



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 06:46 AM UTC

PDB ID : 4UG8 / pdb_00004ug8
Title : Structure of Bacillus subtilis Nitric Oxide Synthase in complex with 6-(5-((3R,4R)-4-((6-azany1-4-methyl-pyridin-2-yl)methyl)pyrrolidin-3-yl)oxypentyl)-4-methyl-pyridin-2-amine
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Deposited on : 2015-03-22
Resolution : 1.89 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)	:	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

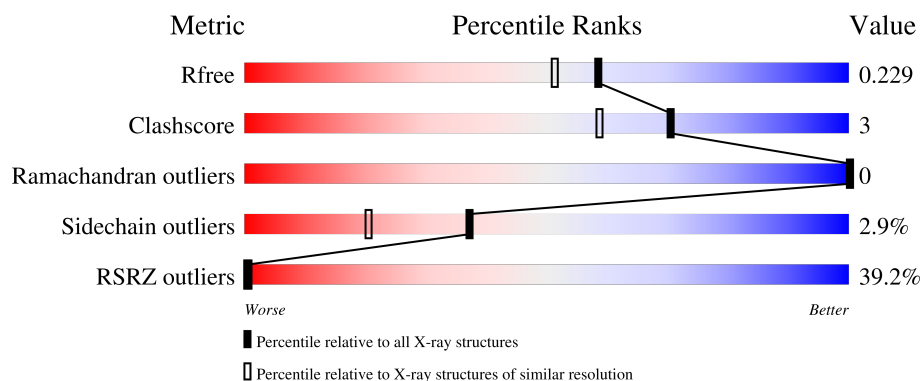
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.89 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	1220 (1.88-1.88)
Clashscore	190562	1234 (1.88-1.88)
Ramachandran outliers	187476	1222 (1.88-1.88)
Sidechain outliers	187428	1222 (1.88-1.88)
RSRZ outliers	180081	1220 (1.88-1.88)

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	363	39% 89% 11%

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 3279 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.

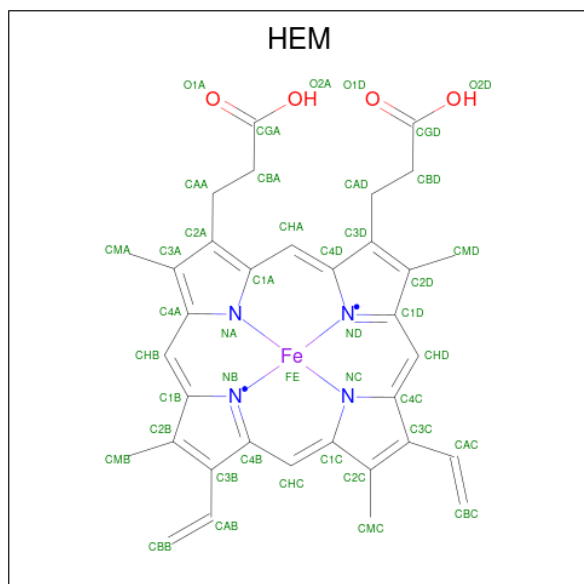
- Molecule 1 is a protein called NITRIC OXIDE SYNTHASE OXYGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	362	Total	C	N	O	S	0	0	0
			2941	1873	507	553	8			

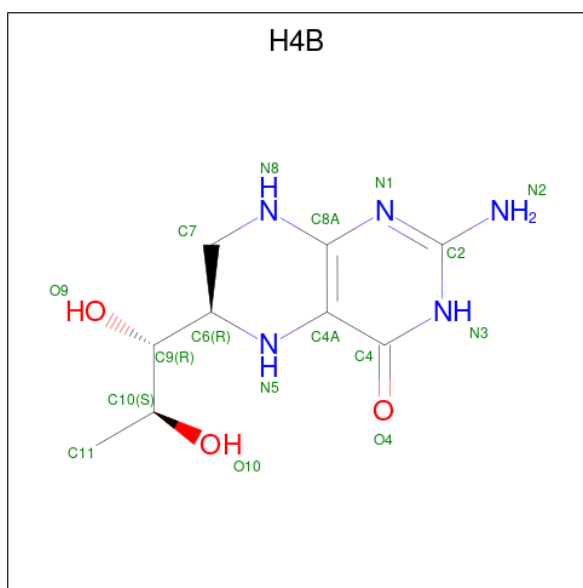
There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	25	ALA	GLU	engineered mutation	UNP O34453
A	26	ALA	GLU	engineered mutation	UNP O34453
A	316	ALA	GLU	engineered mutation	UNP O34453

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



- Molecule 3 is 5,6,7,8-TETRAHYDROBIOPTERIN (CCD ID: H4B) (formula: $C_9H_{15}N_5O_3$).

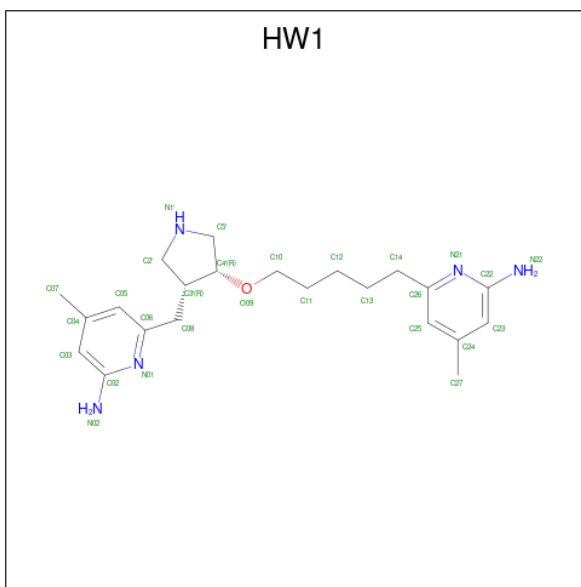


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			17	9	5	3		

- Molecule 4 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

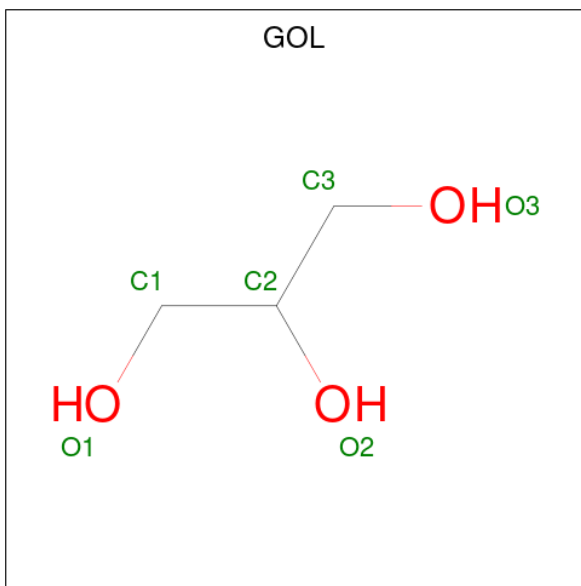
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Cl	0	0
			1	1		

- Molecule 5 is 6-[5-((3R,4R)-4-[(6-amino-4-methylpyridin-2-yl)methyl]pyrrolidin-3-yl)oxy)pentyl]-4-methylpyridin-2-amine (CCD ID: HW1) (formula: $C_{22}H_{33}N_5O$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			28	22	5	1		

- Molecule 6 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			6	3	3		

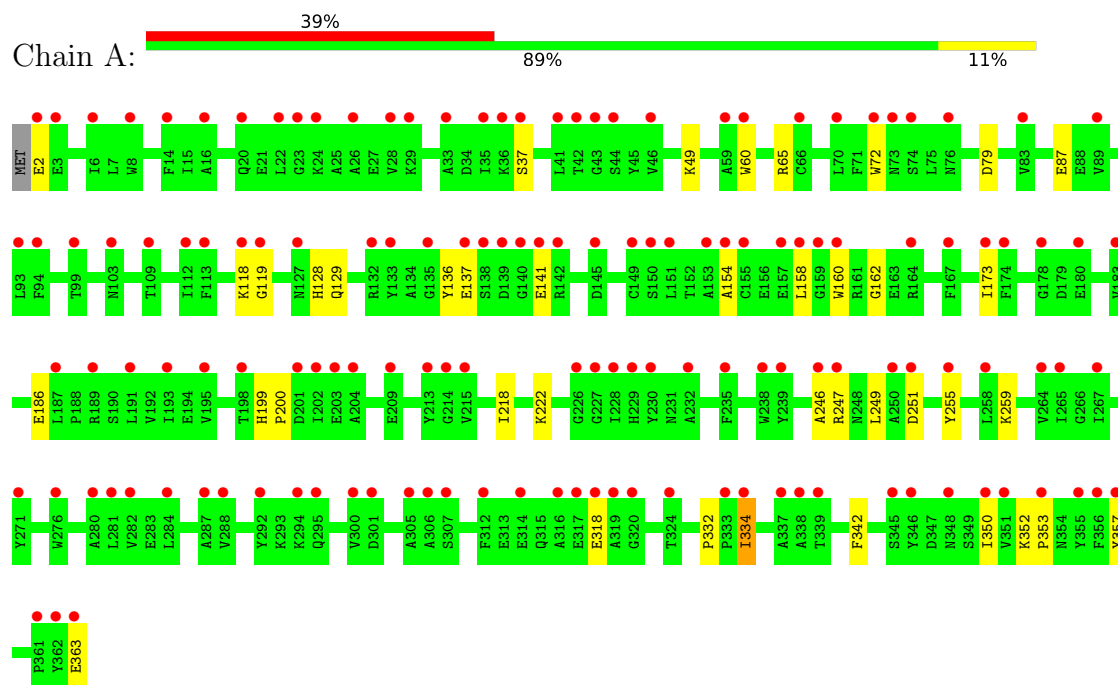
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	243	Total 243	O 243	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 OXYGENASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	80.55Å 95.19Å 62.77Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	37.09 – 1.89 37.09 – 1.89	Depositor EDS
% Data completeness (in resolution range)	80.8 (37.09-1.89) 80.8 (37.09-1.89)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.49 (at 1.89Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.187 , 0.228 0.188 , 0.229	Depositor DCC
R_{free} test set	1615 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	25.8	Xtriage
Anisotropy	0.019	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 42.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	3279	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.43% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, CL, HW1, H4B, 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	0.41	0/3017	0.78	2/4088 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	119	GLY	N-CA-C	-5.50	108.04	115.36
1	A	332	PRO	O-C-N	5.26	123.62	121.15

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2941	0	2856	16	2
2	A	43	0	30	4	0
3	A	17	0	15	2	1
4	A	1	0	0	0	0
5	A	28	0	33	3	0
6	A	6	0	8	0	0
7	A	243	0	0	0	0
All	All	3279	0	2942	20	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 3.

All (20) 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:218:ILE:HD11	5:A:904:HW1:H25	1.76	0.66
1:A:246:ALA:HB2	1:A:334:ILE:HD12	1.83	0.59
1:A:60:TRP:CE3	1:A:72:TRP:HA	2.40	0.56
1:A:247:ARG:NH1	1:A:251:ASP:OD2	2.43	0.52
2:A:901:HEM:O1A	3:A:902:H4B:N3	2.43	0.51
1:A:49:LYS:NZ	1:A:79:ASP:OD2	2.45	0.49
1:A:160:TRP:NE1	1:A:186:GLU:OE2	2.44	0.49
2:A:901:HEM:HBA1	5:A:904:HW1:H27	1.93	0.49
1:A:128:HIS:CE1	5:A:904:HW1:H18	2.49	0.48
1:A:154:ALA:O	1:A:158:LEU:HD12	2.15	0.47
1:A:65:ARG:NH1	1:A:357:TYR:OH	2.48	0.46
2:A:901:HEM:O1A	3:A:902:H4B:N2	2.45	0.46
1:A:136:TYR:OH	1:A:162:GLY:HA3	2.17	0.45
1:A:249:LEU:HA	1:A:255:TYR:HB2	1.98	0.45
2:A:901:HEM:HHC	2:A:901:HEM:HBB2	1.99	0.44
1:A:129:GLN:HG2	1:A:218:ILE:HD13	1.99	0.43
1:A:222:LYS:HE3	1:A:222:LYS:HB2	1.81	0.42
1:A:118:LYS:HE3	1:A:118:LYS:HB2	1.81	0.42
1:A:352:LYS:HA	1:A:353:PRO:C	2.45	0.41
1:A:199:HIS:CG	1:A:200:PRO:HD2	2.56	0.41

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)
1:A:342:PHE:O	3:A:902:H4B:O9[2_555]	2.00	0.20
1:A:87:GLU:OE2	1:A:259:LYS:NZ[3_556]	2.07	0.13

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	360/363 (99%)	350 (97%)	10 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	309/310 (100%)	300 (97%)	9 (3%)	37	20

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	GLU
1	A	37	SER
1	A	137	GLU
1	A	141	GLU
1	A	173	ILE
1	A	318	GLU
1	A	334	ILE
1	A	350	ILE
1	A	363	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	129	GLN
1	A	229	HIS
1	A	290	HIS

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 5 ligands modelled in this entry, 1 is monoatomic - leaving 4 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$
5	HW1	A	904	-	29,30,30	0.45	0	31,40,40	1.63	4 (12%)
2	HEM	A	901	1	50,50,50	1.53	5 (10%)	67,82,82	1.65	15 (22%)
6	GOL	A	905	-	5,5,5	0.21	0	5,5,5	0.48	0
3	H4B	A	902	-	17,18,18	1.80	3 (17%)	14,26,26	1.89	5 (35%)

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
5	HW1	A	904	-	-	4/13/23/23	0/3/3/3
2	HEM	A	901	1	-	2/14/54/54	-
6	GOL	A	905	-	-	0/4/4/4	-
3	H4B	A	902	-	-	3/8/17/17	0/2/2/2

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	902	H4B	C4A-C8A	5.68	1.48	1.38
2	A	901	HEM	FE-NB	5.19	2.10	1.94
2	A	901	HEM	FE-NC	4.18	2.09	1.95
2	A	901	HEM	C4D-ND	-3.12	1.34	1.40
2	A	901	HEM	C1B-NB	-3.07	1.34	1.40
2	A	901	HEM	C3B-C4B	2.67	1.50	1.44
3	A	902	H4B	C4-N3	-2.63	1.33	1.38
3	A	902	H4B	C4A-C4	2.24	1.48	1.41

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	904	HW1	C02-N01-C06	5.36	122.08	118.07
2	A	901	HEM	CHC-C4B-NB	4.78	129.57	124.42
3	A	902	H4B	C2-N1-C8A	4.57	121.42	113.36
2	A	901	HEM	CHD-C1D-ND	4.31	129.06	124.42
5	A	904	HW1	C22-N21-C26	4.18	121.20	118.07
2	A	901	HEM	CHA-C4D-ND	3.49	128.68	124.37
2	A	901	HEM	C1B-NB-C4B	3.27	109.08	105.21
5	A	904	HW1	C05-C06-N01	-3.03	119.28	122.73
2	A	901	HEM	C3B-C4B-NB	-2.89	107.39	109.47
2	A	901	HEM	CHB-C1B-NB	2.85	127.90	124.37
2	A	901	HEM	C4A-CHB-C1B	-2.72	119.85	126.25
5	A	904	HW1	C08-C06-N01	2.68	119.85	116.57
2	A	901	HEM	CHA-C4D-C3D	-2.65	120.34	125.23
2	A	901	HEM	C1A-CHA-C4D	-2.59	120.15	126.25
2	A	901	HEM	CHD-C1D-C2D	-2.58	120.96	125.03
3	A	902	H4B	C4-C4A-N5	2.52	123.13	116.27
3	A	902	H4B	O4-C4-C4A	-2.46	121.32	127.26
2	A	901	HEM	O2D-CGD-CBD	2.45	121.74	114.00
3	A	902	H4B	C4A-C4-N3	2.38	118.67	112.13
2	A	901	HEM	C4C-CHD-C1D	-2.25	121.23	126.02
3	A	902	H4B	O10-C10-C9	2.24	113.48	109.77
2	A	901	HEM	CHA-C1A-NA	2.20	127.84	123.86
2	A	901	HEM	C4C-NC-C1C	2.14	109.30	105.82
2	A	901	HEM	CBD-CAD-C3D	-2.07	106.80	112.53

There are no chirality outliers.

All (9) torsion outliers are listed below:

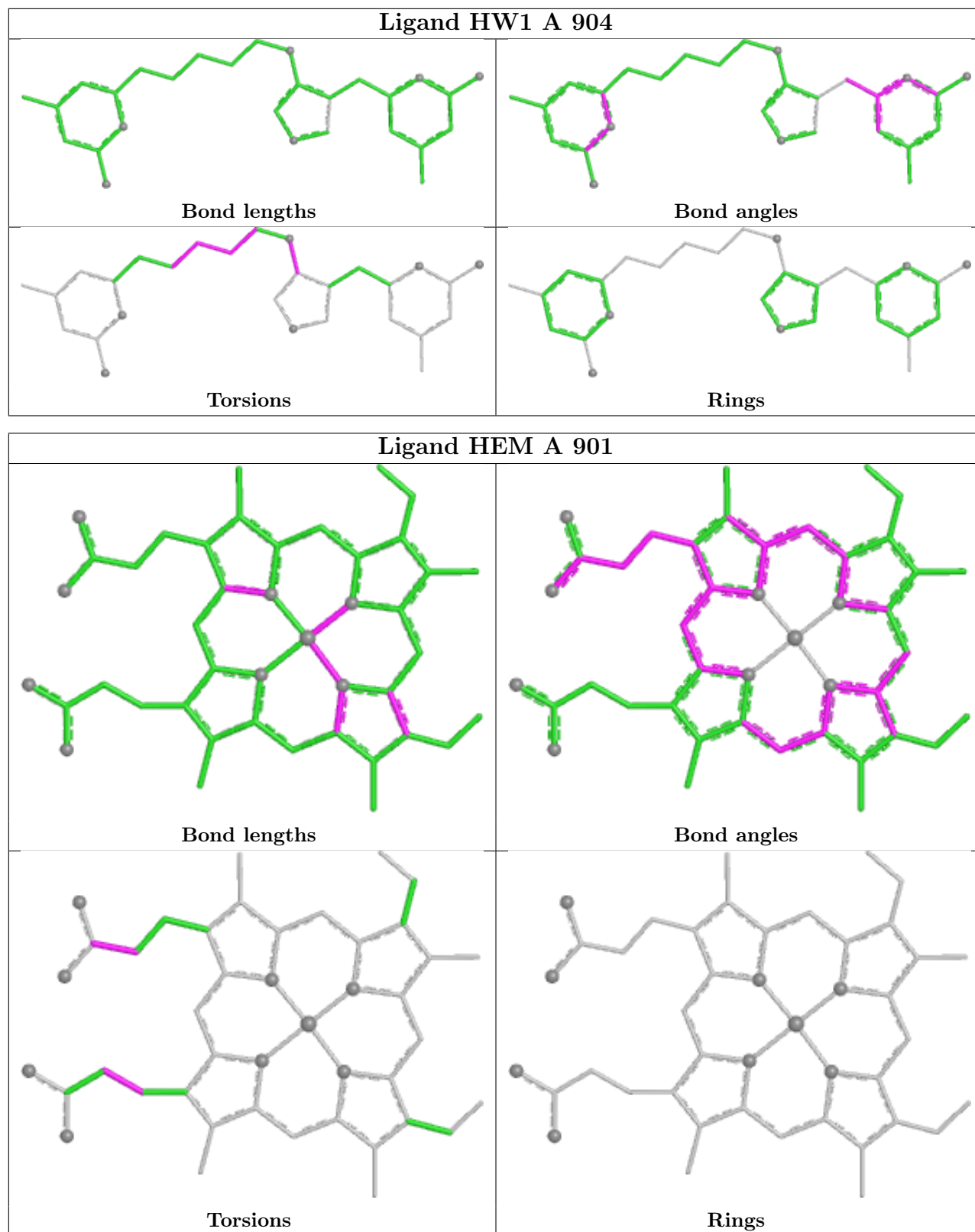
Mol	Chain	Res	Type	Atoms
3	A	902	H4B	O10-C10-C9-O9
5	A	904	HW1	C10-C11-C12-C13
5	A	904	HW1	C5'-C4'-O09-C10
5	A	904	HW1	O09-C10-C11-C12
3	A	902	H4B	C11-C10-C9-C6
3	A	902	H4B	C11-C10-C9-O9
2	A	901	HEM	C2A-CAA-CBA-CGA
5	A	904	HW1	C11-C12-C13-C14
2	A	901	HEM	CAD-CBD-CGD-O2D

There are no ring outliers.

3 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	904	HW1	3	0
2	A	901	HEM	4	0
3	A	902	H4B	2	1

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 ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	362/363 (99%)	1.86	142 (39%) 1 0	21, 33, 67, 92	0

All (142) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	138	SER	4.5
1	A	334	ILE	4.0
1	A	362	TYR	4.0
1	A	158	LEU	3.8
1	A	26	ALA	3.8
1	A	119	GLY	3.6
1	A	8	TRP	3.6
1	A	118	LYS	3.5
1	A	356	PHE	3.5
1	A	193	ILE	3.5
1	A	23	GLY	3.4
1	A	137	GLU	3.4
1	A	247	ARG	3.4
1	A	41	LEU	3.4
1	A	140	GLY	3.3
1	A	350	ILE	3.3
1	A	89	VAL	3.3
1	A	232	ALA	3.3
1	A	307	SER	3.2
1	A	288	VAL	3.2
1	A	141	GLU	3.2
1	A	250	ALA	3.2
1	A	355	TYR	3.1
1	A	43	GLY	3.1
1	A	305	ALA	3.1
1	A	295	GLN	3.0
1	A	180	GLU	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	363	GLU	3.0
1	A	142	ARG	2.9
1	A	312	PHE	2.9
1	A	33	ALA	2.9
1	A	246	ALA	2.9
1	A	160	TRP	2.9
1	A	150	SER	2.9
1	A	280	ALA	2.8
1	A	316	ALA	2.8
1	A	187	LEU	2.8
1	A	284	LEU	2.8
1	A	173	ILE	2.8
1	A	99	THR	2.8
1	A	300	VAL	2.8
1	A	201	ASP	2.8
1	A	154	ALA	2.8
1	A	133	TYR	2.7
1	A	59	ALA	2.7
1	A	76	ASN	2.7
1	A	153	ALA	2.7
1	A	28	VAL	2.7
1	A	346	TYR	2.7
1	A	112	ILE	2.7
1	A	46	VAL	2.7
1	A	159	GLY	2.7
1	A	238	TRP	2.6
1	A	155	CYS	2.6
1	A	178	GLY	2.6
1	A	60	TRP	2.6
1	A	132	ARG	2.6
1	A	251	ASP	2.6
1	A	351	VAL	2.6
1	A	357	TYR	2.6
1	A	72	TRP	2.6
1	A	281	LEU	2.6
1	A	20	GLN	2.6
1	A	228	ILE	2.6
1	A	36	LYS	2.5
1	A	214	GLY	2.5
1	A	320	GLY	2.5
1	A	337	ALA	2.5
1	A	42	THR	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	70	LEU	2.5
1	A	93	LEU	2.5
1	A	324	THR	2.5
1	A	157	GLU	2.5
1	A	306	ALA	2.5
1	A	29	LYS	2.5
1	A	294	LYS	2.4
1	A	139	ASP	2.4
1	A	109	THR	2.4
1	A	151	LEU	2.4
1	A	235	PHE	2.4
1	A	301	ASP	2.4
1	A	203	GLU	2.4
1	A	44	SER	2.4
1	A	239	TYR	2.3
1	A	167	PHE	2.3
1	A	35	ILE	2.3
1	A	3	GLU	2.3
1	A	94	PHE	2.3
1	A	174	PHE	2.3
1	A	16	ALA	2.3
1	A	204	ALA	2.3
1	A	333	PRO	2.3
1	A	282	VAL	2.3
1	A	2	GLU	2.3
1	A	276	TRP	2.3
1	A	189	ARG	2.3
1	A	345	SER	2.3
1	A	317	GLU	2.3
1	A	149	CYS	2.3
1	A	24	LYS	2.3
1	A	191	LEU	2.3
1	A	135	GLY	2.2
1	A	267	ILE	2.2
1	A	215	VAL	2.2
1	A	264	VAL	2.2
1	A	213	TYR	2.2
1	A	271	TYR	2.2
1	A	202	ILE	2.2
1	A	22	LEU	2.2
1	A	14	PHE	2.2
1	A	195	VAL	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	339	THR	2.2
1	A	74	SER	2.2
1	A	361	PRO	2.2
1	A	319	ALA	2.2
1	A	338	ALA	2.2
1	A	292	TYR	2.1
1	A	145	ASP	2.1
1	A	83	VAL	2.1
1	A	227	GLY	2.1
1	A	66	CYS	2.1
1	A	258	LEU	2.1
1	A	103	ASN	2.1
1	A	198	THR	2.1
1	A	287	ALA	2.1
1	A	209	GLU	2.1
1	A	314	GLU	2.1
1	A	226	GLY	2.1
1	A	73	ASN	2.1
1	A	183	VAL	2.1
1	A	6	ILE	2.0
1	A	113	PHE	2.0
1	A	37	SER	2.0
1	A	127	ASN	2.0
1	A	348	ASN	2.0
1	A	318	GLU	2.0
1	A	229	HIS	2.0
1	A	265	ILE	2.0
1	A	230	TYR	2.0
1	A	255	TYR	2.0
1	A	353	PRO	2.0
1	A	164	ARG	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](#)

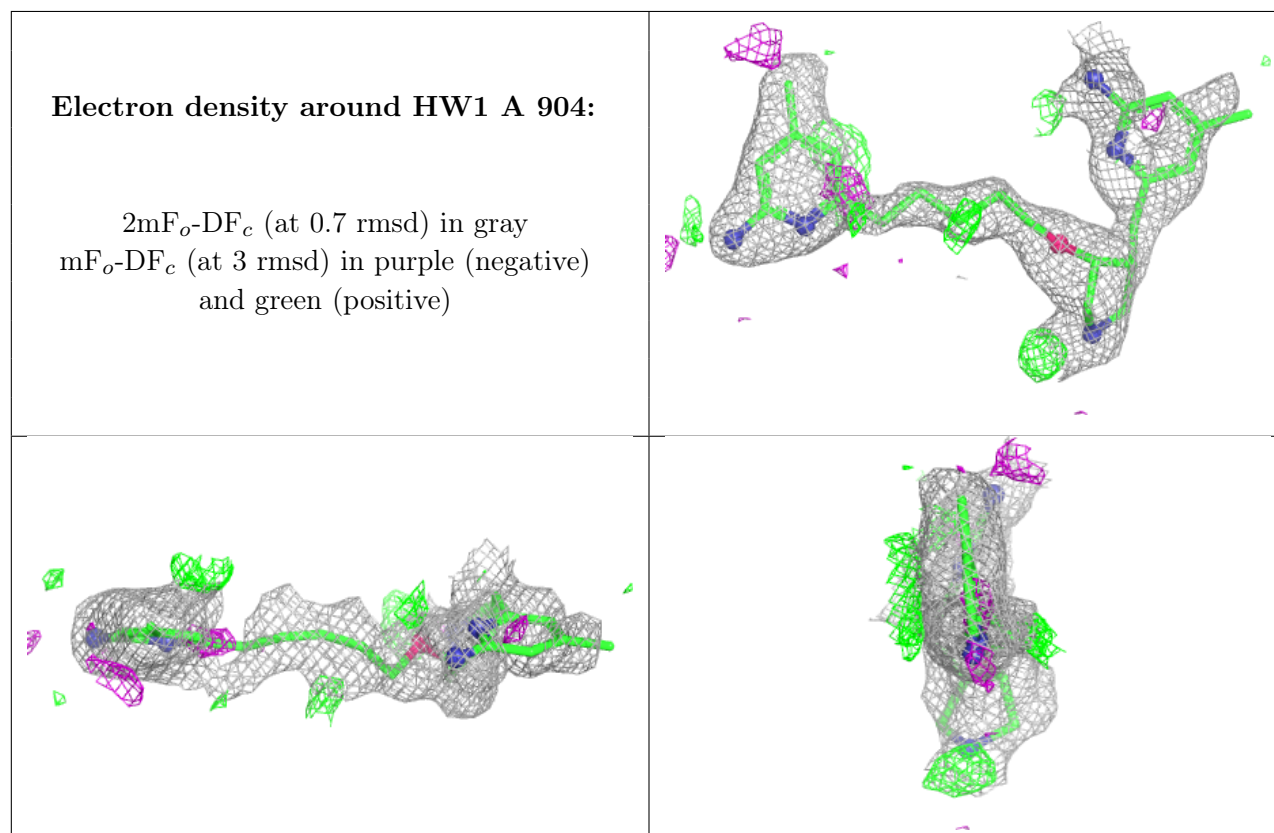
There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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.

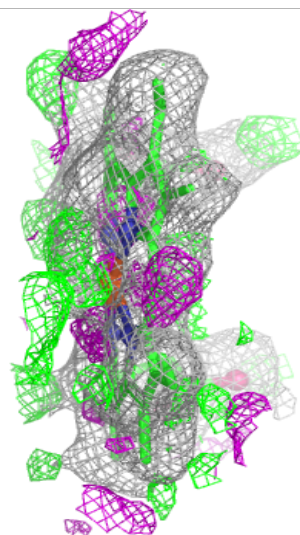
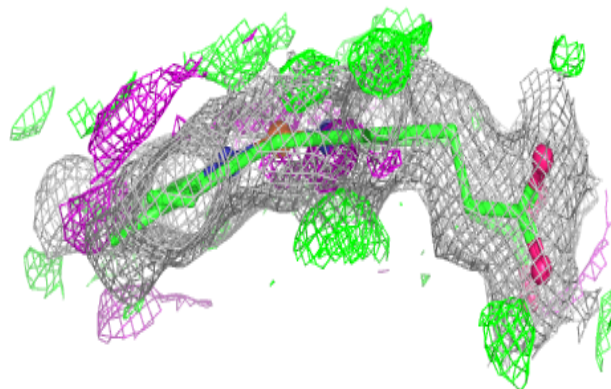
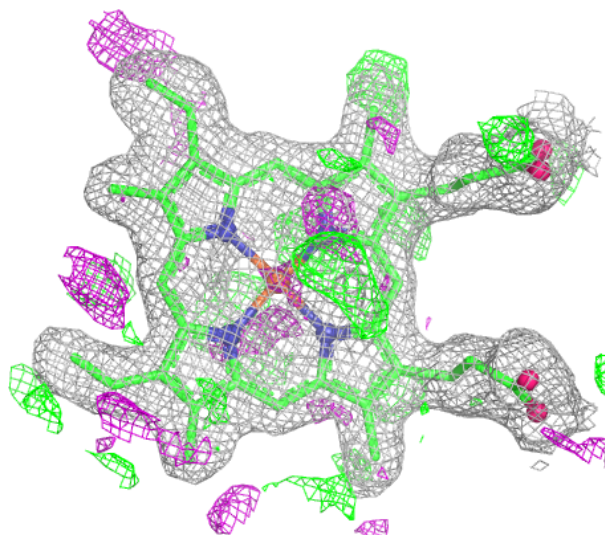
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	GOL	A	905	6/6	0.68	0.25	53,55,58,60	0
3	H4B	A	902	17/17	0.76	0.23	41,50,76,77	0
5	HW1	A	904	28/28	0.78	0.21	24,50,67,68	0
2	HEM	A	901	43/43	0.91	0.13	19,25,40,45	0
4	CL	A	903	1/1	0.97	0.06	30,30,30,30	0

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 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)



6.5 Other polymers [i](#)

There are no such residues in this entry.