

wwPDB X-ray Structure Validation Summary Report (i)

Jan 15, 2024 – 07:35 pm GMT

PDB ID : 6QO9

Title : Crystal structure of ribonucleotide reductase NrdF from Bacillus anthracis

soaked with manganese ions

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Deposited on : 2019-02-12

Resolution : 1.30 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 as541be (2020)

Xtriage (Phenix) : 1.13 EDS : 2.36

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

 $Refmac \quad : \quad 5.8.0158$

CCP4 : 7.0.044 (Gargrove)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

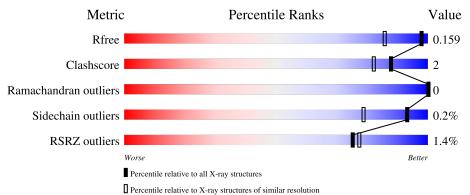
Validation Pipeline (wwPDB-VP) : 2.36

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 1.30 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 $(\#\text{Entries})$	Similar resolution $(\#\text{Entries, resolution range}(\mathring{A}))$
R_{free}	130704	1058 (1.30-1.30)
Clashscore	141614	1101 (1.30-1.30)
Ramachandran outliers	138981	1058 (1.30-1.30)
Sidechain outliers	138945	1058 (1.30-1.30)
RSRZ outliers	127900	1029 (1.30-1.30)

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	A	322	85 %		11%			
1	В	322	85%	5%	11%			



2 Entry composition (i)

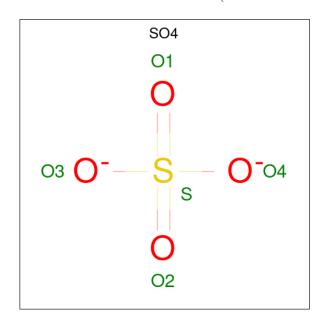
There are 5 unique types of molecules in this entry. The entry contains 10253 atoms, of which 4868 are hydrogens and 0 are deuteriums.

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 Ribonucleoside-diphosphate reductase subunit beta.

Mol	Chain	Residues		Atoms					ZeroOcc	AltConf	Trace
1	A	288	Total 4820	C 1558	H 2406	N 383	O 463	S 10	0	18	0
1	В	288	Total 4846	C 1564	Н	N 386	O 462	S 10	0	19	0

• Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total O S 5 4 1	0	0
2	A	1	Total O S 5 4 1	0	0
2	A	1	Total O S 5 4 1	0	0
2	В	1	Total O S 5 4 1	0	0

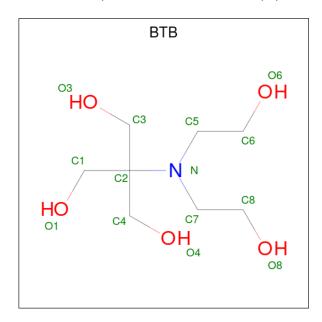
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Mol	Chain	Residues	Atom	5	ZeroOcc	AltConf
2	В	1	Total O 5 4	S 1	0	0

• Molecule 3 is 2-[BIS-(2-HYDROXY-ETHYL)-AMINO]-2-HYDROXYMETHYL-PROPAN E-1,3-DIOL (three-letter code: BTB) (formula: $C_8H_{19}NO_5$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf		
2	Λ	1	Total	С	Н	N	О	0	0	
3	A	1	33	8	19	1	5	0		
2	D	1	Total	С	Н	N	О	0	0	
)	Б	1	33	8	19	1	5	U	U	

• Molecule 4 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	2	Total Mn 2 2	0	0
4	В	2	Total Mn 2 2	0	0

• Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	247	Total O 247 247	0	0

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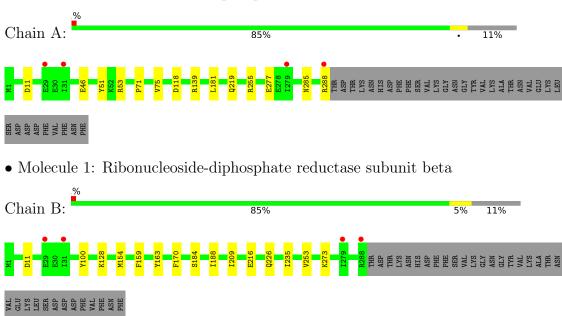
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	В	245	Total O 245 245	0	0



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: Ribonucleoside-diphosphate reductase subunit beta





4 Data and refinement statistics (i)

Property	Value	Source		
Space group	P 1 21 1	Depositor		
Cell constants	57.53Å 61.23Å 95.84Å	Depositor		
a, b, c, α , β , γ	90.00° 106.48° 90.00°	Depositor		
Resolution (Å)	28.76 - 1.30	Depositor		
resolution (A)	42.30 - 1.30	Depositor Depositor EDS Depositor EDS Depositor Depositor Depositor Xtriage Depositor Depositor Depositor Variage Depositor Decc WwPDB-VP Xtriage Xtriage EDS Xtriage Xtriage EDS Xtriage EDS Xtriage EDS WwPDB-VP		
% Data completeness	99.0 (28.76-1.30)	Depositor		
(in resolution range)	99.0 (42.30-1.30)	EDS		
R_{merge}	0.05	Depositor		
R_{sym}	(Not available)	Depositor		
$< I/\sigma(I) > 1$	1.99 (at 1.30Å)	Xtriage		
Refinement program	PHENIX 1.12_2829	Depositor		
R, R_{free}	0.136 , 0.159	Depositor		
it, it free	0.136 , 0.159	DCC		
R_{free} test set	7764 reflections (5.00%)	wwPDB-VP		
Wilson B-factor (Å ²)	10.7	0		
Anisotropy	0.488	Xtriage		
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.42, 46.6	EDS		
L-test for twinning ²	$< L > = 0.48, < L^2> = 0.31$	Xtriage		
Estimated twinning fraction	0.000 for h,-k,-h-l	Xtriage		
F_o, F_c correlation	0.98	EDS		
Total number of atoms	10253	wwPDB-VP		
Average B, all atoms (\mathring{A}^2)	17.0	wwPDB-VP		

Xtriage's analysis on translational NCS is as follows: The analyses of the Patterson function reveals a significant off-origin peak that is 77.57 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 8.2838e-07. The detected translational NCS is most likely also responsible for the elevated intensity ratio.

²Theoretical values of <|L|>, $<L^2>$ 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: BTB, MN, SO4

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		nd lengths	Bond angles		
IVIOI	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	A	0.61	$1/2531 \ (0.0\%)$	0.73	1/3422 (0.0%)	
1	В	0.59	$1/2540 \ (0.0\%)$	0.72	0/3431	
All	All	0.60	$2/5071 \ (0.0\%)$	0.72	1/6853 (0.0%)	

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(A)	Ideal(A)
1	В	216	GLU	CG-CD	5.51	1.60	1.51
1	A	46	GLU	CG-CD	5.49	1.60	1.51

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\mathbf{Observed}(^o)$	$Ideal(^{o})$
1	A	53	ARG	NE-CZ-NH1	-6.67	116.96	120.30

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	2414	2406	2336	11	0
1	В	2422	2424	2353	8	0
2	A	15	0	0	0	0

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Continued	trom	mmoninonic	maaa
COHABABACA		DIEUIUU	DUIUE
0 0 1000100000			

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	В	10	0	0	0	0
3	A	14	19	19	3	0
3	В	14	19	19	0	0
4	A	2	0	0	0	0
4	В	2	0	0	0	0
5	A	247	0	0	4	0
5	В	245	0	0	1	0
All	All	5385	4868	4727	19	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 2.

The worst 5 of 19 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$egin{aligned} & ext{Interatomic} \ & ext{distance} \ & ext{(Å)} \end{aligned}$	Clash overlap (Å)
1:B:100[B]:TYR:OH	5:B:501:HOH:O	1.78	1.00
1:A:285:ASN:OD1	1:A:288:ARG:NH1	2.23	0.71
1:A:277:GLU:OE1	5:A:501:HOH:O	2.11	0.68
1:B:184:SER:O	1:B:188[A]:ILE:HG12	2.02	0.59
1:A:71:PRO:O	1:A:75[B]:VAL:HG13	2.03	0.58

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	Perce	ntiles
1	A	$304/322 \ (94\%)$	301 (99%)	3 (1%)	0	100	100
1	В	305/322~(95%)	303 (99%)	2 (1%)	0	100	100
All	All	609/644 (95%)	604 (99%)	5 (1%)	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	270/283 (95%)	268 (99%)	2 (1%)	84 61		
1	В	271/283 (96%)	271 (100%)	0	100 100		
All	All	541/566 (96%)	539 (100%)	2 (0%)	93 76		

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	219[A]	GLN
1	A	219[B]	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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)

Of 11 ligands modelled in this entry, 4 are monoatomic - leaving 7 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	Tuno	Chain	hain Res Link Bond lengths				ths	Bond angles		
MIOI	Type	Chain	nes	Lilik	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	SO4	A	402	-	4,4,4	0.09	0	6,6,6	0.31	0
2	SO4	A	403	-	4,4,4	0.11	0	6,6,6	0.19	0
2	SO4	A	401	-	4,4,4	0.17	0	6,6,6	0.28	0
3	ВТВ	В	401	-	13,13,13	1.15	1 (7%)	7,16,16	0.44	0
2	SO4	В	402	-	4,4,4	0.17	0	6,6,6	0.23	0
3	BTB	A	404	-	13,13,13	0.54	0	7,16,16	0.75	0
2	SO4	В	403	-	4,4,4	0.17	0	6,6,6	0.15	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
3	BTB	В	401	-	-	1/21/21/21	-
3	BTB	A	404	_	-	7/21/21/21	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(\text{\AA})$	$Ideal(\AA)$
3	В	401	BTB	C4-C2	2.90	1.57	1.53

There are no bond angle outliers.

There are no chirality outliers.

5 of 8 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	404	BTB	C3-C2-C4-O4
3	A	404	BTB	N-C2-C4-O4
3	A	404	BTB	C1-C2-N-C5
3	A	404	BTB	C3-C2-N-C5
3	A	404	BTB	C4-C2-N-C5

There are no ring outliers.

1 monomer is involved in 3 short contacts:



Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	404	BTB	3	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 $>$	$\#\mathrm{RSRZ}{>}2$	$OWAB(Å^2)$	Q<0.9
1	A	288/322 (89%)	-0.25	4 (1%) 75 77	7, 15, 26, 44	0
1	В	288/322 (89%)	-0.27	4 (1%) 75 77	8, 14, 25, 39	0
All	All	576/644 (89%)	-0.26	8 (1%) 75 77	7, 14, 26, 44	0

The worst 5 of 8 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	279	ILE	3.5
1	В	31	ILE	3.2
1	В	29	GLU	2.6
1	A	288	ARG	2.5
1	A	29	GLU	2.3

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	$\operatorname{B-factors}(\mathring{\mathrm{A}}^2)$	Q < 0.9
2	SO4	A	403	5/5	0.68	0.18	53,55,55,55	5
2	SO4	В	403	5/5	0.68	0.24	62,64,64,64	0
2	SO4	В	402	5/5	0.79	0.15	62,62,62,62	0
2	SO4	A	402	5/5	0.89	0.13	61,61,61,61	0
3	BTB	A	404	14/14	0.90	0.12	20,29,39,42	33
3	BTB	В	401	14/14	0.96	0.09	16,20,29,31	0
2	SO4	A	401	5/5	1.00	0.04	12,13,14,16	0
4	MN	A	405	1/1	1.00	0.11	6,6,6,6	0
4	MN	A	406	1/1	1.00	0.12	6,6,6,6	0
4	MN	В	404	1/1	1.00	0.11	5,5,5,5	0
4	MN	В	405	1/1	1.00	0.11	5,5,5,5	0

6.5 Other polymers (i)

There are no such residues in this entry.

