



Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2020 – 05:12 am BST

PDB ID : 4I79
Title : Crystal structure of human NUP43
Authors : Xu, C.; Tempel, W.; Li, Z.; He, H.; Wernimont, A.K.; Bountra, C.; Arrow-smith, C.H.; Edwards, A.M.; Min, J.; Structural Genomics Consortium (SGC)
Deposited on : 2012-11-30
Resolution : 1.75 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
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)	:	2.11

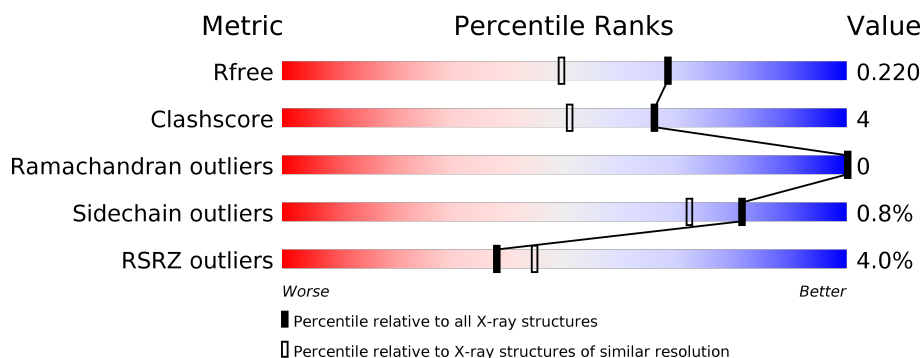
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.75 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	2340 (1.76-1.76)
Clashscore	141614	2466 (1.76-1.76)
Ramachandran outliers	138981	2437 (1.76-1.76)
Sidechain outliers	138945	2437 (1.76-1.76)
RSRZ outliers	127900	2298 (1.76-1.76)

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 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 $\leq 5\%$. 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	399	<div> <div>4%</div> <div> <div></div> <div>74%</div> <div>5%</div> <div>21%</div> </div> </div>
1	B	399	<div> <div>2%</div> <div> <div></div> <div>72%</div> <div>7%</div> <div>21%</div> </div> </div>
2	C	9	<div> <div></div> <div>100%</div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5444 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 Nucleoporin Nup43.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	316	Total	C	N	O	S	0	22	0
			2502	1594	430	466	12			
1	B	316	Total	C	N	O	S	0	33	1
			2587	1639	447	488	13			

There are 38 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-18	MET	-	EXPRESSION TAG	UNP Q8NFH3
A	-17	GLY	-	EXPRESSION TAG	UNP Q8NFH3
A	-16	SER	-	EXPRESSION TAG	UNP Q8NFH3
A	-15	SER	-	EXPRESSION TAG	UNP Q8NFH3
A	-14	HIS	-	EXPRESSION TAG	UNP Q8NFH3
A	-13	HIS	-	EXPRESSION TAG	UNP Q8NFH3
A	-12	HIS	-	EXPRESSION TAG	UNP Q8NFH3
A	-11	HIS	-	EXPRESSION TAG	UNP Q8NFH3
A	-10	HIS	-	EXPRESSION TAG	UNP Q8NFH3
A	-9	HIS	-	EXPRESSION TAG	UNP Q8NFH3
A	-8	SER	-	EXPRESSION TAG	UNP Q8NFH3
A	-7	SER	-	EXPRESSION TAG	UNP Q8NFH3
A	-6	GLY	-	EXPRESSION TAG	UNP Q8NFH3
A	-5	LEU	-	EXPRESSION TAG	UNP Q8NFH3
A	-4	VAL	-	EXPRESSION TAG	UNP Q8NFH3
A	-3	PRO	-	EXPRESSION TAG	UNP Q8NFH3
A	-2	ARG	-	EXPRESSION TAG	UNP Q8NFH3
A	-1	GLY	-	EXPRESSION TAG	UNP Q8NFH3
A	0	SER	-	EXPRESSION TAG	UNP Q8NFH3
B	-18	MET	-	EXPRESSION TAG	UNP Q8NFH3
B	-17	GLY	-	EXPRESSION TAG	UNP Q8NFH3
B	-16	SER	-	EXPRESSION TAG	UNP Q8NFH3
B	-15	SER	-	EXPRESSION TAG	UNP Q8NFH3
B	-14	HIS	-	EXPRESSION TAG	UNP Q8NFH3
B	-13	HIS	-	EXPRESSION TAG	UNP Q8NFH3

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-12	HIS	-	EXPRESSION TAG	UNP Q8NFH3
B	-11	HIS	-	EXPRESSION TAG	UNP Q8NFH3
B	-10	HIS	-	EXPRESSION TAG	UNP Q8NFH3
B	-9	HIS	-	EXPRESSION TAG	UNP Q8NFH3
B	-8	SER	-	EXPRESSION TAG	UNP Q8NFH3
B	-7	SER	-	EXPRESSION TAG	UNP Q8NFH3
B	-6	GLY	-	EXPRESSION TAG	UNP Q8NFH3
B	-5	LEU	-	EXPRESSION TAG	UNP Q8NFH3
B	-4	VAL	-	EXPRESSION TAG	UNP Q8NFH3
B	-3	PRO	-	EXPRESSION TAG	UNP Q8NFH3
B	-2	ARG	-	EXPRESSION TAG	UNP Q8NFH3
B	-1	GLY	-	EXPRESSION TAG	UNP Q8NFH3
B	0	SER	-	EXPRESSION TAG	UNP Q8NFH3

- Molecule 2 is a protein called floating chain, unknown sequence.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	9	Total	C	N	O	0	9	0
			45	27	9	9			

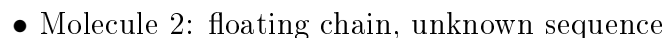
- Molecule 3 is UNKNOWN ATOM OR ION (three-letter code: UNX) (formula: X).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	19	Total	X	0	0
			19	19		
3	A	13	Total	X	0	0
			13	13		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	130	Total	O	0	0
			130	130		
4	B	148	Total	O	0	0
			148	148		

- Molecule 1: Nucleoporin Nup43



100%

WORLDWIDE
PDB
PROTEIN DATA BANK

4 Data and refinement statistics

Property	Value	Source
Space group	P 2 ₁ 2 ₁ 2	Depositor
Cell constants a, b, c, α , β , γ	80.74 Å 161.96 Å 58.20 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.00 – 1.75 45.32 – 1.75	Depositor EDS
% Data completeness (in resolution range)	99.9 (45.00-1.75) 100.0 (45.32-1.75)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.14 (at 1.75 Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.181 , 0.212 0.190 , 0.220	Depositor DCC
R_{free} test set	2490 reflections (3.20%)	wwPDB-VP
Wilson B-factor (Å ²)	22.7	Xtriage
Anisotropy	0.489	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 44.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5444	wwPDB-VP
Average B, all atoms (Å ²)	29.0	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 43.05 % 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 1.8650e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical 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: UNX

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	0.83	0/2622	0.75	1/3584 (0.0%)
1	B	0.78	0/2723	0.78	0/3715
All	All	0.80	0/5345	0.76	1/7299 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	376	ARG	NE-CZ-NH1	5.01	122.81	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	2502	0	2365	20	0
1	B	2587	0	2485	21	0
2	C	45	0	11	0	0
3	A	13	0	0	0	0
3	B	19	0	0	0	0
4	A	130	0	0	0	0
4	B	148	0	0	0	0
All	All	5444	0	4861	40	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 4.

All (40) 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:358[B]:VAL:CG1	1:A:363[B]:LEU:HD12	1.94	0.97
1:B:252[B]:MET:HG2	1:B:253:PRO:HD2	1.55	0.89
1:A:358[B]:VAL:HG13	1:A:363[B]:LEU:HD12	1.67	0.77
1:A:358[B]:VAL:HG12	1:A:363[B]:LEU:HD12	1.67	0.75
1:A:358[B]:VAL:CG1	1:A:363[B]:LEU:CD1	2.71	0.68
1:B:252[B]:MET:CG	1:B:253:PRO:HD2	2.26	0.63
1:B:13:ILE:HG13	1:B:372[B]:ILE:HD11	1.89	0.55
1:B:358[A]:VAL:HG13	1:B:363[A]:LEU:HD23	1.89	0.55
1:B:9:VAL:HG11	1:B:372[A]:ILE:HD12	1.89	0.54
1:B:214[B]:THR:HG21	1:B:252[B]:MET:HE1	1.91	0.53
1:A:358[B]:VAL:HG12	1:A:363[B]:LEU:CD1	2.38	0.53
1:A:286[B]:LEU:HD21	1:A:356:LEU:HD22	1.90	0.52
1:B:231[A]:HIS:CD2	1:B:231[A]:HIS:H	2.28	0.52
1:A:286[B]:LEU:HD12	1:A:345:LEU:HD12	1.92	0.51
1:A:171:ASP:OD2	1:A:196:LYS:NZ	2.40	0.51
1:A:16:THR:HG22	1:A:34:THR:HG22	1.93	0.50
1:B:6:ALA:HB2	1:B:373:TYR:CE2	2.47	0.49
1:A:286[B]:LEU:CD2	1:A:356:LEU:HD22	2.43	0.48
1:B:91:ALA:HB2	1:B:137[A]:VAL:HG13	1.96	0.48
1:A:79:ASP:HB3	1:A:137[A]:VAL:HG12	1.96	0.48
1:B:363[A]:LEU:HD13	1:B:364:VAL:N	2.29	0.48
1:B:240:GLY:O	1:B:259:ALA:HB3	2.15	0.47
1:A:117:THR:HG22	1:B:201:ARG:NH2	2.31	0.46
1:A:151:GLY:O	1:A:169:ASN:HB3	2.16	0.46
1:A:91:ALA:HB2	1:A:137[B]:VAL:HG13	1.98	0.45
1:B:16:THR:HG22	1:B:34:THR:HG22	1.99	0.44
1:B:79:ASP:HB3	1:B:137[B]:VAL:HG12	2.00	0.43
1:A:231[A]:HIS:H	1:A:231[A]:HIS:CD2	2.37	0.42
1:B:231[A]:HIS:CE1	1:B:248:ARG:NH2	2.87	0.42
1:B:44[B]:ILE:HD12	1:B:100:PHE:CE2	2.54	0.42
1:B:365:CYS:SG	1:B:373:TYR:HB2	2.61	0.41
1:B:8:PHE:CE1	1:B:371:ALA:HB2	2.56	0.41
1:A:83:PHE:CZ	1:A:142:PRO:HA	2.55	0.41
1:A:83:PHE:HZ	1:A:142:PRO:HA	1.86	0.41
1:A:286[B]:LEU:CD1	1:A:345:LEU:HD12	2.51	0.40
1:B:286:LEU:HD23	1:B:286:LEU:C	2.42	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	324/399 (81%)	314 (97%)	10 (3%)	0	100	100
1	B	334/399 (84%)	325 (97%)	9 (3%)	0	100	100
All	All	658/798 (82%)	639 (97%)	19 (3%)	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	275/351 (78%)	273 (99%)	2 (1%)	84	75
1	B	293/351 (84%)	288 (98%)	5 (2%)	60	42
All	All	568/702 (81%)	561 (99%)	7 (1%)	81	56

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11[A]	GLN
1	A	11[B]	GLN
1	B	72	ARG
1	B	214[A]	THR
1	B	214[B]	THR
1	B	363[A]	LEU
1	B	363[B]	LEU

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

Mol	Chain	Res	Type
1	B	230	GLN

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](#)

Of 32 ligands modelled in this entry, 32 are unknown - leaving 0 for Mogul analysis.

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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
2	C	4

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	C	335[A]:UNK	C	336[A]:UNK	N	12.49
1	C	330[A]:UNK	C	331[A]:UNK	N	4.08
1	C	331[A]:UNK	C	332[A]:UNK	N	3.85
1	C	337[A]:UNK	C	338[A]:UNK	N	3.55

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	316/399 (79%)	-0.06	16 (5%) 28 34	17, 28, 50, 77	4 (1%)
1	B	316/399 (79%)	-0.03	9 (2%) 53 58	14, 25, 48, 61	5 (1%)
2	C	0/9	-	-	-	-
All	All	632/807 (78%)	-0.04	25 (3%) 38 45	14, 26, 50, 77	9 (1%)

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	215	GLY	4.5
1	A	24	GLY	3.9
1	A	331	LEU	3.8
1	A	170	ALA	3.6
1	B	213[A]	LEU	3.4
1	B	214[A]	THR	3.3
1	B	331[A]	LEU	3.2
1	B	332[A]	SER	3.2
1	B	184	PRO	3.1
1	B	216	ASP	2.9
1	A	105[A]	ASN	2.6
1	A	184	PRO	2.6
1	A	172	SER	2.5
1	A	217	ARG	2.4
1	A	332	SER	2.3
1	A	256	LEU	2.3
1	B	329[A]	SER	2.3
1	B	330[A]	TRP	2.2
1	A	49	ILE	2.2
1	A	26	LEU	2.2
1	B	215	GLY	2.2
1	A	330	TRP	2.2
1	A	62	GLU	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	107[A]	GLN	2.0
1	A	8	PHE	2.0

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, 95th 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(\AA^2)	Q<0.9
3	UNX	A	404	1/1	0.71	0.14	40,40,40,40	0
3	UNX	A	410	1/1	0.74	0.13	36,36,36,36	0
3	UNX	B	415	1/1	0.75	0.14	42,42,42,42	0
3	UNX	B	417	1/1	0.78	0.18	40,40,40,40	0
3	UNX	B	402	1/1	0.80	0.21	38,38,38,38	0
3	UNX	B	405	1/1	0.81	0.21	38,38,38,38	0
3	UNX	B	416	1/1	0.81	0.22	48,48,48,48	0
3	UNX	B	411	1/1	0.81	0.25	39,39,39,39	0
3	UNX	B	407	1/1	0.82	0.30	48,48,48,48	0
3	UNX	B	408	1/1	0.84	0.13	37,37,37,37	0
3	UNX	B	413	1/1	0.87	0.36	43,43,43,43	0
3	UNX	A	411	1/1	0.88	0.16	33,33,33,33	0
3	UNX	B	409	1/1	0.88	0.07	25,25,25,25	0
3	UNX	B	410	1/1	0.88	0.10	36,36,36,36	0
3	UNX	B	401	1/1	0.89	0.20	38,38,38,38	0
3	UNX	A	412	1/1	0.89	0.14	32,32,32,32	0
3	UNX	A	406	1/1	0.90	0.11	38,38,38,38	0
3	UNX	B	412	1/1	0.91	0.16	31,31,31,31	0
3	UNX	A	407	1/1	0.91	0.30	35,35,35,35	0
3	UNX	B	414	1/1	0.91	0.15	32,32,32,32	0
3	UNX	B	406	1/1	0.92	0.20	37,37,37,37	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	UNX	A	405	1/1	0.92	0.11	30,30,30,30	0
3	UNX	A	408	1/1	0.93	0.23	49,49,49,49	0
3	UNX	A	409	1/1	0.94	0.11	32,32,32,32	0
3	UNX	B	403	1/1	0.94	0.08	30,30,30,30	0
3	UNX	A	413	1/1	0.95	0.18	25,25,25,25	0
3	UNX	B	404	1/1	0.95	0.10	27,27,27,27	0
3	UNX	B	418	1/1	0.96	0.13	19,19,19,19	0
3	UNX	A	403	1/1	0.96	0.08	17,17,17,17	0
3	UNX	B	419	1/1	0.97	0.11	34,34,34,34	0
3	UNX	A	401	1/1	0.98	0.08	16,16,16,16	0
3	UNX	A	402	1/1	0.99	0.11	17,17,17,17	0

6.5 Other polymers [i](#)

There are no such residues in this entry.