



## Full wwPDB X-ray Structure Validation Report ⓘ

Aug 15, 2018 – 12:21 PM EDT

PDB ID : 6EDR  
Title : Crystal Structure of Human CD38 in Complex with 4'-Thioribose NAD<sup>+</sup>  
Deposited on : 2018-08-10  
Resolution : 2.40 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report.

This report is produced by the wwPDB biocuration pipeline after annotation of the structure.

We welcome your comments at [validation@mail.wwpdb.org](mailto: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 ⓘ symbol.

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

MolProbity : 4.02b-467  
Mogul : 1.7.3 (157068), CSD as539be (2018)  
Xtriage (Phenix) : 1.13  
EDS : rb-20031172  
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)  
Refmac : 5.8.0158  
CCP4 : 7.0 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : rb-20031172

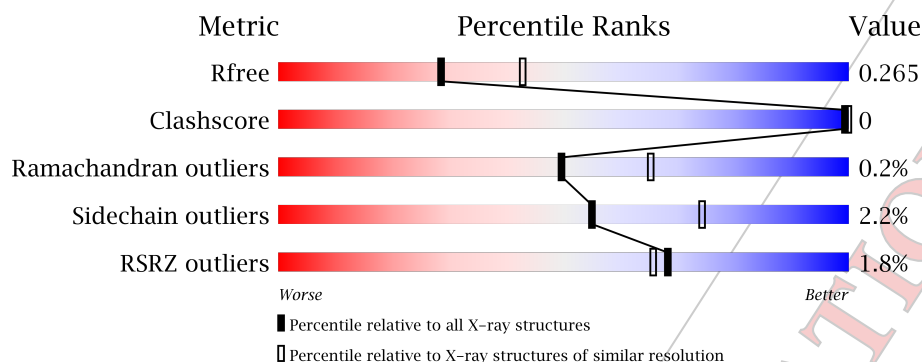
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

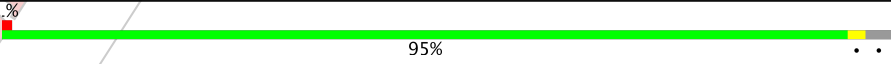
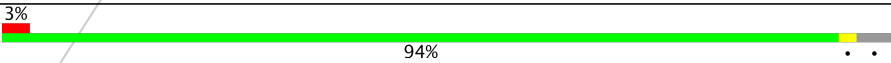
The reported resolution of this entry is 2.40 Å.

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))
$R_{free}$	111664	3481 (2.40-2.40)
Clashscore	122126	3956 (2.40-2.40)
Ramachandran outliers	120053	3897 (2.40-2.40)
Sidechain outliers	120020	3898 (2.40-2.40)
RSRZ outliers	108989	3386 (2.40-2.40)

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  $\geq 3$ , 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	257	 95%
1	B	257	 94%

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 4083 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.

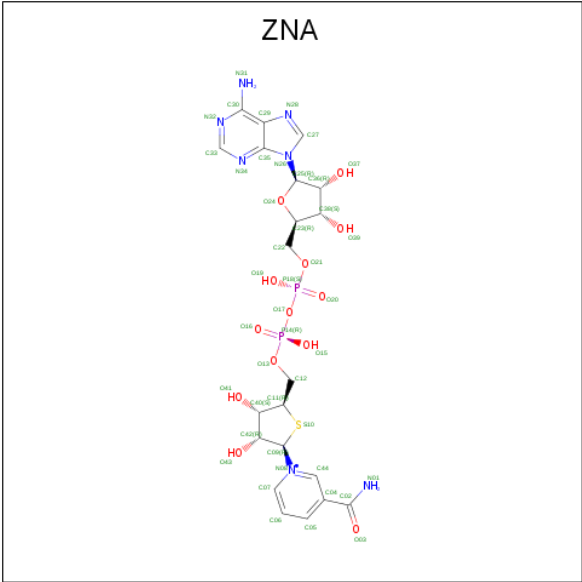
- Molecule 1 is a protein called ADP-ribosyl cyclase/cyclic ADP-ribose hydrolase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	249	Total	C	N	O	S	0	2	0
			1984	1265	342	362	15			
1	B	246	Total	C	N	O	S	0	1	0
			1924	1228	325	356	15			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	44	SER	-	expression tag	UNP P28907
A	100	ASP	ASN	engineered mutation	UNP P28907
A	164	ALA	ASN	engineered mutation	UNP P28907
A	209	ASP	ASN	engineered mutation	UNP P28907
A	219	ASP	ASN	engineered mutation	UNP P28907
B	44	SER	-	expression tag	UNP P28907
B	100	ASP	ASN	engineered mutation	UNP P28907
B	164	ALA	ASN	engineered mutation	UNP P28907
B	209	ASP	ASN	engineered mutation	UNP P28907
B	219	ASP	ASN	engineered mutation	UNP P28907

- Molecule 2 is [(2 {R},3 {S},4 {R},5 {R})-5-(3-aminocarbonylpyridin-1-yl)-3,4-bis(oxidanyl)thiolan-2-yl]methoxy-oxidanyl-phosphoryl [(2 {R},3 {S},4 {R},5 {R})-5-(6-aminopurin-9-yl)-3,4-bis(oxidanyl)oxolan-2-yl]methyl hydrogen phosphate (three-letter code: ZNA) (formula: C<sub>21</sub>H<sub>28</sub>N<sub>7</sub>O<sub>13</sub>P<sub>2</sub>S).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	S	0	0
			44	21	7	13	2	1		
2	B	1	Total	C	N	O	P	S	0	0
			44	21	7	13	2	1		

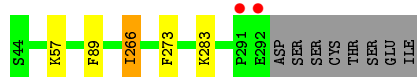
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	49	Total	O	0	0
			49	49		
3	B	38	Total	O	0	0
			38	38		

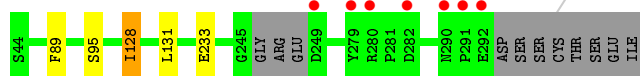
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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: ADP-ribosyl cyclase/cyclic ADP-ribose hydrolase 1



- Molecule 1: ADP-ribosyl cyclase/cyclic ADP-ribose hydrolase 1



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	57.74Å 51.11Å 100.67Å 90.00° 96.89° 90.00°	Depositor
Resolution (Å)	30.00 – 2.40 29.22 – 2.39	Depositor EDS
% Data completeness (in resolution range)	99.7 (30.00-2.40) 99.4 (29.22-2.39)	Depositor EDS
$R_{merge}$	0.37	Depositor
$R_{sym}$	0.37	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.66 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.8.0158	Depositor
R, $R_{free}$	0.216 , 0.262 0.220 , 0.265	Depositor DCC
$R_{free}$ test set	1181 reflections (5.11%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	34.3	Xtriage
Anisotropy	1.101	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 31.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	4083	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.53% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup> Intensities estimated from amplitudes.

<sup>2</sup> Theoretical 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: ZNA

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	0.49	0/2038	0.74	0/2765
1	B	0.49	0/1977	0.66	0/2691
All	All	0.49	0/4015	0.70	0/5456

There are no bond length outliers.

There are no bond angle outliers.

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	1984	0	1895	1	0
1	B	1924	0	1785	0	0
2	A	44	0	0	0	0
2	B	44	0	0	0	0
3	A	49	0	0	0	0
3	B	38	0	0	0	0
All	All	4083	0	3680	1	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 0.

All (1) 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:266:ILE:HD11	1:A:273:PHE:CD1	2.46	0.51

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	A	249/257 (97%)	242 (97%)	7 (3%)	0	100	100
1	B	243/257 (95%)	236 (97%)	6 (2%)	1 (0%)	36	51
All	All	492/514 (96%)	478 (97%)	13 (3%)	1 (0%)	49	65

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	128	ILE

### 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	213/237 (90%)	209 (98%)	4 (2%)	60	78
1	B	201/237 (85%)	196 (98%)	5 (2%)	50	71
All	All	414/474 (87%)	405 (98%)	9 (2%)	55	74



All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	57	LYS
1	A	89	PHE
1	A	266	ILE
1	A	283	LYS
1	B	89	PHE
1	B	95	SER
1	B	128	ILE
1	B	131	LEU
1	B	233	GLU

Some 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 carbohydrates in this entry.

### 5.6 Ligand geometry [i](#)

2 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	ZNA	A	401	-	43,48,48	2.56	14 (32%)	49,73,73	3.90	14 (28%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	ZNA	B	401	-	43,48,48	2.60	13 (30%)	49,73,73	3.75	18 (36%)

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	ZNA	A	401	-	-	0/24/72/72	0/5/5/5
2	ZNA	B	401	-	-	0/24/72/72	0/5/5/5

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	ZNA	C44-C04	-8.49	1.39	1.53
2	A	401	ZNA	C44-C04	-7.56	1.41	1.53
2	B	401	ZNA	C04-C02	-6.96	1.40	1.52
2	A	401	ZNA	C04-C02	-6.00	1.42	1.52
2	B	401	ZNA	C06-C05	-4.94	1.40	1.53
2	A	401	ZNA	C06-C05	-4.90	1.40	1.53
2	B	401	ZNA	C05-C04	-4.79	1.40	1.53
2	A	401	ZNA	C05-C04	-4.50	1.41	1.53
2	B	401	ZNA	C44-N08	-4.18	1.40	1.47
2	A	401	ZNA	C07-N08	-3.55	1.40	1.47
2	A	401	ZNA	C44-N08	-3.43	1.41	1.47
2	A	401	ZNA	C06-C07	-3.16	1.39	1.51
2	B	401	ZNA	C06-C07	-3.09	1.40	1.51
2	A	401	ZNA	C27-N26	-2.80	1.33	1.36
2	B	401	ZNA	C27-N26	-2.67	1.33	1.36
2	B	401	ZNA	C29-C35	-2.58	1.34	1.40
2	B	401	ZNA	C07-N08	-2.54	1.42	1.47
2	A	401	ZNA	C29-C35	-2.37	1.35	1.40
2	A	401	ZNA	C09-C42	2.38	1.57	1.54
2	B	401	ZNA	C09-N08	2.57	1.49	1.46
2	A	401	ZNA	C09-N08	2.66	1.49	1.46
2	B	401	ZNA	O24-C25	3.25	1.45	1.41
2	B	401	ZNA	C33-N32	3.79	1.41	1.33
2	A	401	ZNA	O24-C25	3.82	1.46	1.41
2	A	401	ZNA	C33-N32	3.85	1.41	1.33
2	B	401	ZNA	C33-N34	5.01	1.40	1.32
2	A	401	ZNA	C33-N34	5.69	1.41	1.32

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	ZNA	N34-C33-N32	-12.28	118.35	128.86
2	B	401	ZNA	N34-C33-N32	-11.79	118.77	128.86
2	A	401	ZNA	S10-C09-N08	-8.96	103.03	112.85
2	B	401	ZNA	S10-C09-N08	-7.75	104.36	112.85
2	A	401	ZNA	C23-O24-C25	-4.46	105.18	109.83
2	B	401	ZNA	C42-C09-S10	-3.93	101.42	106.11
2	A	401	ZNA	C42-C09-S10	-3.20	102.29	106.11
2	B	401	ZNA	C23-O24-C25	-2.76	106.95	109.83
2	B	401	ZNA	C35-C29-N28	-2.31	107.18	109.41
2	B	401	ZNA	O43-C42-C09	-2.24	107.05	111.56
2	B	401	ZNA	P18-O17-P14	-2.14	125.44	132.63
2	A	401	ZNA	C35-C29-N28	-2.06	107.42	109.41
2	B	401	ZNA	C25-N26-C35	-2.03	123.12	126.64
2	B	401	ZNA	C07-N08-C09	2.14	125.59	114.87
2	A	401	ZNA	C05-C04-C44	4.21	116.88	109.80
2	A	401	ZNA	C09-S10-C11	4.38	98.79	94.29
2	A	401	ZNA	C06-C07-N08	4.45	119.00	111.21
2	B	401	ZNA	C06-C07-N08	4.48	119.05	111.21
2	B	401	ZNA	C06-C05-C04	5.07	121.04	111.36
2	A	401	ZNA	C05-C04-C02	5.21	119.79	110.76
2	B	401	ZNA	C05-C04-C44	5.28	118.69	109.80
2	B	401	ZNA	C09-S10-C11	5.28	99.71	94.29
2	A	401	ZNA	C06-C05-C04	5.32	121.50	111.36
2	B	401	ZNA	C44-C04-C02	5.64	120.67	110.07
2	B	401	ZNA	C05-C04-C02	5.70	120.64	110.76
2	A	401	ZNA	C44-C04-C02	7.06	123.33	110.07
2	B	401	ZNA	C05-C06-C07	7.31	120.31	110.95
2	B	401	ZNA	C04-C44-N08	7.40	121.18	111.16
2	A	401	ZNA	C05-C06-C07	7.79	120.93	110.95
2	A	401	ZNA	C04-C44-N08	7.89	121.85	111.16
2	B	401	ZNA	C07-N08-C44	12.10	119.73	108.17
2	A	401	ZNA	C07-N08-C44	12.20	119.83	108.17

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 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.

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## 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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	249/257 (96%)	-0.40	2 (0%) 86 84	29, 44, 68, 97	0
1	B	246/257 (95%)	-0.24	7 (2%) 53 50	28, 48, 96, 133	0
All	All	495/514 (96%)	-0.32	9 (1%) 68 66	28, 46, 79, 133	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	249	ASP	5.0
1	B	290	ASN	4.4
1	A	292	GLU	4.3
1	B	282	ASP	3.5
1	B	291	PRO	2.7
1	B	279	TYR	2.4
1	A	291	PRO	2.3
1	B	280	ARG	2.2
1	B	292	GLU	2.2

### 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. The B-factors column lists the minimum,

median, 95<sup>th</sup> 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	B-factors(Å <sup>2</sup> )	Q<0.9
2	ZNA	A	401	44/44	0.59	0.36	73,138,169,172	0
2	ZNA	B	401	44/44	0.83	0.24	61,94,155,163	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.