



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 26, 2014 – 04:15 PM GMT

PDB ID : 3E7M
Title : Structure of murine iNOS oxygenase domain with inhibitor AR-C95791
Authors : Garcin, E.D.; Arvai, A.S.; Rosenfeld, R.J.; Kroeger, M.D.; Crane, B.R.; Andersson, G.; Andrews, G.; Hamley, P.J.; Mallinder, P.R.; Nicholls, D.J.; St-Gallay, S.A.; Tinker, A.C.; Gensmantel, N.P.; Mete, A.; Cheshire, D.R.; Connolly, S.; Stuehr, D.J.; Aberg, A.; Wallace, A.V.; Tainer, J.A.; Getzoff, E.D.
Deposited on : 2008-08-18
Resolution : 2.00 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

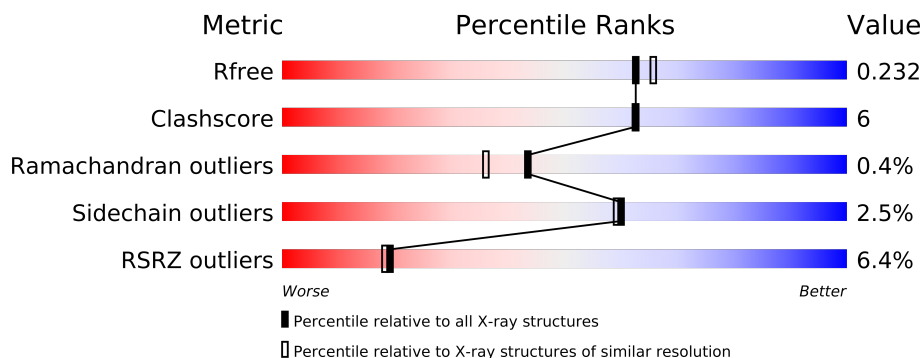
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

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}	66092	4888 (2.00-2.00)
Clashscore	79885	6188 (2.00-2.00)
Ramachandran outliers	78287	6102 (2.00-2.00)
Sidechain outliers	78261	6100 (2.00-2.00)
RSRZ outliers	66119	4890 (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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	433	
1	B	433	

2 Entry composition i

There are 7 unique types of molecules in this entry. The entry contains 7484 atoms, of which 0 are hydrogen and 0 are deuterium.

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.

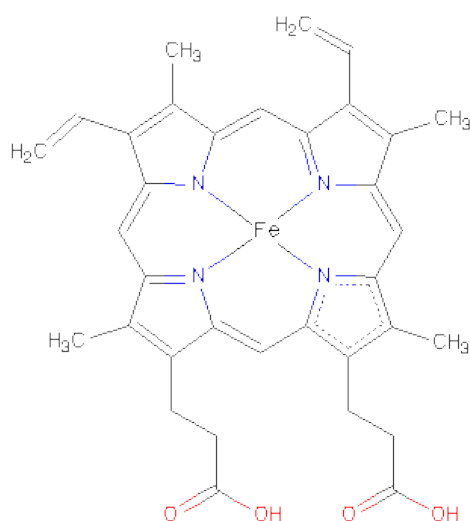
- Molecule 1 is a protein called Nitric oxide synthase, inducible.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	416	Total	C	N	O	S	0	0	0
			3389	2174	583	612	20			
1	B	421	Total	C	N	O	S	0	0	0
			3421	2193	589	618	21			

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

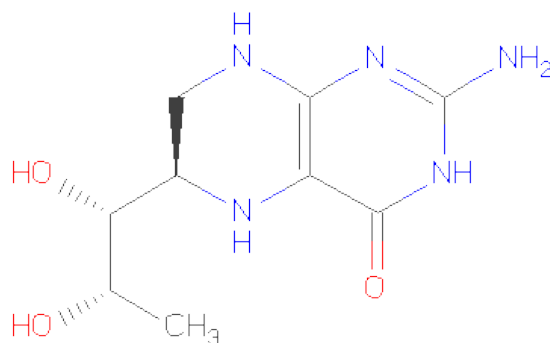
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Zn	0	0
			1	1		
2	A	1	Total	Zn	0	0
			1	1		

- Molecule 3 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



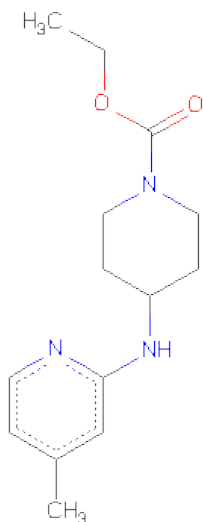
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
3	B	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- Molecule 4 is 5,6,7,8-TETRAHYDROBIOPTERIN (three-letter code: H4B) (formula: $C_9H_{15}N_5O_3$).



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

- Molecule 5 is ETHYL 4-[(4-METHYLPYRIDIN-2-YL)AMINO]PIPERIDINE-1-CARBOXYLATE (three-letter code: AT2) (formula: $C_{14}H_{21}N_3O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			19	14	3	2		
5	B	1	Total	C	N	O	0	0
			19	14	3	2		

- Molecule 6 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



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

- Molecule 7 is water.

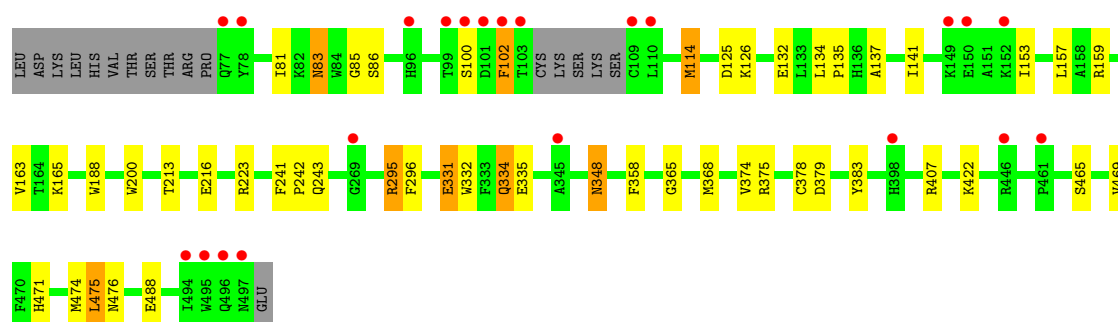
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	228	Total 228	O 228	0	0
7	B	278	Total 278	O 278	0	0

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.

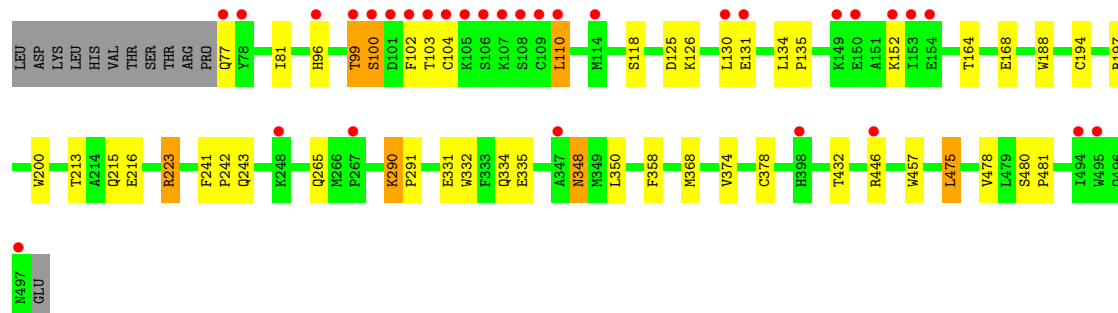
- Molecule 1: Nitric oxide synthase, inducible

Chain A: 



- Molecule 1: Nitric oxide synthase, inducible

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	213.85Å 213.85Å 116.64Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	19.96 – 2.00 39.41 – 2.00	Depositor EDS
% Data completeness (in resolution range)	98.7 (19.96-2.00) 98.7 (39.41-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.03	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.37 (at 2.00Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.223 , 0.237 0.218 , 0.232	Depositor DCC
R_{free} test set	5249 reflections (5.31%)	DCC
Wilson B-factor (Å ²)	30.3	Xtriage
Anisotropy	0.707	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 47.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	0 of 104070 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7484	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 1.90% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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: HEM, ZN, AT2, H4B, EDO

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	0.33	0/3488	0.59	2/4743 (0.0%)
1	B	0.36	0/3521	0.63	1/4787 (0.0%)
All	All	0.34	0/7009	0.61	3/9530 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	365	GLY	N-CA-C	-5.39	99.64	113.10
1	B	368	MET	N-CA-C	-5.19	96.98	111.00
1	A	368	MET	N-CA-C	-5.14	97.13	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3389	0	3276	36	0
1	B	3421	0	3310	41	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	43	0	30	0	0
3	B	43	0	30	0	0
4	A	17	0	15	0	0
4	B	17	0	15	1	0
5	A	19	0	21	0	0
5	B	19	0	21	0	0
6	A	8	0	12	0	0
7	A	228	0	0	5	2
7	B	278	0	0	5	1
All	All	7484	0	6730	77	3

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 6.

All (77) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:103:THR:O	1:B:110:LEU:HD23	1.62	1.00
1:B:81:ILE:HD11	1:B:475:LEU:HD13	1.44	0.97
1:A:81:ILE:HD11	1:A:475:LEU:HD13	1.62	0.79
1:A:83:ASN:HD22	1:A:85:GLY:H	1.35	0.75
1:A:83:ASN:ND2	1:A:85:GLY:H	1.86	0.72
1:A:331:GLU:H	1:A:331:GLU:CD	1.93	0.72
1:B:432:THR:CG2	7:B:1158:HOH:O	2.39	0.70
1:B:290:LYS:HE2	1:B:291:PRO:HD2	1.76	0.68
1:B:223:ARG:HH21	1:B:223:ARG:CG	2.07	0.67
1:B:215:GLN:OE1	7:B:1476:HOH:O	2.12	0.66
1:B:223:ARG:HH21	1:B:223:ARG:HG3	1.61	0.66
1:B:290:LYS:HE2	1:B:291:PRO:CD	2.31	0.61
1:B:81:ILE:HD11	1:B:475:LEU:CD1	2.26	0.60
1:B:102:PHE:HE1	1:B:478:VAL:HG23	1.66	0.60
1:B:194:CYS:HB3	1:B:197:ARG:HD2	1.83	0.60
1:A:465:SER:O	1:A:471:HIS:HE1	1.85	0.59
1:A:469:VAL:HG13	1:A:474:MET:HE3	1.86	0.58
1:B:194:CYS:O	1:B:197:ARG:HD3	2.04	0.58
1:B:432:THR:HG21	7:B:1158:HOH:O	1.99	0.57
1:A:469:VAL:HG13	1:A:474:MET:CE	2.35	0.56
1:A:295:ARG:HD2	1:A:296:PHE:CE2	2.42	0.55
1:A:488:GLU:HG2	7:A:1379:HOH:O	2.08	0.54
1:B:102:PHE:CE1	1:B:478:VAL:HG23	2.42	0.54
1:B:188:TRP:CE3	1:B:200:TRP:HA	2.44	0.52
1:A:134:LEU:HB3	1:A:135:PRO:HD3	1.91	0.52
1:B:103:THR:HG22	1:B:118:SER:OG	2.10	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:223:ARG:HH21	1:A:223:ARG:HG3	1.74	0.52
1:A:348:ASN:HB2	7:A:1221:HOH:O	2.09	0.52
1:A:83:ASN:ND2	1:A:86:SER:H	2.08	0.52
1:B:104:CYS:SG	1:B:110:LEU:N	2.81	0.51
1:A:188:TRP:CE3	1:A:200:TRP:HA	2.45	0.51
1:A:125:ASP:OD1	1:A:126:LYS:HG3	2.10	0.51
1:B:446:ARG:HH21	1:B:446:ARG:HG2	1.77	0.49
1:A:348:ASN:H	1:A:348:ASN:HD22	1.61	0.48
1:A:114:MET:HE3	7:A:1011:HOH:O	2.14	0.47
1:B:348:ASN:H	1:B:348:ASN:HD22	1.63	0.47
1:A:102:PHE:HE2	1:A:476:ASN:O	1.97	0.46
1:A:153:ILE:O	1:A:157:LEU:HD23	2.14	0.46
1:B:125:ASP:OD1	1:B:126:LYS:HG3	2.16	0.46
1:A:407:ARG:HD2	7:A:1364:HOH:O	2.15	0.46
1:B:152:LYS:HB3	7:B:1410:HOH:O	2.15	0.46
1:B:213:THR:OG1	1:B:216:GLU:HG3	2.15	0.46
1:A:295:ARG:HD3	1:A:383:TYR:HE2	1.81	0.45
1:B:194:CYS:O	1:B:197:ARG:CD	2.63	0.45
1:A:331:GLU:N	1:A:331:GLU:CD	2.66	0.45
1:A:332:TRP:O	1:A:335:GLU:HB2	2.16	0.45
1:B:134:LEU:HB3	1:B:135:PRO:HD3	1.99	0.44
1:A:348:ASN:HD22	1:A:348:ASN:N	2.16	0.44
1:B:332:TRP:O	1:B:335:GLU:HB2	2.16	0.44
1:A:348:ASN:ND2	1:A:348:ASN:H	2.16	0.44
1:A:241:PHE:HB3	1:A:242:PRO:CD	2.48	0.44
1:B:243:GLN:HB3	1:B:358:PHE:CE2	2.53	0.44
1:A:243:GLN:HB3	1:A:358:PHE:CE2	2.52	0.43
1:B:241:PHE:HB3	1:B:242:PRO:CD	2.49	0.43
1:B:241:PHE:HB3	1:B:242:PRO:HD2	2.00	0.43
1:B:457:TRP:HA	4:B:1902:H4B:N1	2.34	0.43
1:B:374:VAL:O	1:B:378:CYS:HB2	2.19	0.42
1:B:348:ASN:N	1:B:348:ASN:HD22	2.17	0.42
1:A:165:LYS:HD2	1:A:165:LYS:N	2.34	0.42
1:A:334:GLN:OE1	1:A:335:GLU:N	2.53	0.42
1:A:132:GLU:O	1:A:135:PRO:HD2	2.20	0.42
1:A:374:VAL:O	1:A:378:CYS:HB2	2.20	0.42
1:A:422:LYS:NZ	7:A:1357:HOH:O	2.53	0.41
1:B:99:THR:HG23	1:B:478:VAL:O	2.21	0.41
1:A:213:THR:OG1	1:A:216:GLU:HG3	2.21	0.41
1:B:81:ILE:HA	1:B:81:ILE:HD13	1.92	0.41
1:A:137:ALA:O	1:A:141:ILE:HG12	2.20	0.41
1:B:77:GLN:O	1:B:96:HIS:HE1	2.04	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:350:LEU:C	1:B:350:LEU:HD23	2.40	0.41
1:A:159:ARG:O	1:A:163:VAL:HG23	2.21	0.41
1:B:99:THR:HB	1:B:100:SER:H	1.51	0.40
1:B:223:ARG:HD3	7:B:1189:HOH:O	2.20	0.40
1:A:375:ARG:O	1:A:379:ASP:HB2	2.21	0.40
1:B:102:PHE:CG	1:B:103:THR:N	2.86	0.40
1:B:348:ASN:H	1:B:348:ASN:ND2	2.20	0.40
1:B:480:SER:HA	1:B:481:PRO:C	2.41	0.40
1:B:164:THR:O	1:B:168:GLU:HG3	2.22	0.40

All (3) 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	Distance(Å)	Clash(Å)
7:A:1082:HOH:O	7:A:1082:HOH:O[11_655]	0.98	1.22
7:A:1376:HOH:O	7:A:1376:HOH:O[11_655]	1.19	1.01
7:B:1215:HOH:O	7:B:1303:HOH:O[9_766]	2.00	0.20

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	412/433 (95%)	393 (95%)	18 (4%)	1 (0%)	56	51
1	B	419/433 (97%)	402 (96%)	15 (4%)	2 (0%)	38	29
All	All	831/866 (96%)	795 (96%)	33 (4%)	3 (0%)	43	36

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	100	SER
1	B	100	SER
1	B	99	THR

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	362/381 (95%)	354 (98%)	8 (2%)	64	65
1	B	366/381 (96%)	356 (97%)	10 (3%)	57	56
All	All	728/762 (96%)	710 (98%)	18 (2%)	60	59

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

Mol	Chain	Res	Type
1	A	83	ASN
1	A	102	PHE
1	A	114	MET
1	A	295	ARG
1	A	331	GLU
1	A	334	GLN
1	A	348	ASN
1	A	475	LEU
1	B	110	LEU
1	B	130	LEU
1	B	131	GLU
1	B	223	ARG
1	B	265	GLN
1	B	290	LYS
1	B	331	GLU
1	B	334	GLN
1	B	348	ASN
1	B	475	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	83	ASN
1	A	96	HIS
1	A	215	GLN
1	A	219	GLN
1	A	348	ASN
1	A	421	GLN

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Mol	Chain	Res	Type
1	A	442	GLN
1	A	471	HIS
1	B	96	HIS
1	B	143	GLN
1	B	215	GLN
1	B	219	GLN
1	B	348	ASN
1	B	421	GLN

5.3.3 RNA ⓘ

There are no RNA chains 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 10 ligands modelled in this entry, 2 are monoatomic - leaving 8 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$
6	EDO	A	5001	-	3,3,3	0.67	0	2,2,2	0.61	0
6	EDO	A	5002	-	3,3,3	0.60	0	2,2,2	0.70	0
3	HEM	A	901	-	49,50,50	2.25	17 (34%)	46,82,82	1.20	1 (2%)
4	H4B	A	902	-	18,18,18	1.48	3 (16%)	24,26,26	1.89	7 (29%)
5	AT2	A	906	-	20,20,20	1.81	5 (25%)	26,26,26	1.46	3 (11%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	HEM	B	1901	-	49,50,50	2.36	17 (34%)	46,82,82	1.33	3 (6%)
4	H4B	B	1902	-	18,18,18	1.61	5 (27%)	24,26,26	1.86	7 (29%)
5	AT2	B	1906	-	20,20,20	1.74	4 (20%)	26,26,26	1.44	4 (15%)

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
6	EDO	A	5001	-	-	0/1/1/1	0/0/0/0
6	EDO	A	5002	-	-	0/1/1/1	0/0/0/0
3	HEM	A	901	-	-	0/14/114/114	0/0/8/8
4	H4B	A	902	-	-	0/8/17/17	0/0/2/2
5	AT2	A	906	-	-	0/11/21/21	0/2/2/2
3	HEM	B	1901	-	-	0/14/114/114	0/0/8/8
4	H4B	B	1902	-	-	0/8/17/17	0/0/2/2
5	AT2	B	1906	-	-	0/11/21/21	0/2/2/2

All (51) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	901	HEM	C3D-C4D	7.37	1.46	1.44
3	B	1901	HEM	C2B-C1B	6.24	1.46	1.44
3	B	1901	HEM	C3D-C4D	5.92	1.46	1.44
3	A	901	HEM	C3D-C2D	-5.21	1.34	1.43
3	A	901	HEM	C2B-C1B	5.18	1.45	1.44
3	B	1901	HEM	C3D-C2D	-5.18	1.34	1.43
5	A	906	AT2	O16-C15	4.68	1.28	1.21
5	B	1906	AT2	O16-C15	4.30	1.28	1.21
3	B	1901	HEM	C2D-C1D	4.27	1.45	1.44
3	B	1901	HEM	CHB-C1B	4.14	1.41	1.35
3	B	1901	HEM	C3B-C2B	-4.00	1.36	1.43
5	A	906	AT2	C6-N5	3.96	1.41	1.34
5	B	1906	AT2	C6-N5	3.91	1.41	1.34
3	A	901	HEM	CHB-C1B	3.77	1.41	1.35
3	B	1901	HEM	CHA-C4D	3.71	1.41	1.35
3	A	901	HEM	CHA-C4D	3.58	1.41	1.35
4	A	902	H4B	C7-C6	-3.47	1.48	1.52
3	A	901	HEM	C3B-C2B	-3.46	1.37	1.43
3	B	1901	HEM	C3B-C4B	3.26	1.48	1.44
3	B	1901	HEM	FE-NA	3.22	2.06	1.92

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	901	HEM	C4A-C3A	3.19	1.44	1.40
3	B	1901	HEM	CHC-C1C	3.16	1.42	1.36
3	B	1901	HEM	CHD-C4C	3.13	1.42	1.36
4	A	902	H4B	C4A-C8A	3.10	1.45	1.41
4	B	1902	H4B	C4A-C8A	3.07	1.45	1.41
3	A	901	HEM	C3C-C2C	-2.97	1.38	1.43
3	B	1901	HEM	C4A-C3A	2.81	1.43	1.40
4	B	1902	H4B	C7-C6	-2.80	1.49	1.52
4	B	1902	H4B	C9-C10	2.67	1.58	1.52
3	A	901	HEM	C3B-C4B	2.65	1.47	1.44
5	A	906	AT2	C4-N5	2.63	1.40	1.34
3	A	901	HEM	FE-NA	2.56	2.03	1.92
3	A	901	HEM	FE-NB	2.54	2.07	1.97
3	B	1901	HEM	FE-ND	2.52	2.07	1.97
3	A	901	HEM	FE-NC	2.50	2.07	1.97
3	B	1901	HEM	FE-NB	2.46	2.06	1.97
4	B	1902	H4B	C7-N8	-2.43	1.42	1.46
3	B	1901	HEM	CMC-C2C	2.38	1.54	1.47
3	A	901	HEM	CMC-C2C	2.37	1.54	1.47
5	B	1906	AT2	C4-N5	2.33	1.40	1.34
4	A	902	H4B	C7-N8	-2.33	1.42	1.46
3	B	1901	HEM	C3C-C2C	-2.31	1.39	1.43
3	A	901	HEM	CHC-C1C	2.30	1.40	1.36
4	B	1902	H4B	C6-N5	2.30	1.50	1.46
3	A	901	HEM	FE-ND	2.24	2.06	1.97
5	B	1906	AT2	C3-C4	2.22	1.43	1.38
3	A	901	HEM	CHD-C4C	2.20	1.40	1.36
3	A	901	HEM	CMB-C2B	2.17	1.54	1.47
5	A	906	AT2	C3-C4	2.16	1.43	1.38
3	B	1901	HEM	CMB-C2B	2.16	1.54	1.47
5	A	906	AT2	C7-C6	2.03	1.44	1.39

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	901	HEM	C3B-C4B-NB	-5.70	109.92	114.00
3	B	1901	HEM	C3B-C4B-NB	-5.41	110.13	114.00
4	A	902	H4B	C7-C6-C9	5.19	121.67	113.66
4	B	1902	H4B	C7-C6-C9	5.07	121.47	113.66
5	B	1906	AT2	O17-C15-N12	3.66	114.67	111.60
5	A	906	AT2	O17-C15-N12	3.48	114.51	111.60
4	A	902	H4B	N8-C8A-N1	3.10	120.37	115.82

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1902	H4B	N8-C8A-N1	3.09	120.34	115.82
5	A	906	AT2	C3-C4-N5	-3.05	120.49	123.88
5	B	1906	AT2	C3-C4-N5	-2.89	120.67	123.88
4	B	1902	H4B	C6-C7-N8	2.89	115.39	111.66
4	A	902	H4B	C6-C7-N8	2.78	115.24	111.66
3	B	1901	HEM	C4A-C3A-C2A	2.77	108.92	107.00
4	B	1902	H4B	C4-C4A-N5	2.63	122.92	119.10
4	A	902	H4B	C2-N1-C8A	2.60	121.31	117.61
4	B	1902	H4B	C2-N1-C8A	2.52	121.20	117.61
4	B	1902	H4B	C4A-C4-N3	2.44	120.32	114.06
4	A	902	H4B	C4-C4A-N5	2.40	122.59	119.10
4	B	1902	H4B	C9-C6-N5	2.38	114.13	109.69
4	A	902	H4B	C4A-C4-N3	2.34	120.06	114.06
3	B	1901	HEM	CBD-CAD-C3D	-2.24	109.47	114.37
5	A	906	AT2	C7-C6-N5	-2.24	119.47	122.69
5	B	1906	AT2	C7-C6-N5	-2.24	119.47	122.69
5	B	1906	AT2	O16-C15-N12	-2.09	120.25	124.36
4	A	902	H4B	C4-C4A-C8A	2.02	116.43	114.56

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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	416/433 (96%)	0.20	22 (5%)	25 25	23, 35, 51, 64	2 (0%)
1	B	421/433 (97%)	0.32	31 (7%)	14 14	22, 33, 50, 67	7 (1%)
All	All	837/866 (96%)	0.26	53 (6%)	19 19	22, 34, 50, 67	9 (1%)

All (53) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	102	PHE	12.0
1	B	105	LYS	9.7
1	B	107	LYS	9.2
1	B	100	SER	8.8
1	B	108	SER	8.4
1	A	102	PHE	7.8
1	B	497	ASN	7.0
1	B	103	THR	6.9
1	B	101	ASP	6.4
1	A	103	THR	6.2
1	B	109	CYS	6.0
1	A	109	CYS	5.8
1	A	101	ASP	5.5
1	A	100	SER	5.5
1	A	497	ASN	5.1
1	A	78	TYR	5.1
1	B	494	ILE	5.1
1	A	494	ILE	5.0
1	B	106	SER	4.7
1	B	149	LYS	4.5
1	B	78	TYR	4.4
1	A	496	GLN	4.4
1	B	104	CYS	4.1
1	B	110	LEU	3.8

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Mol	Chain	Res	Type	RSRZ
1	A	99	THR	3.5
1	A	149	LYS	3.4
1	B	150	GLU	3.2
1	A	495	TRP	3.1
1	B	495	TRP	3.1
1	B	99	THR	2.9
1	A	77	GLN	2.8
1	A	110	LEU	2.8
1	B	131	GLU	2.7
1	B	77	GLN	2.7
1	A	446	ARG	2.6
1	A	398	HIS	2.6
1	A	152	LYS	2.6
1	B	446	ARG	2.4
1	B	130	LEU	2.3
1	B	153	ILE	2.3
1	A	96	HIS	2.3
1	B	398	HIS	2.3
1	A	345	ALA	2.2
1	B	267	PRO	2.2
1	B	114	MET	2.2
1	B	347	ALA	2.1
1	A	269	GLY	2.1
1	A	461	PRO	2.1
1	A	150	GLU	2.1
1	B	96	HIS	2.1
1	B	248	LYS	2.1
1	B	154	GLU	2.1
1	B	152	LYS	2.1

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates 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. 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
5	AT2	B	1906	19/19	0.21	1.55	18,22,28,30	0
4	H4B	A	902	17/17	0.20	1.33	23,25,29,29	0
3	HEM	A	901	43/43	0.17	1.08	21,24,26,29	0
5	AT2	A	906	19/19	0.17	0.89	20,22,29,30	0
3	HEM	B	1901	43/43	0.18	0.83	18,22,25,26	0
4	H4B	B	1902	17/17	0.17	0.67	22,25,28,29	0
6	EDO	A	5002	4/4	0.15	-0.61	55,55,55,55	0
2	ZN	B	1900	1/1	0.20	-1.50	34,34,34,34	1
2	ZN	A	900	1/1	0.14	-2.02	45,45,45,45	1
6	EDO	A	5001	4/4	0.10	-2.64	52,54,55,55	0

6.5 Other polymers

There are no such residues in this entry.