



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 09:24 PM UTC

PDB ID : 1M9K / pdb_00001m9k
Title : Human Endothelial Nitric Oxide Synthase with 7-Nitroindazole Bound
Authors : Rosenfeld, R.J.; Garcin, E.D.; Panda, K.; Andersson, G.; Aberg, A.; Wallace, A.V.; Stuehr, D.J.; Tainer, J.A.; Getzoff, E.D.
Deposited on : 2002-07-29
Resolution : 2.01 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

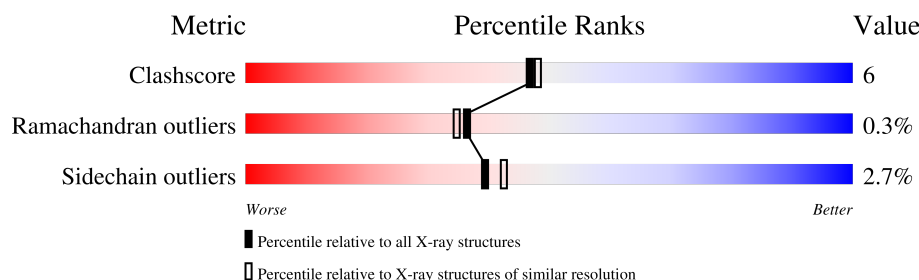
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.01 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	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	415	
1	B	415	

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	7NI	A	906[A]	-	X	-	-
4	7NI	A	906[B]	-	X	-	-
4	7NI	B	907[A]	-	X	-	-

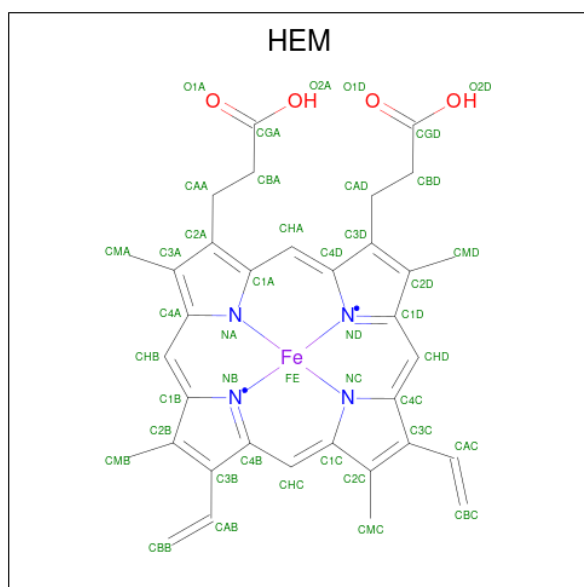
In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	400	Total 3179	C 2028	N 559	O 576	S 16	0	0	0
1	B	401	Total 3189	C 2034	N 559	O 580	S 16	0	0	0

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

- $$\text{C}_{34}\text{H}_{32}\text{FeN}_4\text{O}_4).$$



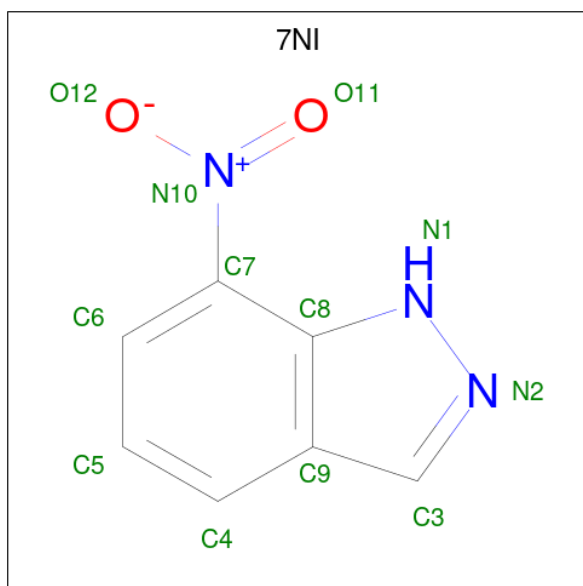
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

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

- Molecule 4 is 7-NITROINDAZOLE (CCD ID: 7NI) (formula: $C_7H_5N_3O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	1
			24	14	6	4		
4	B	1	Total	C	N	O	0	1
			24	14	6	4		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	208	Total	O	0	0
			208	208		
5	B	204	Total	O	0	0
			204	204		

Note EDS was not executed.

- [illegible]

4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	69.79Å 91.57Å 156.10Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.80 – 2.01	Depositor
% Data completeness (in resolution range)	98.6 (19.80-2.01)	Depositor
R_{merge}	0.04	Depositor
R_{sym}	0.04	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.196 , 0.224	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6915	wwPDB-VP
Average B, all atoms (Å ²)	33.0	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: 7NI, HEM, ZN

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	0.48	0/3270	0.96	14/4457 (0.3%)
1	B	0.50	0/3280	0.96	13/4470 (0.3%)
All	All	0.49	0/6550	0.96	27/8927 (0.3%)

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

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

There are no bond length outliers.

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	358	MET	N-CA-C	-7.66	96.79	108.96
1	B	359	SER	N-CA-C	7.55	120.46	111.33
1	A	358	MET	N-CA-C	-7.44	97.10	109.46
1	B	357	TYR	N-CA-C	7.43	121.27	110.28
1	B	437	ALA	N-CA-C	7.00	118.70	111.14
1	A	359	SER	N-CA-C	6.90	119.68	111.33
1	A	121	GLU	N-CA-C	-6.83	103.62	112.23
1	B	333	LEU	N-CA-C	6.68	120.71	108.69
1	A	241	PHE	N-CA-C	-6.61	98.85	109.96
1	A	357	TYR	N-CA-C	6.59	120.62	110.20
1	A	190	TRP	N-CA-C	6.52	119.22	111.33
1	A	333	LEU	N-CA-C	6.51	120.40	108.69
1	B	241	PHE	N-CA-C	-6.47	99.98	110.20
1	A	187	ARG	N-CA-C	6.45	120.75	113.01

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	282	GLY	N-CA-C	-6.14	103.82	112.60
1	B	190	TRP	N-CA-C	5.97	119.39	111.75
1	A	73	ASN	N-CA-C	-5.65	99.53	108.73
1	A	449	VAL	CA-C-N	5.42	123.61	119.66
1	A	449	VAL	C-N-CA	5.42	123.61	119.66
1	A	348	PHE	CA-C-N	5.41	125.08	119.56
1	A	348	PHE	C-N-CA	5.41	125.08	119.56
1	B	280	THR	N-CA-C	-5.41	101.42	109.42
1	B	187	ARG	N-CA-C	5.36	119.30	112.87
1	B	450	PRO	CA-C-N	5.05	124.72	119.56
1	B	450	PRO	C-N-CA	5.05	124.72	119.56
1	A	278	GLY	N-CA-C	5.03	122.36	115.27
1	B	121	GLU	N-CA-C	-5.01	106.01	111.82

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	475	TYR	Sidechain

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	3179	0	3077	31	0
1	B	3189	0	3087	43	0
2	A	1	0	0	0	0
3	A	43	0	30	0	0
3	B	43	0	30	0	0
4	A	24	0	10	0	0
4	B	24	0	10	0	0
5	A	208	0	0	4	0
5	B	204	0	0	3	0
All	All	6915	0	6244	72	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 6.

All (72) 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:292:LEU:HD13	1:B:302:LEU:HD13	1.51	0.91
1:B:250:ARG:HH11	1:B:250:ARG:HG2	1.36	0.91
1:A:250:ARG:HH11	1:A:250:ARG:HG3	1.41	0.83
1:B:250:ARG:HH11	1:B:250:ARG:CG	1.94	0.81
1:B:453:SER:HB3	1:B:456:LEU:HD12	1.65	0.79
1:A:247:GLN:HB2	1:A:250:ARG:HD3	1.69	0.75
1:B:124:LEU:HD11	1:B:154:GLU:HG2	1.71	0.72
1:B:250:ARG:HG2	1:B:250:ARG:NH1	1.96	0.70
1:A:149:ARG:HD2	1:A:169:GLU:OE1	1.92	0.70
1:A:250:ARG:HG3	1:A:250:ARG:NH1	2.11	0.65
1:B:216:LYS:HG3	1:B:309:LEU:HD22	1.80	0.63
1:B:250:ARG:HH11	1:B:250:ARG:HA	1.62	0.63
1:A:67:LYS:HA	1:A:67:LYS:HE2	1.79	0.62
1:B:73:ASN:HD22	1:B:462:GLN:NE2	1.98	0.61
1:B:90:GLN:HB3	1:B:468:PHE:CD2	2.36	0.60
1:A:453:SER:HB3	1:A:456:LEU:HD12	1.86	0.57
1:B:176:GLN:HE22	1:B:179:ARG:HH11	1.52	0.57
1:A:368:CYS:SG	1:A:376:LEU:HD13	2.45	0.56
1:B:368:CYS:SG	1:B:376:LEU:HD13	2.46	0.56
1:A:176:GLN:HE22	1:A:179:ARG:HH11	1.54	0.55
1:B:73:ASN:HD22	1:B:462:GLN:HE21	1.54	0.55
1:A:376:LEU:HB2	5:A:946:HOH:O	2.06	0.54
1:A:246:SER:OG	1:A:250:ARG:NH1	2.40	0.54
1:A:235:CYS:HB3	1:A:238:ARG:HE	1.73	0.54
1:B:67:LYS:HE3	1:B:68:PHE:CZ	2.44	0.53
1:A:428:MET:HE1	1:A:462:GLN:HE21	1.73	0.53
1:B:376:LEU:HB2	5:B:947:HOH:O	2.09	0.52
1:B:474:ARG:HD2	5:B:1074:HOH:O	2.08	0.52
1:A:90:GLN:HB3	1:A:468:PHE:CD2	2.45	0.52
1:B:250:ARG:NH1	1:B:331:TYR:OH	2.44	0.51
1:A:358:MET:HE1	5:A:982:HOH:O	2.11	0.50
1:B:255:ARG:HG3	5:B:1024:HOH:O	2.11	0.50
1:B:135:TYR:CE1	1:B:149:ARG:HG3	2.46	0.50
1:A:156:GLU:HG3	5:A:1083:HOH:O	2.11	0.50
1:A:124:LEU:HD23	1:A:128:ARG:NH2	2.27	0.49
1:A:470:SER:HA	1:A:471:PRO:C	2.38	0.49
1:A:183:ARG:HD3	1:A:447:TRP:CD2	2.48	0.49
1:B:292:LEU:HD13	1:B:302:LEU:CD1	2.34	0.48
1:B:250:ARG:CG	1:B:250:ARG:NH1	2.61	0.48
1:B:149:ARG:HD2	1:B:169:GLU:OE1	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:250:ARG:NH1	1:B:250:ARG:HA	2.27	0.47
1:B:244:TRP:HB2	1:B:292:LEU:HB3	1.96	0.46
1:B:150:LEU:O	1:B:154:GLU:HG3	2.16	0.46
1:B:257:GLN:O	1:B:258:ASP:HB3	2.14	0.46
1:B:480:TRP:HE1	1:B:481:LYS:HE3	1.80	0.46
1:A:364:THR:HG21	1:A:452:ILE:HG23	1.98	0.46
1:B:135:TYR:HE1	1:B:149:ARG:HG3	1.81	0.46
1:B:470:SER:HA	1:B:471:PRO:C	2.42	0.45
1:A:238:ARG:NH2	5:A:1086:HOH:O	2.49	0.45
1:B:428:MET:HE1	1:B:462:GLN:HE21	1.82	0.44
1:A:257:GLN:NE2	1:A:258:ASP:N	2.66	0.44
1:B:478:ASP:HB2	1:B:481:LYS:HD2	2.00	0.44
1:A:247:GLN:HB2	1:A:250:ARG:CD	2.45	0.44
1:B:290:PRO:HB3	1:B:304:LEU:CD2	2.48	0.43
1:B:219:THR:HA	1:B:224:LEU:HD22	2.00	0.43
1:A:75:GLU:HG3	1:B:370:PRO:HG2	1.99	0.43
1:B:364:THR:O	1:B:368:CYS:HB2	2.18	0.43
1:A:257:GLN:NE2	1:A:258:ASP:H	2.17	0.42
1:B:124:LEU:O	1:B:128:ARG:HG3	2.19	0.42
1:B:250:ARG:HH11	1:B:250:ARG:CB	2.31	0.42
1:B:90:GLN:HG2	1:B:91:ASP:N	2.34	0.42
1:A:364:THR:O	1:A:368:CYS:HB2	2.20	0.42
1:B:204:ALA:HA	1:B:207:MET:HE3	2.02	0.41
1:A:103:LEU:HD22	1:B:463:GLU:HB3	2.01	0.41
1:A:150:LEU:O	1:A:153:VAL:HG22	2.19	0.41
1:B:207:MET:HE2	1:B:241:PHE:HD2	1.85	0.41
1:B:207:MET:HG3	1:B:231:PHE:CE2	2.55	0.41
1:A:257:GLN:HE21	1:A:258:ASP:N	2.19	0.41
1:A:216:LYS:HG2	1:A:309:LEU:HD22	2.02	0.41
1:B:250:ARG:HH11	1:B:250:ARG:CA	2.29	0.41
1:A:73:ASN:HD22	1:A:462:GLN:NE2	2.19	0.40
1:A:247:GLN:CB	1:A:250:ARG:HD3	2.47	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	396/415 (95%)	386 (98%)	9 (2%)	1 (0%)	36	35
1	B	397/415 (96%)	388 (98%)	8 (2%)	1 (0%)	36	35
All	All	793/830 (96%)	774 (98%)	17 (2%)	2 (0%)	36	35

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	258	ASP
1	B	258	ASP

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	336/352 (96%)	325 (97%)	11 (3%)	33	34
1	B	338/352 (96%)	331 (98%)	7 (2%)	47	52
All	All	674/704 (96%)	656 (97%)	18 (3%)	39	42

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

Mol	Chain	Res	Type
1	A	67	LYS
1	A	87	GLN
1	A	103	LEU
1	A	148	GLN
1	A	149	ARG
1	A	207	MET
1	A	224	LEU
1	A	250	ARG
1	A	255	ARG
1	A	262	ARG

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Mol	Chain	Res	Type
1	A	417	ILE
1	B	103	LEU
1	B	124	LEU
1	B	148	GLN
1	B	149	ARG
1	B	250	ARG
1	B	302	LEU
1	B	376	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	90	GLN
1	A	133	GLN
1	A	164	GLN
1	A	176	GLN
1	A	233	GLN
1	A	256	GLN
1	A	257	GLN
1	A	435	GLN
1	A	462	GLN
1	A	466	ASN
1	B	126	GLN
1	B	133	GLN
1	B	176	GLN
1	B	256	GLN
1	B	462	GLN
1	B	466	ASN
1	B	476	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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 1 is monoatomic - leaving 6 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	$\# Z > 2$
4	7NI	B	907[B]	-	13,13,13	3.80	6 (46%)	13,18,18	5.88	8 (61%)
3	HEM	B	901	1	50,50,50	1.36	6 (12%)	67,82,82	0.95	3 (4%)
4	7NI	A	906[B]	-	13,13,13	3.52	8 (61%)	13,18,18	6.09	7 (53%)
3	HEM	A	901	1	50,50,50	1.23	6 (12%)	67,82,82	0.90	0
4	7NI	A	906[A]	-	13,13,13	3.58	9 (69%)	13,18,18	6.12	7 (53%)
4	7NI	B	907[A]	-	13,13,13	3.41	8 (61%)	13,18,18	6.24	7 (53%)

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	7NI	B	907[B]	-	-	0/2/4/4	0/2/2/2
3	HEM	B	901	1	-	0/14/54/54	-
4	7NI	A	906[B]	-	-	0/2/4/4	0/2/2/2
3	HEM	A	901	1	-	2/14/54/54	-
4	7NI	A	906[A]	-	-	0/2/4/4	0/2/2/2
4	7NI	B	907[A]	-	-	0/2/4/4	0/2/2/2

All (43) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	907[B]	7NI	C5-C4	9.49	1.55	1.38
4	A	906[A]	7NI	C5-C4	7.63	1.51	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	907[A]	7NI	C5-C4	7.58	1.51	1.38
4	A	906[B]	7NI	C5-C4	7.57	1.51	1.38
4	B	907[B]	7NI	C9-C3	5.27	1.48	1.42
4	A	906[A]	7NI	C8-C7	-4.53	1.34	1.42
4	B	907[A]	7NI	C9-C8	4.50	1.46	1.41
4	B	907[B]	7NI	C8-C7	-4.43	1.34	1.42
4	A	906[B]	7NI	C8-C7	-4.39	1.34	1.42
4	A	906[A]	7NI	C9-C8	4.37	1.46	1.41
4	B	907[A]	7NI	C8-C7	-4.28	1.34	1.42
4	A	906[A]	7NI	C4-C9	-4.28	1.33	1.41
4	A	906[B]	7NI	C4-C9	-4.25	1.33	1.41
4	A	906[B]	7NI	C9-C8	4.20	1.46	1.41
4	A	906[A]	7NI	C9-C3	4.19	1.47	1.42
4	B	907[A]	7NI	C4-C9	-4.11	1.33	1.41
4	B	907[B]	7NI	C9-C8	4.09	1.46	1.41
4	B	907[B]	7NI	C3-N2	3.85	1.34	1.32
4	A	906[B]	7NI	C9-C3	3.80	1.46	1.42
4	A	906[A]	7NI	N1-N2	-3.72	1.33	1.36
4	A	906[B]	7NI	N1-N2	-3.63	1.33	1.36
4	B	907[A]	7NI	C9-C3	3.61	1.46	1.42
3	B	901	HEM	FE-ND	3.54	2.05	1.94
4	B	907[A]	7NI	N1-N2	-3.35	1.34	1.36
3	B	901	HEM	FE-NC	3.35	2.06	1.95
3	A	901	HEM	FE-NA	3.05	2.05	1.95
3	B	901	HEM	FE-NA	3.03	2.05	1.95
4	A	906[B]	7NI	C3-N2	-2.90	1.30	1.32
3	A	901	HEM	FE-NC	2.89	2.04	1.95
4	A	906[B]	7NI	C8-N1	2.86	1.38	1.36
4	A	906[A]	7NI	C3-N2	-2.71	1.30	1.32
3	A	901	HEM	FE-NB	2.64	2.03	1.94
4	A	906[A]	7NI	C8-N1	2.62	1.38	1.36
3	A	901	HEM	FE-ND	2.57	2.02	1.94
4	B	907[B]	7NI	N1-N2	-2.55	1.34	1.36
3	B	901	HEM	FE-NB	2.48	2.02	1.94
3	A	901	HEM	CMC-C2C	2.32	1.55	1.50
4	B	907[A]	7NI	C3-N2	-2.29	1.30	1.32
3	B	901	HEM	CBC-CAC	2.22	1.41	1.30
3	B	901	HEM	CMC-C2C	2.14	1.55	1.50
3	A	901	HEM	CBC-CAC	2.11	1.40	1.30
4	A	906[A]	7NI	O12-N10	-2.08	1.21	1.35
4	B	907[A]	7NI	O12-N10	-2.04	1.21	1.35

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	907[A]	7NI	C3-N2-N1	15.66	114.20	106.02
4	A	906[A]	7NI	C3-N2-N1	15.51	114.12	106.02
4	A	906[B]	7NI	C3-N2-N1	15.24	113.97	106.02
4	B	907[B]	7NI	C9-C3-N2	-11.27	103.48	111.68
4	B	907[B]	7NI	C3-N2-N1	10.99	111.76	106.02
4	B	907[A]	7NI	C9-C3-N2	-10.32	104.17	111.68
4	A	906[A]	7NI	C9-C3-N2	-9.92	104.46	111.68
4	A	906[B]	7NI	C9-C3-N2	-9.70	104.62	111.68
4	B	907[B]	7NI	C8-C9-C3	8.64	108.34	103.94
4	A	906[B]	7NI	C9-C8-N1	-8.43	102.73	106.90
4	B	907[B]	7NI	C9-C8-N1	-8.35	102.77	106.90
4	A	906[A]	7NI	C9-C8-N1	-8.18	102.86	106.90
4	B	907[A]	7NI	C9-C8-N1	-8.01	102.94	106.90
4	B	907[A]	7NI	C8-C9-C3	7.40	107.70	103.94
4	A	906[B]	7NI	C8-C9-C3	7.19	107.60	103.94
4	A	906[A]	7NI	C8-C9-C3	6.89	107.45	103.94
4	B	907[B]	7NI	C8-N1-N2	4.65	113.66	111.42
4	B	907[B]	7NI	C5-C4-C9	-4.39	114.31	120.95
4	A	906[A]	7NI	C5-C4-C9	-3.75	115.29	120.95
4	B	907[A]	7NI	C5-C4-C9	-3.65	115.43	120.95
4	A	906[B]	7NI	C5-C4-C9	-3.62	115.47	120.95
4	A	906[A]	7NI	C4-C9-C8	3.32	121.75	119.04
4	B	907[A]	7NI	C4-C9-C8	3.28	121.71	119.04
4	A	906[B]	7NI	C4-C9-C8	3.16	121.61	119.04
4	B	907[B]	7NI	C4-C9-C8	2.41	121.00	119.04
3	B	901	HEM	CMD-C2D-C1D	2.39	128.77	125.03
3	B	901	HEM	CBD-CAD-C3D	-2.38	105.95	112.53
4	B	907[A]	7NI	C4-C9-C3	-2.25	130.59	135.71
4	B	907[B]	7NI	C4-C9-C3	-2.22	130.66	135.71
4	A	906[B]	7NI	C4-C9-C3	-2.17	130.78	135.71
4	A	906[A]	7NI	C4-C9-C3	-2.16	130.80	135.71
3	B	901	HEM	C1D-C2D-C3D	-2.02	104.85	106.98

There are no chirality outliers.

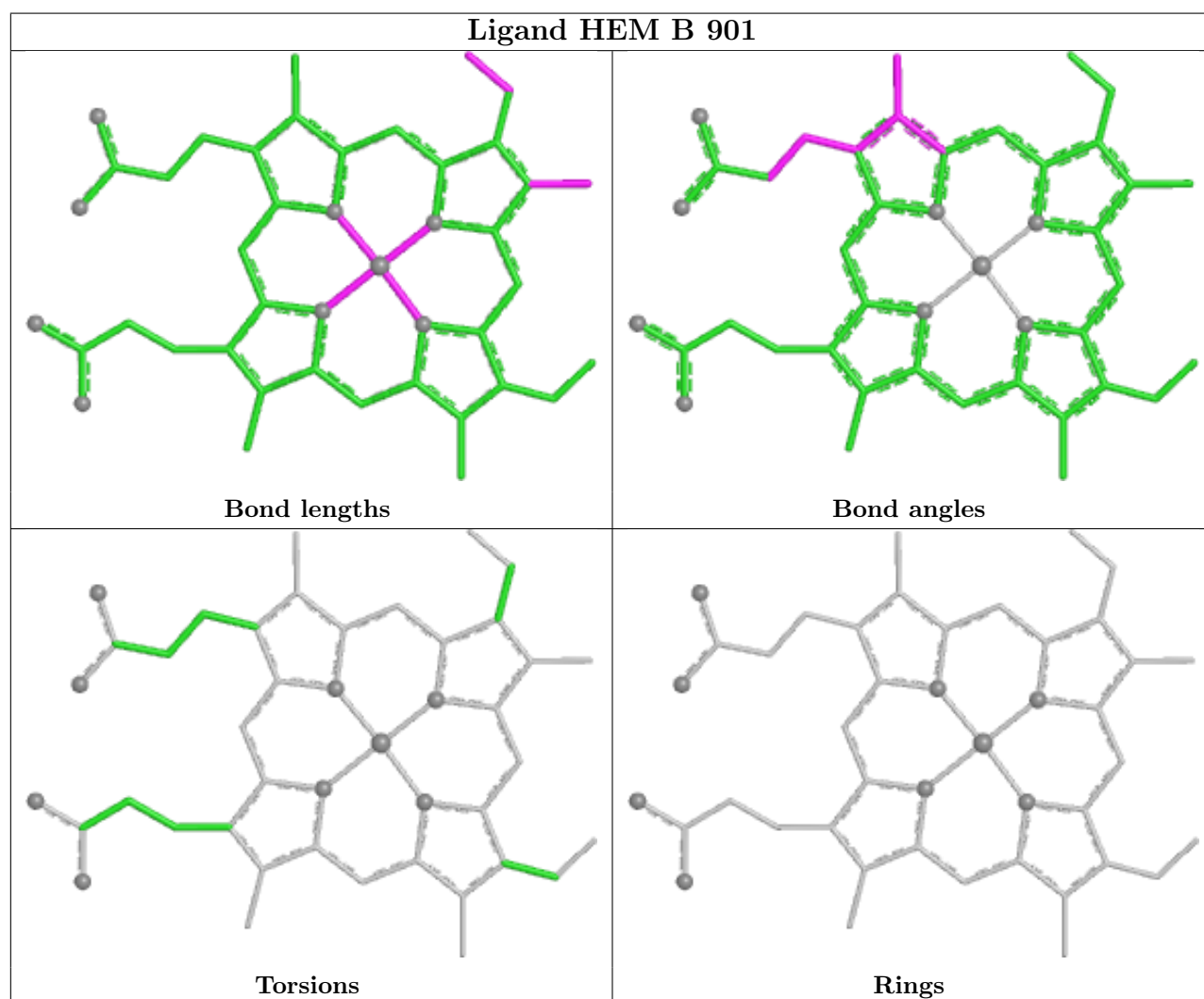
All (2) torsion outliers are listed below:

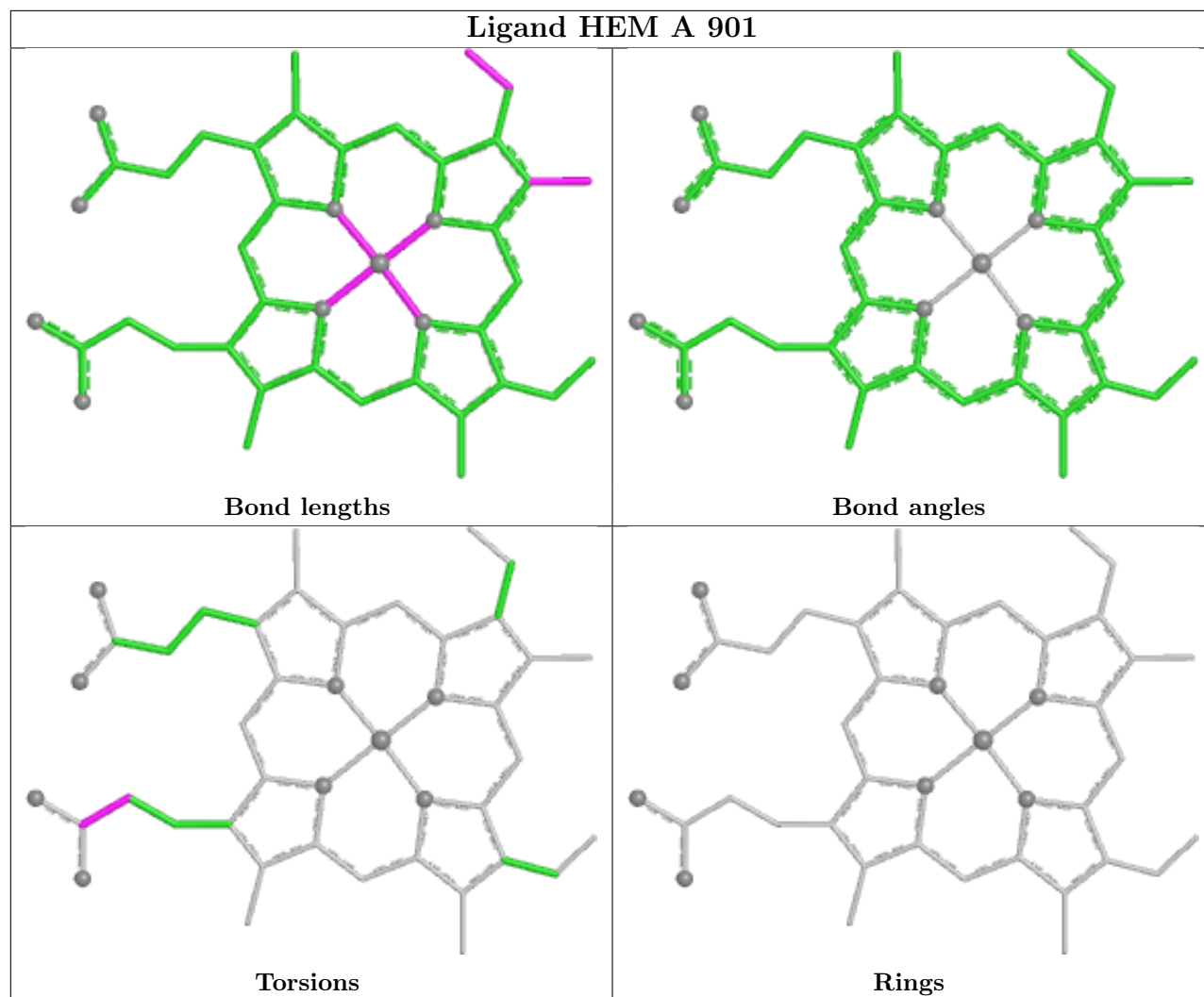
Mol	Chain	Res	Type	Atoms
3	A	901	HEM	CAA-CBA-CGA-O1A
3	A	901	HEM	CAA-CBA-CGA-O2A

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.