



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 12:19 PM UTC

PDB ID : 3E7M / pdb_00003e7m
Title : Structure of murine iNOS oxygenase domain with inhibitor AR-C95791
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 : 2.00 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 ⓘ 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	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

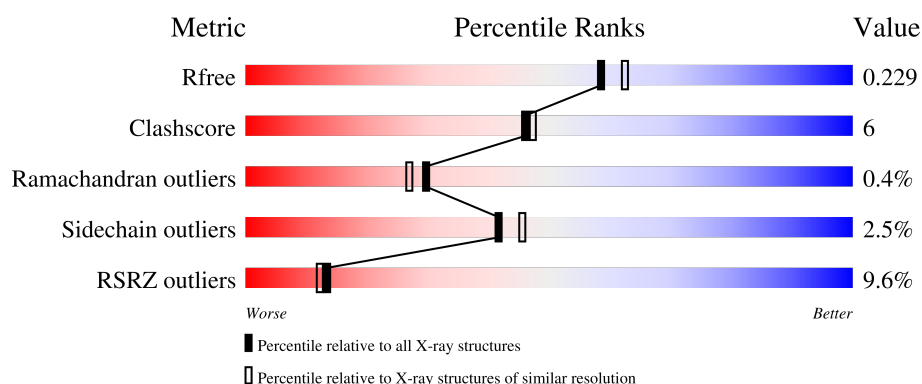
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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))
R_{free}	180053	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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 $\leq 5\%$. 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	
1	B	433	

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 7484 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.

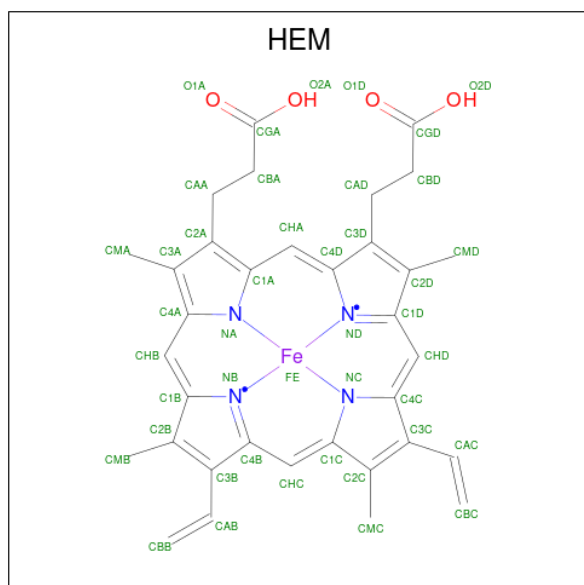
- Molecule 1 is a protein called Nitric oxide synthase, inducible.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	416	Total	C	N	O	S	0	0	0
			3389	2174	583	612	20			
1	B	421	Total	C	N	O	S	0	0	0
			3421	2193	589	618	21			

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

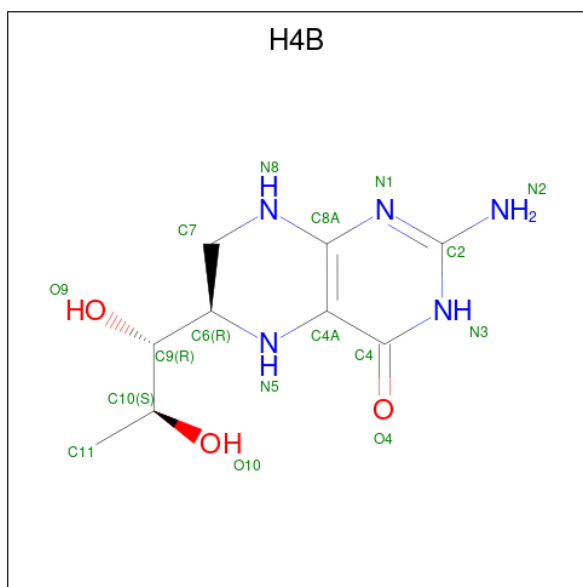
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		
2	B	1	Total	Zn	0	0
			1	1		

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



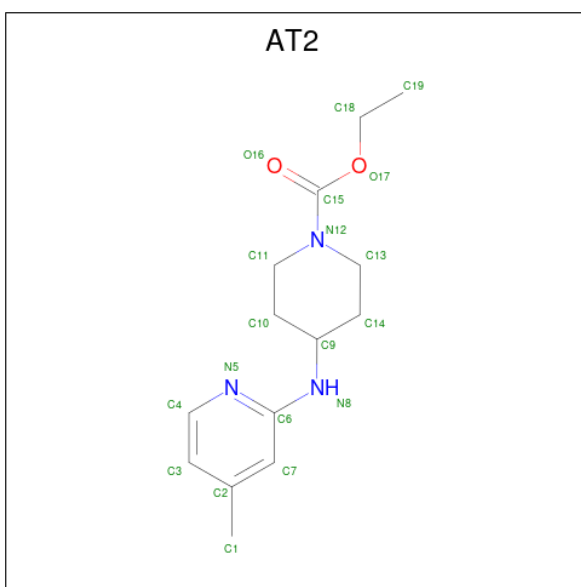
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
3	B	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 4 is 5,6,7,8-TETRAHYDROBIOPTERIN (CCD ID: H4B) (formula: $C_9H_{15}N_5O_3$).



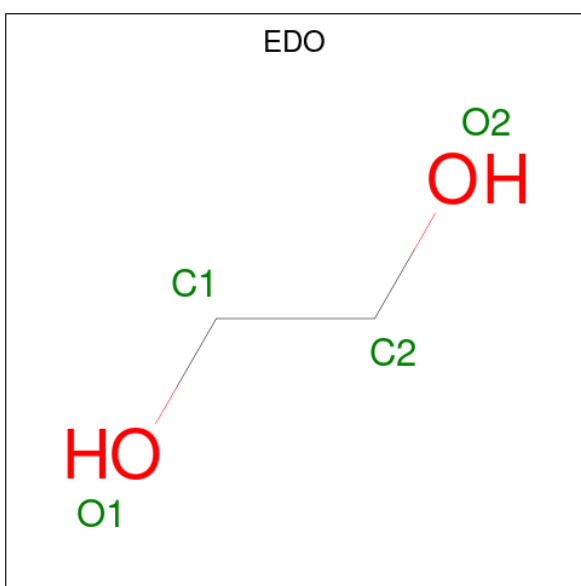
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			17	9	5	3		
4	B	1	Total	C	N	O	0	0
			17	9	5	3		

- Molecule 5 is ETHYL 4-[(4-METHYLPYRIDIN-2-YL)AMINO]PIPERIDINE-1-CARBOXYLATE (CCD ID: AT2) (formula: $C_{14}H_{21}N_3O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			19	14	3	2		
5	B	1	Total	C	N	O	0	0
			19	14	3	2		

- Molecule 6 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			4	2	2		
6	A	1	Total	C	O	0	0
			4	2	2		

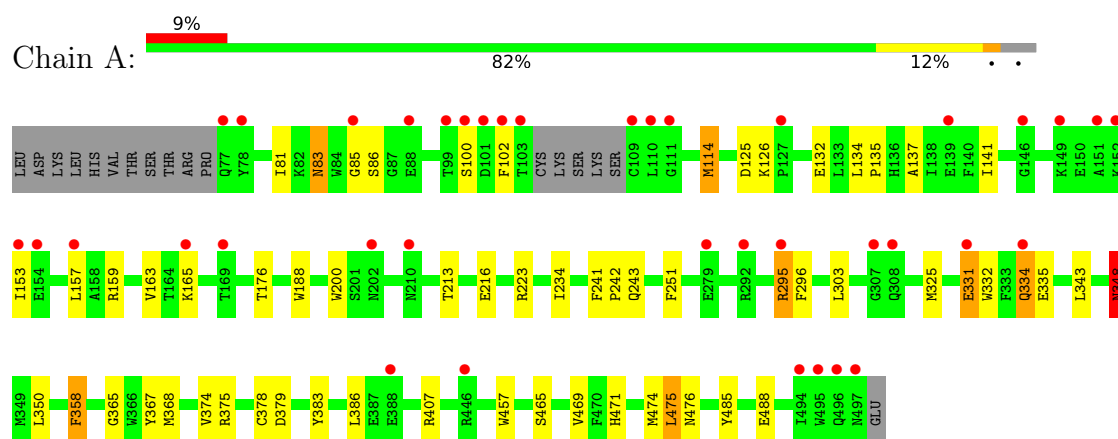
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	228	Total 228	O 228	0	0
7	B	278	Total 278	O 278	0	0

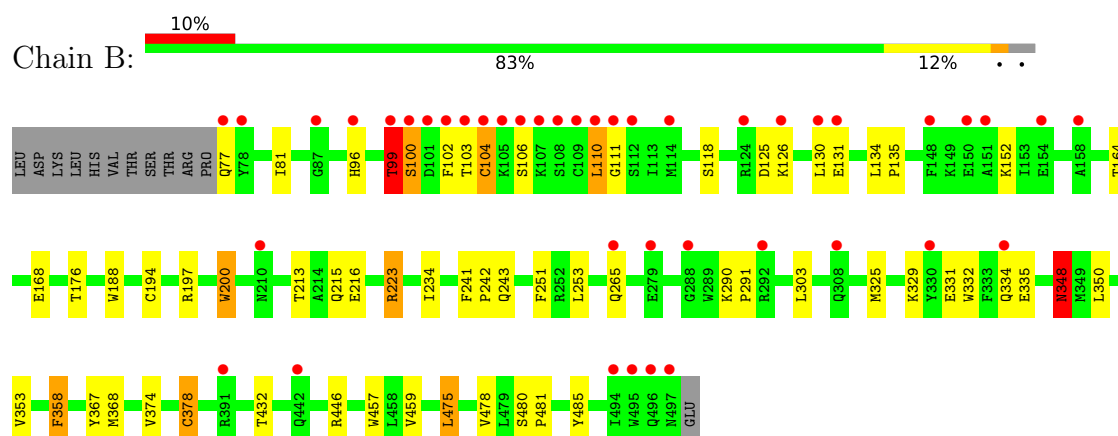
3 Residue-property plots [i](#)

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 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: Nitric oxide synthase, inducible



- Molecule 1: Nitric oxide synthase, inducible



4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	213.85Å 213.85Å 116.64Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	19.96 – 2.00 19.96 – 2.00	Depositor EDS
% Data completeness (in resolution range)	98.7 (19.96-2.00) 98.6 (19.96-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.03	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.37 (at 2.00Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.223 , 0.237 0.215 , 0.229	Depositor DCC
R_{free} test set	5249 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	30.3	Xtriage
Anisotropy	0.707	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 40.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7484	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.90% 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: ZN, AT2, H4B, HEM, EDO

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	0.36	0/3488	0.90	14/4743 (0.3%)
1	B	0.41	0/3521	0.94	20/4787 (0.4%)
All	All	0.38	0/7009	0.92	34/9530 (0.4%)

There are no bond length outliers.

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	106	SER	N-CA-C	7.63	120.67	111.82
1	B	368	MET	N-CA-C	-7.51	96.98	109.07
1	A	368	MET	N-CA-C	-7.42	97.13	109.07
1	B	111	GLY	N-CA-C	7.41	121.87	112.83
1	B	348	ASN	N-CA-C	7.29	121.48	112.59
1	A	365	GLY	N-CA-C	-7.00	99.64	110.77
1	A	348	ASN	N-CA-C	6.66	120.72	112.59
1	B	200	TRP	N-CA-C	6.61	120.21	111.75
1	A	485	TYR	N-CA-C	-6.03	102.75	110.53
1	B	303	LEU	N-CA-C	6.02	118.76	109.07
1	B	485	TYR	N-CA-C	-5.92	102.90	110.53
1	A	176	THR	N-CA-C	-5.77	103.09	110.53
1	B	104	CYS	N-CA-C	5.75	119.63	109.96
1	A	367	TYR	N-CA-C	5.70	118.54	110.50
1	B	325	MET	N-CA-C	5.56	117.76	109.25
1	B	459	VAL	CA-C-N	5.55	123.71	119.66
1	B	459	VAL	C-N-CA	5.55	123.71	119.66
1	B	358	PHE	CA-C-N	5.53	125.20	119.56
1	B	358	PHE	C-N-CA	5.53	125.20	119.56
1	B	367	TYR	N-CA-C	5.51	118.27	110.50
1	B	234	ILE	N-CA-C	5.47	116.47	109.30
1	A	358	PHE	CA-C-N	5.39	125.01	119.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	358	PHE	C-N-CA	5.39	125.01	119.56
1	A	251	PHE	N-CA-C	-5.37	100.55	109.46
1	A	234	ILE	N-CA-C	5.35	116.31	109.30
1	B	253	LEU	N-CA-C	-5.31	100.07	108.73
1	A	303	LEU	N-CA-C	5.30	117.61	109.07
1	B	99	THR	N-CA-C	5.29	122.06	110.80
1	B	176	THR	N-CA-C	-5.27	102.94	110.59
1	A	343	LEU	N-CA-C	5.27	118.17	108.69
1	B	378	CYS	N-CA-C	5.24	119.82	113.38
1	A	386	LEU	N-CA-C	5.11	117.25	111.02
1	A	325	MET	N-CA-C	5.04	116.97	109.25
1	B	251	PHE	N-CA-C	-5.04	101.03	109.24

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	3389	0	3276	39	0
1	B	3421	0	3310	41	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	43	0	30	0	0
3	B	43	0	30	0	0
4	A	17	0	15	0	0
4	B	17	0	15	1	0
5	A	19	0	21	0	0
5	B	19	0	21	0	0
6	A	8	0	12	0	0
7	A	228	0	0	4	2
7	B	278	0	0	5	1
All	All	7484	0	6730	80	3

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 6.

All (80) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:103:THR:O	1:B:110:LEU:HD23	1.62	1.00
1:B:81:ILE:HD11	1:B:475:LEU:HD13	1.44	0.97
1:A:81:ILE:HD11	1:A:475:LEU:HD13	1.62	0.79
1:A:331:GLU:H	1:A:331:GLU:CD	1.93	0.76
1:A:83:ASN:HD22	1:A:85:GLY:H	1.35	0.75
1:A:83:ASN:ND2	1:A:85:GLY:H	1.86	0.72
1:B:432:THR:CG2	7:B:1158:HOH:O	2.39	0.70
1:A:469:VAL:HG13	1:A:474:MET:HE1	1.76	0.68
1:B:290:LYS:HE2	1:B:291:PRO:HD2	1.76	0.68
1:B:223:ARG:HH21	1:B:223:ARG:CG	2.07	0.67
1:B:215:GLN:OE1	7:B:1476:HOH:O	2.12	0.66
1:B:223:ARG:HH21	1:B:223:ARG:HG3	1.61	0.66
1:B:290:LYS:HE2	1:B:291:PRO:CD	2.31	0.61
1:B:81:ILE:HD11	1:B:475:LEU:CD1	2.26	0.60
1:B:102:PHE:HE1	1:B:478:VAL:HG23	1.66	0.60
1:B:194:CYS:HB3	1:B:197:ARG:HD2	1.83	0.60
1:A:465:SER:O	1:A:471:HIS:HE1	1.85	0.59
1:B:194:CYS:O	1:B:197:ARG:HD3	2.04	0.58
1:B:432:THR:HG21	7:B:1158:HOH:O	1.99	0.57
1:A:469:VAL:HG13	1:A:474:MET:CE	2.35	0.56
1:A:295:ARG:HD2	1:A:296:PHE:CE2	2.42	0.55
1:A:114:MET:HE3	7:A:1011:HOH:O	2.07	0.54
1:A:488:GLU:HG2	7:A:1379:HOH:O	2.08	0.54
1:B:102:PHE:CE1	1:B:478:VAL:HG23	2.42	0.54
1:A:134:LEU:HB3	1:A:135:PRO:HD3	1.91	0.52
1:B:188:TRP:CE3	1:B:200:TRP:HA	2.44	0.52
1:A:223:ARG:HG3	1:A:223:ARG:HH21	1.74	0.52
1:B:103:THR:HG22	1:B:118:SER:OG	2.10	0.52
1:A:348:ASN:HB2	7:A:1221:HOH:O	2.09	0.52
1:A:83:ASN:ND2	1:A:86:SER:H	2.08	0.52
1:B:104:CYS:SG	1:B:110:LEU:N	2.81	0.51
1:A:188:TRP:CE3	1:A:200:TRP:HA	2.45	0.51
1:A:125:ASP:OD1	1:A:126:LYS:HG3	2.10	0.51
1:B:446:ARG:HG2	1:B:446:ARG:HH21	1.77	0.49
1:A:114:MET:HE2	1:A:457:TRP:CZ2	2.48	0.49
1:A:348:ASN:HD22	1:A:348:ASN:H	1.61	0.48
1:A:331:GLU:CD	1:A:331:GLU:N	2.66	0.47
1:B:348:ASN:HD22	1:B:348:ASN:H	1.63	0.47
1:A:102:PHE:HE2	1:A:476:ASN:O	1.97	0.46
1:A:153:ILE:O	1:A:157:LEU:HD23	2.14	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:407:ARG:HD2	7:A:1364:HOH:O	2.15	0.46
1:B:125:ASP:OD1	1:B:126:LYS:HG3	2.16	0.46
1:B:152:LYS:HB3	7:B:1410:HOH:O	2.15	0.46
1:B:350:LEU:C	1:B:350:LEU:HD23	2.40	0.46
1:B:213:THR:OG1	1:B:216:GLU:HG3	2.15	0.46
1:A:295:ARG:HD3	1:A:383:TYR:HE2	1.81	0.45
1:B:194:CYS:O	1:B:197:ARG:CD	2.63	0.45
1:A:114:MET:HE2	1:A:457:TRP:HZ2	1.80	0.45
1:A:332:TRP:O	1:A:335:GLU:HB2	2.16	0.45
1:A:83:ASN:HD22	1:A:83:ASN:C	2.24	0.44
1:B:480:SER:HA	1:B:481:PRO:C	2.41	0.44
1:B:134:LEU:HB3	1:B:135:PRO:HD3	1.99	0.44
1:A:348:ASN:HD22	1:A:348:ASN:N	2.16	0.44
1:B:332:TRP:O	1:B:335:GLU:HB2	2.16	0.44
1:A:348:ASN:H	1:A:348:ASN:ND2	2.16	0.44
1:A:241:PHE:HB3	1:A:242:PRO:CD	2.48	0.44
1:B:243:GLN:HB3	1:B:358:PHE:CE2	2.53	0.44
1:A:350:LEU:C	1:A:350:LEU:HD23	2.43	0.44
1:A:243:GLN:HB3	1:A:358:PHE:CE2	2.52	0.43
1:B:241:PHE:HB3	1:B:242:PRO:CD	2.49	0.43
1:B:241:PHE:HB3	1:B:242:PRO:HD2	2.00	0.43
1:B:457:TRP:HA	4:B:1902:H4B:N1	2.34	0.43
1:B:374:VAL:O	1:B:378:CYS:HB2	2.19	0.42
1:B:348:ASN:HD22	1:B:348:ASN:N	2.17	0.42
1:A:165:LYS:N	1:A:165:LYS:HD2	2.34	0.42
1:A:334:GLN:OE1	1:A:335:GLU:N	2.53	0.42
1:A:132:GLU:O	1:A:135:PRO:HD2	2.20	0.42
1:A:374:VAL:O	1:A:378:CYS:HB2	2.20	0.42
1:B:99:THR:HG23	1:B:478:VAL:O	2.21	0.41
1:A:213:THR:OG1	1:A:216:GLU:HG3	2.21	0.41
1:B:81:ILE:HD13	1:B:81:ILE:HA	1.92	0.41
1:A:137:ALA:O	1:A:141:ILE:HG12	2.20	0.41
1:B:77:GLN:O	1:B:96:HIS:HE1	2.04	0.41
1:A:159:ARG:O	1:A:163:VAL:HG23	2.21	0.41
1:A:375:ARG:O	1:A:379:ASP:HB2	2.21	0.40
1:B:99:THR:HB	1:B:100:SER:H	1.51	0.40
1:B:223:ARG:HD3	7:B:1189:HOH:O	2.20	0.40
1:B:102:PHE:CG	1:B:103:THR:N	2.86	0.40
1:B:348:ASN:H	1:B:348:ASN:ND2	2.20	0.40
1:B:164:THR:O	1:B:168:GLU:HG3	2.22	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the sym-

metry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:A:1082:HOH:O	7:A:1082:HOH:O[11_655]	0.98	1.22
7:A:1376:HOH:O	7:A:1376:HOH:O[11_655]	1.19	1.01
7:B:1215:HOH:O	7:B:1303:HOH:O[9_766]	2.00	0.20

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	412/433 (95%)	393 (95%)	18 (4%)	1 (0%)	43	42
1	B	419/433 (97%)	402 (96%)	15 (4%)	2 (0%)	24	21
All	All	831/866 (96%)	795 (96%)	33 (4%)	3 (0%)	30	27

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	100	SER
1	B	100	SER
1	B	99	THR

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	362/381 (95%)	355 (98%)	7 (2%)	50	56
1	B	366/381 (96%)	355 (97%)	11 (3%)	36	38

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	728/762 (96%)	710 (98%)	18 (2%)	42	45

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	83	ASN
1	A	114	MET
1	A	295	ARG
1	A	331	GLU
1	A	334	GLN
1	A	348	ASN
1	A	475	LEU
1	B	110	LEU
1	B	130	LEU
1	B	131	GLU
1	B	223	ARG
1	B	265	GLN
1	B	329	LYS
1	B	331	GLU
1	B	334	GLN
1	B	348	ASN
1	B	353	VAL
1	B	475	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	83	ASN
1	A	96	HIS
1	A	115	ASN
1	A	348	ASN
1	A	418	HIS
1	A	421	GLN
1	A	431	HIS
1	A	442	GLN
1	A	471	HIS
1	A	486	GLN
1	B	96	HIS
1	B	115	ASN
1	B	348	ASN
1	B	418	HIS

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Mol	Chain	Res	Type
1	B	421	GLN
1	B	486	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 10 ligands modelled in this entry, 2 are monoatomic - leaving 8 for Mogul analysis.

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	$\# Z > 2$
6	EDO	A	5002	-	3,3,3	0.47	0	2,2,2	0.67	0
3	HEM	A	901	-	50,50,50	1.44	6 (12%)	67,82,82	0.74	0
3	HEM	B	1901	-	50,50,50	1.50	6 (12%)	67,82,82	0.89	1 (1%)
4	H4B	B	1902	-	17,18,18	1.93	4 (23%)	14,26,26	2.06	5 (35%)
5	AT2	B	1906	-	20,20,20	1.85	5 (25%)	25,26,26	1.32	3 (12%)
4	H4B	A	902	-	17,18,18	1.85	3 (17%)	14,26,26	2.09	5 (35%)
5	AT2	A	906	-	20,20,20	1.93	5 (25%)	25,26,26	1.34	3 (12%)
6	EDO	A	5001	-	3,3,3	0.54	0	2,2,2	0.60	0

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
6	EDO	A	5002	-	-	0/1/1/1	-
3	HEM	A	901	-	-	0/14/54/54	-
3	HEM	B	1901	-	-	0/14/54/54	-
4	H4B	B	1902	-	-	0/8/17/17	0/2/2/2
5	AT2	B	1906	-	-	0/11/21/21	0/2/2/2
4	H4B	A	902	-	-	0/8/17/17	0/2/2/2
5	AT2	A	906	-	-	0/11/21/21	0/2/2/2
6	EDO	A	5001	-	-	0/1/1/1	-

All (29) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	906	AT2	O16-C15	4.99	1.28	1.21
5	B	1906	AT2	O16-C15	4.59	1.28	1.21
5	A	906	AT2	C6-N5	4.25	1.41	1.34
5	B	1906	AT2	C6-N5	4.19	1.41	1.34
4	A	902	H4B	C4A-C8A	4.13	1.45	1.38
4	B	1902	H4B	C4A-C8A	4.10	1.45	1.38
4	A	902	H4B	C7-C6	-4.00	1.48	1.52
3	A	901	HEM	FE-NB	3.96	2.07	1.94
3	B	1901	HEM	FE-ND	3.94	2.07	1.94
4	B	1902	H4B	C7-N8	-3.89	1.42	1.46
3	B	1901	HEM	FE-NB	3.88	2.06	1.94
4	A	902	H4B	C7-N8	-3.73	1.42	1.46
3	A	901	HEM	FE-NC	3.65	2.07	1.95
3	A	901	HEM	FE-ND	3.61	2.06	1.94
3	B	1901	HEM	FE-NA	3.36	2.06	1.95
4	B	1902	H4B	C7-C6	-3.22	1.49	1.52
5	A	906	AT2	C4-N5	2.91	1.40	1.34
3	B	1901	HEM	FE-NC	2.90	2.04	1.95
4	B	1902	H4B	C9-C10	2.89	1.58	1.53
3	A	901	HEM	CAB-C3B	-2.63	1.40	1.47
5	B	1906	AT2	C4-N5	2.58	1.40	1.34
3	B	1901	HEM	CAB-C3B	-2.52	1.40	1.47
3	A	901	HEM	FE-NA	2.49	2.03	1.95
5	B	1906	AT2	C3-C4	2.46	1.43	1.38
5	A	906	AT2	C3-C4	2.40	1.43	1.38
3	B	1901	HEM	CAC-C3C	-2.22	1.41	1.47
5	A	906	AT2	C10-C9	2.15	1.57	1.52
5	B	1906	AT2	C10-C9	2.11	1.56	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	901	HEM	CAC-C3C	-2.02	1.42	1.47

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	902	H4B	C2-N1-C8A	4.50	121.31	113.36
4	B	1902	H4B	C2-N1-C8A	4.44	121.20	113.36
4	A	902	H4B	O4-C4-C4A	-4.11	117.35	127.26
4	B	1902	H4B	O4-C4-C4A	-3.99	117.65	127.26
4	B	1902	H4B	C4A-C4-N3	2.98	120.32	112.13
4	A	902	H4B	C4A-C4-N3	2.89	120.06	112.13
5	A	906	AT2	C3-C4-N5	-2.84	120.49	123.97
5	B	1906	AT2	C3-C4-N5	-2.69	120.67	123.97
5	A	906	AT2	C7-C6-N5	-2.63	119.47	122.92
5	B	1906	AT2	C7-C6-N5	-2.63	119.47	122.92
3	B	1901	HEM	CMD-C2D-C1D	2.61	129.11	125.03
4	B	1902	H4B	C4-C4A-N5	2.44	122.92	116.27
4	A	902	H4B	C4-C4A-N5	2.32	122.59	116.27
4	B	1902	H4B	C2-N3-C4	-2.29	120.95	125.11
4	A	902	H4B	C2-N3-C4	-2.23	121.07	125.11
5	B	1906	AT2	O16-C15-N12	-2.19	120.25	124.30
5	A	906	AT2	O16-C15-N12	-2.05	120.51	124.30

There are no chirality outliers.

There are no torsion outliers.

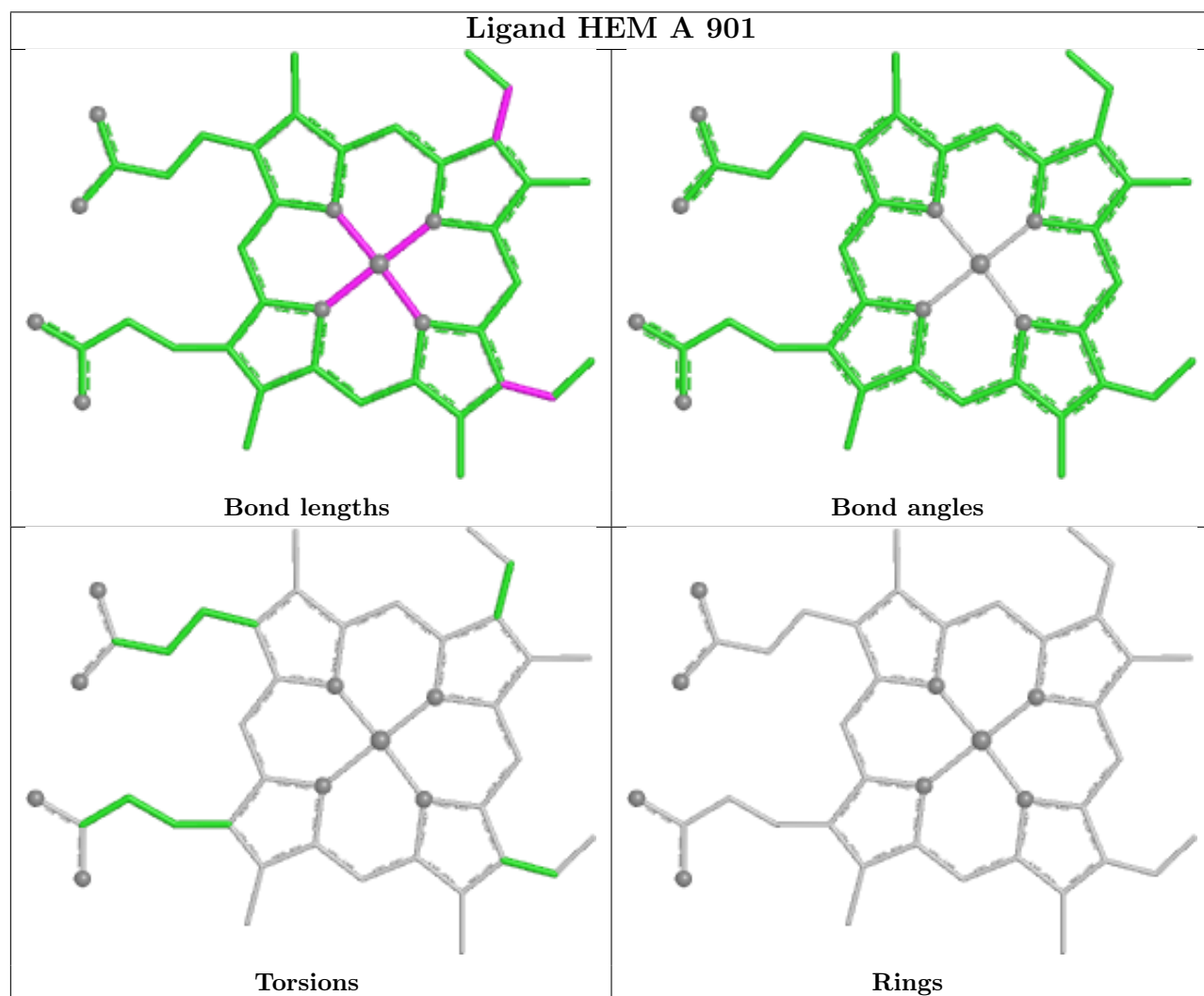
There are no ring outliers.

1 monomer is involved in 1 short contact:

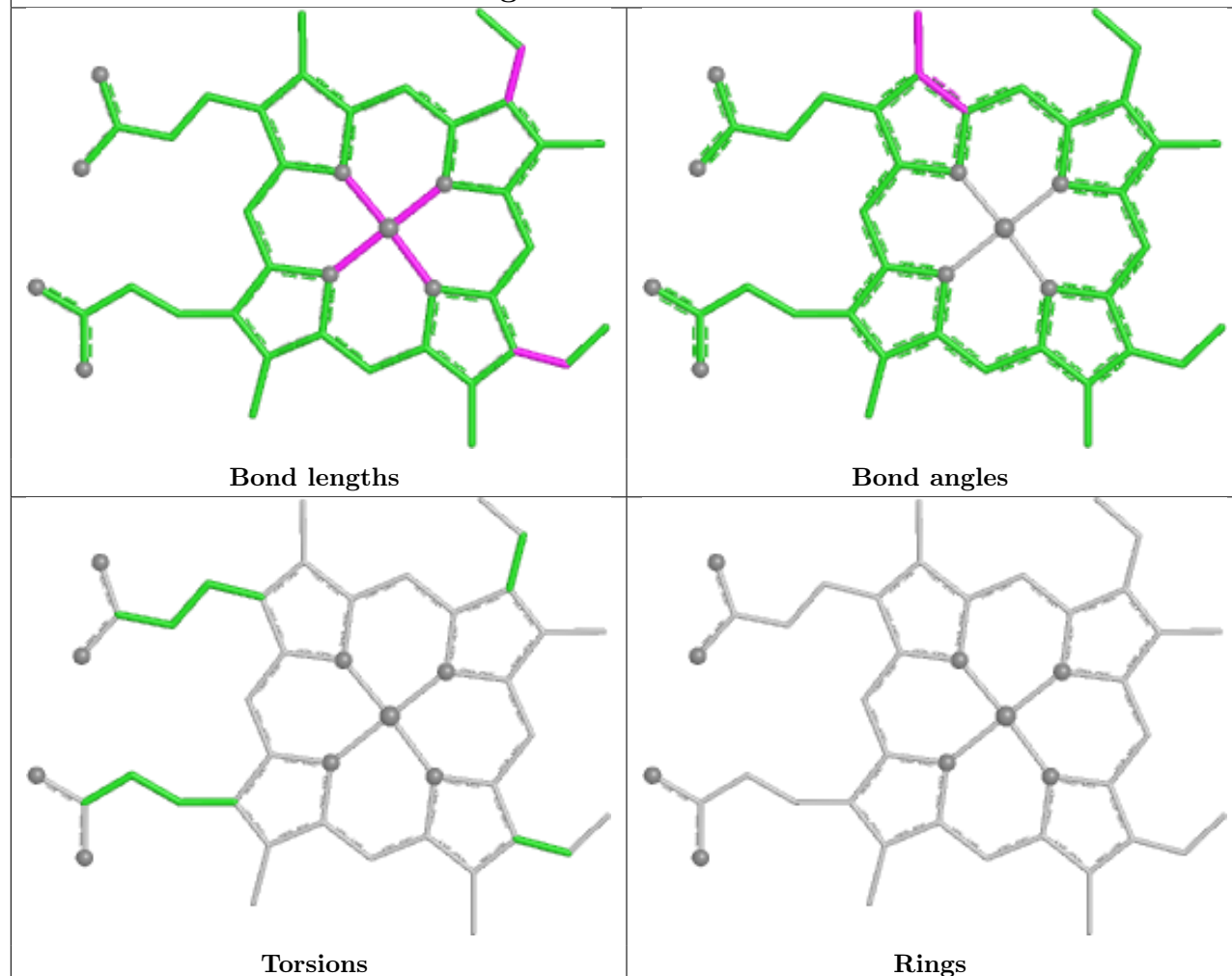
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	1902	H4B	1	0

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.

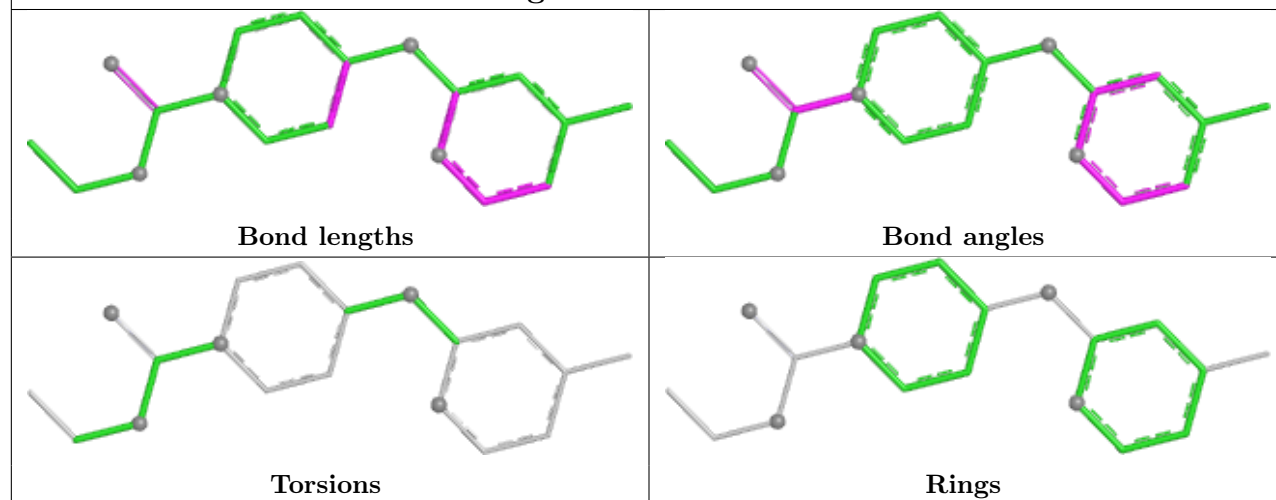
The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

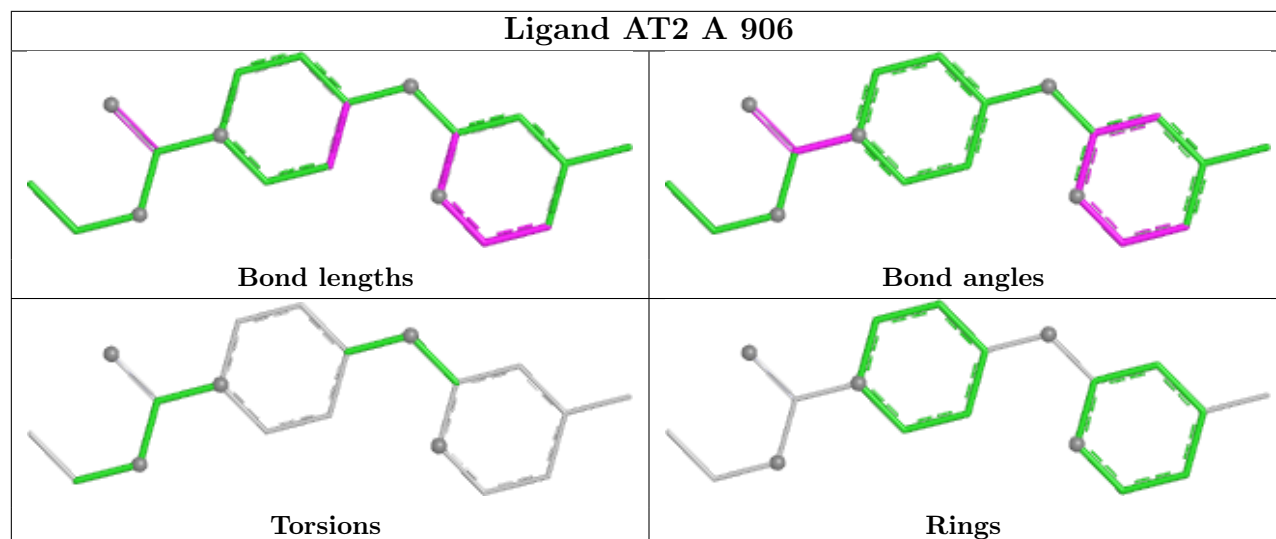


Ligand HEM B 1901



Ligand AT2 B 1906





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	416/433 (96%)	0.57	38 (9%) 15 13	21, 35, 51, 64	2 (0%)
1	B	421/433 (97%)	0.51	42 (9%) 12 11	15, 33, 50, 67	7 (1%)
All	All	837/866 (96%)	0.54	80 (9%) 13 12	15, 34, 50, 67	9 (1%)

All (80) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	109	CYS	9.4
1	A	109	CYS	8.5
1	B	100	SER	8.3
1	A	103	THR	8.0
1	B	102	PHE	8.0
1	B	105	LYS	6.2
1	A	102	PHE	5.9
1	B	107	LYS	5.5
1	B	104	CYS	5.3
1	B	101	ASP	5.2
1	A	497	ASN	4.8
1	B	497	ASN	4.8
1	B	114	MET	4.6
1	B	103	THR	4.2
1	B	131	GLU	4.1
1	A	496	GLN	4.1
1	B	99	THR	4.1
1	A	77	GLN	4.0
1	A	78	TYR	4.0
1	A	100	SER	3.8
1	A	101	ASP	3.8
1	A	111	GLY	3.6
1	A	99	THR	3.5
1	A	279	GLU	3.5

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Mol	Chain	Res	Type	RSRZ
1	B	108	SER	3.4
1	B	496	GLN	3.3
1	A	146	GLY	3.2
1	A	210	ASN	3.2
1	B	494	ILE	3.2
1	A	139	GLU	3.1
1	B	111	GLY	3.1
1	B	106	SER	3.1
1	B	442	GLN	3.1
1	A	157	LEU	3.1
1	B	334	GLN	3.1
1	A	110	LEU	3.0
1	A	202	ASN	2.9
1	B	130	LEU	2.9
1	A	292	ARG	2.9
1	B	78	TYR	2.9
1	A	165	LYS	2.9
1	A	153	ILE	2.9
1	B	210	ASN	2.9
1	A	334	GLN	2.8
1	B	112	SER	2.6
1	B	288	GLY	2.6
1	B	110	LEU	2.6
1	A	149	LYS	2.6
1	B	391	ARG	2.6
1	B	148	PHE	2.5
1	A	307	GLY	2.5
1	A	446	ARG	2.5
1	A	169	THR	2.5
1	A	85	GLY	2.5
1	B	126	LYS	2.5
1	A	154	GLU	2.4
1	B	96	HIS	2.4
1	B	292	ARG	2.4
1	A	88	GLU	2.4
1	A	308	GLN	2.4
1	A	494	ILE	2.4
1	B	77	GLN	2.3
1	A	151	ALA	2.3
1	B	158	ALA	2.3
1	A	295	ARG	2.3
1	A	331	GLU	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	151	ALA	2.3
1	A	495	TRP	2.2
1	B	279	GLU	2.2
1	B	124	ARG	2.2
1	B	150	GLU	2.1
1	B	308	GLN	2.1
1	A	152	LYS	2.1
1	A	388	GLU	2.1
1	B	265	GLN	2.1
1	A	127	PRO	2.1
1	B	495	TRP	2.1
1	B	154	GLU	2.0
1	B	87	GLY	2.0
1	B	330	TYR	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
6	EDO	A	5002	4/4	0.61	0.26	55,55,55,55	0
6	EDO	A	5001	4/4	0.77	0.24	52,54,55,55	0
4	H4B	B	1902	17/17	0.93	0.08	22,25,28,29	0
2	ZN	A	900	1/1	0.93	0.15	45,45,45,45	1
4	H4B	A	902	17/17	0.93	0.08	23,25,29,29	0
5	AT2	A	906	19/19	0.95	0.07	20,22,29,30	0
5	AT2	B	1906	19/19	0.95	0.07	18,22,28,30	0
2	ZN	B	1900	1/1	0.96	0.22	34,34,34,34	1
3	HEM	A	901	43/43	0.98	0.07	21,24,26,29	0

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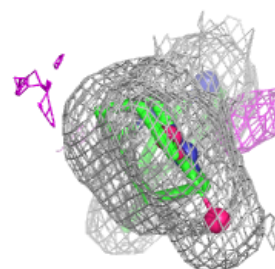
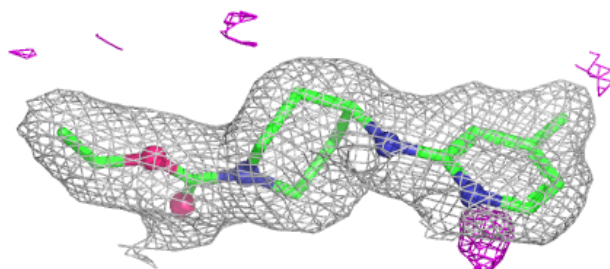
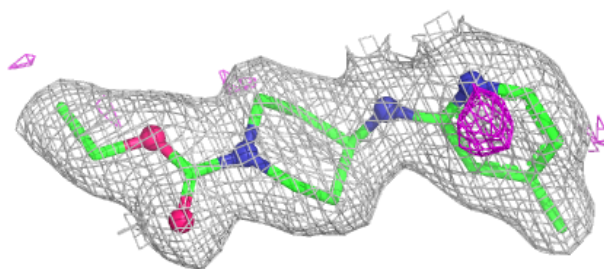
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	HEM	B	1901	43/43	0.98	0.06	18,22,25,26	0

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.

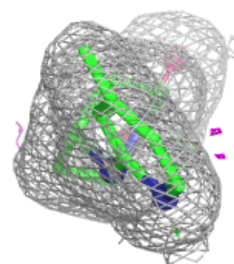
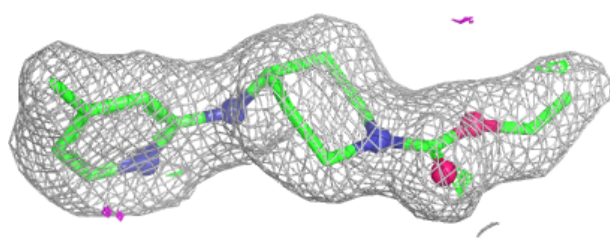
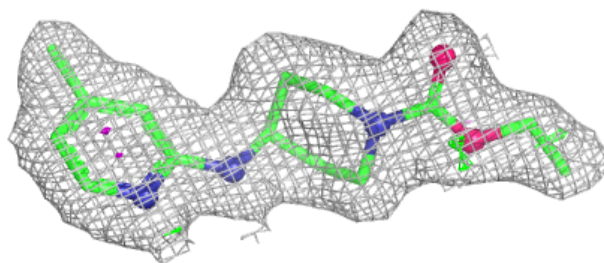
Electron density around AT2 A 906:

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



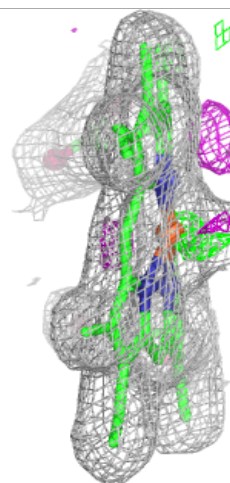
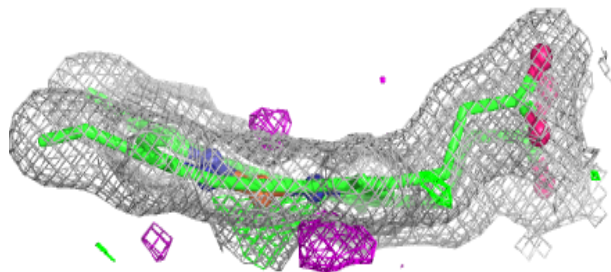
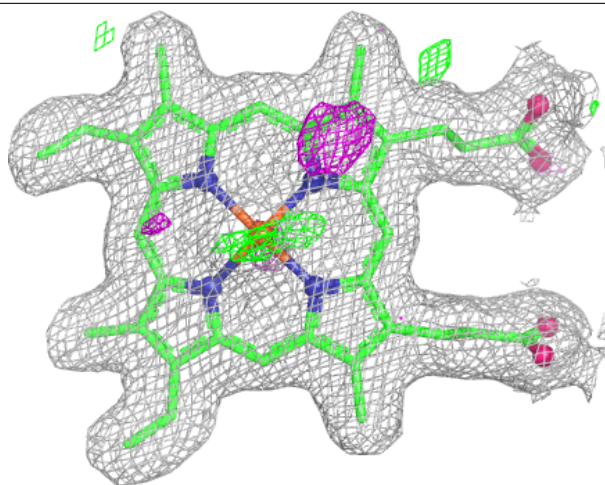
Electron density around AT2 B 1906:

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



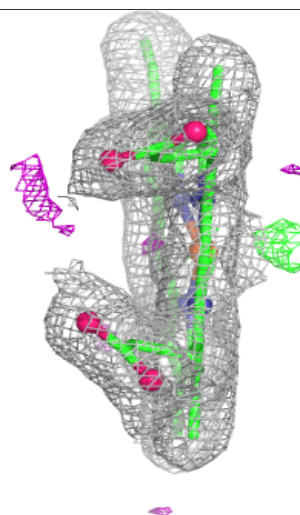
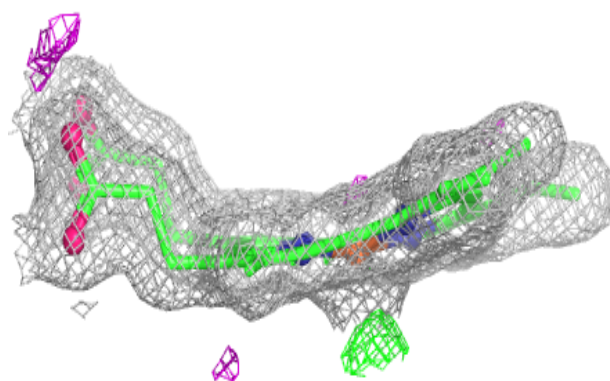
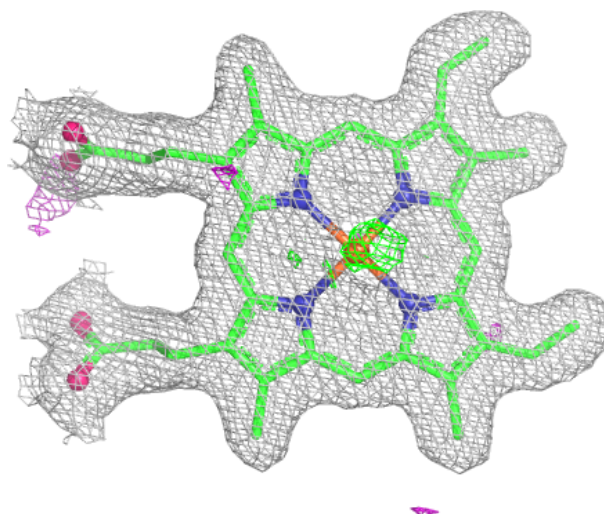
Electron density around HEM A 901:

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



Electron density around HEM B 1901:

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



6.5 Other polymers ⓘ

There are no such residues in this entry.