



Full wwPDB X-ray Structure Validation Report

Mar 1, 2014 – 02:41 AM GMT

PDB ID : 1F20
Title : CRYSTAL STRUCTURE OF RAT NEURONAL NITRIC-OXIDE SYNTHASE FAD/NADP+ DOMAIN AT 1.9A RESOLUTION.
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Deposited on : 2000-05-22
Resolution : 1.90 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>

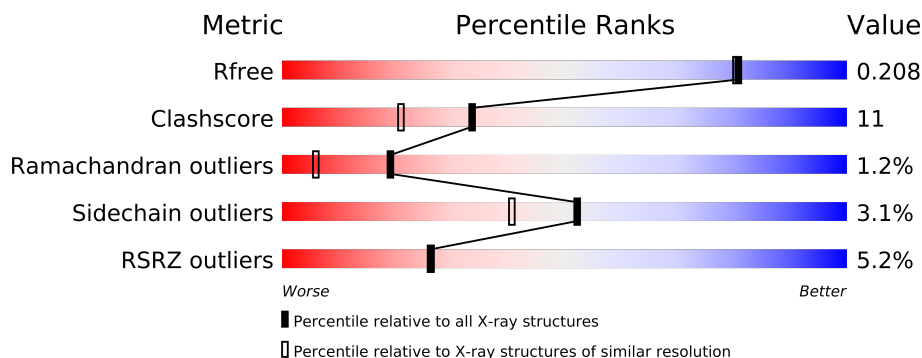
The following versions of software and data (see [references](#)) were used in the production of this report:

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 1.90 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	3684 (1.90-1.90)
Clashscore	79885	4465 (1.90-1.90)
Ramachandran outliers	78287	4413 (1.90-1.90)
Sidechain outliers	78261	4414 (1.90-1.90)
RSRZ outliers	66119	3686 (1.90-1.90)

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	435	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
4	GOL	A	2081	-	X
4	GOL	A	2082	-	X
4	GOL	A	2084	-	X
5	FMT	A	2085	-	X
5	FMT	A	2086	-	X
5	FMT	A	2087	-	X

2 Entry composition i

There are 6 unique types of molecules in this entry. The entry contains 4120 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.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	435	3501	2217	622	647	15	0	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1008	SER	PHE	CONFLICT	UNP P29476

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	A	1	53	27	9	15	2	0	0

- Molecule 3 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDEPHOSPHATE (three-letter code: NAP) (formula: $C_{21}H_{28}N_7O_{17}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

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



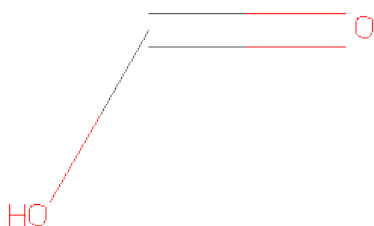
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		

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

- Molecule 5 is FORMIC ACID (three-letter code: FMT) (formula: CH₂O₂).



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

- Molecule 6 is water.

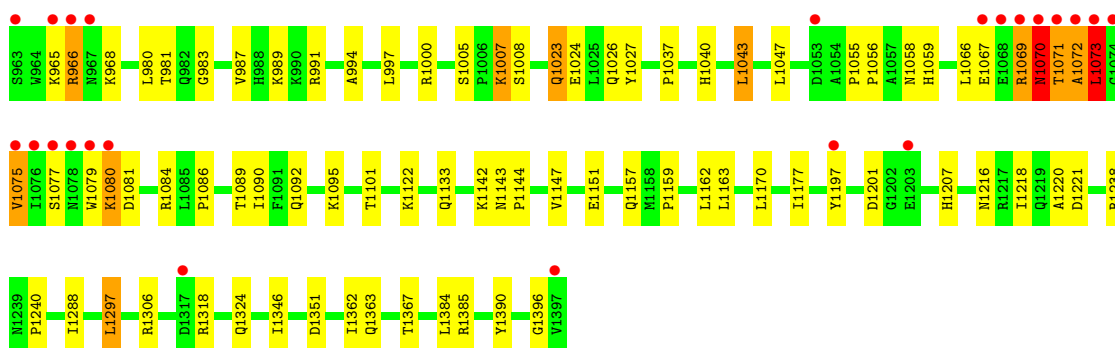
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	479	Total	O	0	0
			479	479		

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

Chain A: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	59.07Å 64.37Å 159.43Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	32.00 – 1.90 31.55 – 1.90	Depositor EDS
% Data completeness (in resolution range)	84.6 (32.00-1.90) 84.6 (31.55-1.90)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.00 (at 1.89Å)	Xtriage
Refinement program	CNS 0.5	Depositor
R, R_{free}	0.186 , 0.206 0.186 , 0.208	Depositor DCC
R_{free} test set	3346 reflections (8.11%)	DCC
Wilson B-factor (Å ²)	20.6	Xtriage
Anisotropy	0.315	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 37.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 41269 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4120	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.88% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, FMT, NAP, FAD

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	0/3587	0.64	3/4867 (0.1%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1070	ASN	N-CA-C	7.01	129.92	111.00
1	A	1072	ALA	N-CA-C	-6.63	93.11	111.00
1	A	1069	ARG	C-N-CA	5.52	135.50	121.70

There are no chirality outliers.

There are no planarity outliers.

5.2 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 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	3501	0	3465	74	0
2	A	53	0	31	0	0
3	A	48	0	25	6	0
4	A	30	0	40	6	0
5	A	9	0	3	0	0
6	A	479	0	0	14	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	4120	0	3564	79	0

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 11.

All (79) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:1073:LEU:HD13	1:A:1073:LEU:H	0.96	1.08
1:A:1073:LEU:H	1:A:1073:LEU:CD1	1.62	1.06
1:A:1073:LEU:HD13	1:A:1073:LEU:N	1.78	0.98
1:A:1056:PRO:HG2	1:A:1059:HIS:HD2	1.43	0.84
1:A:1101:THR:HG21	6:A:2024:HOH:O	1.79	0.82
1:A:1095:LYS:HZ1	4:A:2084:GOL:H12	1.51	0.75
1:A:1070:ASN:CG	1:A:1071:THR:H	1.89	0.74
1:A:1071:THR:HG22	1:A:1071:THR:O	1.88	0.73
1:A:1220:ALA:O	1:A:1221:ASP:HB2	1.92	0.69
1:A:1056:PRO:HG2	1:A:1059:HIS:CD2	2.27	0.69
1:A:1095:LYS:NZ	4:A:2084:GOL:H12	2.09	0.67
3:A:1502:NAP:H6N	6:A:1903:HOH:O	1.95	0.66
1:A:1396:GLY:HA3	6:A:1968:HOH:O	1.95	0.65
1:A:1070:ASN:C	1:A:1072:ALA:H	2.01	0.65
1:A:1007:LYS:HD3	6:A:1929:HOH:O	2.00	0.61
1:A:1069:ARG:O	1:A:1075:VAL:HG13	2.01	0.61
1:A:1073:LEU:N	1:A:1073:LEU:CD1	2.47	0.61
1:A:981:THR:HG21	1:A:994:ALA:HB2	1.85	0.58
1:A:1023:GLN:O	1:A:1026:GLN:HG3	2.03	0.58
1:A:991:ARG:HG2	1:A:1024:GLU:OE1	2.03	0.58
1:A:1101:THR:CG2	6:A:1992:HOH:O	2.52	0.57
1:A:980:LEU:HB2	4:A:2081:GOL:H32	1.85	0.57
1:A:1147:VAL:O	1:A:1151:GLU:HG3	2.05	0.56
1:A:1056:PRO:HB2	1:A:1058:ASN:OD1	2.06	0.55
1:A:1089:THR:OG1	1:A:1092:GLN:HG3	2.06	0.55
1:A:1005:SER:OG	1:A:1007:LYS:HG2	2.08	0.54
1:A:966:ARG:HD2	1:A:966:ARG:H	1.72	0.54
1:A:997:LEU:HD22	1:A:1218:ILE:HG13	1.89	0.54
1:A:1197:TYR:CE2	1:A:1207:HIS:HB2	2.43	0.53
4:A:2084:GOL:H11	6:A:2043:HOH:O	2.08	0.53
1:A:1066:LEU:HD11	1:A:1077:SER:HB2	1.89	0.53
1:A:1157:GLN:O	1:A:1159:PRO:HD3	2.10	0.52
1:A:987:VAL:CG2	1:A:1086:PRO:HD3	2.41	0.50
3:A:1502:NAP:N7N	4:A:2082:GOL:H11	2.27	0.50
1:A:1133:GLN:HB2	6:A:1946:HOH:O	2.12	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:1040:HIS:HB3	1:A:1043:LEU:HD22	1.94	0.50
1:A:1081:ASP:HB3	6:A:1758:HOH:O	2.11	0.49
1:A:1220:ALA:O	1:A:1221:ASP:CB	2.59	0.49
1:A:989:LYS:HE3	6:A:1923:HOH:O	2.12	0.49
1:A:966:ARG:HG2	1:A:966:ARG:HH11	1.77	0.48
1:A:989:LYS:HD2	1:A:989:LYS:N	2.28	0.48
1:A:1070:ASN:CG	1:A:1071:THR:N	2.63	0.48
1:A:1324:GLN:OE1	3:A:1502:NAP:H2A	2.13	0.48
1:A:1055:PRO:HD2	1:A:1090:ILE:HD11	1.95	0.48
1:A:1071:THR:O	1:A:1071:THR:CG2	2.60	0.47
1:A:1027:TYR:CE1	1:A:1177:ILE:HG21	2.48	0.47
1:A:1008:SER:HB2	1:A:1288:ILE:HG23	1.97	0.47
1:A:1122:LYS:NZ	6:A:1761:HOH:O	2.48	0.47
1:A:991:ARG:NH1	6:A:1750:HOH:O	2.49	0.46
3:A:1502:NAP:H71N	4:A:2082:GOL:H11	1.80	0.45
1:A:1000:ARG:HD2	1:A:1216:ASN:HD22	1.81	0.45
1:A:1072:ALA:HA	1:A:1073:LEU:HD13	1.98	0.45
1:A:965:LYS:HG2	1:A:968:LYS:HD2	1.98	0.45
3:A:1502:NAP:N7N	3:A:1502:NAP:PA	2.90	0.44
1:A:965:LYS:HG3	1:A:968:LYS:HB2	1.98	0.44
1:A:1101:THR:CG2	6:A:2024:HOH:O	2.50	0.44
1:A:1346:ILE:HD13	1:A:1362:ILE:CD1	2.47	0.44
1:A:1143:ASN:N	1:A:1144:PRO:CD	2.80	0.44
1:A:1238:ARG:O	1:A:1240:PRO:HD3	2.17	0.44
1:A:1346:ILE:HD13	1:A:1362:ILE:HD13	2.00	0.43
1:A:1007:LYS:HB2	1:A:1007:LYS:NZ	2.34	0.43
1:A:1363:GLN:O	1:A:1367:THR:HG23	2.18	0.43
1:A:1385:ARG:NH1	6:A:1657:HOH:O	2.51	0.43
1:A:1162:LEU:HD23	1:A:1162:LEU:C	2.39	0.43
1:A:1072:ALA:CA	1:A:1073:LEU:HD13	2.49	0.43
1:A:1384:LEU:HB3	1:A:1390:TYR:HB2	2.02	0.42
1:A:1297:LEU:HA	1:A:1297:LEU:HD22	1.91	0.42
1:A:991:ARG:HG2	1:A:1024:GLU:CD	2.40	0.42
1:A:1306:ARG:HD3	6:A:2061:HOH:O	2.19	0.42
1:A:1067:GLU:OE2	1:A:1080:LYS:HE3	2.20	0.42
1:A:966:ARG:NH1	1:A:966:ARG:HG2	2.35	0.42
1:A:1351:ASP:CB	3:A:1502:NAP:O7N	2.68	0.41
1:A:1023:GLN:HG3	1:A:1026:GLN:NE2	2.35	0.41
1:A:1047:LEU:HA	1:A:1147:VAL:HG23	2.02	0.41
1:A:1070:ASN:C	1:A:1072:ALA:N	2.69	0.41
1:A:965:LYS:CG	1:A:968:LYS:HD2	2.50	0.41
1:A:1079:TRP:C	1:A:1080:LYS:HG2	2.40	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:983:GLY:O	1:A:987:VAL:HG23	2.21	0.40
1:A:1142:LYS:O	1:A:1143:ASN:C	2.58	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	433/435 (100%)	418 (96%)	10 (2%)	5 (1%)	19 6

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1070	ASN
1	A	1073	LEU
1	A	1075	VAL
1	A	1201	ASP
1	A	1071	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	387/387 (100%)	375 (97%)	12 (3%)	52 41

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	966	ARG

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Mol	Chain	Res	Type
1	A	1007	LYS
1	A	1023	GLN
1	A	1037	PRO
1	A	1043	LEU
1	A	1073	LEU
1	A	1080	LYS
1	A	1084	ARG
1	A	1163	LEU
1	A	1170	LEU
1	A	1297	LEU
1	A	1318	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1004	GLN
1	A	1026	GLN
1	A	1045	ASN
1	A	1059	HIS
1	A	1216	ASN
1	A	1264	GLN
1	A	1285	GLN
1	A	1330	GLN
1	A	1363	GLN
1	A	1369	GLN
1	A	1388	ASN

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 ⓘ

10 ligands are modelled in this entry.

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	FAD	A	1501	-	58,58,58	3.62	32 (55%)	85,89,89	2.47	22 (25%)
3	NAP	A	1502	-	52,52,52	3.45	21 (40%)	80,80,80	1.82	15 (18%)
4	GOL	A	2080	-	5,5,5	0.25	0	5,5,5	0.68	0
4	GOL	A	2081	-	5,5,5	0.39	0	5,5,5	0.99	0
4	GOL	A	2082	-	5,5,5	0.34	0	5,5,5	0.67	0
4	GOL	A	2083	-	5,5,5	0.56	0	5,5,5	0.69	0
4	GOL	A	2084	-	5,5,5	0.57	0	5,5,5	0.91	0
5	FMT	A	2085	-	2,2,2	0.83	0	1,1,1	0.29	0
5	FMT	A	2086	-	2,2,2	0.70	0	1,1,1	0.36	0
5	FMT	A	2087	-	2,2,2	0.64	0	1,1,1	0.36	0

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	FAD	A	1501	-	-	0/34/50/50	0/1/6/6
3	NAP	A	1502	-	-	0/35/67/67	0/3/5/5
4	GOL	A	2080	-	-	0/4/4/4	0/0/0/0
4	GOL	A	2081	-	-	0/4/4/4	0/0/0/0
4	GOL	A	2082	-	-	0/4/4/4	0/0/0/0
4	GOL	A	2083	-	-	0/4/4/4	0/0/0/0
4	GOL	A	2084	-	-	0/4/4/4	0/0/0/0
5	FMT	A	2085	-	-	0/0/0/0	0/0/0/0
5	FMT	A	2086	-	-	0/0/0/0	0/0/0/0
5	FMT	A	2087	-	-	0/0/0/0	0/0/0/0

All (53) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1502	NAP	C2N-C3N	11.92	1.55	1.38
3	A	1502	NAP	C2N-N1N	10.64	1.49	1.35
2	A	1501	FAD	C9A-N10	8.57	1.51	1.38
2	A	1501	FAD	C4A-N3A	8.34	1.48	1.35
3	A	1502	NAP	C4N-C3N	7.62	1.52	1.39
3	A	1502	NAP	C5N-C4N	6.91	1.55	1.39
2	A	1501	FAD	C4X-C10	6.86	1.52	1.40
2	A	1501	FAD	C2-N3	6.67	1.50	1.37
3	A	1502	NAP	C6N-C5N	6.38	1.53	1.38
2	A	1501	FAD	P-O3P	-6.18	1.48	1.59
2	A	1501	FAD	C9A-C5X	6.09	1.55	1.42
2	A	1501	FAD	C1'-C2'	5.73	1.57	1.51
2	A	1501	FAD	C8-C7	5.72	1.57	1.40
2	A	1501	FAD	C6-C5X	5.66	1.48	1.41
2	A	1501	FAD	C4-N3	5.66	1.46	1.37
3	A	1502	NAP	PA-O3	-5.31	1.50	1.59
2	A	1501	FAD	PA-O3P	-5.29	1.50	1.59
2	A	1501	FAD	C2A-N1A	5.20	1.44	1.33
2	A	1501	FAD	C9-C9A	5.03	1.50	1.40
3	A	1502	NAP	C4A-N3A	4.86	1.43	1.35
2	A	1501	FAD	C5X-N5	4.85	1.42	1.35
2	A	1501	FAD	C2A-N3A	4.78	1.41	1.32
2	A	1501	FAD	C6-C7	4.33	1.49	1.37
3	A	1502	NAP	PN-O3	-4.33	1.51	1.60
2	A	1501	FAD	C8A-N9A	4.22	1.42	1.36
2	A	1501	FAD	C4-C4X	4.20	1.48	1.41
3	A	1502	NAP	P2B-O2B	-4.17	1.47	1.59
2	A	1501	FAD	O4B-C1B	4.06	1.47	1.41
3	A	1502	NAP	O4D-C1D	3.91	1.47	1.41
3	A	1502	NAP	C6N-N1N	3.77	1.46	1.35
3	A	1502	NAP	PN-O5D	-3.63	1.48	1.60
2	A	1501	FAD	C10-N10	3.58	1.46	1.38
2	A	1501	FAD	C10-N1	3.51	1.41	1.35
3	A	1502	NAP	C2A-N3A	3.35	1.38	1.32
2	A	1501	FAD	C5A-C4A	3.29	1.47	1.40
3	A	1502	NAP	C2A-N1A	3.27	1.40	1.33
3	A	1502	NAP	PA-O5B	-2.98	1.45	1.59
2	A	1501	FAD	PA-O5B	-2.89	1.46	1.59
2	A	1501	FAD	C5'-C4'	2.88	1.56	1.51
2	A	1501	FAD	C9-C8	2.76	1.45	1.37
2	A	1501	FAD	P-O5'	-2.72	1.46	1.59
3	A	1502	NAP	O2B-C2B	2.58	1.52	1.44
3	A	1502	NAP	O4B-C1B	2.55	1.45	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1502	NAP	P2B-O3X	-2.53	1.45	1.54
3	A	1502	NAP	O2D-C2D	-2.46	1.37	1.43
3	A	1502	NAP	P2B-O2X	-2.38	1.46	1.54
3	A	1502	NAP	O3D-C3D	-2.20	1.37	1.43
2	A	1501	FAD	PA-O2A	-2.20	1.45	1.55
2	A	1501	FAD	C6A-N1A	2.17	1.47	1.37
2	A	1501	FAD	P-O2P	-2.15	1.45	1.55
2	A	1501	FAD	O3'-C3'	2.12	1.48	1.43
2	A	1501	FAD	C6A-C5A	2.08	1.54	1.42
2	A	1501	FAD	C2B-C1B	2.04	1.56	1.53

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1501	FAD	C2-N1-C10	8.23	123.28	114.98
2	A	1501	FAD	N3A-C2A-N1A	-6.98	122.88	128.71
3	A	1502	NAP	O4B-C1B-C2B	-6.90	100.50	106.95
2	A	1501	FAD	C4X-C10-N10	-6.70	117.17	120.51
2	A	1501	FAD	C1B-N9A-C4A	-6.64	115.16	126.64
2	A	1501	FAD	O4B-C1B-N9A	-6.38	102.51	108.44
2	A	1501	FAD	C8A-N9A-C1B	6.11	138.42	126.38
2	A	1501	FAD	C4B-O4B-C1B	-6.07	103.15	109.75
3	A	1502	NAP	P2B-O2B-C2B	6.01	134.61	121.96
2	A	1501	FAD	C1'-N10-C10	-5.83	110.89	119.17
2	A	1501	FAD	C4X-N5-C5X	4.96	122.27	116.69
3	A	1502	NAP	O4B-C1B-N9A	4.62	112.74	108.44
2	A	1501	FAD	C1'-N10-C9A	4.18	122.94	118.87
3	A	1502	NAP	C6N-C5N-C4N	-3.66	113.62	119.44
3	A	1502	NAP	N3A-C2A-N1A	-3.57	125.72	128.71
3	A	1502	NAP	O7N-C7N-C3N	3.37	123.37	119.58
3	A	1502	NAP	C4N-C3N-C7N	3.15	129.49	121.10
3	A	1502	NAP	C2N-C3N-C7N	-2.95	110.58	119.35
2	A	1501	FAD	C4A-C5A-N7A	2.73	111.86	109.52
2	A	1501	FAD	C5A-C4A-N9A	-2.66	103.31	107.16
3	A	1502	NAP	O4D-C1D-N1N	2.64	110.65	107.95
3	A	1502	NAP	C1B-N9A-C4A	-2.61	122.13	126.64
2	A	1501	FAD	N3A-C4A-N9A	2.57	130.07	125.43
2	A	1501	FAD	O3B-C3B-C4B	2.55	118.59	111.08
3	A	1502	NAP	C5N-C6N-N1N	2.53	124.69	120.43
2	A	1501	FAD	C2A-N1A-C6A	2.47	123.24	118.77
2	A	1501	FAD	O4'-C4'-C3'	-2.42	103.01	109.05
2	A	1501	FAD	O5B-C5B-C4B	-2.36	100.28	108.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1502	NAP	O2D-C2D-C1D	2.32	118.23	111.23
3	A	1502	NAP	O4D-C1D-C2D	2.29	110.28	106.77
2	A	1501	FAD	C8M-C8-C9	-2.24	114.98	120.38
3	A	1502	NAP	C8A-N9A-C1B	2.23	130.77	126.38
2	A	1501	FAD	O4B-C4B-C5B	-2.11	101.83	109.36
2	A	1501	FAD	N1-C10-N10	2.08	121.44	115.97
2	A	1501	FAD	O4B-C4B-C3B	-2.08	100.96	105.17
3	A	1502	NAP	C2D-C3D-C4D	2.03	106.71	102.65
2	A	1501	FAD	C2B-C3B-C4B	-2.01	98.65	102.65

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	435/435 (100%)	0.05	23 (5%) 25 26	9, 21, 50, 62	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1071	THR	10.0
1	A	1072	ALA	8.3
1	A	1070	ASN	8.3
1	A	1074	GLY	7.8
1	A	1075	VAL	7.2
1	A	1076	ILE	6.8
1	A	1073	LEU	5.0
1	A	1069	ARG	4.6
1	A	1397	VAL	4.2
1	A	1077	SER	4.0
1	A	1078	ASN	3.5
1	A	1067	GLU	3.2
1	A	1068	GLU	3.1
1	A	966	ARG	3.0
1	A	1053	ASP	2.7
1	A	967	ASN	2.7
1	A	1080	LYS	2.5
1	A	1203	GLU	2.5
1	A	965	LYS	2.5
1	A	1317	ASP	2.2
1	A	1079	TRP	2.1
1	A	963	SER	2.0
1	A	1197	TYR	2.0

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
4	GOL	A	2081	6/6	0.18	5.08	30,34,37,46	0
4	GOL	A	2082	6/6	0.13	4.74	10,17,26,32	0
5	FMT	A	2086	3/3	0.15	4.56	42,42,44,47	0
4	GOL	A	2084	6/6	0.22	4.22	47,48,52,54	0
5	FMT	A	2087	3/3	0.14	2.89	18,18,35,42	0
5	FMT	A	2085	3/3	0.17	2.49	29,29,40,41	0
4	GOL	A	2083	6/6	0.11	1.63	22,23,24,29	0
4	GOL	A	2080	6/6	0.08	1.36	12,17,20,20	0
2	FAD	A	1501	53/53	0.11	0.08	5,15,49,54	0
3	NAP	A	1502	48/48	0.09	-0.21	8,13,30,55	0

6.5 Other polymers ⓘ

There are no such residues in this entry.