



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 10:50 AM GMT

PDB ID : 3N63
Title : Structure of neuronal nitric oxide synthase D597N/M336V mutant heme domain in complex with 6,6'-(2,2'-(5-amino-1,3-phenylene)bis(ethane-2,1-diyl))bis(4-methylpyridin-2-amine)
Authors : Li, H.; Delker, S.L.; Poulos, T.L.
Deposited on : 2010-05-25
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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

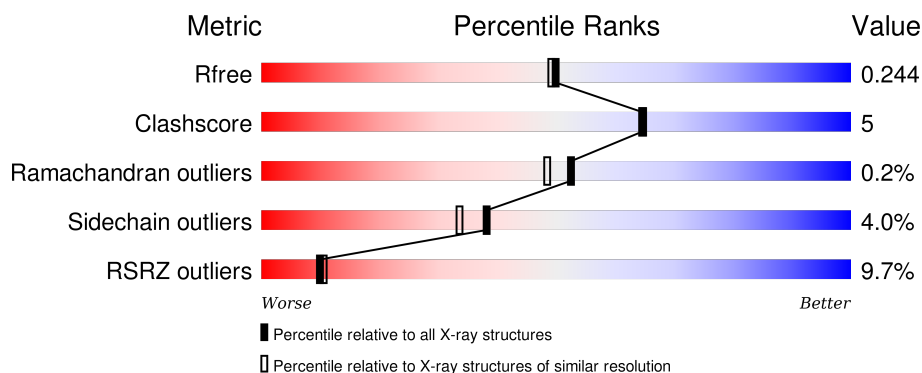
1 Overall quality at a glance

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}	91344	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	422	<div> <div>12%</div> <div>84%</div> <div>10%</div> <div>• •</div> </div>
1	B	422	<div> <div>6%</div> <div>88%</div> <div>9%</div> <div>•</div> </div>

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
3	XFN	A	800	-	-	-	X
4	ACT	A	860	-	-	-	X
4	ACT	B	860	-	-	-	X

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 7089 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.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	407	Total	C	N	O	S	0	2	0
			3318	2125	567	605	21			
1	B	411	Total	C	N	O	S	0	1	0
			3347	2142	575	609	21			

There are 4 discrepancies between the modelled and reference sequences:

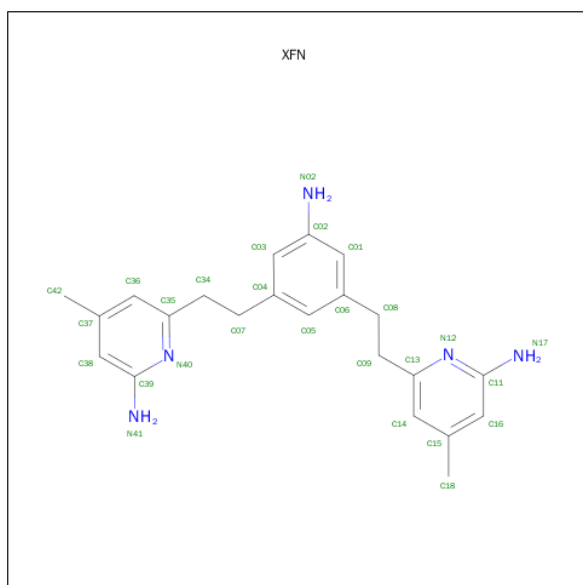
Chain	Residue	Modelled	Actual	Comment	Reference
A	336	VAL	MET	ENGINEERED MUTATION	UNP P29476
A	597	ASN	ASP	ENGINEERED MUTATION	UNP P29476
B	336	VAL	MET	ENGINEERED MUTATION	UNP P29476
B	597	ASN	ASP	ENGINEERED MUTATION	UNP P29476

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



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

- Molecule 3 is 6,6'-[(5-AMINOBENZENE-1,3-DIYL)DIETHANE-2,1-DIYL]BIS(4-METHYL PYRIDIN-2-AMINE) (three-letter code: XFN) (formula: C₂₂H₂₇N₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	N		
			27	22	5		
3	B	1	Total	C	N		
			27	22	5		

- Molecule 4 is ACETATE ION (three-letter code: ACT) (formula: C₂H₃O₂).



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

- Molecule 5 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	3	Total	Zn	0	0
			3	3		

- Molecule 6 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	2	Total	Cl	0	0
			2	2		
6	A	2	Total	Cl	0	0
			2	2		

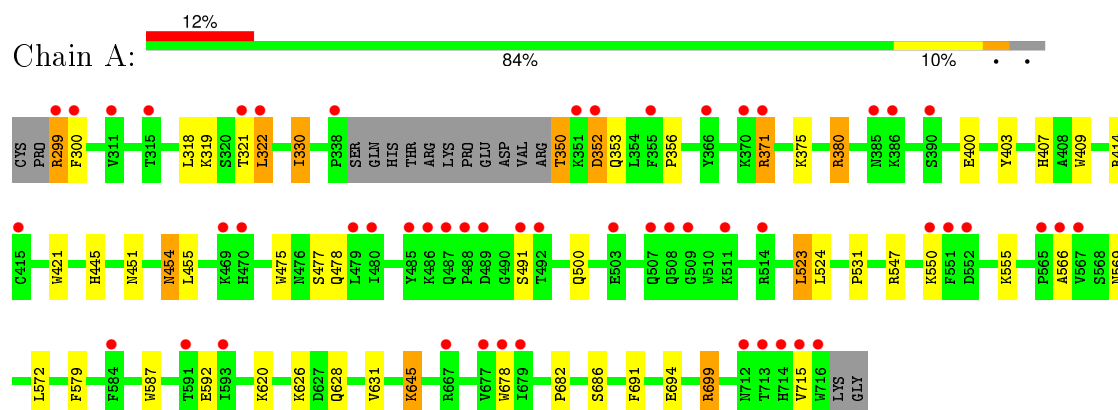
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	113	Total	O	0	0
			113	113		
7	B	156	Total	O	0	0
			156	156		

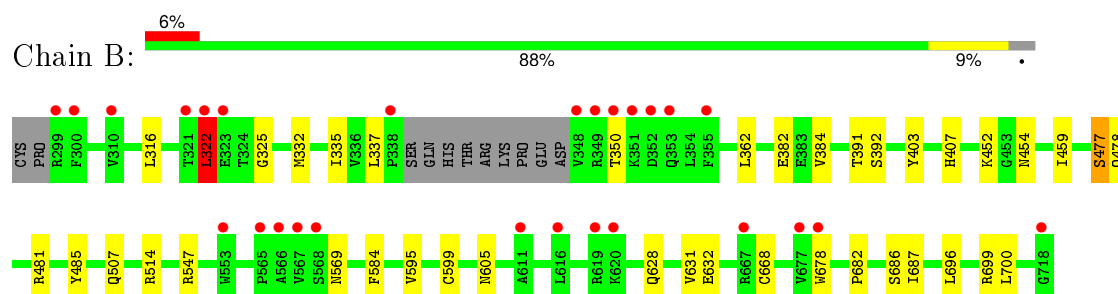
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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 ($\text{RSRZ} > 2$). 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



• Molecule 1: Nitric oxide synthase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	51.80Å 110.98Å 164.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.06 – 2.00 40.74 – 2.00	Depositor EDS
% Data completeness (in resolution range)	96.9 (41.06-2.00) 96.9 (40.74-2.00)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.28 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.5.0089	Depositor
R, R_{free}	0.195 , 0.236 0.206 , 0.244	Depositor DCC
R_{free} test set	3093 reflections (5.19%)	DCC
Wilson B-factor (Å ²)	37.5	Xtriage
Anisotropy	0.162	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 41.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 62725 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7089	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.12% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: ACT, HEM, ZN, CL, XFN

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	0.61	0/3417	0.65	0/4637
1	B	0.65	0/3443	0.66	0/4669
All	All	0.63	0/6860	0.66	0/9306

There are no bond length outliers.

There are no bond angle outliers.

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	3318	0	3233	38	0
1	B	3347	0	3266	30	0
2	A	43	0	30	4	0
2	B	43	0	30	5	0
3	A	27	0	27	2	0
3	B	27	0	27	1	0
4	A	4	0	3	0	0
4	B	4	0	3	1	0
5	A	3	0	0	0	0
6	A	2	0	0	0	0
6	B	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	A	113	0	0	2	0
7	B	156	0	0	7	0
All	All	7089	0	6619	68	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 5.

All (68) 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:371:ARG:CG	1:A:371:ARG:HH21	1.96	0.79
2:B:750:HEM:HH2	2:B:750:HEM:HBB2	1.64	0.79
1:A:523:LEU:HD22	1:A:531:PRO:HB2	1.73	0.70
1:A:371:ARG:HG3	1:A:371:ARG:HH21	1.58	0.68
1:A:699:ARG:HD2	7:A:1035:HOH:O	1.93	0.66
2:A:750:HEM:CMC	2:A:750:HEM:HBC2	2.26	0.66
1:B:382:GLU:HG3	7:B:1216:HOH:O	1.99	0.61
2:A:750:HEM:HMC1	2:A:750:HEM:HBC2	1.82	0.61
1:A:631:VAL:HG11	1:B:628:GLN:HG3	1.85	0.57
1:A:353:GLN:O	1:A:356:PRO:HD2	2.04	0.57
1:A:330:ILE:HD11	1:B:696:LEU:HD22	1.85	0.56
1:B:322:LEU:HB2	1:B:699:ARG:HE	1.70	0.56
1:B:452:LYS:HE3	7:B:1180:HOH:O	2.06	0.56
1:A:682:PRO:HB2	1:B:686:SER:HB2	1.88	0.56
1:A:350:THR:N	1:A:353:GLN:NE2	2.53	0.55
1:B:678:TRP:CZ3	3:B:800:XFN:H08	2.42	0.55
1:B:699:ARG:NH1	7:B:1224:HOH:O	2.40	0.55
1:B:699:ARG:CZ	7:B:1224:HOH:O	2.56	0.54
1:A:491:SER:HB2	7:A:1107:HOH:O	2.09	0.52
1:A:686:SER:HB2	1:B:682:PRO:HB2	1.90	0.52
1:A:380:ARG:HD3	1:A:400:GLU:OE1	2.10	0.52
1:A:321:THR:HG23	1:A:322:LEU:H	1.75	0.52
1:A:592:GLU:OE1	3:A:800:XFN:N40	2.43	0.51
1:A:475:TRP:HB2	1:A:523:LEU:HB3	1.93	0.50
1:A:628:GLN:NE2	1:B:632:GLU:OE2	2.45	0.50
1:A:455:LEU:HD12	1:A:587:TRP:HB3	1.93	0.50
1:A:371:ARG:CG	1:A:371:ARG:NH2	2.62	0.50
1:A:478:GLN:HA	1:A:566:ALA:O	2.12	0.49
1:A:628:GLN:HG2	1:B:631:VAL:HG11	1.94	0.49
1:A:352:ASP:N	1:A:352:ASP:OD2	2.46	0.49
1:A:299:ARG:HH11	1:A:318:LEU:HD22	1.79	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:595:VAL:O	1:B:599:CYS:HB2	2.14	0.47
1:B:403:TYR:CE1	1:B:407:HIS:CE1	3.02	0.47
1:A:475:TRP:CZ2	1:A:531:PRO:HG3	2.50	0.47
2:B:750:HEM:HMB3	4:B:860:ACT:H2	1.97	0.46
2:A:750:HEM:HMC1	2:A:750:HEM:CBC	2.44	0.46
1:A:371:ARG:HG2	1:A:371:ARG:HH21	1.78	0.46
1:B:325:GLY:O	1:B:332:MET:HG3	2.16	0.46
1:B:605:ASN:ND2	7:B:1121:HOH:O	2.44	0.46
1:B:485:TYR:CZ	1:B:514:ARG:HA	2.52	0.45
1:B:322:LEU:HD12	1:B:699:ARG:HB3	1.98	0.45
1:A:371:ARG:HG2	1:A:371:ARG:NH2	2.31	0.45
1:B:485:TYR:CE1	1:B:514:ARG:HA	2.51	0.44
1:A:403:TYR:CE1	1:A:407:HIS:CE1	3.05	0.44
2:B:750:HEM:HBC2	2:B:750:HEM:CMC	2.48	0.44
1:A:626:LYS:HB3	1:B:687:ILE:HD12	1.99	0.44
1:A:350:THR:N	1:A:353:GLN:HE21	2.16	0.43
1:A:686:SER:HA	1:A:691:PHE:CG	2.53	0.43
2:A:750:HEM:HMA2	2:A:750:HEM:CGA	2.48	0.43
2:B:750:HEM:HBC2	2:B:750:HEM:HMC1	2.01	0.42
1:B:316:LEU:HD21	1:B:700:LEU:HD11	2.01	0.42
1:B:459:ILE:HD12	1:B:584:PHE:HB3	2.01	0.42
1:A:524:LEU:O	1:A:531:PRO:HA	2.19	0.42
1:A:694:GLU:HB3	1:B:335:ILE:HD13	2.02	0.42
1:B:391:THR:O	1:B:392:SER:HB2	2.20	0.42
1:B:699:ARG:CD	7:B:1224:HOH:O	2.68	0.42
1:B:478:GLN:HB2	1:B:481:ARG:HG3	2.01	0.42
1:B:477:SER:HB3	1:B:569:ASN:ND2	2.34	0.42
1:A:645:LYS:HB2	1:A:645:LYS:HE3	1.91	0.41
1:B:699:ARG:HG2	7:B:1135:HOH:O	2.20	0.41
1:B:362:LEU:HD11	1:B:384:VAL:HG21	2.02	0.41
1:A:678:TRP:CZ3	3:A:800:XFN:H08	2.55	0.41
1:A:445:HIS:C	1:A:445:HIS:CD2	2.94	0.41
1:A:409:TRP:CE3	1:A:421:TRP:HA	2.56	0.41
1:B:584:PHE:CD1	2:B:750:HEM:CAC	3.04	0.41
1:A:451:ASN:HB3	1:A:454:ASN:O	2.21	0.41
1:A:414:ARG:HD3	1:A:678:TRP:CD2	2.56	0.40
1:A:572:LEU:HB3	1:A:579:PHE:HB2	2.03	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	405/422 (96%)	392 (97%)	12 (3%)	1 (0%)	52	48
1	B	408/422 (97%)	398 (98%)	9 (2%)	1 (0%)	52	48
All	All	813/844 (96%)	790 (97%)	21 (3%)	2 (0%)	52	48

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	322	LEU
1	B	322	LEU

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	365/377 (97%)	344 (94%)	21 (6%)	25	19
1	B	367/377 (97%)	358 (98%)	9 (2%)	55	55
All	All	732/754 (97%)	702 (96%)	30 (4%)	38	32

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	299	ARG
1	A	300	PHE
1	A	319	LYS
1	A	330	ILE

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Mol	Chain	Res	Type
1	A	350	THR
1	A	352	ASP
1	A	371	ARG
1	A	375	LYS
1	A	380	ARG
1	A	454	ASN
1	A	477	SER
1	A	500	GLN
1	A	523	LEU
1	A	547	ARG
1	A	550	LYS
1	A	555	LYS
1	A	569	ASN
1	A	620	LYS
1	A	645	LYS
1	A	699	ARG
1	A	715	VAL
1	B	322	LEU
1	B	337	LEU
1	B	350	THR
1	B	454	ASN
1	B	477	SER
1	B	507	GLN
1	B	547	ARG
1	B	668[A]	CYS
1	B	668[B]	CYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	353	GLN
1	A	454	ASN
1	A	527	ASN
1	A	569	ASN
1	A	605	ASN
1	A	628	GLN
1	A	642	GLN
1	A	697	ASN
1	B	364	GLN
1	B	454	ASN
1	B	507	GLN
1	B	508	GLN

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Mol	Chain	Res	Type
1	B	527	ASN
1	B	605	ASN
1	B	697	ASN

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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 13 ligands modelled in this entry, 7 are 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	$\# Z > 2$
2	HEM	A	750	1	30,50,50	2.24	10 (33%)	24,82,82	2.84	13 (54%)
3	XFN	A	800	-	29,29,29	0.96	1 (3%)	40,40,40	2.16	14 (35%)
4	ACT	A	860	-	1,3,3	1.37	0	0,3,3	0.00	-
2	HEM	B	750	1	30,50,50	1.99	9 (30%)	24,82,82	2.80	14 (58%)
3	XFN	B	800	-	29,29,29	1.00	1 (3%)	40,40,40	2.07	12 (30%)
4	ACT	B	860	-	1,3,3	1.61	0	0,3,3	0.00	-

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
2	HEM	A	750	1	-	0/10/54/54	0/0/8/8
3	XFN	A	800	-	-	0/10/10/10	0/3/3/3
4	ACT	A	860	-	-	0/0/0/0	0/0/0/0
2	HEM	B	750	1	-	0/10/54/54	0/0/8/8
3	XFN	B	800	-	-	0/10/10/10	0/3/3/3
4	ACT	B	860	-	-	0/0/0/0	0/0/0/0

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	750	HEM	C3D-C4D	-6.11	1.43	1.51
2	A	750	HEM	C3B-C4B	-5.59	1.46	1.51
2	B	750	HEM	C3D-C4D	-5.53	1.44	1.51
2	B	750	HEM	C3B-C4B	-5.33	1.47	1.51
2	A	750	HEM	C2C-C1C	-3.81	1.45	1.52
2	B	750	HEM	C2C-C1C	-3.34	1.46	1.52
2	A	750	HEM	C2B-C1B	-2.19	1.44	1.51
2	B	750	HEM	C2D-C1D	-2.10	1.45	1.51
2	B	750	HEM	C2B-C1B	-2.07	1.45	1.51
2	A	750	HEM	C2D-C1D	-2.00	1.45	1.51
3	B	800	XFN	C35-N40	2.00	1.38	1.34
2	B	750	HEM	FE-NC	2.06	2.04	1.95
2	B	750	HEM	C3C-CAC	2.07	1.55	1.51
2	A	750	HEM	CMA-C3A	2.11	1.56	1.51
2	B	750	HEM	CAA-C2A	2.30	1.56	1.52
3	A	800	XFN	C05-C04	2.39	1.43	1.39
2	B	750	HEM	CMA-C3A	2.44	1.56	1.51
2	A	750	HEM	C1C-NC	2.54	1.39	1.36
2	A	750	HEM	CAA-C2A	2.89	1.57	1.52
2	A	750	HEM	FE-NB	2.96	2.13	1.97
2	A	750	HEM	FE-NC	3.14	2.08	1.95

All (53) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	750	HEM	CAA-C2A-C1A	-5.90	120.60	127.01
2	A	750	HEM	CAA-C2A-C1A	-5.29	121.26	127.01
3	A	800	XFN	C14-C13-N12	-4.21	118.21	122.96
2	A	750	HEM	C3C-CAC-CBC	-4.06	118.23	124.46
3	B	800	XFN	C14-C13-N12	-3.93	118.53	122.96
3	B	800	XFN	C36-C35-N40	-3.79	118.68	122.96

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	750	HEM	CMA-C3A-C4A	-3.54	122.51	128.36
3	A	800	XFN	C36-C35-N40	-3.51	119.00	122.96
2	B	750	HEM	CMA-C3A-C4A	-3.47	122.62	128.36
3	A	800	XFN	C18-C15-C16	-3.17	116.13	120.95
3	A	800	XFN	C15-C14-C13	-2.83	118.51	120.28
2	B	750	HEM	CBD-CAD-C3D	-2.82	105.35	113.55
2	A	750	HEM	CBD-CAD-C3D	-2.62	105.92	113.55
3	B	800	XFN	C07-C04-C03	-2.44	116.47	120.56
2	B	750	HEM	C3C-CAC-CBC	-2.42	120.74	124.46
2	A	750	HEM	C3B-C4B-NB	-2.34	107.16	111.63
3	A	800	XFN	C07-C04-C03	-2.14	116.97	120.56
3	B	800	XFN	C16-C15-C14	2.03	120.67	118.08
3	A	800	XFN	N17-C11-N12	2.14	120.40	116.50
3	A	800	XFN	C07-C04-C05	2.16	124.18	120.56
2	B	750	HEM	CBA-CAA-C2A	2.21	116.49	112.53
2	A	750	HEM	CBA-CAA-C2A	2.22	116.52	112.53
3	B	800	XFN	C09-C13-C14	2.23	124.26	121.13
2	B	750	HEM	CMA-C3A-C2A	2.36	130.17	125.24
3	B	800	XFN	N17-C11-N12	2.40	120.88	116.50
2	B	750	HEM	C3B-C4B-CHC	2.40	126.55	123.16
3	A	800	XFN	C08-C09-C13	2.54	117.42	112.53
2	A	750	HEM	CMD-C2D-C3D	2.55	125.64	114.35
3	B	800	XFN	C07-C04-C05	2.65	125.00	120.56
2	B	750	HEM	C2C-C1C-CHC	2.69	127.77	123.68
2	B	750	HEM	C2D-C3D-C4D	2.70	106.07	101.50
3	A	800	XFN	C16-C15-C14	2.71	121.54	118.08
3	A	800	XFN	N41-C39-N40	2.75	121.52	116.50
2	B	750	HEM	CMD-C2D-C3D	2.79	126.69	114.35
3	B	800	XFN	C08-C09-C13	2.80	117.91	112.53
3	B	800	XFN	N41-C39-N40	2.86	121.72	116.50
2	A	750	HEM	C2D-C3D-C4D	2.97	106.54	101.50
3	A	800	XFN	C09-C13-C14	3.20	125.62	121.13
3	A	800	XFN	C34-C35-N40	3.75	121.22	115.69
2	B	750	HEM	CAD-C3D-C4D	3.87	126.11	112.47
2	A	750	HEM	CAD-C3D-C4D	3.89	126.19	112.47
2	A	750	HEM	C3B-C4B-CHC	3.91	128.67	123.16
3	B	800	XFN	C39-N40-C35	4.06	121.11	118.23
2	A	750	HEM	CMB-C2B-C3B	4.17	126.93	116.53
2	A	750	HEM	CMC-C2C-C3C	4.36	127.41	116.53
2	B	750	HEM	CMB-C2B-C3B	4.36	127.41	116.53
2	B	750	HEM	CMC-C2C-C3C	4.37	127.44	116.53
3	B	800	XFN	C34-C35-N40	4.62	122.51	115.69

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	800	XFN	C39-N40-C35	4.85	121.67	118.23
2	A	750	HEM	CAD-C3D-C2D	4.87	127.21	113.22
2	B	750	HEM	CAD-C3D-C2D	5.04	127.70	113.22
3	B	800	XFN	C11-N12-C13	6.08	122.55	118.23
3	A	800	XFN	C11-N12-C13	6.75	123.03	118.23

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	750	HEM	4	0
3	A	800	XFN	2	0
2	B	750	HEM	5	0
3	B	800	XFN	1	0
4	B	860	ACT	1	0

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	407/422 (96%)	0.73	52 (12%) 5 5	27, 51, 87, 110	0
1	B	411/422 (97%)	0.35	27 (6%) 22 22	27, 43, 67, 89	0
All	All	818/844 (96%)	0.54	79 (9%) 10 10	27, 46, 82, 110	0

All (79) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	716	TRP	6.9
1	B	718	GLY	6.6
1	B	300	PHE	6.6
1	A	715	VAL	5.6
1	B	348	VAL	5.1
1	A	300	PHE	5.0
1	B	619	ARG	4.7
1	B	352	ASP	4.5
1	B	351	LYS	4.1
1	B	350	THR	4.1
1	A	486	LYS	4.0
1	A	355	PHE	3.9
1	A	713	THR	3.9
1	A	299	ARG	3.8
1	A	488	PRO	3.6
1	A	352	ASP	3.6
1	A	489	ASP	3.4
1	B	322	LEU	3.4
1	A	667	ARG	3.4
1	A	514	ARG	3.3
1	B	321	THR	3.1
1	A	371	ARG	3.1
1	A	567	VAL	3.1
1	B	611	ALA	3.0

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Mol	Chain	Res	Type	RSRZ
1	B	667	ARG	3.0
1	A	511	LYS	3.0
1	A	678	TRP	2.9
1	A	322	LEU	2.9
1	A	509	GLY	2.9
1	B	355	PHE	2.8
1	B	567	VAL	2.8
1	A	714	HIS	2.8
1	A	566	ALA	2.8
1	A	315	THR	2.8
1	A	487	GLN	2.8
1	A	479	LEU	2.8
1	A	351	LYS	2.7
1	A	507	GLN	2.7
1	A	712	ASN	2.7
1	A	386	LYS	2.7
1	A	311	VAL	2.6
1	A	677	VAL	2.6
1	B	566	ALA	2.6
1	A	551	PHE	2.5
1	B	616	LEU	2.5
1	A	366	TYR	2.5
1	A	584	PHE	2.5
1	A	593	ILE	2.5
1	A	485	TYR	2.5
1	A	491	SER	2.5
1	B	553	TRP	2.5
1	A	338	PRO	2.4
1	A	415	CYS	2.4
1	A	508	GLN	2.4
1	B	338	PRO	2.4
1	B	678	TRP	2.4
1	A	480	ILE	2.4
1	B	620	LYS	2.4
1	A	503	GLU	2.3
1	A	321	THR	2.3
1	A	565	PRO	2.3
1	B	677	VAL	2.3
1	A	591	THR	2.2
1	A	390	SER	2.2
1	A	470	HIS	2.2
1	A	370	LYS	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	492	THR	2.2
1	B	568	SER	2.1
1	B	565	PRO	2.1
1	B	299	ARG	2.1
1	B	310	VAL	2.1
1	A	385	ASN	2.1
1	B	323	GLU	2.1
1	A	550	LYS	2.1
1	B	349	ARG	2.1
1	B	353	GLN	2.1
1	A	469	LYS	2.0
1	A	552	ASP	2.0
1	A	679	ILE	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
4	ACT	A	860	4/4	0.70	0.39	7.57	80,82,82,82	0
4	ACT	B	860	4/4	0.94	0.24	7.20	56,56,58,59	0
3	XFN	A	800	27/27	0.84	0.32	3.03	36,52,56,57	0
6	CL	A	930	1/1	0.98	0.21	1.38	48,48,48,48	0
3	XFN	B	800	27/27	0.89	0.21	0.95	35,45,52,52	0
2	HEM	B	750	43/43	0.97	0.17	0.77	30,35,47,50	0
2	HEM	A	750	43/43	0.97	0.20	0.60	31,34,49,55	0
5	ZN	A	900	1/1	0.99	0.10	-1.13	38,38,38,38	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
6	CL	B	930	1/1	0.99	0.12	-1.41	43,43,43,43	0
6	CL	A	910	1/1	0.98	0.08	-	42,42,42,42	0
5	ZN	A	719	1/1	0.99	0.10	-	36,36,36,36	1
6	CL	B	910	1/1	0.97	0.11	-	45,45,45,45	0
5	ZN	A	901	1/1	0.99	0.11	-	35,35,35,35	1

6.5 Other polymers [i](#)

There are no such residues in this entry.