



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 10:50 AM GMT

PDB ID : 3N69  
Title : Structure of endothelial nitric oxide synthase N368D/V106M double mutant heme domain complexed with 6,6'-(2,2'-(pyridine-3,5-diyl)bis(ethane-2,1-diyl))bis(4-methylpyridin-2-amine)  
Authors : Delker, S.L.; Li, H.; Poulos, T.L.  
Deposited on : 2010-05-25  
Resolution : 2.65 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](mailto: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.

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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

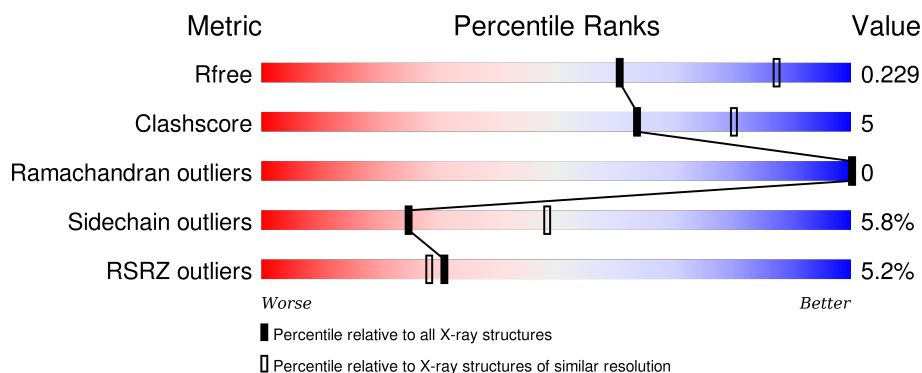
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.65 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	3152 (2.70-2.62)
Clashscore	102246	3524 (2.70-2.62)
Ramachandran outliers	100387	3469 (2.70-2.62)
Sidechain outliers	100360	3469 (2.70-2.62)
RSRZ outliers	91569	3161 (2.70-2.62)

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  $\geq 3$ , 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	444	
1	B	444	

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	ACT	B	860	-	-	-	X
4	CAD	A	950	-	-	-	X

## 2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 7116 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	404	Total	C	N	O	S	0	3	0
			3240	2057	572	594	17			
1	B	403	Total	C	N	O	S	0	4	0
			3249	2061	576	595	17			

There are 6 discrepancies between the modelled and reference sequences:

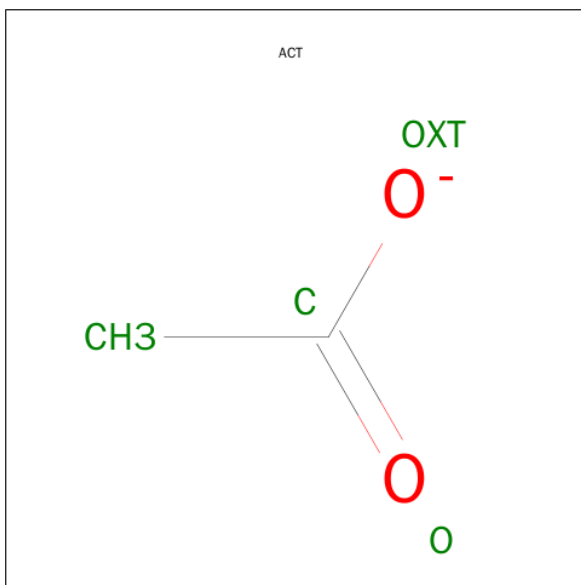
Chain	Residue	Modelled	Actual	Comment	Reference
A	100	ARG	CYS	SEE REMARK 999	UNP P29476
A	106	MET	VAL	ENGINEERED MUTATION	UNP P29476
A	368	ASP	ASN	ENGINEERED MUTATION	UNP P29476
B	100	ARG	CYS	SEE REMARK 999	UNP P29476
B	106	MET	VAL	ENGINEERED MUTATION	UNP P29476
B	368	ASP	ASN	ENGINEERED MUTATION	UNP P29476

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



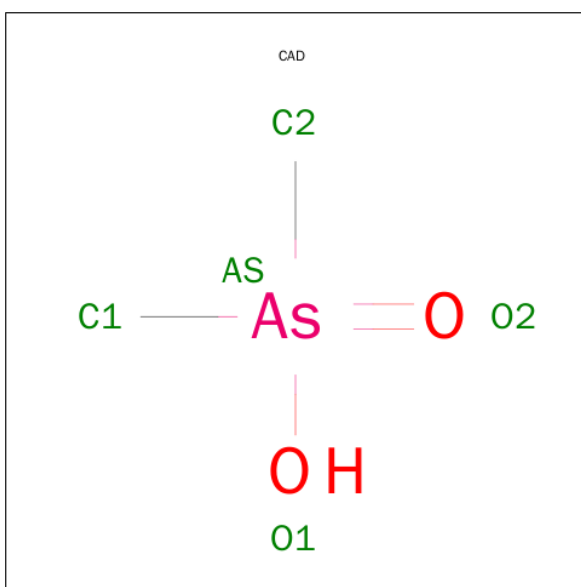
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	B	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- Molecule 3 is ACETATE ION (three-letter code: ACT) (formula:  $C_2H_3O_2$ ).



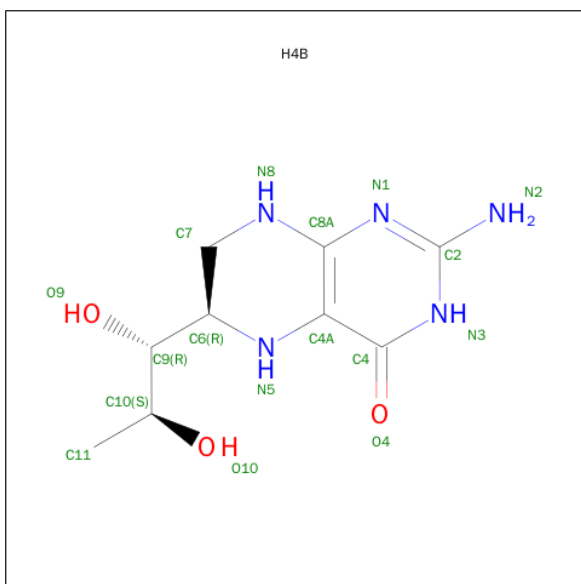
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 4 is CACODYLIC ACID (three-letter code: CAD) (formula:  $C_2H_7AsO_2$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	As	C	0	0
			3	1	2		
4	B	1	Total	As	C	0	0
			3	1	2		

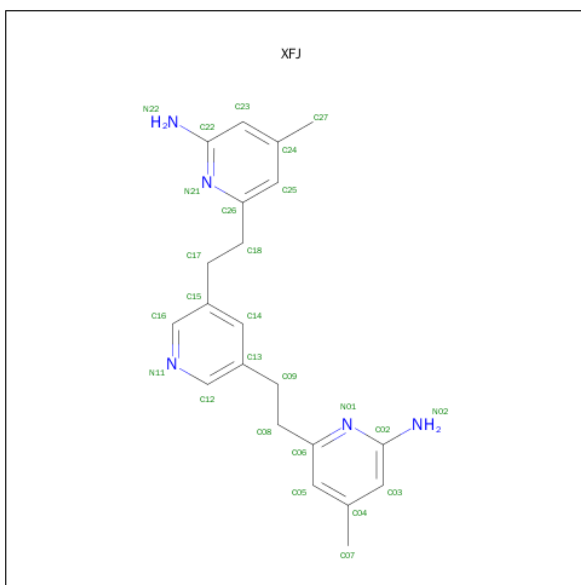
- Molecule 5 is 5,6,7,8-TETRAHYDROBIOPTERIN (three-letter code: H4B) (formula:  $C_9H_{15}N_5O_3$ ).



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

- Molecule 6 is 6,6'-(PYRIDINE-3,5-DIYLDIETHANE-2,1-DIYL)BIS(4-METHYLPYRIDIN-2-AMINE) (three-letter code: XFJ) (formula: C<sub>21</sub>H<sub>25</sub>N<sub>5</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	N		0	0
			26	21	5			
6	B	1	Total	C	N		0	0
			26	21	5			

- Molecule 7 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

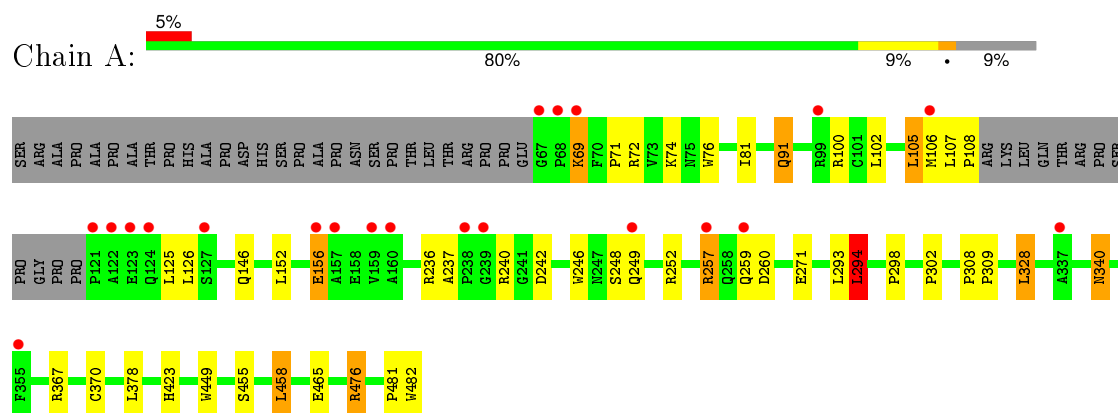
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	224	Total	O	0	0
			224	224		
8	B	216	Total	O	0	0
			216	216		

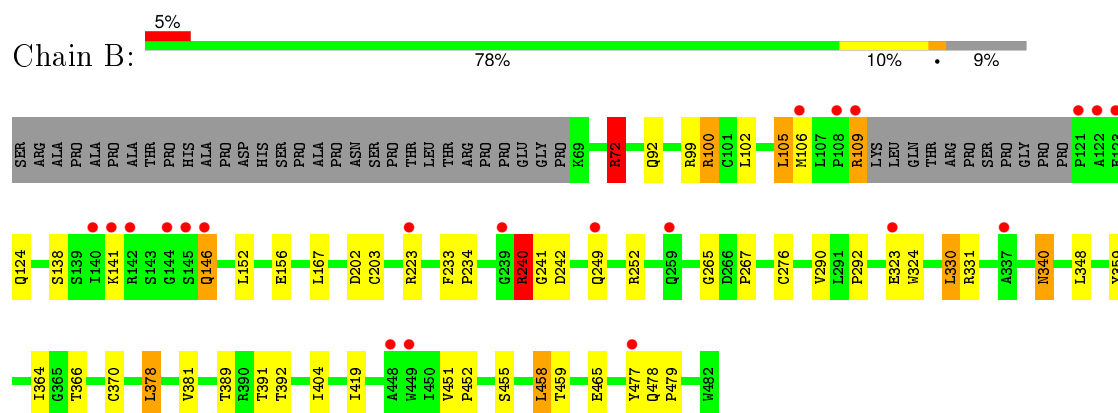
### 3 Residue-property plots

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 ( $RSRZ > 2$ ). 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, $\alpha$ , $\beta$ , $\gamma$	58.15Å 106.63Å 156.96Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.30 – 2.65 39.24 – 2.65	Depositor EDS
% Data completeness (in resolution range)	99.4 (39.30-2.65) 97.9 (39.24-2.65)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	0.07	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.20 (at 2.65Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, $R_{free}$	0.167 , 0.215 0.198 , 0.229	Depositor DCC
$R_{free}$ test set	1396 reflections (5.14%)	DCC
Wilson B-factor (Å <sup>2</sup> )	50.0	Xtriage
Anisotropy	0.405	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 37.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 28592 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	7116	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.66% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: ZN, XFJ, H4B, ACT, HEM, CAD

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	0.75	0/3330	0.75	3/4533 (0.1%)
1	B	0.77	1/3337 (0.0%)	0.76	5/4539 (0.1%)
All	All	0.76	1/6667 (0.0%)	0.76	8/9072 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	203	CYS	CB-SG	-6.47	1.71	1.82

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	240	ARG	NE-CZ-NH1	8.56	124.58	120.30
1	B	72	ARG	NE-CZ-NH1	7.80	124.20	120.30
1	B	240	ARG	NE-CZ-NH2	-7.42	116.59	120.30
1	A	367	ARG	NE-CZ-NH1	-7.03	116.78	120.30
1	B	72	ARG	NE-CZ-NH2	-6.08	117.26	120.30
1	A	328	LEU	CA-CB-CG	6.06	129.24	115.30
1	B	242	ASP	CB-CG-OD1	5.17	122.95	118.30
1	A	294	LEU	CA-CB-CG	5.10	127.02	115.30

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	3240	0	3139	34	0
1	B	3249	0	3150	35	0
2	A	43	0	30	1	0
2	B	43	0	30	5	0
3	A	4	0	3	0	0
3	B	4	0	3	0	0
4	A	3	0	0	0	0
4	B	3	0	0	1	0
5	A	17	0	15	1	0
5	B	17	0	15	0	0
6	A	26	0	25	2	0
6	B	26	0	25	2	0
7	A	1	0	0	0	0
8	A	224	0	0	4	0
8	B	216	0	0	3	0
All	All	7116	0	6435	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:107:LEU:HB3	1:A:108:PRO:HD2	1.61	0.83
1:B:249[A]:GLN:HB2	1:B:252:ARG:HG3	1.64	0.79
1:A:240:ARG:HD3	1:A:298:PRO:HB3	1.69	0.75
1:B:381:VAL:HG21	1:B:404:ILE:HD11	1.72	0.70
1:B:324:TRP:HB2	4:B:950:CAD:C1	2.22	0.69
1:A:249:GLN:HB2	1:A:252:ARG:HG3	1.78	0.66
1:B:366:THR:O	1:B:370:CYS:HB2	1.96	0.65
1:A:259:GLN:H	1:A:259:GLN:CD	1.99	0.65
1:A:340:ASN:HD22	1:A:340:ASN:H	1.44	0.63
2:A:500:HEM:O2D	6:A:800:XFJ:H17	1.98	0.63
1:B:72:ARG:HH11	1:B:72:ARG:HG2	1.64	0.62
1:B:109[B]:ARG:NH1	1:B:109[B]:ARG:HB2	2.14	0.62
1:A:237:ALA:O	8:A:1206:HOH:O	2.16	0.62
1:A:76:TRP:CE2	1:B:106:MET:HE3	2.36	0.60
1:A:71:PRO:HA	1:B:109[B]:ARG:HH21	1.65	0.60
1:A:69:LYS:HE2	1:A:69:LYS:HA	1.83	0.59
1:B:370:CYS:SG	1:B:378:LEU:HD13	2.43	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:152:LEU:O	1:A:156:GLU:HG2	2.04	0.58
1:A:370:CYS:SG	1:A:378:LEU:HD13	2.44	0.58
1:B:249[A]:GLN:HB2	1:B:252:ARG:CG	2.33	0.57
1:B:109[B]:ARG:CZ	1:B:109[B]:ARG:HB2	2.35	0.56
1:B:478:GLN:HB2	1:B:479:PRO:HD2	1.88	0.56
1:A:246:TRP:HB2	1:A:294:LEU:HB3	1.88	0.54
1:A:71:PRO:HA	1:B:109[B]:ARG:NH2	2.24	0.53
1:A:476:ARG:HD2	8:A:1148:HOH:O	2.09	0.53
1:B:378:LEU:HB2	8:B:1232:HOH:O	2.09	0.52
1:B:330:LEU:HB2	8:B:1233:HOH:O	2.08	0.51
2:B:500:HEM:HBC2	2:B:500:HEM:CMC	2.41	0.51
1:B:72:ARG:NH1	1:B:72:ARG:HG2	2.27	0.49
1:B:359:TYR:CD2	1:B:364:ILE:HD11	2.47	0.49
1:A:106:MET:HE1	6:A:800:XFJ:C24	2.43	0.48
2:B:500:HEM:HBB2	2:B:500:HEM:HHC	1.95	0.48
1:A:465:GLU:HB3	1:B:105:LEU:HD22	1.96	0.48
1:B:340:ASN:HD22	1:B:340:ASN:H	1.61	0.48
1:A:126:LEU:HD11	1:A:156:GLU:HA	1.96	0.47
1:B:152:LEU:O	1:B:156:GLU:HG2	2.15	0.47
1:B:290:VAL:HG23	8:B:1388:HOH:O	2.14	0.47
2:B:500:HEM:HBC2	2:B:500:HEM:HMC1	1.97	0.47
1:A:248:SER:HA	1:A:340:ASN:HB3	1.97	0.46
1:A:476:ARG:NE	8:A:1148:HOH:O	2.49	0.46
1:A:105:LEU:HD22	1:B:465:GLU:HB3	1.98	0.46
1:B:455:SER:O	1:B:458:LEU:HB2	2.16	0.45
1:A:72[B]:ARG:H	1:B:109[B]:ARG:CZ	2.30	0.45
1:A:455:SER:HB3	1:A:458:LEU:HD22	1.99	0.45
1:B:451:VAL:HA	1:B:452:PRO:HD3	1.85	0.45
1:B:233:PHE:HB3	1:B:234:PRO:CD	2.47	0.44
1:A:246:TRP:CZ2	1:A:302:PRO:HG3	2.52	0.43
1:A:91:GLN:HE21	1:A:91:GLN:HA	1.84	0.43
1:A:308:PRO:HA	1:A:309:PRO:HD3	1.93	0.43
1:A:74:LYS:HD2	1:A:81:ILE:HD11	2.00	0.43
1:B:167:LEU:HG	1:B:348:LEU:HD12	2.01	0.43
2:B:500:HEM:O2D	6:B:800:XFJ:H17	2.18	0.43
1:A:423:HIS:HB2	1:B:392:THR:HB	2.00	0.42
1:A:378:LEU:HB2	8:A:1005:HOH:O	2.18	0.42
1:B:452:PRO:HG2	1:B:459:THR:HG21	2.01	0.42
1:B:240:ARG:HD2	1:B:241:GLY:O	2.20	0.42
1:B:265:GLY:O	1:B:267:PRO:HD3	2.20	0.42
1:A:449:TRP:HA	5:A:600:H4B:N1	2.35	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:236:ARG:HD3	1:A:242:ASP:OD1	2.19	0.42
1:A:240:ARG:HD3	1:A:298:PRO:CB	2.45	0.41
1:B:72:ARG:CG	1:B:72:ARG:HH11	2.31	0.41
1:A:126:LEU:HD21	1:A:156:GLU:HB3	2.02	0.41
1:A:481:PRO:HD2	1:A:482:TRP:CE3	2.56	0.41
1:B:99:ARG:HG2	1:B:100:ARG:HD2	2.02	0.40
1:B:276:CYS:SG	1:B:292:PRO:HG2	2.61	0.40
1:B:146:GLN:HA	1:B:146:GLN:HE21	1.87	0.40
1:A:257:ARG:HG3	1:A:257:ARG:HH11	1.86	0.40
2:B:500:HEM:C1C	6:B:800:XFJ:H07B	2.56	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	403/444 (91%)	394 (98%)	9 (2%)	0	100	100
1	B	402/444 (90%)	391 (97%)	11 (3%)	0	100	100
All	All	805/888 (91%)	785 (98%)	20 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 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	346/377 (92%)	329 (95%)	17 (5%)	31	58
1	B	347/377 (92%)	323 (93%)	24 (7%)	19	39
All	All	693/754 (92%)	652 (94%)	41 (6%)	25	48

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	69	LYS
1	A	91	GLN
1	A	100	ARG
1	A	102	LEU
1	A	105	LEU
1	A	125	LEU
1	A	146	GLN
1	A	156	GLU
1	A	257	ARG
1	A	260	ASP
1	A	271	GLU
1	A	293	LEU
1	A	294	LEU
1	A	328	LEU
1	A	340	ASN
1	A	458	LEU
1	A	476	ARG
1	B	72	ARG
1	B	92	GLN
1	B	100	ARG
1	B	102	LEU
1	B	105	LEU
1	B	109[A]	ARG
1	B	109[B]	ARG
1	B	124	GLN
1	B	138	SER
1	B	141	LYS
1	B	146	GLN
1	B	202	ASP
1	B	223	ARG
1	B	240	ARG
1	B	323	GLU
1	B	330	LEU
1	B	331	ARG
1	B	340	ASN

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Mol	Chain	Res	Type
1	B	378	LEU
1	B	389	THR
1	B	391	THR
1	B	419	ILE
1	B	458	LEU
1	B	477	TYR

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	91	GLN
1	A	146	GLN
1	A	191	GLN
1	A	278	GLN
1	A	340	ASN
1	A	376	ASN
1	A	413	GLN
1	A	468	ASN
1	B	146	GLN
1	B	191	GLN
1	B	222	ASN
1	B	225	ASN
1	B	340	ASN
1	B	376	ASN
1	B	405	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 11 ligands modelled in this entry, 1 is monoatomic - leaving 10 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	500	1	30,50,50	1.97	9 (30%)	24,82,82	2.64	12 (50%)
5	H4B	A	600	-	13,18,18	0.97	0	11,26,26	2.71	6 (54%)
6	XFJ	A	800	-	28,28,28	0.93	1 (3%)	38,38,38	2.08	13 (34%)
3	ACT	A	860	-	1,3,3	1.17	0	0,3,3	0.00	-
4	CAD	A	950	-	0,2,4	0.00	-	0,1,6	0.00	-
2	HEM	B	500	1	30,50,50	2.21	8 (26%)	24,82,82	2.66	12 (50%)
5	H4B	B	600	-	13,18,18	0.84	0	11,26,26	2.49	5 (45%)
6	XFJ	B	800	-	28,28,28	0.88	0	38,38,38	2.15	11 (28%)
3	ACT	B	860	-	1,3,3	1.47	0	0,3,3	0.00	-
4	CAD	B	950	-	0,2,4	0.00	-	0,1,6	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	500	1	-	0/10/54/54	0/0/8/8
5	H4B	A	600	-	-	0/8/17/17	0/2/2/2
6	XFJ	A	800	-	-	0/10/10/10	0/3/3/3
3	ACT	A	860	-	-	0/0/0/0	0/0/0/0
4	CAD	A	950	-	-	0/0/0/0	0/0/0/0
2	HEM	B	500	1	-	0/10/54/54	0/0/8/8
5	H4B	B	600	-	-	0/8/17/17	0/2/2/2
6	XFJ	B	800	-	-	0/10/10/10	0/3/3/3
3	ACT	B	860	-	-	0/0/0/0	0/0/0/0
4	CAD	B	950	-	-	0/0/0/0	0/0/0/0

All (18) bond length outliers are listed below:



Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	500	HEM	C3B-C4B	-7.37	1.45	1.51
2	A	500	HEM	C3D-C4D	-5.37	1.44	1.51
2	B	500	HEM	C3D-C4D	-5.30	1.44	1.51
2	A	500	HEM	C3B-C4B	-3.52	1.48	1.51
2	B	500	HEM	C2C-C1C	-3.38	1.46	1.52
2	A	500	HEM	C2C-C1C	-2.88	1.47	1.52
2	B	500	HEM	C2D-C1D	-2.35	1.44	1.51
2	A	500	HEM	C2D-C1D	-2.30	1.44	1.51
2	B	500	HEM	C2B-C1B	-2.10	1.45	1.51
2	A	500	HEM	C2B-C1B	-2.03	1.45	1.51
6	A	800	XFJ	C16-C15	2.05	1.41	1.38
2	A	500	HEM	C4C-NC	2.07	1.38	1.36
2	B	500	HEM	C3C-CAC	2.10	1.55	1.51
2	A	500	HEM	FE-NB	2.37	2.10	1.97
2	B	500	HEM	C3B-CAB	2.43	1.55	1.51
2	A	500	HEM	C3C-CAC	2.56	1.56	1.51
2	B	500	HEM	FE-NC	2.86	2.07	1.95
2	A	500	HEM	C1C-NC	3.65	1.40	1.36

All (59) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	500	HEM	CBA-CAA-C2A	-5.44	102.77	112.53
2	B	500	HEM	C3C-CAC-CBC	-4.22	117.98	124.46
2	A	500	HEM	CAA-C2A-C1A	-4.12	122.54	127.01
5	A	600	H4B	N3-C2-N1	-4.08	118.85	125.53
5	B	600	H4B	N3-C2-N1	-3.64	119.56	125.53
2	A	500	HEM	CBA-CAA-C2A	-3.64	106.01	112.53
6	A	800	XFJ	C18-C26-C25	-3.24	116.59	121.13
6	B	800	XFJ	C18-C26-C25	-3.03	116.88	121.13
6	A	800	XFJ	C24-C25-C26	-2.94	118.44	120.28
6	A	800	XFJ	C15-C14-C13	-2.89	117.50	121.25
2	A	500	HEM	CBD-CAD-C3D	-2.74	105.59	113.55
6	B	800	XFJ	C05-C06-N01	-2.72	119.89	122.96
6	B	800	XFJ	C25-C26-N21	-2.56	120.07	122.96
6	A	800	XFJ	C15-C16-N11	-2.45	119.09	123.66
6	B	800	XFJ	C15-C16-N11	-2.36	119.26	123.66
2	A	500	HEM	C3B-C4B-NB	-2.33	107.17	111.63
6	B	800	XFJ	C15-C14-C13	-2.18	118.43	121.25
2	B	500	HEM	CAA-C2A-C1A	-2.15	124.67	127.01
2	A	500	HEM	C3C-CAC-CBC	-2.10	121.24	124.46
6	A	800	XFJ	C25-C26-N21	-2.07	120.62	122.96
2	B	500	HEM	C3B-C4B-NB	-2.04	107.73	111.63

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	800	XFJ	C25-C24-C23	2.06	120.71	118.08
5	B	600	H4B	N2-C2-N1	2.09	120.67	117.20
2	B	500	HEM	C2C-C1C-CHC	2.17	126.97	123.68
6	A	800	XFJ	C12-N11-C16	2.46	121.20	117.50
2	B	500	HEM	C3B-C4B-CHC	2.48	126.66	123.16
5	A	600	H4B	N2-C2-N3	2.60	121.50	117.20
6	B	800	XFJ	C12-N11-C16	2.69	121.55	117.50
5	B	600	H4B	C2-N1-C8A	2.73	120.69	114.54
2	A	500	HEM	CMD-C2D-C3D	2.82	126.84	114.35
2	B	500	HEM	C2D-C3D-C4D	2.85	106.34	101.50
6	A	800	XFJ	C02-N01-C06	2.92	120.30	118.23
6	A	800	XFJ	N22-C22-N21	2.95	121.87	116.50
6	A	800	XFJ	C14-C13-C12	2.99	119.83	116.57
5	A	600	H4B	C4A-C8A-N8	3.01	121.98	118.43
2	B	500	HEM	CMD-C2D-C3D	3.02	127.72	114.35
2	A	500	HEM	C2D-C3D-C4D	3.03	106.64	101.50
6	B	800	XFJ	N22-C22-N21	3.05	122.06	116.50
6	B	800	XFJ	C02-N01-C06	3.09	120.43	118.23
5	A	600	H4B	C2-N1-C8A	3.63	122.71	114.54
2	B	500	HEM	CMB-C2B-C3B	3.64	125.62	116.53
6	B	800	XFJ	C14-C15-C16	3.76	120.68	116.57
5	A	600	H4B	C4-N3-C2	3.83	121.26	115.94
2	B	500	HEM	CAD-C3D-C4D	3.91	126.26	112.47
5	B	600	H4B	C4-C4A-C8A	3.92	118.11	114.56
2	A	500	HEM	C3B-C4B-CHC	3.92	128.68	123.16
6	A	800	XFJ	C14-C15-C16	3.99	120.93	116.57
2	A	500	HEM	CAD-C3D-C4D	4.19	127.25	112.47
2	A	500	HEM	CMC-C2C-C3C	4.19	127.00	116.53
5	A	600	H4B	C4-C4A-C8A	4.21	118.37	114.56
2	A	500	HEM	CMB-C2B-C3B	4.30	127.27	116.53
2	A	500	HEM	CAD-C3D-C2D	4.48	126.09	113.22
2	B	500	HEM	CMC-C2C-C3C	4.53	127.83	116.53
5	B	600	H4B	C4-N3-C2	4.54	122.25	115.94
6	A	800	XFJ	C18-C26-N21	4.80	122.78	115.69
2	B	500	HEM	CAD-C3D-C2D	4.92	127.35	113.22
6	B	800	XFJ	C18-C26-N21	4.98	123.04	115.69
6	A	800	XFJ	C22-N21-C26	5.33	122.02	118.23
6	B	800	XFJ	C22-N21-C26	6.51	122.85	118.23

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	500	HEM	1	0
5	A	600	H4B	1	0
6	A	800	XFJ	2	0
2	B	500	HEM	5	0
6	B	800	XFJ	2	0
4	B	950	CAD	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	404/444 (90%)	0.26	21 (5%) 31 28	24, 39, 63, 82	0
1	B	403/444 (90%)	0.15	21 (5%) 31 28	25, 41, 65, 90	0
All	All	807/888 (90%)	0.20	42 (5%) 31 28	24, 40, 64, 90	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	121	PRO	5.4
1	A	239	GLY	5.3
1	A	160	ALA	4.9
1	B	259	GLN	4.8
1	A	159	VAL	4.8
1	B	122	ALA	4.4
1	A	156	GLU	4.2
1	A	123	GLU	4.1
1	A	122	ALA	3.8
1	A	238	PRO	3.7
1	A	259	GLN	3.6
1	A	69	LYS	3.5
1	A	124	GLN	3.4
1	A	68	PRO	3.2
1	B	477	TYR	3.1
1	B	123	GLU	3.1
1	A	99	ARG	2.9
1	B	142	ARG	2.9
1	B	106	MET	2.9
1	B	223	ARG	2.9
1	B	249[A]	GLN	2.8
1	A	106	MET	2.7
1	B	121	PRO	2.7
1	A	337	ALA	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	239	GLY	2.6
1	A	67	GLY	2.4
1	B	144	GLY	2.4
1	A	127	SER	2.4
1	A	257	ARG	2.4
1	B	323	GLU	2.3
1	B	448	ALA	2.3
1	B	146	GLN	2.2
1	B	145	SER	2.2
1	B	108	PRO	2.2
1	B	141	LYS	2.2
1	A	249	GLN	2.2
1	B	140	ILE	2.2
1	A	157	ALA	2.2
1	B	337	ALA	2.1
1	A	355	PHE	2.1
1	B	449	TRP	2.0
1	B	109[A]	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 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
3	ACT	B	860	4/4	0.95	0.25	5.75	48,50,50,50	0
4	CAD	A	950	3/5	0.94	0.19	3.25	56,56,59,60	0
6	XFJ	B	800	26/26	0.81	0.40	1.59	33,49,55,55	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
3	ACT	A	860	4/4	0.95	0.20	1.19	43,45,45,48	0
6	XFJ	A	800	26/26	0.87	0.31	0.78	24,40,54,55	0
2	HEM	A	500	43/43	0.96	0.27	0.78	27,29,37,41	0
5	H4B	A	600	17/17	0.98	0.22	-0.02	27,31,32,34	0
2	HEM	B	500	43/43	0.98	0.20	-0.17	25,30,34,37	0
5	H4B	B	600	17/17	0.97	0.21	-0.45	25,30,31,32	0
4	CAD	B	950	3/5	0.99	0.10	-1.75	60,60,61,62	0
7	ZN	A	900	1/1	0.99	0.07	-3.16	32,32,32,32	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.