



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 01:34 AM UTC

PDB ID : 1DV4 / pdb_00001dv4
Title : PARTIAL STRUCTURE OF 16S RNA OF THE SMALL RIBOSOMAL SUB-UNIT FROM THERMUS THERMOPHILUS
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Deposited on : 2000-01-19
Resolution : 4.50 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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

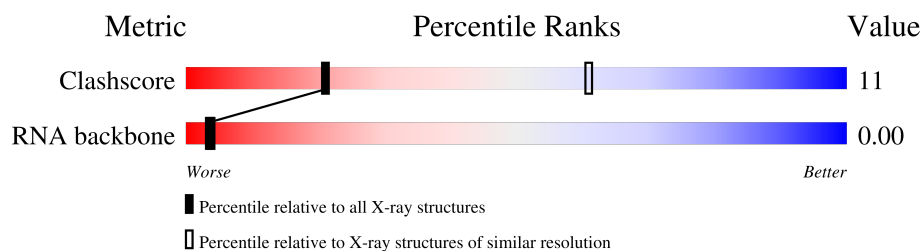
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 4.50 Å.

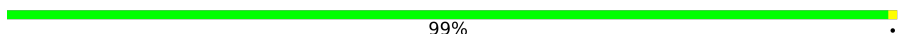
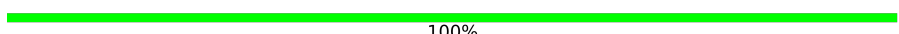
Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	1002 (5.06-3.94)
RNA backbone	3983	1039 (5.90-3.10)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	362	 85% 15%
2	E	145	 99% .
3	G	135	 100%



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	E	1	Total	O	P	W	0	0
			82	62	2	18		

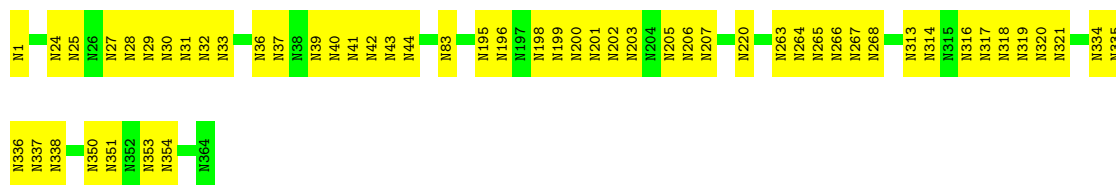
3 Residue-property plots

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. 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. 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.

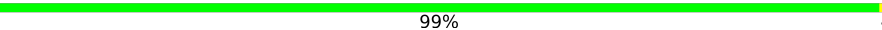
Note EDS was not executed.

• Molecule 1: 16S RIBOSOMAL RNA

Chain A:  85% 15%

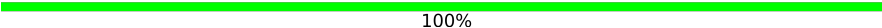


• Molecule 2: RIBOSOMAL PROTEIN S5

Chain E:  99% .



• Molecule 3: RIBOSOMAL PROTEIN S7

Chain G:  100%

There are no outlier residues recorded for this chain.

4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	407.10Å 407.10Å 176.01Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 4.50	Depositor
% Data completeness (in resolution range)	(Not available) (30.00-4.50)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	12.50	Depositor
Refinement program		Depositor
R, R_{free}	(Not available) , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3956	wwPDB-VP
Average B, all atoms (Å ²)	3.0	wwPDB-VP

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: WO2

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).

There are no protein, RNA or DNA chains available to summarize Z scores of covalent bonds and angles.

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	3594	0	2934	75	0
2	E	145	0	0	1	0
3	G	135	0	0	0	0
4	E	82	0	0	1	0
All	All	3956	0	2934	76	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 (76) 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:83:N:H4'	1:A:264:N:P	1.80	1.20
1:A:83:N:H4'	1:A:263:N:O3'	0.98	1.15

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:83:N:C4'	1:A:264:N:P	2.35	1.15
1:A:83:N:C5'	1:A:264:N:P	2.35	1.15
1:A:83:N:C4'	1:A:263:N:O3'	1.95	1.13
1:A:195:N:H2'	1:A:196:N:O4'	1.55	1.07
1:A:313:N:H2'	1:A:314:N:O4'	1.55	1.06
1:A:24:N:H2'	1:A:25:N:O4'	1.55	1.05
1:A:36:N:H2'	1:A:37:N:O4'	1.55	1.02
1:A:83:N:H5'	1:A:264:N:P	1.95	1.02
1:A:206:N:P	1:A:220:N:O3'	2.35	0.85
1:A:1:N:P	1:A:33:N:H3'	2.19	0.82
1:A:265:N:H1'2	1:A:266:N:O4'	1.83	0.79
1:A:200:N:H1'2	1:A:201:N:O4'	1.83	0.79
1:A:41:N:H1'2	1:A:42:N:O4'	1.83	0.79
1:A:318:N:H1'2	1:A:319:N:O4'	1.83	0.79
1:A:29:N:H1'2	1:A:30:N:O4'	1.83	0.78
1:A:1:N:H5'	1:A:33:N:C2'	2.20	0.71
2:E:49:ARG:CA	4:E:1:WO2:O44	2.38	0.71
1:A:31:N:H2'	1:A:32:N:O4'	1.91	0.70
1:A:43:N:H2'	1:A:44:N:O4'	1.91	0.70
1:A:202:N:H2'	1:A:203:N:O4'	1.91	0.70
1:A:334:N:H2'	1:A:335:N:O4'	1.92	0.70
1:A:267:N:H2'	1:A:268:N:O4'	1.91	0.70
1:A:320:N:H2'	1:A:321:N:O4'	1.91	0.70
1:A:1:N:H5'	1:A:33:N:H2'	1.74	0.69
1:A:195:N:H2'	1:A:196:N:C1'	2.25	0.66
1:A:313:N:H2'	1:A:314:N:C1'	2.26	0.66
1:A:195:N:C2'	1:A:196:N:O4'	2.39	0.66
1:A:24:N:H2'	1:A:25:N:C1'	2.26	0.65
1:A:36:N:H2'	1:A:37:N:C1'	2.26	0.65
1:A:24:N:C2'	1:A:25:N:O4'	2.39	0.64
1:A:205:N:O3'	1:A:207:N:P	2.57	0.63
1:A:36:N:C2'	1:A:37:N:O4'	2.39	0.62
1:A:313:N:C2'	1:A:314:N:O4'	2.39	0.61
1:A:335:N:H2'	1:A:336:N:O4'	2.01	0.60
1:A:353:N:H2'	1:A:354:N:O4'	2.02	0.59
1:A:200:N:O2'	1:A:201:N:P	2.64	0.56
1:A:41:N:O2'	1:A:42:N:P	2.64	0.56
1:A:29:N:O2'	1:A:30:N:P	2.64	0.56
1:A:318:N:O2'	1:A:319:N:P	2.64	0.55
1:A:350:N:H2'	1:A:351:N:O4'	2.06	0.55
1:A:265:N:O2'	1:A:266:N:P	2.64	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:40:N:H1'2	1:A:41:N:O5'	2.11	0.51
1:A:199:N:H1'2	1:A:200:N:O5'	2.11	0.51
1:A:264:N:H1'2	1:A:265:N:O5'	2.11	0.51
1:A:317:N:H1'2	1:A:318:N:O5'	2.11	0.51
1:A:28:N:H1'2	1:A:29:N:O5'	2.11	0.50
1:A:1:N:H5'	1:A:33:N:C3'	2.42	0.49
1:A:337:N:H2'	1:A:338:N:O4'	2.15	0.47
1:A:202:N:C2'	1:A:203:N:O4'	2.63	0.46
1:A:267:N:C2'	1:A:268:N:O4'	2.63	0.45
1:A:31:N:C2'	1:A:32:N:O4'	2.63	0.45
1:A:320:N:C2'	1:A:321:N:O4'	2.63	0.45
1:A:198:N:O2'	1:A:199:N:C5'	2.65	0.45
1:A:316:N:O2'	1:A:317:N:C5'	2.65	0.45
1:A:27:N:O2'	1:A:28:N:C5'	2.65	0.45
1:A:263:N:O2'	1:A:264:N:C5'	2.65	0.44
1:A:39:N:O2'	1:A:40:N:C5'	2.65	0.43
1:A:42:N:O2'	1:A:43:N:O5'	2.36	0.43
1:A:264:N:H1'2	1:A:265:N:C5'	2.48	0.43
1:A:40:N:H1'2	1:A:41:N:C5'	2.48	0.43
1:A:28:N:H1'2	1:A:29:N:C5'	2.48	0.43
1:A:199:N:H1'2	1:A:200:N:C5'	2.48	0.43
1:A:266:N:O2'	1:A:267:N:O5'	2.35	0.43
1:A:317:N:H1'2	1:A:318:N:C5'	2.48	0.42
1:A:202:N:H2'	1:A:203:N:C1'	2.49	0.42
1:A:30:N:O2'	1:A:31:N:O5'	2.35	0.42
1:A:83:N:H5''	1:A:264:N:P	2.47	0.42
1:A:267:N:H2'	1:A:268:N:C1'	2.49	0.42
1:A:43:N:C2'	1:A:44:N:O4'	2.63	0.41
1:A:31:N:H2'	1:A:32:N:C1'	2.49	0.41
1:A:320:N:H2'	1:A:321:N:C1'	2.50	0.41
1:A:43:N:H2'	1:A:44:N:C1'	2.49	0.41
1:A:319:N:O2'	1:A:320:N:O5'	2.35	0.41
1:A:200:N:O2'	1:A:201:N:O5'	2.38	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

There are no protein backbone outliers to report in this entry.

5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	A	0/362	-	-

There are no RNA backbone outliers to report.

There are no RNA pucker outliers to report.

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](#)

1 ligand is 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	$\# Z > 2$
4	WO2	E	1	-	60,116,116	1.16	9 (15%)	6,348,348	1.07	0

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	E	1	WO2	P1-OP1	2.64	1.58	1.53
4	E	1	WO2	P1-OP3	2.45	1.58	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	E	1	WO2	W17-O27	-2.42	1.59	2.16
4	E	1	WO2	W11-O16	-2.37	1.83	1.94
4	E	1	WO2	W13-O51	-2.34	1.60	2.16
4	E	1	WO2	W5-O16	2.34	2.05	1.94
4	E	1	WO2	P2-OP6	2.33	1.57	1.53
4	E	1	WO2	W15-O49	-2.07	1.67	2.16
4	E	1	WO2	W7-O15	-2.05	1.84	1.94

There are no bond angle outliers.

There are no chirality outliers.

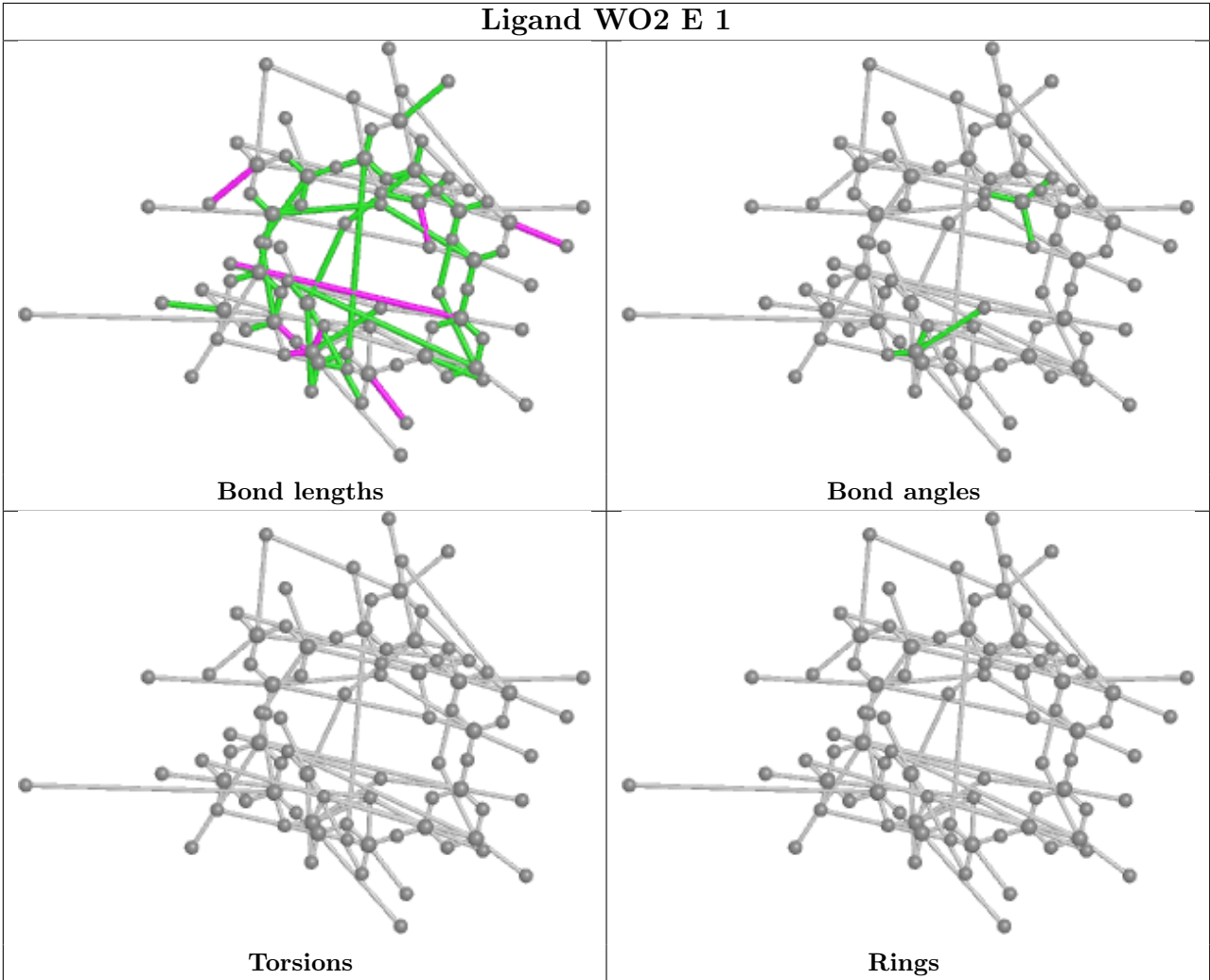
There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	E	1	WO2	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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
1	A	27

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	204:N	O3'	205:N	P	137.12
1	A	332:N	O3'	333:N	P	128.07

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Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	150:N	O3'	151:N	P	111.77
1	A	85:N	O3'	86:N	P	59.85
1	A	74:N	O3'	75:N	P	18.32
1	A	95:N	O3'	96:N	P	18.32
1	A	117:N	O3'	118:N	P	18.32
1	A	139:N	O3'	140:N	P	18.32
1	A	158:N	O3'	159:N	P	18.32
1	A	180:N	O3'	181:N	P	18.32
1	A	342:N	O3'	343:N	P	17.51
1	A	205:N	O3'	206:N	P	17.35
1	A	213:N	O3'	214:N	P	17.35
1	A	11:N	O3'	12:N	P	17.34
1	A	53:N	O3'	54:N	P	17.34
1	A	225:N	O3'	226:N	P	17.33
1	A	235:N	O3'	236:N	P	17.33
1	A	245:N	O3'	246:N	P	17.33
1	A	255:N	O3'	256:N	P	17.33
1	A	274:N	O3'	275:N	P	17.33
1	A	285:N	O3'	286:N	P	17.33
1	A	295:N	O3'	296:N	P	17.33
1	A	305:N	O3'	306:N	P	17.33
1	A	327:N	O3'	328:N	P	17.33
1	A	359:N	O3'	360:N	P	17.33
1	A	206:N	O3'	207:N	P	16.56
1	A	45:N	O3'	46:N	P	9.56

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.