Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:06 PM GMT

PDB ID	:	1NKD
Title	:	ATOMIC RESOLUTION (1.07 ANGSTROMS) STRUCTURE OF THE ROP MUTANT <2AA>
Authors	:	Vlassi, M.; Kokkinidis, M.
Deposited on	:	1997-09-23
Resolution	:	1.09 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
http://wwpdb.org/validation/2016/XrayValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

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

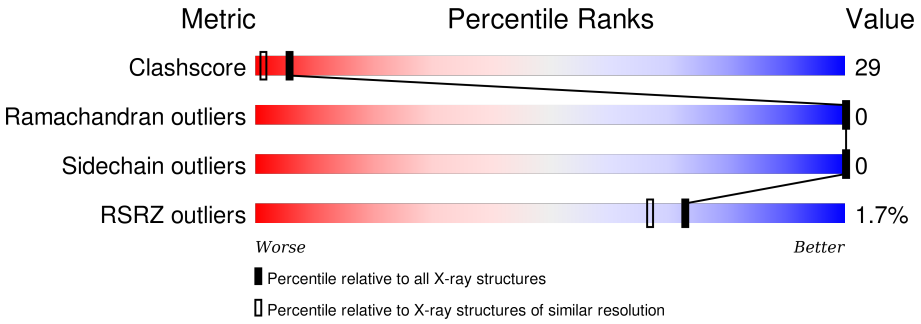
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.09 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))
Clashscore	102246	1055 (1.14-1.06)
Ramachandran outliers	100387	1016 (1.14-1.06)
Sidechain outliers	100360	1014 (1.14-1.06)
RSRZ outliers	91569	1009 (1.14-1.06)

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. 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	65	2% 52% 23% 15% 9%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 1063 atoms, of which 439 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 ROP.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	59	949	315	439	87	103	5	0	10	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	30	ALA	-	INSERTION	UNP P03051
A	32	ALA	-	INSERTION	UNP P03051

- Molecule 2 is water.

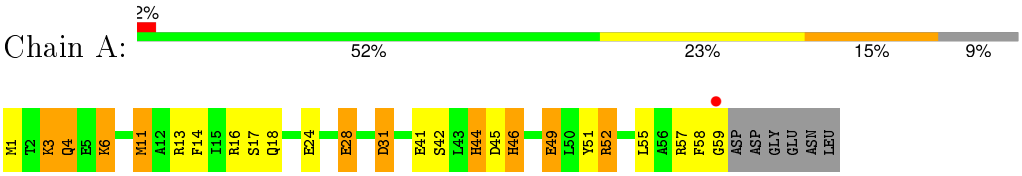
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	114	114	114	0	0

3 Residue-property plots (i)

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: ROP

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	47.06Å 37.88Å 31.65Å 90.00° 100.80° 90.00°	Depositor
Resolution (Å)	8.00 – 1.09 7.88 – 1.10	Depositor EDS
% Data completeness (in resolution range)	98.2 (8.00-1.09) 98.4 (7.88-1.10)	Depositor EDS
R_{merge}	0.45	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.73 (at 1.10Å)	Xtriage
Refinement program	SHELXL-93	Depositor
R, R_{free}	0.101 , 0.134 0.160 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å^2)	11.8	Xtriage
Anisotropy	0.456	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.30 , 67.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.51$, $< L^2 > = 0.34$	Xtriage
Outliers	0 of 21972 reflections	Xtriage
F_o,F_c correlation	0.98	EDS
Total number of atoms	1063	wwPDB-VP
Average B, all atoms (Å^2)	21.0	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 23.47 % 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 4.6360e-03.The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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	3.94	56/573 (9.8%)	2.56	46/765 (6.0%)

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	A	0	1

All (56) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	49[A]	GLU	CD-OE1	35.33	1.64	1.25
1	A	49[B]	GLU	CD-OE1	35.33	1.64	1.25
1	A	52	ARG	CZ-NH1	25.10	1.65	1.33
1	A	41[A]	GLU	CG-CD	-19.50	1.22	1.51
1	A	41[B]	GLU	CG-CD	-19.50	1.22	1.51
1	A	4	GLN	CD-NE2	17.88	1.77	1.32
1	A	59	GLY	C-O	15.53	1.48	1.23
1	A	13	ARG	CZ-NH2	15.24	1.52	1.33
1	A	6[A]	LYS	CD-CE	-13.16	1.18	1.51
1	A	6[B]	LYS	CD-CE	-13.16	1.18	1.51
1	A	46[A]	HIS	CG-CD2	12.21	1.56	1.35
1	A	46[B]	HIS	CG-CD2	12.21	1.56	1.35
1	A	59	GLY	CA-C	12.11	1.71	1.51
1	A	41[A]	GLU	CD-OE1	11.53	1.38	1.25
1	A	41[B]	GLU	CD-OE1	11.53	1.38	1.25
1	A	57	ARG	CD-NE	10.39	1.64	1.46
1	A	18	GLN	CD-NE2	-8.98	1.10	1.32
1	A	3[A]	LYS	CE-NZ	8.85	1.71	1.49
1	A	3[B]	LYS	CE-NZ	8.85	1.71	1.49
1	A	3[C]	LYS	CE-NZ	8.85	1.71	1.49

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	41[A]	GLU	CD-OE2	-8.81	1.16	1.25
1	A	41[B]	GLU	CD-OE2	-8.81	1.16	1.25
1	A	4	GLN	CD-OE1	-8.52	1.05	1.24
1	A	52	ARG	CD-NE	-8.31	1.32	1.46
1	A	16	ARG	CD-NE	-7.93	1.32	1.46
1	A	49[A]	GLU	CG-CD	-7.86	1.40	1.51
1	A	49[B]	GLU	CG-CD	-7.86	1.40	1.51
1	A	1	MET	CA-CB	-7.75	1.36	1.53
1	A	28[A]	GLU	CB-CG	-7.42	1.38	1.52
1	A	28[B]	GLU	CB-CG	-7.42	1.38	1.52
1	A	51	TYR	CE1-CZ	-7.29	1.29	1.38
1	A	31[A]	ASP	CB-CG	-6.87	1.37	1.51
1	A	31[B]	ASP	CB-CG	-6.87	1.37	1.51
1	A	1	MET	CB-CG	-6.56	1.30	1.51
1	A	44	HIS	CD2-NE2	-6.49	1.23	1.38
1	A	46[A]	HIS	CD2-NE2	-6.37	1.24	1.38
1	A	46[B]	HIS	CD2-NE2	-6.37	1.24	1.38
1	A	51	TYR	CE2-CZ	6.33	1.46	1.38
1	A	17[A]	SER	CB-OG	-6.24	1.34	1.42
1	A	17[B]	SER	CB-OG	-6.24	1.34	1.42
1	A	17[C]	SER	CB-OG	-6.24	1.34	1.42
1	A	24	GLU	CD-OE1	5.94	1.32	1.25
1	A	13	ARG	CD-NE	-5.93	1.36	1.46
1	A	55	LEU	CB-CG	-5.63	1.36	1.52
1	A	13	ARG	NE-CZ	-5.62	1.25	1.33
1	A	44	HIS	CE1-NE2	-5.45	1.20	1.32
1	A	44	HIS	ND1-CE1	5.40	1.48	1.34
1	A	42	SER	CA-CB	-5.36	1.45	1.52
1	A	11[A]	MET	SD-CE	5.34	2.07	1.77
1	A	11[B]	MET	SD-CE	5.34	2.07	1.77
1	A	28[A]	GLU	CD-OE1	5.32	1.31	1.25
1	A	28[B]	GLU	CD-OE1	5.32	1.31	1.25
1	A	6[A]	LYS	CE-NZ	-5.29	1.35	1.49
1	A	6[B]	LYS	CE-NZ	-5.29	1.35	1.49
1	A	59	GLY	N-CA	-5.19	1.38	1.46
1	A	14	PHE	CD2-CE2	5.05	1.49	1.39

All (46) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	13	ARG	NE-CZ-NH2	-21.53	109.54	120.30
1	A	52	ARG	NH1-CZ-NH2	-16.96	100.75	119.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	52	ARG	NE-CZ-NH2	16.15	128.38	120.30
1	A	41[A]	GLU	OE1-CD-OE2	-14.62	105.76	123.30
1	A	41[B]	GLU	OE1-CD-OE2	-14.62	105.76	123.30
1	A	41[A]	GLU	CG-CD-OE2	12.13	142.55	118.30
1	A	41[B]	GLU	CG-CD-OE2	12.13	142.55	118.30
1	A	6[A]	LYS	CD-CE-NZ	11.30	137.68	111.70
1	A	6[B]	LYS	CD-CE-NZ	11.30	137.68	111.70
1	A	41[A]	GLU	CB-CG-CD	10.28	141.95	114.20
1	A	41[B]	GLU	CB-CG-CD	10.28	141.95	114.20
1	A	45	ASP	CB-CG-OD2	9.73	127.06	118.30
1	A	31[A]	ASP	CB-CG-OD1	-9.16	110.05	118.30
1	A	31[B]	ASP	CB-CG-OD1	-9.16	110.05	118.30
1	A	13	ARG	NE-CZ-NH1	9.08	124.84	120.30
1	A	1	MET	CA-CB-CG	8.88	128.39	113.30
1	A	6[A]	LYS	CG-CD-CE	8.39	137.06	111.90
1	A	6[B]	LYS	CG-CD-CE	8.39	137.06	111.90
1	A	45	ASP	OD1-CG-OD2	-7.96	108.17	123.30
1	A	57	ARG	CA-CB-CG	-7.40	97.12	113.40
1	A	45	ASP	CB-CG-OD1	7.17	124.75	118.30
1	A	57	ARG	NE-CZ-NH1	6.93	123.77	120.30
1	A	49[A]	GLU	OE1-CD-OE2	-6.80	115.14	123.30
1	A	49[B]	GLU	OE1-CD-OE2	-6.80	115.14	123.30
1	A	31[A]	ASP	CB-CG-OD2	6.79	124.41	118.30
1	A	31[B]	ASP	CB-CG-OD2	6.79	124.41	118.30
1	A	59	GLY	N-CA-C	-6.68	96.41	113.10
1	A	46[A]	HIS	N-CA-CB	-6.66	98.60	110.60
1	A	46[B]	HIS	N-CA-CB	-6.66	98.60	110.60
1	A	3[A]	LYS	CD-CE-NZ	-6.64	96.42	111.70
1	A	3[B]	LYS	CD-CE-NZ	-6.64	96.42	111.70
1	A	3[C]	LYS	CD-CE-NZ	-6.64	96.42	111.70
1	A	52	ARG	NE-CZ-NH1	5.93	123.27	120.30
1	A	18	GLN	CG-CD-OE1	-5.75	110.09	121.60
1	A	41[A]	GLU	CG-CD-OE1	-5.75	106.81	118.30
1	A	41[B]	GLU	CG-CD-OE1	-5.75	106.81	118.30
1	A	1	MET	CB-CG-SD	5.67	129.40	112.40
1	A	13	ARG	NH1-CZ-NH2	5.50	125.45	119.40
1	A	58	PHE	CB-CG-CD2	5.20	124.44	120.80
1	A	44	HIS	CE1-NE2-CD2	5.06	119.25	106.60
1	A	28[A]	GLU	CA-CB-CG	5.04	124.49	113.40
1	A	28[B]	GLU	CA-CB-CG	5.04	124.49	113.40
1	A	46[A]	HIS	CB-CG-CD2	-5.02	115.24	130.80
1	A	46[B]	HIS	CB-CG-CD2	-5.02	115.24	130.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	28[A]	GLU	CG-CD-OE1	-5.01	108.29	118.30
1	A	28[B]	GLU	CG-CD-OE1	-5.01	108.29	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	52	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	A	510	439	507	30	0
2	A	114	0	0	18	0
All	All	624	439	507	30	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 29.

All (30) 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:3[B]:LYS:NZ	1:A:3[B]:LYS:CE	1.71	1.51
1:A:11[B]:MET:SD	1:A:11[B]:MET:CE	2.08	1.41
1:A:11[A]:MET:SD	1:A:11[A]:MET:CE	2.07	1.41
1:A:4:GLN:CD	1:A:4:GLN:NE2	1.77	1.37
1:A:49[A]:GLU:OE1	1:A:49[A]:GLU:CD	1.64	1.34
1:A:49[B]:GLU:CD	1:A:49[B]:GLU:OE2	1.64	1.34
1:A:3[C]:LYS:NZ	2:A:110:HOH:O	1.61	1.27
1:A:3[B]:LYS:CE	2:A:127:HOH:O	1.87	1.18
1:A:31[A]:ASP:OD1	2:A:172:HOH:O	1.65	1.13
1:A:28[A]:GLU:OE1	2:A:174:HOH:O	1.67	1.12
1:A:31[A]:ASP:OD1	2:A:171:HOH:O	1.69	1.10
1:A:3[B]:LYS:HD3	2:A:149:HOH:O	1.59	1.03

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:46[B]:HIS:CG	2:A:134:HOH:O	2.13	1.00
1:A:3[B]:LYS:CE	2:A:110:HOH:O	2.24	0.85
1:A:6[B]:LYS:HE3	2:A:110:HOH:O	1.83	0.78
1:A:3[B]:LYS:NZ	1:A:3[B]:LYS:CD	2.49	0.75
1:A:3[A]:LYS:NZ	2:A:137:HOH:O	2.24	0.69
1:A:4:GLN:OE1	2:A:143:HOH:O	2.11	0.68
1:A:44:HIS:HE1	2:A:71:HOH:O	1.78	0.66
1:A:3[B]:LYS:HE3	2:A:110:HOH:O	1.91	0.64
1:A:3[C]:LYS:HD3	2:A:149:HOH:O	1.98	0.62
1:A:11[A]:MET:CE	1:A:11[A]:MET:HB2	2.34	0.57
1:A:3[B]:LYS:HE2	2:A:110:HOH:O	1.96	0.56
1:A:3[B]:LYS:CD	2:A:149:HOH:O	2.34	0.53
1:A:6[B]:LYS:CE	2:A:110:HOH:O	2.49	0.51
1:A:11[A]:MET:CE	1:A:11[A]:MET:CB	2.91	0.49
1:A:44:HIS:HD2	2:A:81:HOH:O	1.98	0.47
1:A:3[B]:LYS:HE3	1:A:6[B]:LYS:HE2	1.99	0.44
1:A:4:GLN:CG	1:A:4:GLN:NE2	2.73	0.40
1:A:3[B]:LYS:HE3	1:A:6[B]:LYS:CE	2.51	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	69/65 (106%)	68 (99%)	1 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	62/55 (113%)	62 (100%)	0	100 100

There are no protein residues with a non-rotameric sidechain to report.

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

Mol	Chain	Res	Type
1	A	44	HIS

5.3.3 RNA

There are no RNA molecules 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

There are no ligands in this entry.

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, 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)	Q<0.9
1	A	59/65 (90%)	-0.01	1 (1%) 73 67	10, 16, 23, 37	2 (3%)

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	59	GLY	3.0

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

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.