



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 11:08 AM UTC

PDB ID : 5HAZ / pdb_00005haz
Title : Structure of Chaetomium thermophilum Nup170 CTD
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Deposited on : 2015-12-31
Resolution : 2.10 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
Mogul	:	2022.3.0, CSD as543be (2022)
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

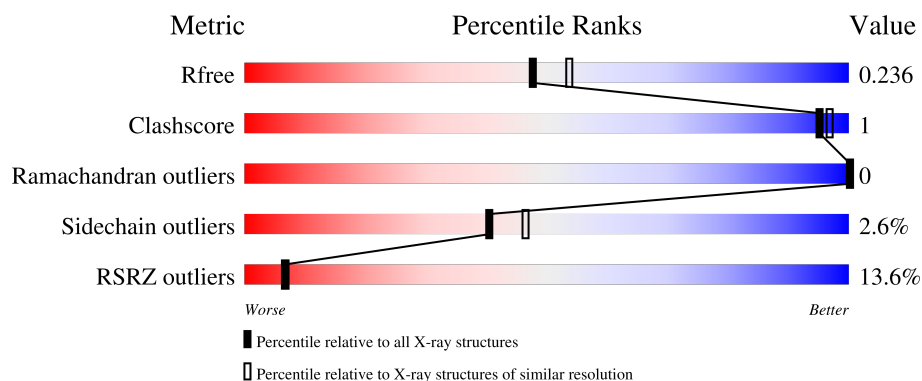
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	567	13% 92% 5%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 9405 atoms, of which 4486 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.

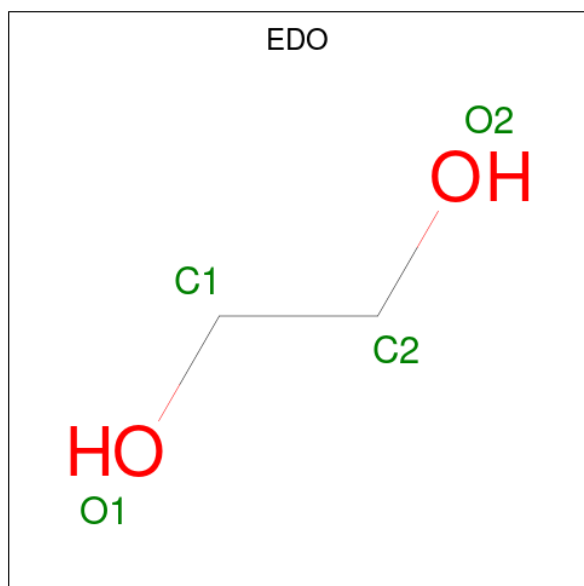
- Molecule 1 is a protein called Nucleoporin NUP170.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	551	Total	C	H	N	O	S	0	23	0
			8974	2844	4457	791	866	16			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	850	SER	-	expression tag	UNP G0S7B6

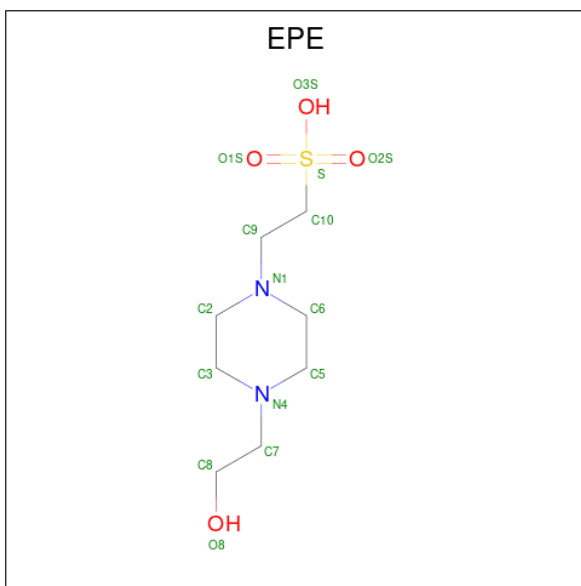
- Molecule 2 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	H	O	0	0
			10	2	6	2		
2	A	1	Total	C	H	O	0	0
			10	2	6	2		

- Molecule 3 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (CCD

ID: EPE) (formula: C₈H₁₈N₂O₄S).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	H	N	O	S	0	1
			32	8	17	2	4	1		

- Molecule 4 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Na	0	0
			1	1		

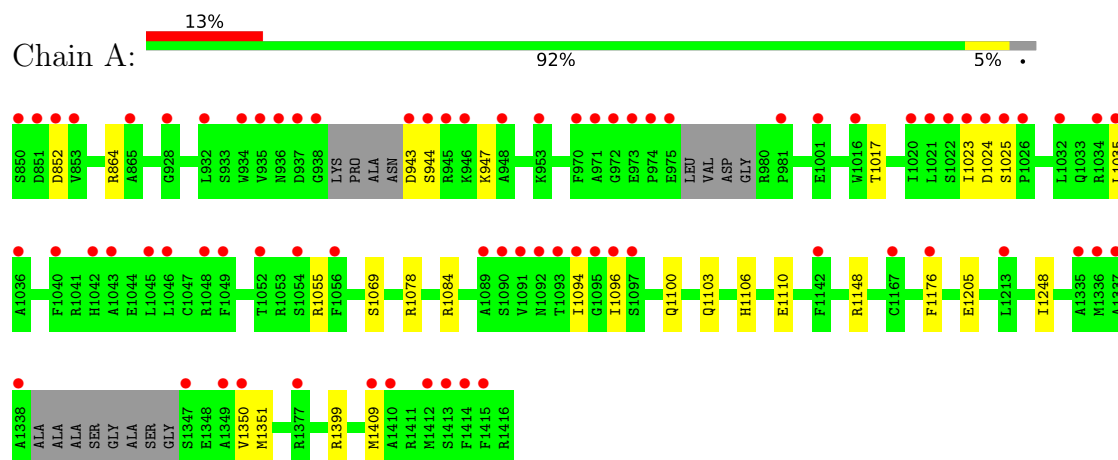
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	378	Total	O	0	2
			378	378		

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: Nucleoporin NUP170



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	141.26Å 54.77Å 108.20Å 90.00° 91.05° 90.00°	Depositor
Resolution (Å)	37.26 – 2.10 37.26 – 2.10	Depositor EDS
% Data completeness (in resolution range)	97.9 (37.26-2.10) 98.2 (37.26-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.70 (at 2.10Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155: ???)	Depositor
R, R_{free}	0.206 , 0.231 0.212 , 0.236	Depositor DCC
R_{free} test set	2396 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	43.5	Xtriage
Anisotropy	0.327	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 48.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.014 for -h,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9405	wwPDB-VP
Average B, all atoms (Å ²)	75.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.03% 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: EPE, NA, EDO

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	0.19	0/4636	0.36	0/6290

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 [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	4517	4457	4438	11	0
2	A	8	12	12	0	0
3	A	15	17	17	0	0
4	A	1	0	0	0	0
5	A	378	0	0	2	0
All	All	4919	4486	4467	11	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 1.

All (11) 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:1069:SER:O	1:A:1078:ARG:NH1	2.27	0.67
1:A:1205[A]:GLU:OE1	5:A:1601:HOH:O	2.13	0.66
1:A:1399:ARG:NH1	5:A:1603:HOH:O	2.35	0.57
1:A:1106:HIS:NE2	1:A:1110[A]:GLU:OE2	2.39	0.56
1:A:1350:VAL:HG12	1:A:1351:MET:N	2.27	0.49
1:A:943:ASP:OD2	1:A:944:SER:N	2.45	0.49
1:A:1055:ARG:NH1	1:A:1094:ILE:O	2.48	0.45
1:A:1084[B]:ARG:HH11	1:A:1084[B]:ARG:HG2	1.82	0.44
1:A:1084[B]:ARG:HH11	1:A:1084[B]:ARG:CG	2.30	0.44
1:A:1023:ILE:HG22	1:A:1025:SER:H	1.85	0.41

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	566/567 (100%)	551 (97%)	15 (3%)	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	489/475 (103%)	477 (98%)	12 (2%)	42	48

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	852	ASP
1	A	864	ARG
1	A	947	LYS
1	A	1017	THR
1	A	1024	ASP
1	A	1035	LEU
1	A	1096	ILE
1	A	1100	GLN
1	A	1103	GLN
1	A	1148	ARG
1	A	1248	ILE
1	A	1409	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1092	ASN
1	A	1155	ASN
1	A	1294	GLN
1	A	1317	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 1 is monoatomic - leaving 3 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
2	EDO	A	1501	-	3,3,3	0.43	0	2,2,2	0.35	0
3	EPE	A	1503[A]	3	15,15,15	0.82	1 (6%)	19,20,20	1.79	5 (26%)
2	EDO	A	1502	-	3,3,3	0.39	0	2,2,2	0.39	0

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
2	EDO	A	1501	-	-	0/1/1/1	-
3	EPE	A	1503[A]	3	-	3/9/19/19	0/1/1/1
2	EDO	A	1502	-	-	0/1/1/1	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1503[A]	EPE	C10-S	2.50	1.81	1.77

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1503[A]	EPE	C5-N4-C3	3.82	117.08	108.84
3	A	1503[A]	EPE	C7-N4-C5	3.43	120.38	111.24
3	A	1503[A]	EPE	O1S-S-C10	2.91	111.13	106.73
3	A	1503[A]	EPE	C6-N1-C2	2.46	114.15	108.84
3	A	1503[A]	EPE	C7-N4-C3	2.35	117.49	111.24

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1503[A]	EPE	S-C10-C9-N1
3	A	1503[A]	EPE	C8-C7-N4-C3
3	A	1503[A]	EPE	N4-C7-C8-O8

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 ⓘ

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	551/567 (97%)	0.64	75 (13%) 7 7	23, 64, 149, 215	21 (3%)

All (75) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1338	ALA	7.9
1	A	938	GLY	7.8
1	A	1337	ALA	6.3
1	A	1094	ILE	6.2
1	A	1167	CYS	5.3
1	A	1095	GLY	5.1
1	A	1096	ILE	4.9
1	A	970	PHE	4.8
1	A	974	PRO	4.8
1	A	945	ARG	4.6
1	A	1413	SER	4.5
1	A	1042	HIS	4.5
1	A	1347	SER	4.3
1	A	1024	ASP	4.2
1	A	981	PRO	4.1
1	A	1349	ALA	3.9
1	A	971	ALA	3.9
1	A	972	GLY	3.8
1	A	1409	MET	3.8
1	A	1025	SER	3.8
1	A	946	LYS	3.6
1	A	853[A]	VAL	3.5
1	A	1410	ALA	3.5
1	A	1049	PHE	3.4
1	A	1023	ILE	3.3
1	A	975	GLU	3.3
1	A	1335	ALA	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	1414	PHE	3.2
1	A	1036	ALA	3.1
1	A	1091	VAL	3.1
1	A	1336	MET	3.0
1	A	1035	LEU	3.0
1	A	973	GLU	3.0
1	A	1097	SER	2.9
1	A	934	TRP	2.8
1	A	1415	PHE	2.8
1	A	1412	MET	2.7
1	A	1040	PHE	2.7
1	A	1052	THR	2.6
1	A	1142	PHE	2.6
1	A	851	ASP	2.6
1	A	944	SER	2.5
1	A	1176	PHE	2.5
1	A	1089	ALA	2.4
1	A	1034	ARG	2.4
1	A	1093	THR	2.4
1	A	865	ALA	2.4
1	A	1020	ILE	2.4
1	A	1001	GLU	2.3
1	A	1016	TRP	2.3
1	A	948	ALA	2.3
1	A	1022	SER	2.3
1	A	1092	ASN	2.3
1	A	943	ASP	2.3
1	A	1032	LEU	2.3
1	A	935	VAL	2.3
1	A	1021	LEU	2.2
1	A	936	ASN	2.2
1	A	852	ASP	2.2
1	A	937	ASP	2.2
1	A	1090	SER	2.2
1	A	1048	ARG	2.2
1	A	1043	ALA	2.2
1	A	1046	LEU	2.2
1	A	1056	PHE	2.1
1	A	1045	LEU	2.1
1	A	932	LEU	2.1
1	A	1377[A]	ARG	2.1
1	A	1054	SER	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	953	LYS	2.0
1	A	928	GLY	2.0
1	A	1350	VAL	2.0
1	A	850	SER	2.0
1	A	1213	LEU	2.0
1	A	1026	PRO	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 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
2	EDO	A	1501	4/4	0.74	0.18	95,114,114,114	0
3	EPE	A	1503[A]	15/15	0.76	0.38	112,135,153,175	32
4	NA	A	1504	1/1	0.82	0.34	245,245,245,245	0
2	EDO	A	1502	4/4	0.84	0.19	86,103,104,104	0

6.5 Other polymers [i](#)

There are no such residues in this entry.