wwPDB X-ray Structure Validation Summary Report (i)

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PDB ID	:	3NED
Title	:	mRouge
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Deposited on	:	2010-06-08
Resolution	:	0.95 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:

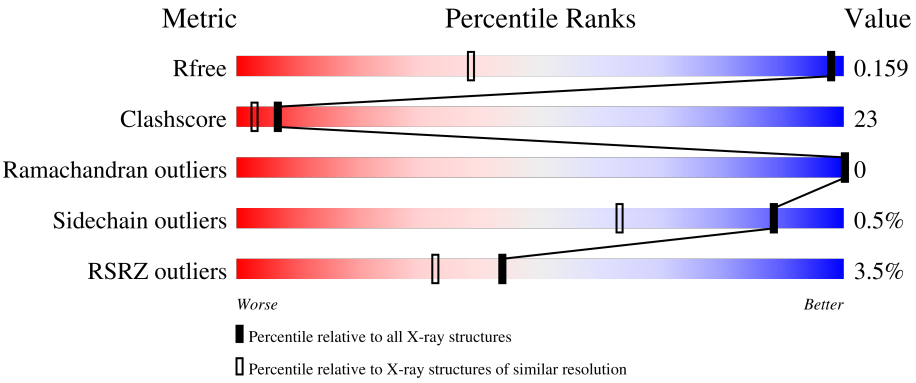
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, 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-RAY DIFFRACTION

The reported resolution of this entry is 0.95 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}	130704	1243 (1.06-0.86)
Clashscore	141614	1321 (1.06-0.86)
Ramachandran outliers	138981	1233 (1.06-0.86)
Sidechain outliers	138945	1235 (1.06-0.86)
RSRZ outliers	127900	1209 (1.06-0.86)

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	244	3% 73% 17% 5% 5%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
1	EYG	A	67[C]	X	-	-	-

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 2617 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 PAmCherry1 protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	229	2140	1367	346	411	16	11	45	0

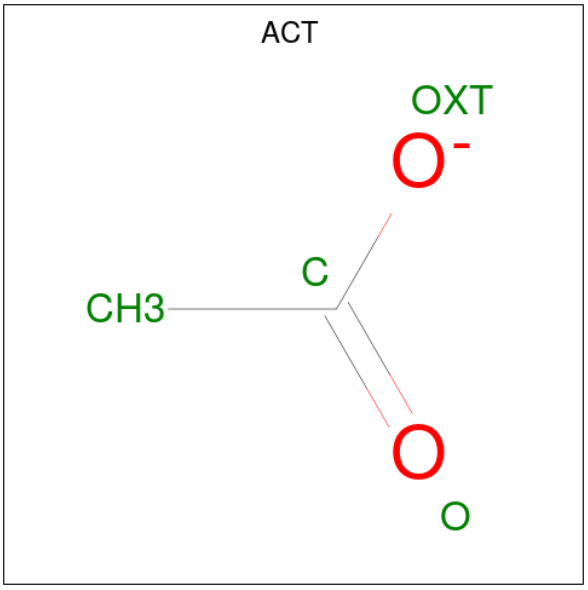
There are 25 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-12	MET	-	expression tag	UNP D1MPT3
A	-11	GLY	-	expression tag	UNP D1MPT3
A	-10	HIS	-	expression tag	UNP D1MPT3
A	-9	HIS	-	expression tag	UNP D1MPT3
A	-8	HIS	-	expression tag	UNP D1MPT3
A	-7	HIS	-	expression tag	UNP D1MPT3
A	-6	HIS	-	expression tag	UNP D1MPT3
A	-5	HIS	-	expression tag	UNP D1MPT3
A	-4	GLY	-	expression tag	UNP D1MPT3
A	16	THR	VAL	engineered mutation	UNP D1MPT3
A	26	GLU	VAL	engineered mutation	UNP D1MPT3
A	57	ALA	THR	engineered mutation	UNP D1MPT3
A	67	NRQ	MET	chromophore	UNP D1MPT3
A	67	NRQ	TYR	chromophore	UNP D1MPT3
A	67	NRQ	GLY	chromophore	UNP D1MPT3
A	70	LYS	ASN	engineered mutation	UNP D1MPT3
A	83	LEU	PHE	engineered mutation	UNP D1MPT3
A	98	ASN	LYS	engineered mutation	UNP D1MPT3
A	146	CYS	LEU	engineered mutation	UNP D1MPT3
A	161	MET	VAL	engineered mutation	UNP D1MPT3
A	163	MET	PRO	engineered mutation	UNP D1MPT3
A	165	LEU	VAL	engineered mutation	UNP D1MPT3
A	195	THR	VAL	engineered mutation	UNP D1MPT3
A	197	THR	ARG	engineered mutation	UNP D1MPT3
A	217	ASN	ALA	engineered mutation	UNP D1MPT3

- Molecule 2 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	3	Total Na 3 3	0	0

- Molecule 3 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 4 2 2	0	0

- Molecule 4 is water.

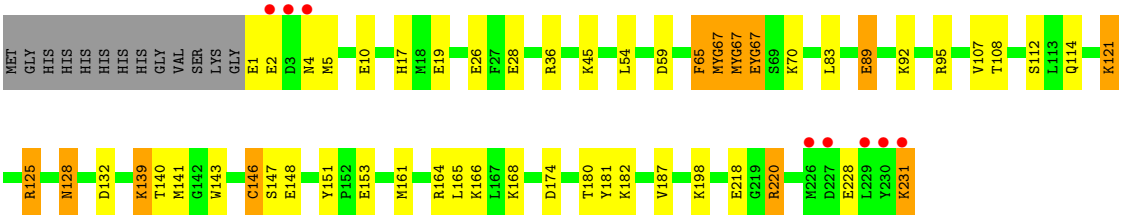
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	414	Total O 470 470	0	46

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: PAmCherry1 protein

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	48.81Å 42.76Å 61.20Å 90.00° 112.35° 90.00°	Depositor
Resolution (Å)	56.60 – 0.95 56.61 – 0.95	Depositor EDS
% Data completeness (in resolution range)	91.9 (56.60-0.95) 92.0 (56.61-0.95)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.65 (at 0.95Å)	Xtriage
Refinement program	REFMAC 5.5.0088	Depositor
R, R_{free}	0.133 , 0.157 0.135 , 0.159	Depositor DCC
R_{free} test set	6732 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å^2)	5.8	Xtriage
Anisotropy	1.569	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 53.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	0.023 for h,-k,-h-l	Xtriage
F_o,F_c correlation	0.98	EDS
Total number of atoms	2617	wwPDB-VP
Average B, all atoms (Å^2)	12.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.04% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical 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.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CH6, NA, EYG, NRQ, ACT

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	1.39	21/2253 (0.9%)	1.19	8/3021 (0.3%)

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	1	0

The worst 5 of 21 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	65[A]	PHE	CA-C	13.57	1.88	1.52
1	A	65[B]	PHE	CA-C	13.57	1.88	1.52
1	A	65[C]	PHE	CA-C	13.57	1.88	1.52
1	A	146[A]	CYS	CB-SG	8.09	1.96	1.82
1	A	146[B]	CYS	CB-SG	8.09	1.96	1.82

The worst 5 of 8 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	121[A]	LYS	CD-CE-NZ	10.55	135.97	111.70
1	A	121[B]	LYS	CD-CE-NZ	10.55	135.97	111.70
1	A	220	ARG	NE-CZ-NH1	8.31	124.46	120.30
1	A	36	ARG	NE-CZ-NH1	8.30	124.45	120.30
1	A	92	LYS	CD-CE-NZ	-7.57	94.29	111.70

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	67[C]	EYG	CA1

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	2140	0	2055	83	0
2	A	3	0	0	0	0
3	A	4	0	3	0	0
4	A	470	0	0	42	0
All	All	2617	0	2058	83	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 23.

The worst 5 of 83 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:168:LYS:HD2	4:A:425:HOH:O	1.31	1.29
1:A:114[A]:GLN:HG2	4:A:365:HOH:O	1.17	1.27
1:A:128[B]:ASN:ND2	4:A:261:HOH:O	1.69	1.22
1:A:161:MET:SD	4:A:420[A]:HOH:O	1.99	1.18
1:A:141[A]:MET:SD	1:A:168:LYS:HG2	1.85	1.17

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	266/244 (109%)	261 (98%)	5 (2%)	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	239/205 (117%)	238 (100%)	1 (0%)	91 68

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	231	LYS

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)

3 non-standard protein/DNA/RNA residues 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	$\#\|Z\|>2$
1	NRQ	A	67[A]	1	23,24,25	1.52	3 (13%)	23,32,34	2.69	8 (34%)
1	CH6	A	67[B]	1	24,24,25	1.70	3 (12%)	28,32,34	2.95	11 (39%)
1	EYG	A	67[C]	1	19,23,26	1.62	3 (15%)	21,33,37	2.92	5 (23%)

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
1	NRQ	A	67[A]	1	-	3/9/31/32	0/2/2/2
1	CH6	A	67[B]	1	-	5/12/31/32	0/2/2/2
1	EYG	A	67[C]	1	1/1/5/8	3/8/34/37	0/2/2/2

The worst 5 of 9 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	67[A]	NRQ	CA2-C2	-4.33	1.44	1.48
1	A	67[B]	CH6	CA2-C2	-4.33	1.44	1.48
1	A	67[B]	CH6	C1-N2	4.22	1.38	1.32
1	A	67[C]	EYG	CA2-C2	-3.75	1.44	1.49
1	A	67[C]	EYG	CB-CA1	-3.27	1.50	1.53

The worst 5 of 24 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	67[C]	EYG	C2-CA2-N2	10.37	113.06	105.12
1	A	67[A]	NRQ	CA2-C2-N3	9.69	107.95	103.37
1	A	67[B]	CH6	CA2-C2-N3	9.69	107.95	103.37
1	A	67[B]	CH6	CE-SD-CG1	-6.07	79.56	100.40
1	A	67[B]	CH6	CB1-CA1-N1	5.45	124.46	110.17

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	67[C]	EYG	CA1

5 of 11 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	67[A]	NRQ	C1-CA1-CB1-CG1
1	A	67[B]	CH6	CA1-CB1-CG1-SD
1	A	67[B]	CH6	N1-CA1-CB1-CG1
1	A	67[C]	EYG	N2-CA2-CB2-CG2
1	A	67[A]	NRQ	N2-CA2-CB2-CG2

There are no ring outliers.

1 monomer is involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	67[A]	NRQ	8	0

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 4 ligands modelled in this entry, 3 are monoatomic - leaving 1 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	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
3	ACT	A	233	-	3,3,3	0.72	0	3,3,3	0.40	0

There are no bond length outliers.

There are no bond angle outliers.

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.

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	228/244 (93%)	-0.01	8 (3%) 44 33	5, 8, 16, 31	3 (1%)

The worst 5 of 8 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	3	ASP	5.8
1	A	230	TYR	5.0
1	A	2	GLU	4.6
1	A	229	LEU	4.1
1	A	227	ASP	3.3

6.2 Non-standard residues in protein, DNA, RNA chains

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	B-factors(Å^2)	Q<0.9
1	CH6	A	67[B]	23/24	0.97	0.16	6,9,28,34	23
1	EYG	A	67[C]	22/25	0.97	0.17	6,14,28,34	22
1	NRQ	A	67[A]	23/24	0.98	0.15	6,10,28,34	23

6.3 Carbohydrates

There are no monosaccharides 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. 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	B-factors(Å^2)	Q<0.9
3	ACT	A	233	4/4	0.98	0.06	8,8,8,9	0
2	NA	A	232	1/1	0.99	0.09	8,8,8,8	0
2	NA	A	235	1/1	1.00	0.05	8,8,8,8	0
2	NA	A	234	1/1	1.00	0.03	9,9,9,9	0

6.5 Other polymers

There are no such residues in this entry.