

Full wwPDB X-ray Structure Validation Report (i)

Nov 5, 2023 – 09:43 pm GMT

PDB ID	:	1W3R
Title	:	NimA from D. radiodurans with Metronidazole and Pyruvate
Authors	:	Leiros, HK.S.; Kozielski-Stuhrmann, S.; Kapp, U.; Terradot, L.; Leonard,
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Deposited on	:	2004-07-17
Resolution	:	1.90 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 https://www.wwpdb.org/validation/2017/XrayValidationReportHelp with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (1)) were used in the production of this report:

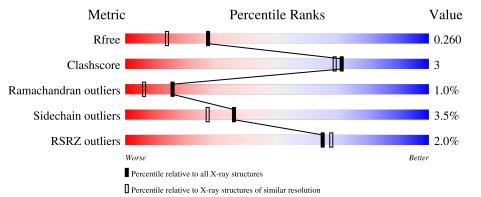
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as 541 be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\text{-}RAY\;DIFFRACTION$

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	$\begin{array}{c} \textbf{Whole archive} \\ (\#\textbf{Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
R_{free}	130704	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (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 of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. 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 <=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	А	216	83%	9%	•• 6%			



1W3R

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 1952 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 NIMA-RELATED PROTEIN.

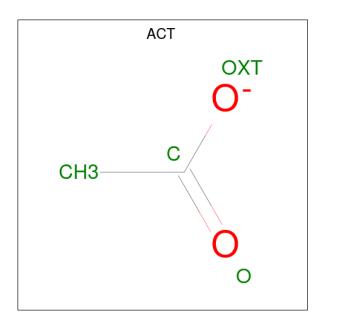
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	А	204	Total 1637	C 1030	N 297	O 309	S 1	18	0	0

Chain	Residue	Modelled	Actual	Comment	Reference
А	-21	MET	-	expression tag	UNP Q9RW27
А	-20	SER	-	expression tag	UNP Q9RW27
A	-19	TYR	-	expression tag	UNP Q9RW27
А	-18	TYR	-	expression tag	UNP Q9RW27
A	-17	HIS	-	expression tag	UNP Q9RW27
A	-16	HIS	-	expression tag	UNP Q9RW27
А	-15	HIS	-	expression tag	UNP Q9RW27
A	-14	HIS	-	expression tag	UNP Q9RW27
А	-13	HIS	-	expression tag	UNP Q9RW27
A	-12	HIS	-	expression tag	UNP Q9RW27
А	-11	LEU	-	expression tag	UNP Q9RW27
A	-10	GLU	-	expression tag	UNP Q9RW27
А	-9	SER	-	expression tag	UNP Q9RW27
А	-8	THR	-	expression tag	UNP Q9RW27
А	-7	SER	-	expression tag	UNP Q9RW27
А	-6	LEU	-	expression tag	UNP Q9RW27
А	-5	TYR	-	expression tag	UNP Q9RW27
А	-4	LYS	-	expression tag	UNP Q9RW27
А	-3	LYS	-	expression tag	UNP Q9RW27
А	-2	ALA	-	expression tag	UNP Q9RW27
А	-1	GLY	-	expression tag	UNP Q9RW27

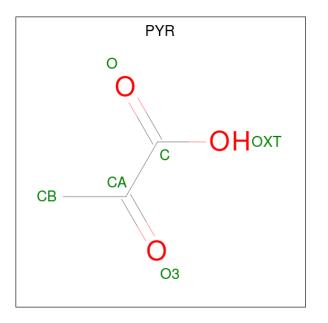
There are 21 discrepancies between the modelled and reference sequences:

• Molecule 2 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).





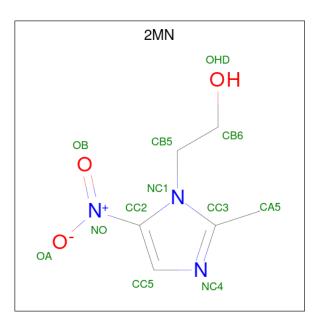
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	А	1	Total 4	$\begin{array}{c} \mathrm{C} \\ \mathrm{2} \end{array}$	O 2	0	0



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	А	1	Total 6	${ m C} { m 3}$	O 3	0	0

• Molecule 4 is Metronidazole (three-letter code: 2MN) (formula: $C_6H_9N_3O_3$).





Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	А	1	Total C N O 12 6 3 3	0	0

• Molecule 5 is water.

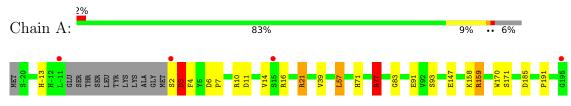
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	А	293	Total O 293 293	0	0



3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: NIMA-RELATED PROTEIN





4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants	99.59Å 39.17 Å 59.84 Å	Depositor
a, b, c, α , β , γ	90.00° 114.09° 90.00°	Depositor
Resolution (Å)	20.00 - 1.90	Depositor
Resolution (A)	19.94 - 1.90	EDS
% Data completeness	100.0 (20.00-1.90)	Depositor
(in resolution range)	$100.0\ (19.94\text{-}1.90)$	EDS
R _{merge}	0.14	Depositor
R _{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	2.54 (at 1.90\AA)	Xtriage
Refinement program	REFMAC 5.2.0003	Depositor
D D.	0.189 , 0.256	Depositor
R, R_{free}	0.192 , 0.260	DCC
R_{free} test set	852 reflections $(5.06%)$	wwPDB-VP
Wilson B-factor $(Å^2)$	15.0	Xtriage
Anisotropy	0.165	Xtriage
Bulk solvent $k_{sol}(e/Å^3), B_{sol}(Å^2)$	0.30 , 47.5	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	0.94	EDS
Total number of atoms	1952	wwPDB-VP
Average B, all atoms $(Å^2)$	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 9.20% of the height of the origin peak. No significant pseudotranslation is detected.

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.



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: ACT, PYR, $2\mathrm{MN}$

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).

Mal	Mol Chain		nd lengths	Bond angles		
NIOI	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	А	2.22	6/1682~(0.4%)	2.96	10/2290~(0.4%)	

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	А	0	2

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	$\operatorname{Ideal}(\operatorname{\AA})$
1	А	21	ARG	CZ-NH1	71.58	2.26	1.33
1	А	21	ARG	CZ-NH2	44.66	1.91	1.33
1	А	159	ARG	NE-CZ	-13.21	1.15	1.33
1	А	158	LYS	CE-NZ	-8.70	1.27	1.49
1	А	-13	HIS	CA-CB	-6.87	1.38	1.53
1	А	147	GLU	CB-CG	5.98	1.63	1.52

All (6) bond length outliers are listed below:

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(^{o})$	$Ideal(^{o})$
1	А	21	ARG	NE-CZ-NH1	-70.77	84.92	120.30
1	А	159	ARG	NE-CZ-NH1	-70.05	85.27	120.30
1	А	159	ARG	NE-CZ-NH2	55.11	147.86	120.30
1	А	21	ARG	NE-CZ-NH2	-51.18	94.71	120.30
1	А	21	ARG	NH1-CZ-NH2	-48.23	66.34	119.40
1	А	159	ARG	CD-NE-CZ	18.20	149.07	123.60
1	A	158	LYS	CB-CG-CD	11.83	142.36	111.60

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Mol	Chain	Res	Type	Atoms	Ζ	$Observed(^{o})$	$Ideal(^{o})$
1	А	185	ASP	CB-CG-OD2	6.59	124.24	118.30
1	А	3	ASP	CB-CG-OD2	5.59	123.33	118.30
1	А	158	LYS	CD-CE-NZ	5.31	123.92	111.70

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There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	А	159	ARG	Sidechain
1	А	21	ARG	Sidechain

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	А	1637	0	1565	11	0
2	А	4	0	3	0	0
3	А	6	0	0	1	0
4	А	12	0	4	0	0
5	А	293	0	0	3	0
All	All	1952	0	1572	11	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 3.

All (11) 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:71:HIS:NE2	3:A:1197:PYR:O	2.12	0.76
1:A:83:GLY:HA2	5:A:2159:HOH:O	1.96	0.64
1:A:2:SER:O	1:A:3:ASP:C	2.40	0.59
1:A:77:ARG:HB3	5:A:2154:HOH:O	2.04	0.57
1:A:91:GLU:OE2	1:A:93:SER:OG	2.16	0.52
1:A:2:SER:O	1:A:4:PHE:N	2.47	0.48
1:A:11:ASP:HB3	1:A:14:VAL:HG23	1.94	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:39:VAL:HG23	1:A:57:LEU:HD11	1.96	0.47
1:A:191:PRO:HD2	5:A:2285:HOH:O	2.15	0.46
1:A:77:ARG:HD3	1:A:77:ARG:HA	1.56	0.45
1:A:6:ASP:HB3	1:A:7:PRO:CD	2.52	0.40

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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	А	200/216~(93%)	193~(96%)	5(2%)	2(1%)	15 6

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	А	3	ASP
1	А	77	ARG

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	А	172/182~(94%)	166~(96%)	6 (4%)	36 27

All (6) residues with a non-rotameric sidechain are listed below:



Mol	Chain	Res	Type
1	А	10	ARG
1	А	16	ARG
1	А	57	LEU
1	А	77	ARG
1	А	170	TRP
1	А	171	SER

Sometimes side chains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such side chains are listed below:

Mol	Chain	Res	Type	
1	А	81	ASN	

5.3.3 RNA (i)

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

3 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	Turne	ype Chain	Res	Link	Bond lengths			Bond angles		
NIOI	Type	Unam	nes	LIIIK	Counts	RMSZ	# Z >2	Counts	RMSZ	# Z > 2
4	2MN	А	1198	-	9,12,12	3.30	3 (33%)	6,16,16	<mark>3.31</mark>	4 (66%)
2	ACT	А	1196	-	3,3,3	0.69	0	3,3,3	0.76	0
3	PYR	А	1197	-	$5,\!5,\!5$	3.07	3 (60%)	$3,\!6,\!6$	2.01	1 (33%)



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
4	2MN	А	1198	-	-	1/3/7/7	0/1/1/1
3	PYR	А	1197	-	-	0/4/4/4	-

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	$\operatorname{Ideal}(\operatorname{\AA})$
4	А	1198	2MN	CB5-CB6	-6.75	1.26	1.51
4	А	1198	2MN	CC2-NO	-5.25	1.36	1.45
4	А	1198	2MN	CB5-NC1	-4.49	1.29	1.47
3	А	1197	PYR	O3-CA	4.28	1.32	1.23
3	А	1197	PYR	CA-C	-4.07	1.39	1.54
3	А	1197	PYR	O-C	3.40	1.31	1.22

All (6) bond length outliers are listed below:

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(^{o})$	$Ideal(^{o})$
4	А	1198	2MN	CC5-CC2-NO	-4.76	123.24	126.98
4	А	1198	2MN	CB6-CB5-NC1	4.67	121.95	110.70
4	А	1198	2MN	CA5-CC3-NC4	-3.35	112.19	121.77
3	А	1197	PYR	OXT-C-CA	3.06	122.34	113.97
4	А	1198	2MN	OHD-CB6-CB5	2.38	119.89	110.82

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	А	1198	2MN	CB6-CB5-NC1-CC3

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	А	1197	PYR	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 (i)

6.1 Protein, DNA and RNA chains (i)

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^{th} 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		$OWAB(Å^2)$	$\mathbf{Q}{<}0.9$	
1	А	204/216~(94%)	-0.17	4 (1%)	65	68	9, 17, 32, 43	12~(5%)

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	А	-11	LEU	3.6
1	А	15	SER	3.1
1	А	195	GLY	2.6
1	А	2	SER	2.1

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 monosaccharides 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. The B-factors column lists the minimum, median, 95^{th} 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	$\mathbf{B} ext{-factors}(\mathbf{A}^2)$	Q<0.9
4	2MN	А	1198	12/12	0.79	0.19	$60,\!61,\!61,\!61$	0
3	PYR	А	1197	6/6	0.86	0.14	36,38,39,41	0
2	ACT	А	1196	4/4	0.96	0.13	23,23,24,25	0



6.5 Other polymers (i)

There are no such residues in this entry.

