WORLDWIDE PDB
PROTEIN DATA BANK

wwPDB X-ray Structure Validation Summary Report ⓘ

Jun 12, 2024 – 05:05 PM EDT

PDB ID	:	3Q0O
Title	:	Crystal structure of the PUMILIO-homology domain from Human PUMILIO1 in complex with erk2 NRE
Authors	:	Lu, G.; Hall, T.M.T.
Deposited on	:	2010-12-15
Resolution	:	2.80 Å(reported)

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

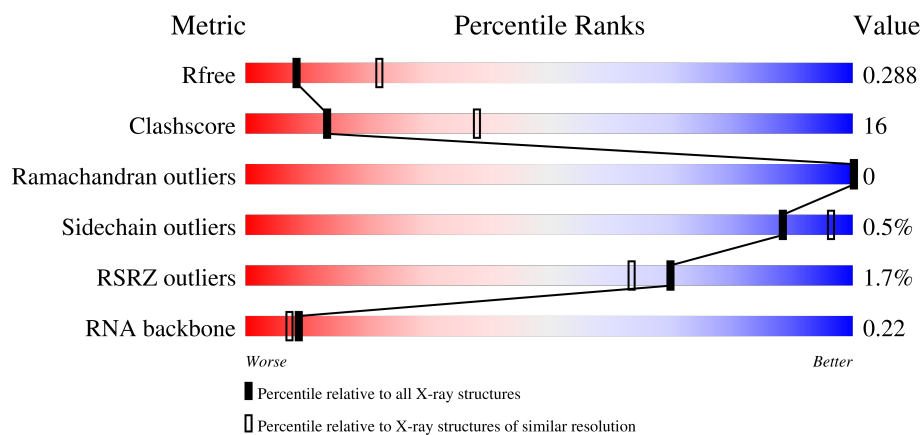|		
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.20.1
EDS	:	2.36.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

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 2.80 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	Percentile Ranks	Value
Rfree		0.288
Clashscore		16
Ramachandran outliers		0
Sidechain outliers		0.5%
RSRZ outliers		1.7%
RNA backbone		0.22

Worse ← → *Better*

■ Percentile relative to all X-ray structures
▯ Percentile relative to X-ray structures of similar resolution

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R$_{free}$	130704	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)
RNA backbone	3102	1227 (3.10-2.50)

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	349	69% 28% ··
1	B	349	67% 28% ·
2	C	8	12% 38% 38% 12%
2	D	8	25% 38% 25% 12%

2 Entry composition ⓘ

There are 3 unique types of molecules in this entry. The entry contains 5881 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 Pumilio homolog 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	341	Total 2768	C 1752	N 501	O 498	S 17	0	0	0
1	B	335	Total 2723	C 1726	N 492	O 489	S 16	0	0	0

- Molecule 2 is a RNA chain called 5'-R(UP*GP*UP*AP*CP*AP*UP*C)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	7	Total 144	C 66	N 24	O 48	P 6	0	0	0
2	D	7	Total 144	C 66	N 24	O 48	P 6	0	0	0

- Molecule 3 is water.

| Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	60	Total 60	O 60	0	0
3	B	32	Total 32	O 32	0	0
3	C	2	Total 2	O 2	0	0
3	D	8	Total 8	O 8	0	0

3 Residue-property plots ⓘ

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: Pumilio homolog 1
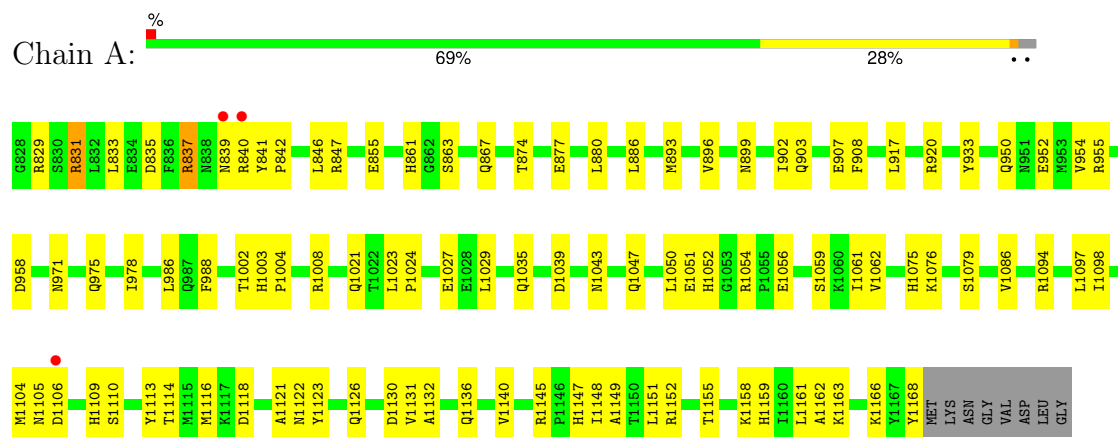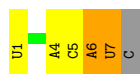
Chain A:

- Molecule 1: Pumilio homolog 1
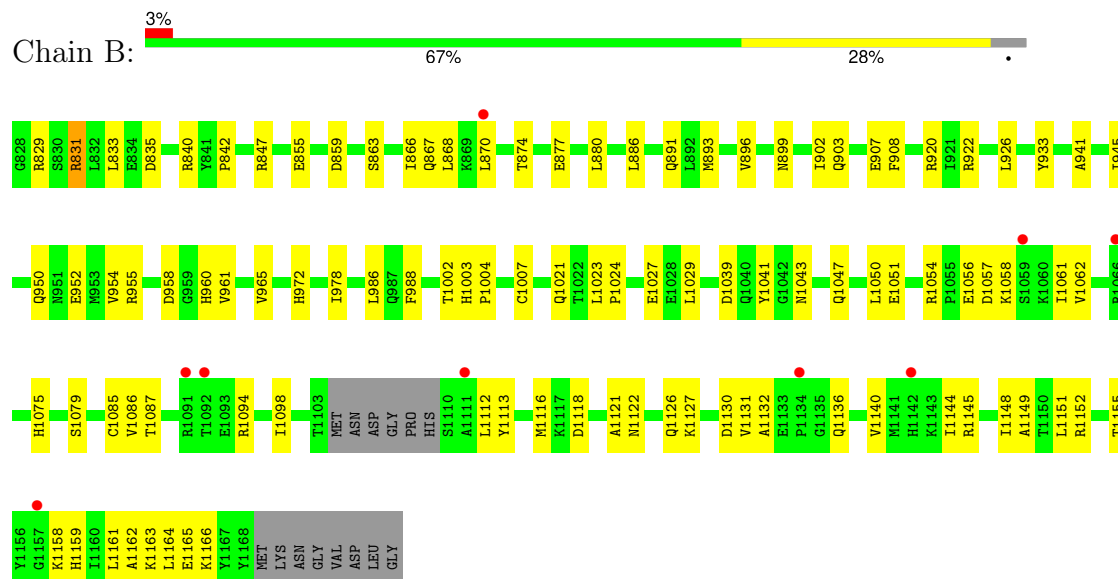
Chain B:

- Molecule 2: 5'-R(UP*GP*UP*AP*CP*AP*UP*C)-3'

Chain C:

• Molecule 2: 5'-R(UP*GP*UP*AP*CP*AP*UP*C)-3'

Chain D:

| 25% | 38% | 25% | 12% |

U1 A4 C5 A6 U7 C

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	36.00Å 59.50Å 341.56Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.10 – 2.80 41.13 – 2.80	Depositor EDS
% Data completeness (in resolution range)	94.7 (41.10-2.80) 94.7 (41.13-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.57 (at 2.81Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.4_6)	Depositor
R, R_{free}	0.224 , 0.297 0.230 , 0.288	Depositor DCC
R_{free} test set	1048 reflections (5.81%)	wwPDB-VP
Wilson B-factor (Å^2)	41.7	Xtriage
Anisotropy	0.745	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 46.7	EDS
L-test for twinning[2]	$< |L| > = 0.46$, $< L^2 > = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.91	EDS
Total number of atoms	5881	wwPDB-VP
Average B, all atoms (Å^2)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.27% of the height of the origin peak. No significant pseudotranslation is detected.*

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[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 ⓘ

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	0.33	0/2821	0.47	0/3804
1	B	0.34	0/2773	0.47	0/3737
2	C	0.38	0/160	0.69	0/247
2	D	0.39	0/160	0.68	0/247
All	All	0.34	0/5914	0.48	0/8035

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 ⓘ

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	2768	0	2783	86	0
1	B	2723	0	2746	90	0
2	C	144	0	76	7	0
2	D	144	0	76	6	0
3	A	60	0	0	7	0
3	B	32	0	0	2	0
3	C	2	0	0	0	0
3	D	8	0	0	0	0
All	All	5881	0	5681	184	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 16.

The worst 5 of 184 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:B:866:ILE:O	1:B:870:LEU:CD1	1.70	1.37
1:A:839:ASN:O	1:A:840:ARG:HD2	1.39	1.19
1:A:831:ARG:HD3	1:A:831:ARG:H	1.00	1.14
1:B:866:ILE:C	1:B:870:LEU:HD13	1.67	1.13
1:B:866:ILE:O	1:B:870:LEU:HD13	0.93	1.11

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	339/349 (97%)	325 (96%)	14 (4%)	0	100	100
1	B	331/349 (95%)	317 (96%)	14 (4%)	0	100	100
All	All	670/698 (96%)	642 (96%)	28 (4%)	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	303/310 (98%)	301 (99%)	2 (1%)	84	95

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	298/310 (96%)	297 (100%)	1 (0%)	92	98
All	All	601/620 (97%)	598 (100%)	3 (0%)	88	96

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	831	ARG
1	A	837	ARG
1	B	831	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 16 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	1088	HIS
1	B	1052	HIS
1	B	845	GLN
1	B	1031	GLN
1	A	1119	GLN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	C	6/8 (75%)	3 (50%)	0
2	D	6/8 (75%)	4 (66%)	0
All	All	12/16 (75%)	7 (58%)	0

5 of 7 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	C	5	C
2	C	6	A
2	C	7	U
2	D	4	A
2	D	5	C

There are no RNA pucker outliers to report.

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 monosaccharides 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, 95th 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	341/349 (97%)	-0.09	3 (0%)	84	80	21, 41, 72, 87	0
1	B	335/349 (95%)	-0.04	9 (2%)	54	44	21, 43, 71, 89	0
2	C	7/8 (87%)	0.30	0	100	100	30, 33, 52, 53	0
2	D	7/8 (87%)	0.21	0	100	100	47, 48, 62, 66	0
All	All	690/714 (96%)	-0.06	12 (1%)	70	63	21, 43, 71, 89	0

The worst 5 of 12 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	1142	HIS	4.2
1	A	1106	ASP	3.7
1	B	1092	THR	2.9
1	B	1059	SER	2.8
1	B	1091	ARG	2.8

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 monosaccharides 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.