



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 07:47 am GMT

PDB ID : 3N5S
Title : Structure of endothelial nitric oxide synthase heme domain complexed with 4-(2-(5-(2-(6-amino-4-methylpyridin-2-yl)ethyl)pyridin-3-yl)ethyl)-6-methylpyridin-2-amine
Authors : Delker, S.L.; Li, H.; Poulos, T.L.
Deposited on : 2010-05-25
Resolution : 2.18 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	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
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)	:	recalc28949

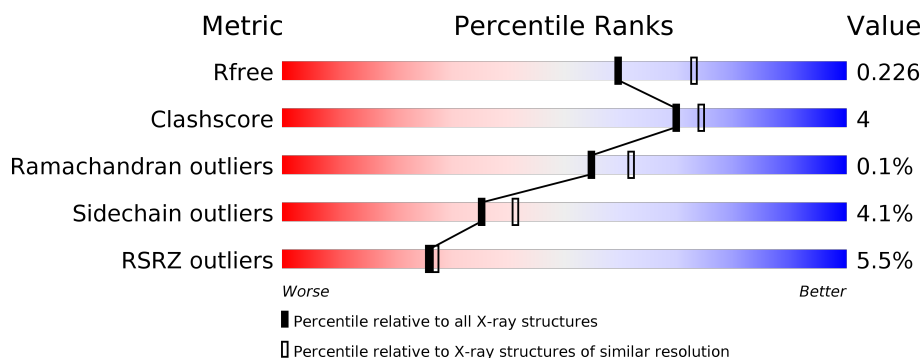
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.18 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}	100719	5526 (2.20-2.16)
Clashscore	112137	6386 (2.20-2.16)
Ramachandran outliers	110173	6282 (2.20-2.16)
Sidechain outliers	110143	6282 (2.20-2.16)
RSRZ outliers	101464	5562 (2.20-2.16)

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	444	
1	B	444	

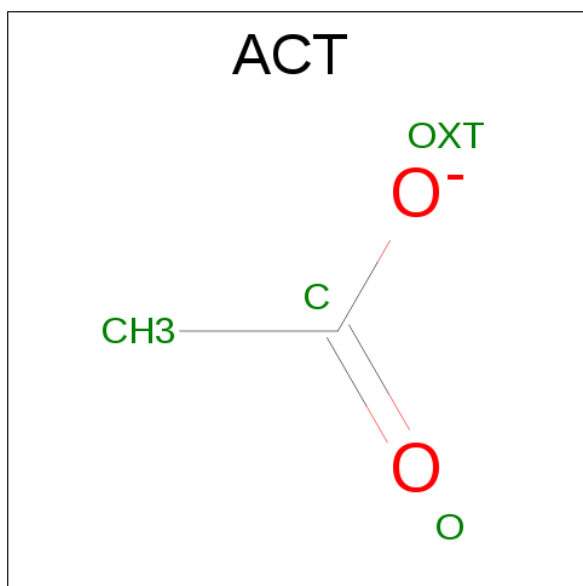
The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	ACT	B	860	-	-	-	X
4	GOL	A	880	-	-	-	X
4	GOL	B	880	-	-	-	X

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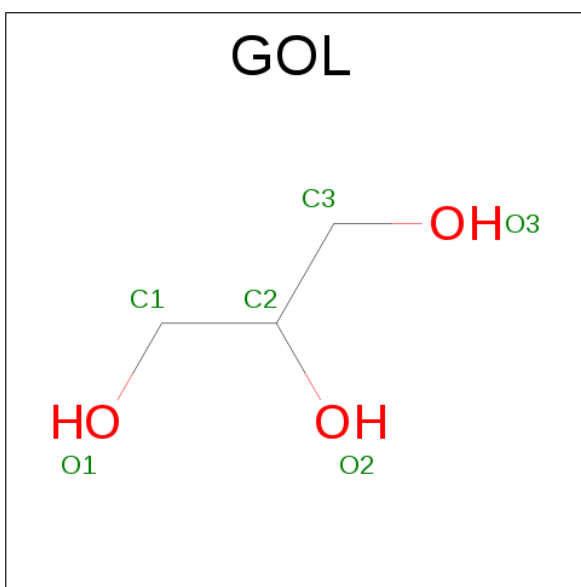
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

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



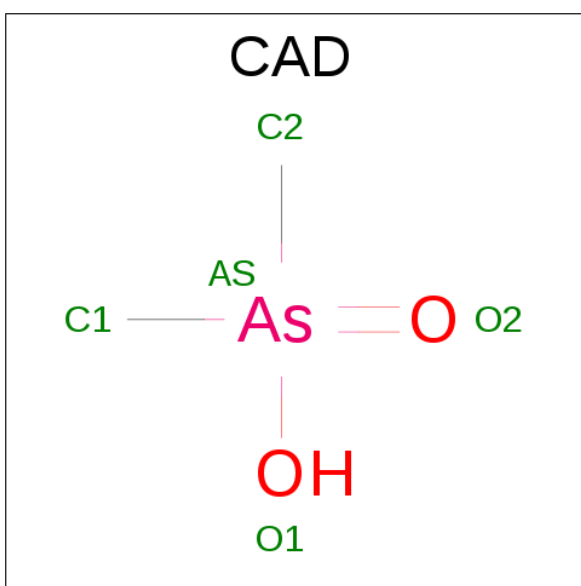
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



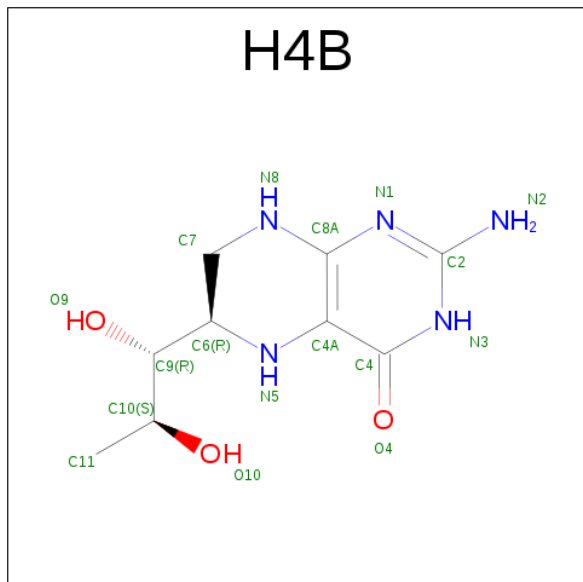
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is CACODYLIC ACID (three-letter code: CAD) (formula: $C_2H_7AsO_2$).



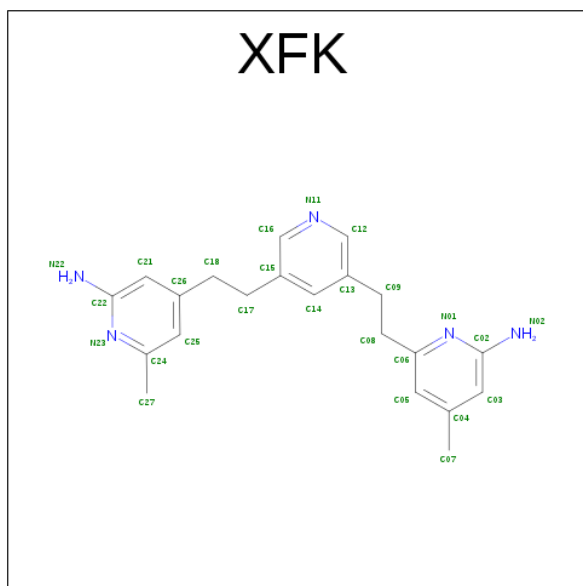
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	As	C	0	0
			3	1	2		
5	B	1	Total	As	C	0	0
			3	1	2		

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



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

- Molecule 7 is 6-(2-{5-[2-(2-AMINO-6-METHYLPYRIDIN-4-YL)ETHYL]PYRIDIN-3-YL}ETHYL)-4-METHYLPYRIDIN-2-AMINE (three-letter code: XFK) (formula: $C_{21}H_{25}N_5$).



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

- Molecule 8 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

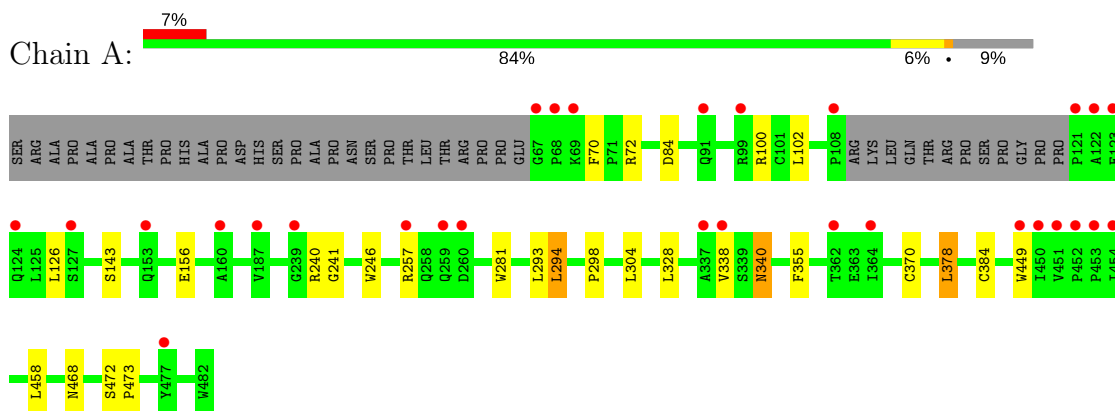
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	194	Total	O	0	0
			194	194		
9	B	189	Total	O	0	0
			189	189		

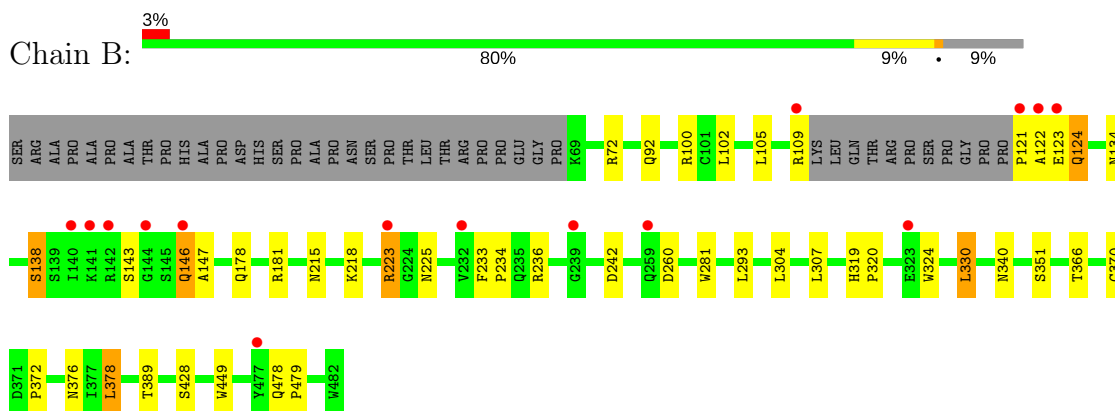
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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



• 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	57.95Å 106.80Å 157.03Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.10 – 2.18 38.10 – 2.18	Depositor EDS
% Data completeness (in resolution range)	99.7 (38.10-2.18) 99.7 (38.10-2.18)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.13 (at 2.18Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.177 , 0.220 0.183 , 0.226	Depositor DCC
R_{free} test set	2563 reflections (5.23%)	DCC
Wilson B-factor (Å ²)	34.6	Xtriage
Anisotropy	0.171	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 37.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7008	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.14% 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, ZN, H4B, XFK, 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	0.54	0/3299	0.61	0/4494
1	B	0.55	0/3306	0.61	1/4502 (0.0%)
All	All	0.55	0/6605	0.61	1/8996 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	378	LEU	CA-CB-CG	5.05	126.92	115.30

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	3209	0	3114	19	0
1	B	3217	0	3120	27	0
2	A	43	0	30	5	0
2	B	43	0	30	4	0
3	A	4	0	3	0	0
3	B	4	0	3	0	0
4	A	6	0	8	0	0
4	B	6	0	8	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	3	0	0	2	0
5	B	3	0	0	1	0
6	A	17	0	15	1	0
6	B	17	0	15	1	0
7	A	26	0	25	2	0
7	B	26	0	25	2	0
8	A	1	0	0	0	0
9	A	194	0	0	2	0
9	B	189	0	0	4	0
All	All	7008	0	6396	54	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 4.

All (54) 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:384:CYS:SG	5:A:950:CAD:AS	2.48	1.32
2:A:500:HEM:O2D	7:A:800:XFK:H21	1.55	1.06
1:B:215:ASN:HB2	9:B:1219:HOH:O	1.61	1.01
2:B:500:HEM:O2D	7:B:800:XFK:H21	1.72	0.89
1:B:146:GLN:NE2	1:B:147:ALA:H	1.73	0.87
1:A:257:ARG:HG3	1:A:257:ARG:HH11	1.48	0.78
1:B:146:GLN:HE21	1:B:147:ALA:H	1.29	0.78
1:A:240:ARG:HD3	1:A:298:PRO:HB3	1.68	0.75
1:B:236:ARG:HD2	1:B:242:ASP:OD1	1.89	0.71
1:B:223:ARG:HB2	1:B:223:ARG:HH11	1.62	0.65
1:A:257:ARG:NH1	1:A:257:ARG:HG3	2.13	0.65
1:B:236:ARG:HD3	1:B:351:SER:HB3	1.79	0.63
2:B:500:HEM:HBA2	7:B:800:XFK:H08	1.84	0.59
1:B:366:THR:O	1:B:370:CYS:HB2	2.02	0.58
1:A:340:ASN:HD22	1:A:340:ASN:H	1.51	0.58
1:A:384:CYS:CB	5:A:950:CAD:AS	3.13	0.57
1:B:72:ARG:NH1	9:B:1232:HOH:O	2.38	0.57
1:B:324:TRP:HB2	5:B:950:CAD:C1	2.35	0.56
1:B:378:LEU:HB2	9:B:1201:HOH:O	2.08	0.53
9:A:1052:HOH:O	4:B:880:GOL:H2	2.09	0.53
1:A:449:TRP:HA	6:A:600:H4B:N1	2.25	0.52
1:A:370:CYS:SG	1:A:378:LEU:HD13	2.50	0.52
1:B:281:TRP:HB2	1:B:304:LEU:HD21	1.92	0.52
1:B:146:GLN:N	1:B:146:GLN:HE21	2.08	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:281:TRP:HB2	1:A:304:LEU:HD21	1.94	0.50
1:A:240:ARG:HD3	1:A:298:PRO:CB	2.41	0.49
1:B:178:GLN:HE22	1:B:181:ARG:HH11	1.59	0.49
1:B:146:GLN:HE21	1:B:147:ALA:N	2.03	0.49
1:A:246:TRP:HB2	1:A:294:LEU:HB3	1.93	0.49
2:A:500:HEM:HBB2	2:A:500:HEM:HHC	1.95	0.49
1:B:449:TRP:HA	6:B:600:H4B:N1	2.27	0.49
1:B:233:PHE:HB3	1:B:234:PRO:CD	2.43	0.48
1:B:370:CYS:SG	1:B:378:LEU:HD13	2.54	0.47
1:A:355:PHE:CD1	2:A:500:HEM:HAC	2.50	0.47
1:A:240:ARG:HD2	1:A:241:GLY:O	2.15	0.46
1:B:293:LEU:HD11	1:B:307:LEU:HD21	1.98	0.45
1:A:378:LEU:HB2	9:A:1002:HOH:O	2.17	0.45
1:B:121:PRO:HB2	1:B:124:GLN:HB3	1.99	0.44
1:A:338:VAL:CG2	7:A:800:XFK:H09A	2.48	0.44
1:B:319:HIS:CG	1:B:320:PRO:HD2	2.53	0.44
1:B:134:ASN:O	1:B:138:SER:HB2	2.18	0.44
1:B:389:THR:O	1:B:389:THR:HG22	2.18	0.43
2:B:500:HEM:HBC2	2:B:500:HEM:CMC	2.49	0.43
1:B:146:GLN:NE2	1:B:147:ALA:N	2.55	0.42
1:B:478:GLN:HB2	1:B:479:PRO:HD2	2.03	0.41
1:A:126:LEU:HD11	1:A:156:GLU:HA	2.03	0.41
1:B:372:PRO:HA	1:B:376:ASN:ND2	2.36	0.41
1:B:389:THR:O	1:B:389:THR:CG2	2.69	0.41
2:A:500:HEM:CBB	2:A:500:HEM:HHC	2.51	0.41
2:B:500:HEM:HBB2	2:B:500:HEM:HHC	2.02	0.41
1:B:330:LEU:HB2	9:B:1205:HOH:O	2.19	0.40
1:A:355:PHE:CD1	2:A:500:HEM:CAC	3.05	0.40
1:A:472:SER:HA	1:A:473:PRO:C	2.42	0.40
1:A:70:PHE:HB3	1:A:84:ASP:O	2.21	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	400/444 (90%)	389 (97%)	11 (3%)	0	100	100
1	B	400/444 (90%)	393 (98%)	6 (2%)	1 (0%)	44	47
All	All	800/888 (90%)	782 (98%)	17 (2%)	1 (0%)	55	62

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	122	ALA

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	343/377 (91%)	332 (97%)	11 (3%)	44	53
1	B	344/377 (91%)	327 (95%)	17 (5%)	29	32
All	All	687/754 (91%)	659 (96%)	28 (4%)	35	41

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	72	ARG
1	A	100	ARG
1	A	102	LEU
1	A	143	SER
1	A	293	LEU
1	A	294	LEU
1	A	328	LEU
1	A	340	ASN
1	A	378	LEU
1	A	458	LEU
1	A	468	ASN
1	B	92	GLN

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Mol	Chain	Res	Type
1	B	100	ARG
1	B	102	LEU
1	B	105	LEU
1	B	109	ARG
1	B	123	GLU
1	B	124	GLN
1	B	138	SER
1	B	143	SER
1	B	146	GLN
1	B	218	LYS
1	B	223	ARG
1	B	225	ASN
1	B	260	ASP
1	B	330	LEU
1	B	340	ASN
1	B	428	SER

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

Mol	Chain	Res	Type
1	A	91	GLN
1	A	153	GLN
1	A	166	HIS
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	124	GLN
1	B	146	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 [i](#)

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 13 ligands modelled in this entry, 1 is monoatomic - leaving 12 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	28,50,50	2.26	7 (25%)	17,82,82	1.87	4 (23%)
6	H4B	A	600	-	14,18,18	0.78	0	12,26,26	2.48	6 (50%)
7	XFK	A	800	-	28,28,28	0.69	0	38,38,38	1.93	13 (34%)
3	ACT	A	860	-	1,3,3	1.55	0	0,3,3	0.00	-
4	GOL	A	880	-	5,5,5	0.44	0	5,5,5	0.24	0
5	CAD	A	950	-	0,2,4	0.00	-	0,1,6	0.00	-
2	HEM	B	500	1	28,50,50	2.24	8 (28%)	17,82,82	1.93	4 (23%)
6	H4B	B	600	-	14,18,18	0.91	1 (7%)	12,26,26	2.31	6 (50%)
7	XFK	B	800	-	28,28,28	0.56	0	38,38,38	2.09	12 (31%)
3	ACT	B	860	-	1,3,3	1.44	0	0,3,3	0.00	-
4	GOL	B	880	-	5,5,5	0.32	0	5,5,5	0.54	0
5	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	-	0/6/54/54	0/0/8/8
6	H4B	A	600	-	-	0/8/17/17	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	XFK	A	800	-	-	0/10/10/10	0/3/3/3
3	ACT	A	860	-	-	0/0/0/0	0/0/0/0
4	GOL	A	880	-	-	0/4/4/4	0/0/0/0
5	CAD	A	950	-	-	0/0/0/0	0/0/0/0
2	HEM	B	500	1	-	0/6/54/54	0/0/8/8
6	H4B	B	600	-	-	0/8/17/17	0/2/2/2
7	XFK	B	800	-	-	0/10/10/10	0/3/3/3
3	ACT	B	860	-	-	0/0/0/0	0/0/0/0
4	GOL	B	880	-	-	0/4/4/4	0/0/0/0
5	CAD	B	950	-	-	0/0/0/0	0/0/0/0

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	500	HEM	C3B-C2B	-5.55	1.33	1.40
2	B	500	HEM	C3B-C2B	-4.59	1.34	1.40
2	A	500	HEM	C3C-C2C	-4.37	1.34	1.40
2	B	500	HEM	C3C-C2C	-4.34	1.34	1.40
6	B	600	H4B	C4-C4A	-2.39	1.38	1.41
2	B	500	HEM	C1D-ND	2.02	1.40	1.36
2	A	500	HEM	CMB-C2B	2.22	1.56	1.51
2	B	500	HEM	CMA-C3A	2.48	1.56	1.51
2	B	500	HEM	C4D-ND	2.68	1.39	1.36
2	A	500	HEM	C4D-ND	2.96	1.40	1.36
2	A	500	HEM	C3B-CAB	3.55	1.54	1.47
2	A	500	HEM	C3C-CAC	3.68	1.55	1.47
2	B	500	HEM	C3C-CAC	3.84	1.55	1.47
2	B	500	HEM	C3B-CAB	3.93	1.55	1.47
2	A	500	HEM	C3D-C2D	4.62	1.51	1.37
2	B	500	HEM	C3D-C2D	5.30	1.53	1.37

All (45) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	500	HEM	CBA-CAA-C2A	-4.70	103.49	112.48
2	A	500	HEM	CBA-CAA-C2A	-4.69	103.51	112.48
2	A	500	HEM	CBD-CAD-C3D	-3.64	105.52	112.47
6	B	600	H4B	N3-C2-N1	-3.62	119.58	125.45
2	B	500	HEM	C1D-C2D-C3D	-3.58	104.50	107.00
7	B	800	XFK	C05-C06-N01	-3.47	119.15	122.91
2	B	500	HEM	CBD-CAD-C3D	-3.43	105.93	112.47
7	B	800	XFK	C26-C25-C24	-3.28	118.18	120.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	A	800	XFK	C15-C14-C13	-3.26	117.14	121.26
6	A	600	H4B	N3-C2-N1	-3.18	120.30	125.45
7	A	800	XFK	C26-C25-C24	-2.84	118.45	120.25
7	B	800	XFK	C15-C14-C13	-2.76	117.78	121.26
7	B	800	XFK	C27-C24-C25	-2.71	117.50	121.82
7	A	800	XFK	C17-C18-C26	-2.62	103.90	113.31
7	A	800	XFK	C05-C06-N01	-2.51	120.19	122.91
7	A	800	XFK	C27-C24-C25	-2.39	118.01	121.82
6	A	600	H4B	C6-C7-N8	-2.31	107.34	111.01
2	A	500	HEM	C1D-C2D-C3D	-2.28	105.41	107.00
6	B	600	H4B	C4A-N5-C6	-2.20	115.17	121.16
7	A	800	XFK	C17-C15-C16	-2.19	118.24	121.82
2	A	500	HEM	CMA-C3A-C4A	-2.15	125.16	128.46
7	A	800	XFK	C09-C13-C12	-2.13	118.34	121.82
7	B	800	XFK	C15-C16-N11	-2.05	119.95	123.70
6	A	600	H4B	N2-C2-N3	2.06	120.53	117.24
2	B	500	HEM	CAA-CBA-CGA	2.15	116.34	112.66
6	B	600	H4B	N2-C2-N3	2.16	120.69	117.24
7	B	800	XFK	C14-C13-C12	2.42	119.05	116.71
7	B	800	XFK	C12-N11-C16	2.43	120.88	117.45
7	A	800	XFK	C27-C24-N23	2.46	120.76	116.55
6	B	600	H4B	C2-N1-C8A	2.50	120.15	114.51
6	A	600	H4B	C2-N1-C8A	2.79	120.78	114.51
7	B	800	XFK	C27-C24-N23	2.86	121.44	116.55
7	A	800	XFK	C14-C13-C12	3.08	119.68	116.71
7	A	800	XFK	N22-C22-N23	3.44	122.44	116.64
7	B	800	XFK	N22-C22-N23	3.45	122.46	116.64
6	A	600	H4B	C4-N3-C2	3.59	121.22	116.06
6	B	600	H4B	C4-C4A-C8A	3.65	117.86	114.56
6	B	600	H4B	C4-N3-C2	3.86	121.61	116.06
7	B	800	XFK	C26-C21-C22	3.88	120.85	117.97
7	A	800	XFK	C02-N01-C06	3.94	120.96	118.17
7	A	800	XFK	C26-C21-C22	4.05	120.98	117.97
7	B	800	XFK	C14-C15-C16	4.06	120.63	116.71
7	A	800	XFK	C14-C15-C16	4.27	120.83	116.71
6	A	600	H4B	C4-C4A-C8A	5.02	119.11	114.56
7	B	800	XFK	C02-N01-C06	5.80	122.27	118.17

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

9 monomers are involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	500	HEM	5	0
6	A	600	H4B	1	0
7	A	800	XFK	2	0
5	A	950	CAD	2	0
2	B	500	HEM	4	0
6	B	600	H4B	1	0
7	B	800	XFK	2	0
4	B	880	GOL	1	0
5	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, 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	404/444 (90%)	0.20	29 (7%) 16 17	23, 34, 57, 72	0
1	B	403/444 (90%)	0.01	15 (3%) 42 42	22, 36, 59, 80	0
All	All	807/888 (90%)	0.11	44 (5%) 26 27	22, 35, 58, 80	0

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	259	GLN	6.2
1	A	69	LYS	5.4
1	A	239	GLY	5.0
1	A	160	ALA	4.6
1	B	121	PRO	4.5
1	A	67	GLY	4.2
1	A	121	PRO	4.2
1	B	477	TYR	4.0
1	A	122	ALA	3.7
1	B	259	GLN	3.6
1	B	142	ARG	3.5
1	B	146	GLN	3.4
1	A	124	GLN	3.4
1	A	123	GLU	3.3
1	B	122	ALA	3.3
1	A	91	GLN	3.0
1	B	239	GLY	2.9
1	A	451	VAL	2.8
1	A	477	TYR	2.8
1	A	338	VAL	2.8
1	A	68	PRO	2.8
1	B	141	LYS	2.7
1	B	123	GLU	2.7
1	A	127	SER	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	449	TRP	2.6
1	A	257	ARG	2.6
1	A	450	ILE	2.6
1	A	99	ARG	2.5
1	A	337	ALA	2.5
1	B	223	ARG	2.5
1	A	362	THR	2.4
1	B	140	ILE	2.4
1	A	364	ILE	2.4
1	A	260	ASP	2.3
1	B	323	GLU	2.2
1	B	144	GLY	2.2
1	A	153	GLN	2.1
1	A	108	PRO	2.1
1	A	454	ILE	2.1
1	A	452	PRO	2.1
1	A	187	VAL	2.1
1	A	453	PRO	2.1
1	B	232	VAL	2.1
1	B	109	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, 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
3	ACT	B	860	4/4	0.97	0.19	5.96	40,40,40,41	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	GOL	A	880	6/6	0.75	0.30	2.78	56,59,60,60	0
4	GOL	B	880	6/6	0.89	0.21	2.25	52,54,55,55	0
7	XFK	B	800	26/26	0.89	0.19	1.01	32,48,57,57	0
2	HEM	A	500	43/43	0.97	0.20	0.78	22,26,38,41	0
7	XFK	A	800	26/26	0.90	0.21	0.67	22,40,45,45	0
6	H4B	A	600	17/17	0.97	0.17	0.46	28,30,31,32	0
3	ACT	A	860	4/4	0.94	0.12	0.32	35,35,36,38	0
2	HEM	B	500	43/43	0.98	0.12	-0.02	23,27,38,40	0
5	CAD	A	950	3/5	0.99	0.08	-0.36	48,48,49,50	0
6	H4B	B	600	17/17	0.97	0.12	-0.73	34,36,37,37	0
8	ZN	A	900	1/1	1.00	0.09	-0.97	26,26,26,26	0
5	CAD	B	950	3/5	0.99	0.05	-	50,50,51,52	0

6.5 Other polymers [i](#)

There are no such residues in this entry.