	B	131	GLU
2	B	132	GLU
2	B	139	GLU
2	B	149	LYS
2	B	152	LYS
2	B	153	ILE
2	B	154	GLU
2	B	157	LEU
2	B	161	GLU
2	B	165	LYS
2	B	169	THR
2	B	170	THR
2	B	177	LEU

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Mol	Chain	Res	Type
2	B	195	ILE
2	B	201	SER
2	B	202	ASN
2	B	210	ASN
2	B	212	SER
2	B	223	ARG
2	B	226	LEU
2	B	233	ASN
2	B	253	LEU
2	B	256	SER
2	B	258	LEU
2	B	260	ARG
2	B	265	GLN
2	B	266	MET
2	B	283	LEU
2	B	290	LYS
2	B	295	ARG
2	B	301	LEU
2	B	324	THR
2	B	329	LYS
2	B	348	ASN
2	B	386	LEU
2	B	399	THR
2	B	403	LEU
2	B	406	ASP
2	B	409	VAL
2	B	421	GLN
2	B	423	GLN
2	B	439	LYS
2	B	444	GLU
2	B	446	ARG
2	B	462	VAL
2	B	467	THR
2	B	475	LEU
2	B	487	ILE
2	B	496	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	77	GLN
1	A	95	HIS

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Mol	Chain	Res	Type
1	A	96	HIS
2	B	202	ASN
2	B	210	ASN
2	B	215	GLN
2	B	219	GLN
2	B	224	HIS
2	B	231	ASN
2	B	233	ASN
2	B	348	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

3 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	H4B	B	1498	-	16,18,18	4.11	7 (43%)	11,26,26	4.50	6 (54%)
5	YWO	B	1499	-	23,24,24	2.21	10 (43%)	30,33,33	2.24	6 (20%)
3	HEM	B	1497	2	41,50,50	1.71	6 (14%)	45,82,82	2.23	9 (20%)

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	H4B	B	1498	-	1/1/3/5	3/8/17/17	0/2/2/2
5	YWO	B	1499	-	-	2/14/16/16	0/2/2/2
3	HEM	B	1497	2	-	2/12/54/54	-

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	1498	H4B	C7-C6	-11.07	1.41	1.52
4	B	1498	H4B	C2-N1	7.05	1.47	1.35
4	B	1498	H4B	C4-N3	5.46	1.42	1.33
3	B	1497	HEM	CBB-CAB	4.78	1.54	1.30
4	B	1498	H4B	C7-N8	-4.39	1.37	1.44
5	B	1499	YWO	C2-N1	4.11	1.53	1.45
3	B	1497	HEM	CBC-CAC	4.00	1.55	1.29
5	B	1499	YWO	C12-C7	3.78	1.46	1.38
5	B	1499	YWO	C4-C3	3.65	1.45	1.38
4	B	1498	H4B	C6-N5	-3.60	1.38	1.45
5	B	1499	YWO	C9-C8	3.51	1.45	1.38
5	B	1499	YWO	C6-C1	3.43	1.45	1.38
3	B	1497	HEM	CAB-C3B	3.38	1.56	1.47
4	B	1498	H4B	O9-C9	-3.24	1.35	1.43
4	B	1498	H4B	C8A-N1	3.16	1.40	1.34
5	B	1499	YWO	C9-C10	2.68	1.44	1.38
5	B	1499	YWO	C11-C10	2.58	1.44	1.38
3	B	1497	HEM	CMC-C2C	2.51	1.57	1.51
5	B	1499	YWO	C16-N3	2.46	1.39	1.32
3	B	1497	HEM	CBA-CGA	2.38	1.56	1.50
5	B	1499	YWO	O2-N1	2.32	1.26	1.22
3	B	1497	HEM	C3C-C2C	-2.30	1.37	1.40
5	B	1499	YWO	C12-C11	2.20	1.42	1.38

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	B	1498	H4B	C8A-C4A-C4	9.59	123.08	114.57
3	B	1497	HEM	CBA-CAA-C2A	-8.66	97.84	112.62
5	B	1499	YWO	O4-C16-C15	-8.33	108.08	120.30
4	B	1498	H4B	C2-N3-C4	8.33	129.17	115.93

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	B	1497	HEM	C4A-C3A-C2A	7.01	111.88	107.00
4	B	1498	H4B	N1-C2-N3	-5.17	117.31	125.42
3	B	1497	HEM	CAA-CBA-CGA	4.63	126.72	113.76
5	B	1499	YWO	C15-C16-N3	-4.35	109.22	116.68
4	B	1498	H4B	C4A-C4-N3	-4.26	111.90	124.01
5	B	1499	YWO	C1-C2-N1	3.89	120.63	116.47
3	B	1497	HEM	CMA-C3A-C4A	-3.87	122.51	128.46
5	B	1499	YWO	C3-O1-C7	3.65	127.10	118.00
3	B	1497	HEM	CMC-C2C-C3C	3.36	130.96	124.68
5	B	1499	YWO	O4-C16-N3	-2.92	117.92	123.00
4	B	1498	H4B	N2-C2-N3	2.86	121.71	117.25
3	B	1497	HEM	C4B-CHC-C1C	2.84	126.30	122.56
3	B	1497	HEM	C4C-CHD-C1D	2.84	126.30	122.56
4	B	1498	H4B	N2-C2-N1	2.40	120.98	117.25
3	B	1497	HEM	CHC-C4B-NB	-2.40	121.83	124.43
3	B	1497	HEM	CBB-CAB-C3B	-2.37	115.85	127.62
5	B	1499	YWO	O2-N1-C2	-2.09	115.45	119.03

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
4	B	1498	H4B	C6

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	1497	HEM	C2B-C3B-CAB-CBB
3	B	1497	HEM	C4B-C3B-CAB-CBB
4	B	1498	H4B	N5-C6-C9-O9
4	B	1498	H4B	N5-C6-C9-C10
5	B	1499	YWO	C11-C10-C14-C15
5	B	1499	YWO	C9-C10-C14-C15
4	B	1498	H4B	C7-C6-C9-O9

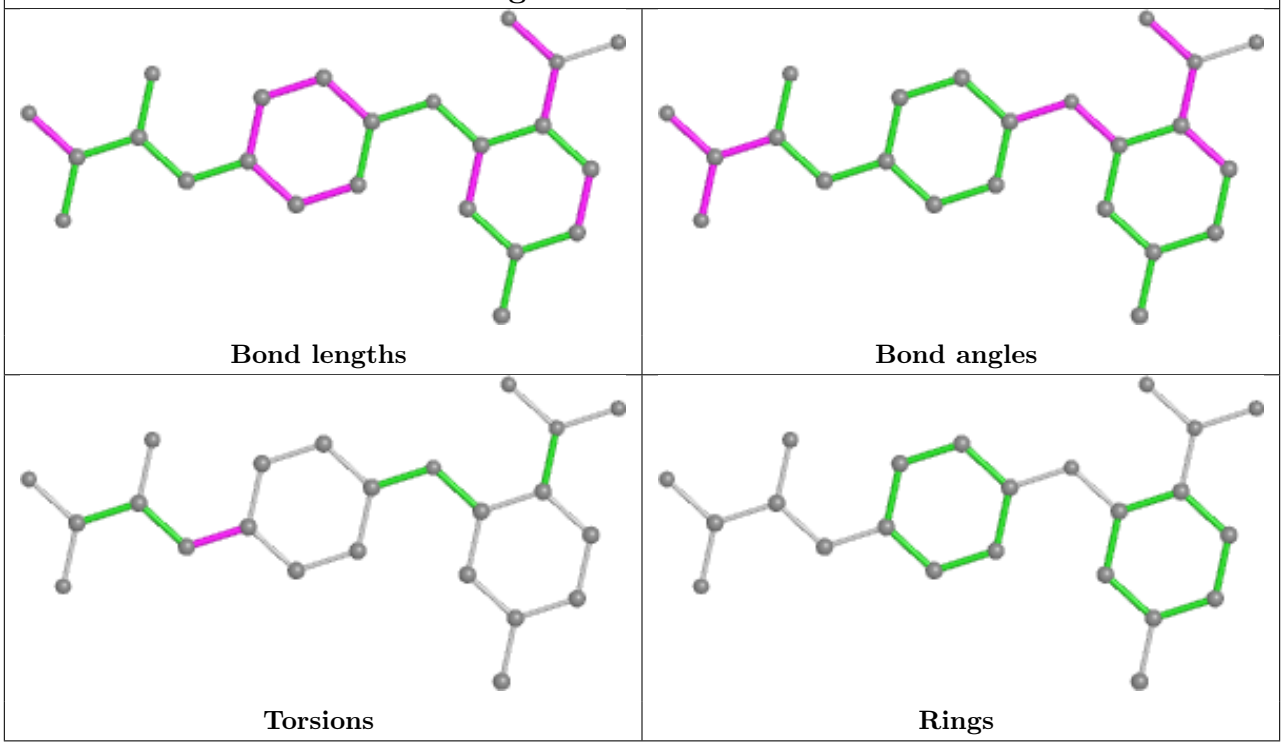
There are no ring outliers.

3 monomers are involved in 14 short contacts:

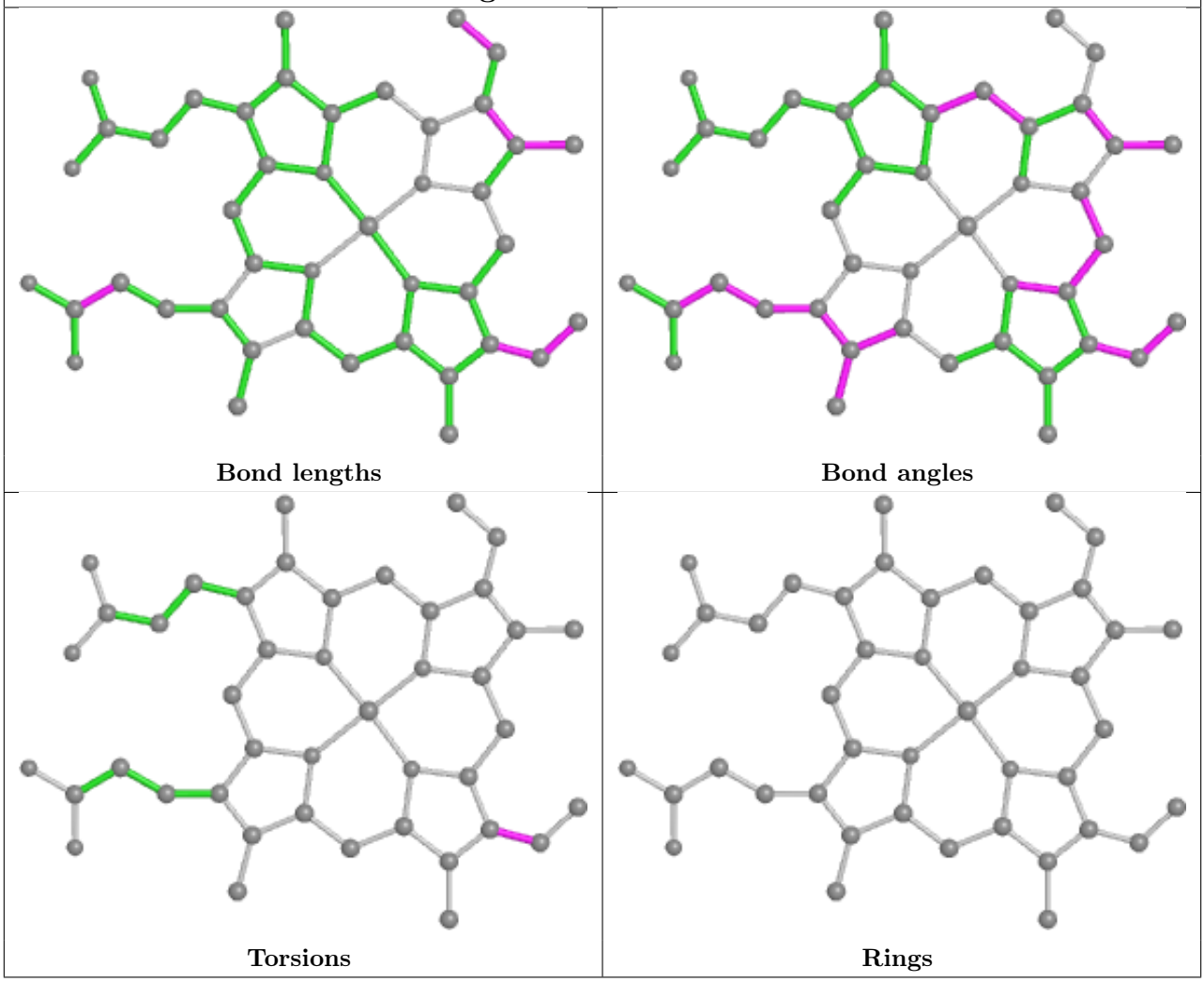
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	1498	H4B	4	0
5	B	1499	YWO	8	0
3	B	1497	HEM	4	0

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 then 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 YWO B 1499



Ligand HEM B 1497



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)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.