



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 03:29 PM UTC

PDB ID : 3NGO / pdb_00003ngo
Title : Crystal structure of the human CNOT6L nuclease domain in complex with poly(A) DNA
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Deposited on : 2010-06-12
Resolution : 2.20 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

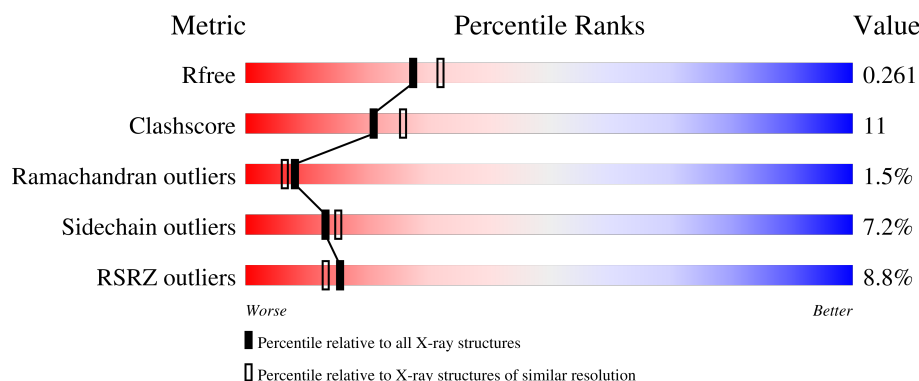
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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}	180053	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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 $\leq 5\%$. 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	398	7% 65% 17% • 15%
2	B	4	50% 75% 25%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 2890 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 CCR4-NOT transcription complex subunit 6-like.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	337	Total	C	N	O	S	0	0	0
			2722	1753	446	503	20			

- Molecule 2 is a DNA chain called 5'-D(*AP*AP*AP*A)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	4	Total	C	N	O	P	0	0	0
			54	25	10	16	3			

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

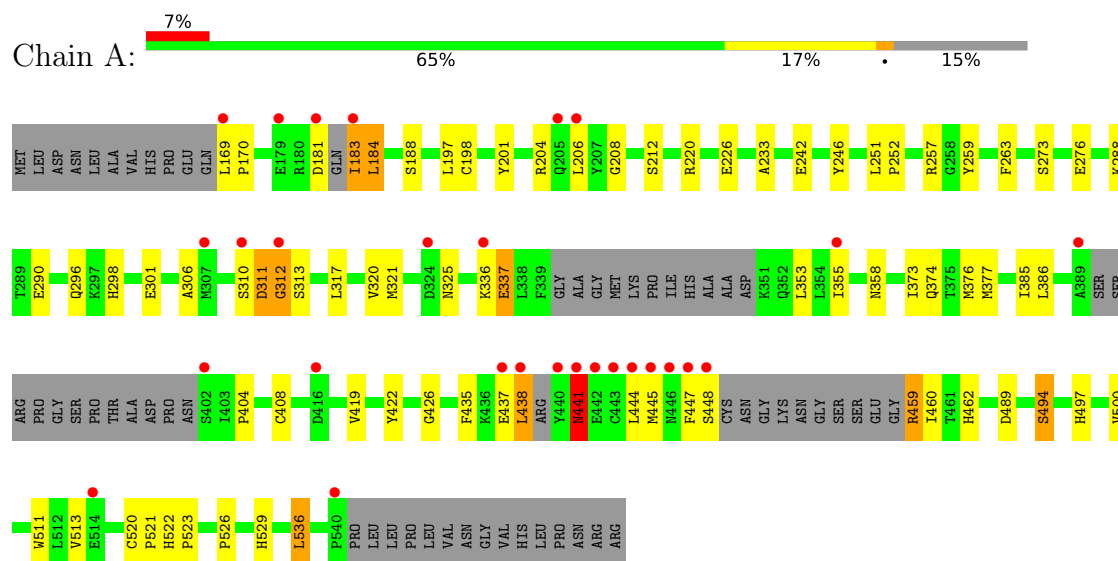
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	109	Total	O	0	0
			109	109		
4	B	3	Total	O	0	0
			3	3		

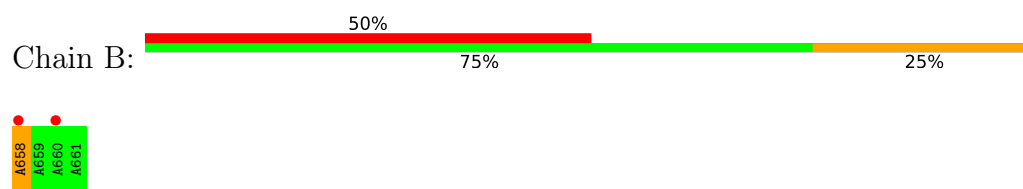
3 Residue-property plots [i](#)

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: CCR4-NOT transcription complex subunit 6-like



- Molecule 2: 5'-D(*AP*AP*AP*A)-3'



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	77.37Å 77.37Å 167.18Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 2.20 50.00 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.6 (50.00-2.20) 99.6 (50.00-2.20)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.11 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.5.0044	Depositor
R, R_{free}	0.215 , 0.263 0.212 , 0.261	Depositor DCC
R_{free} test set	1522 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	39.8	Xtriage
Anisotropy	0.335	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 34.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.028 for -h,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	2890	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 5.41% of the height of the origin peak. No significant pseudotranslation is detected.*

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: MG

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	0.94	0/2788	0.99	2/3774 (0.1%)
2	B	0.47	0/60	1.67	2/91 (2.2%)
All	All	0.93	0/2848	1.01	4/3865 (0.1%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	500	VAL	CB-CA-C	-7.12	102.35	110.96
2	B	658	DA	P-O3'-C3'	6.85	130.48	120.20
1	A	312	GLY	N-CA-C	-6.20	100.97	112.37
2	B	658	DA	O3'-P-O5'	-5.62	95.58	104.00

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	2722	0	2665	59	0
2	B	54	0	29	1	0
3	A	2	0	0	0	0
4	A	109	0	0	5	0
4	B	3	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	2890	0	2694	60	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 11.

All (60) 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:183:ILE:HG23	1:A:184:LEU:CD1	1.58	1.31
1:A:183:ILE:CG2	1:A:184:LEU:HD13	1.62	1.26
1:A:183:ILE:CG2	1:A:184:LEU:H	1.58	1.16
1:A:183:ILE:HG23	1:A:184:LEU:HD12	1.36	1.07
1:A:183:ILE:HG23	1:A:184:LEU:H	1.19	1.03
1:A:183:ILE:CG2	1:A:184:LEU:CD1	2.30	0.96
1:A:184:LEU:CD1	1:A:184:LEU:H	1.88	0.87
1:A:459:ARG:HH11	1:A:459:ARG:HG3	1.40	0.85
1:A:183:ILE:HG23	1:A:184:LEU:N	1.84	0.85
1:A:459:ARG:HH11	1:A:459:ARG:CG	1.88	0.85
1:A:183:ILE:CG2	1:A:184:LEU:N	2.29	0.83
1:A:183:ILE:HG22	1:A:184:LEU:H	1.43	0.81
1:A:184:LEU:CD1	1:A:184:LEU:N	2.46	0.77
1:A:184:LEU:HD12	1:A:184:LEU:N	1.99	0.77
1:A:184:LEU:HD13	1:A:184:LEU:H	1.50	0.75
1:A:204:ARG:NH2	1:A:208:GLY:O	2.21	0.73
1:A:183:ILE:HG22	1:A:184:LEU:HD13	1.67	0.71
1:A:489:ASP:OD2	1:A:529:HIS:HD2	1.74	0.70
1:A:459:ARG:HG3	1:A:459:ARG:NH1	2.07	0.68
1:A:325:ASN:HD22	1:A:374:GLN:HE22	1.40	0.68
1:A:181:ASP:OD1	1:A:183:ILE:HG22	1.96	0.65
1:A:441:ASN:N	1:A:441:ASN:OD1	2.30	0.65
1:A:310:SER:O	1:A:312:GLY:O	2.16	0.64
1:A:242:GLU:HG3	4:A:665:HOH:O	1.98	0.63
1:A:310:SER:O	1:A:311:ASP:C	2.40	0.62
1:A:317:LEU:HD23	1:A:321:MET:HE3	1.84	0.59
1:A:273:SER:OG	1:A:276:GLU:HG3	2.05	0.57
1:A:198:CYS:SG	1:A:201:TYR:CD1	2.98	0.56
1:A:459:ARG:CG	1:A:459:ARG:NH1	2.58	0.55
1:A:181:ASP:OD1	1:A:181:ASP:C	2.47	0.54
1:A:337:GLU:HB3	4:A:636:HOH:O	2.07	0.53
1:A:426:GLY:HA2	1:A:462:HIS:CE1	2.44	0.53
1:A:494:SER:HB2	1:A:497:HIS:HD2	1.75	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:325:ASN:ND2	1:A:374:GLN:HE22	2.06	0.52
1:A:321:MET:C	1:A:321:MET:SD	2.93	0.51
1:A:170:PRO:HD3	1:A:511:TRP:CD1	2.46	0.50
1:A:188:SER:HA	1:A:536:LEU:O	2.11	0.50
1:A:259:TYR:CZ	1:A:288:LYS:HE3	2.47	0.50
1:A:374:GLN:HA	1:A:377:MET:HE3	1.95	0.49
1:A:373:ILE:HG22	1:A:377:MET:HE2	1.95	0.48
1:A:353:LEU:O	1:A:404:PRO:HD2	2.15	0.47
1:A:220:ARG:HD2	1:A:526:PRO:O	2.15	0.46
1:A:257:ARG:CZ	4:A:619:HOH:O	2.63	0.46
1:A:233:ALA:O	1:A:259:TYR:OH	2.25	0.46
1:A:529:HIS:HE1	4:A:587:HOH:O	1.99	0.46
1:A:441:ASN:O	1:A:444:LEU:HB2	2.16	0.46
1:A:298:HIS:CE1	1:A:385:ILE:HG23	2.52	0.45
4:A:560:HOH:O	2:B:658:DA:H3'	2.17	0.45
1:A:522:HIS:CG	1:A:523:PRO:HD2	2.52	0.44
1:A:251:LEU:N	1:A:252:PRO:HD2	2.33	0.43
1:A:296:GLN:OE1	1:A:298:HIS:HE1	2.01	0.43
1:A:520:CYS:HB3	1:A:521:PRO:HA	2.01	0.43
1:A:422:TYR:O	1:A:462:HIS:HE1	2.01	0.42
1:A:358:ASN:HA	1:A:408:CYS:O	2.19	0.42
1:A:376:MET:HE1	1:A:444:LEU:HD22	2.01	0.42
1:A:306:ALA:HB2	1:A:320:VAL:HG12	2.01	0.42
1:A:290:GLU:HG2	1:A:290:GLU:O	2.20	0.41
1:A:435:PHE:O	1:A:438:LEU:HB2	2.21	0.41
1:A:246:TYR:CE1	1:A:263:PHE:HB2	2.55	0.41
1:A:422:TYR:O	1:A:462:HIS:CE1	2.75	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	325/398 (82%)	315 (97%)	5 (2%)	5 (2%)	8 6

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	441	ASN
1	A	313	SER
1	A	437	GLU
1	A	311	ASP
1	A	197	LEU

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	305/354 (86%)	283 (93%)	22 (7%)	13 15

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	169	LEU
1	A	183	ILE
1	A	184	LEU
1	A	206	LEU
1	A	212	SER
1	A	226	GLU
1	A	301	GLU
1	A	336	LYS
1	A	337	GLU
1	A	355	ILE
1	A	386	LEU
1	A	419	VAL
1	A	438	LEU
1	A	441	ASN
1	A	445	MET
1	A	447	PHE

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Mol	Chain	Res	Type
1	A	448	SER
1	A	459	ARG
1	A	460	ILE
1	A	494	SER
1	A	513	VAL
1	A	536	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	216	ASN
1	A	298	HIS
1	A	325	ASN
1	A	362	HIS
1	A	412	ASN
1	A	462	HIS
1	A	473	ASN
1	A	497	HIS
1	A	499	ASN
1	A	529	HIS

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 2 are monoatomic - 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

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 [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, 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	337/398 (84%)	0.54	28 (8%) 17 15	26, 41, 68, 90	0
2	B	4/4 (100%)	1.77	2 (50%) 0 0	49, 64, 94, 101	0
All	All	341/402 (84%)	0.55	30 (8%) 15 13	26, 41, 70, 101	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	438	LEU	5.1
1	A	444	LEU	4.8
1	A	445	MET	4.6
1	A	448	SER	4.3
1	A	447	PHE	4.2
1	A	441	ASN	3.9
1	A	442	GLU	3.4
1	A	402	SER	3.3
2	B	658	DA	3.2
1	A	389	ALA	3.2
1	A	312	GLY	3.2
2	B	660	DA	3.2
1	A	179	GLU	3.1
1	A	169	LEU	3.1
1	A	440	TYR	3.1
1	A	183	ILE	3.0
1	A	310	SER	2.8
1	A	206	LEU	2.7
1	A	443	CYS	2.7
1	A	437	GLU	2.6
1	A	307	MET	2.6
1	A	540	PRO	2.3
1	A	181	ASP	2.3
1	A	324	ASP	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	205	GLN	2.2
1	A	514	GLU	2.2
1	A	416	ASP	2.2
1	A	355	ILE	2.1
1	A	446	ASN	2.1
1	A	336	LYS	2.1

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 oligosaccharides 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(Å ²)	Q<0.9
3	MG	A	557	1/1	0.92	0.16	50,50,50,50	0
3	MG	A	556	1/1	0.98	0.09	39,39,39,39	0

6.5 Other polymers [i](#)

There are no such residues in this entry.