



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 03:16 AM UTC

PDB ID : 1RTA / pdb_00001rta
Title : CRYSTAL STRUCTURE DISPOSITION OF THYMIDYLIC ACID
TETRAMER IN COMPLEX WITH RIBONUCLEASE A
Authors : Birdsall, D.L.; McPherson, A.
Deposited on : 1992-08-28
Resolution : 2.50 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
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

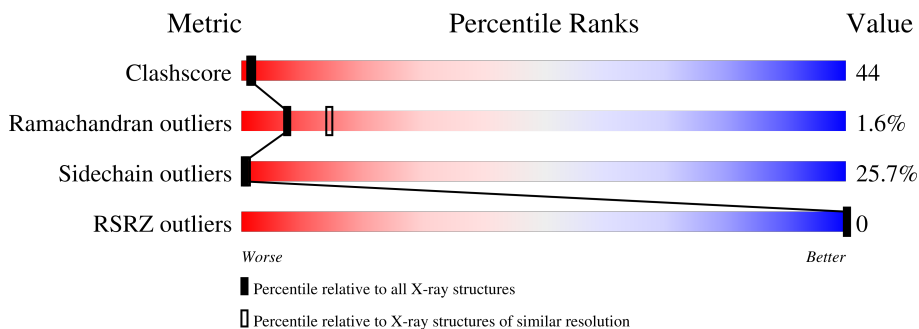
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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 $\leq 5\%$. 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	C	4	
2	E	124	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 1031 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 DNA chain called DNA (5'-D(*TP*TP*TP*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	4	Total	C	N	O	P	0	0	0
			81	40	8	29	4			

- Molecule 2 is a protein called PROTEIN (RIBONUCLEASE A (E.C.3.1.27.5)).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	124	Total	C	N	O	S	0	0	0
			950	575	171	192	12			

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: DNA (5'-D(*TP*TP*TP*T)-3')

Chain C:  75% 25%

T300
T301
T302
T303

- Molecule 2: PROTEIN (RIBONUCLEASE A (E.C.3.1.27.5))

Chain E:  25% 32% 35% 8%

K1 E2 T3 K7 F8 E9 R10 Q11 H12 M13 D14 S15 S16 S23 S24 Y25 C26 N27 Q28 M29 M30 K31 S32 R33 N34 L35 T36 K37 D38 R39 C40 K41 P42 V43 M44 T45 F46 V47 H48 S50 E49 L51 A52 D53 V54 Q55 A56 V57 C58 S59 Q60 K61 N62 V63 A64 C65 K66 N67

G68 Q69 T70 N71 C72 Y73 Q74 S75 Y76 S77 T78 M79 S80 I81 T82 D83 C84 R85 E86 S89 S90 K91 Y92 P93 N94 C95 A96 Y97 K98 T99 T100 Q101 A102 N103 K104 H105 I106 I107 Y108 C110 E111 G112 N113 P114 Y115 V116 P117 V118 H119 F120 V124

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	44.90Å 75.00Å 43.60Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	(Not available) – 2.50 38.52 – 2.50	Depositor EDS
% Data completeness (in resolution range)	(Not available) ((Not available)-2.50) 32.4 (38.52-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1228.42 (at 2.51Å)	Xtriage
Refinement program	X-PLOR, CORELS	Depositor
R, R_{free}	0.235 , (Not available) 0.233 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	7.6	Xtriage
Anisotropy	0.295	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 79.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.25$, $\langle L^2 \rangle = 0.10$	Xtriage
Estimated twinning fraction	0.189 for l,-k,h	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	1031	wwPDB-VP
Average B, all atoms (Å ²)	8.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.01% 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

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	C	2.49	5/88 (5.7%)	3.07	15/132 (11.4%)
2	E	1.73	18/966 (1.9%)	2.63	86/1304 (6.6%)
All	All	1.81	23/1054 (2.2%)	2.68	101/1436 (7.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	1
2	E	0	3
All	All	0	4

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	300	DT	OP3-P	11.01	1.70	1.48
1	C	301	DT	P-O5'	8.50	1.77	1.60
2	E	12	HIS	CG-ND1	-8.14	1.29	1.38
1	C	300	DT	P-O5'	8.12	1.76	1.60
1	C	302	DT	P-O5'	7.55	1.75	1.60
2	E	48	HIS	CD2-NE2	-7.22	1.29	1.37
2	E	39	ARG	NE-CZ	7.03	1.40	1.33
2	E	48	HIS	CG-ND1	-7.00	1.30	1.38
2	E	105	HIS	CG-ND1	-6.75	1.30	1.38
2	E	12	HIS	CD2-NE2	-6.74	1.30	1.37
2	E	116	VAL	CA-CB	6.37	1.62	1.54
2	E	43	VAL	CA-CB	6.36	1.63	1.54
2	E	12	HIS	CA-C	-6.16	1.44	1.52
2	E	119	HIS	CD2-NE2	-5.91	1.31	1.37
2	E	62	ASN	CA-CB	5.71	1.59	1.52
2	E	109	ALA	CA-CB	-5.59	1.45	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	105	HIS	CD2-NE2	-5.48	1.31	1.37
2	E	115	TYR	CA-CB	-5.47	1.45	1.53
2	E	99	THR	CA-CB	5.42	1.59	1.53
2	E	42	PRO	CA-CB	5.33	1.61	1.53
2	E	39	ARG	CA-CB	5.29	1.62	1.53
2	E	119	HIS	CB-CG	5.28	1.57	1.50
1	C	301	DT	C2'-C1'	5.02	1.62	1.52

All (101) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	302	DT	C5'-C4'-C3'	-10.33	99.41	114.90
1	C	302	DT	C2-N1-C1'	-9.64	104.89	119.35
1	C	302	DT	O4'-C1'-N1	9.62	122.83	108.40
2	E	35	LEU	CA-C-O	9.26	130.80	119.78
2	E	111	GLU	O-C-N	9.19	132.48	123.29
2	E	94	ASN	OD1-CG-ND2	-8.77	113.83	122.60
1	C	302	DT	C4'-C3'-O3'	8.70	123.06	110.00
2	E	12	HIS	CA-CB-CG	-8.50	105.30	113.80
2	E	26	CYS	CA-CB-SG	-8.39	95.09	114.40
2	E	34	ASN	CA-CB-CG	8.34	120.94	112.60
2	E	83	ASP	CA-CB-CG	8.18	120.78	112.60
2	E	74	GLN	CA-CB-CG	8.14	130.38	114.10
1	C	303	DT	C4'-C3'-O3'	8.07	122.10	110.00
2	E	33	ARG	CA-CB-CG	-7.90	98.30	114.10
2	E	40	CYS	CA-CB-SG	-7.80	96.46	114.40
2	E	38	ASP	N-CA-C	-7.69	103.86	113.55
2	E	70	THR	N-CA-C	7.62	122.61	113.16
2	E	104	LYS	O-C-N	-7.41	115.06	122.99
1	C	303	DT	O4'-C1'-N1	7.38	119.46	108.40
1	C	301	DT	P-O5'-C5'	-7.29	109.06	120.00
2	E	11	GLN	CA-CB-CG	7.28	128.66	114.10
1	C	300	DT	P-O5'-C5'	-7.23	109.16	120.00
2	E	73	TYR	N-CA-C	7.18	120.22	107.80
2	E	105	HIS	CB-CG-CD2	-7.12	121.94	131.20
2	E	13	MET	N-CA-C	6.99	121.38	107.62
2	E	115	TYR	O-C-N	6.96	131.18	123.04
2	E	11	GLN	CB-CG-CD	-6.95	100.79	112.60
2	E	47	VAL	CA-C-O	6.88	128.49	120.71
2	E	39	ARG	O-C-N	6.75	131.46	122.82
2	E	69	GLN	CA-CB-CG	6.74	127.58	114.10
1	C	300	DT	C4'-C3'-O3'	6.70	120.05	110.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	119	HIS	CB-CG-ND1	6.64	132.66	122.70
2	E	54	VAL	O-C-N	-6.64	114.99	121.83
2	E	94	ASN	CB-CG-ND2	6.61	126.31	116.40
2	E	66	LYS	CA-CB-CG	6.58	127.25	114.10
2	E	77	SER	N-CA-C	6.56	120.17	110.46
2	E	95	CYS	O-C-N	6.50	130.34	123.06
2	E	33	ARG	NE-CZ-NH2	-6.42	113.43	119.20
2	E	106	ILE	CA-CB-CG2	-6.37	99.67	110.50
1	C	303	DT	C2'-C3'-O3'	6.31	120.96	111.50
1	C	302	DT	P-O5'-C5'	-6.29	110.57	120.00
2	E	27	ASN	CB-CG-ND2	6.27	125.81	116.40
2	E	50	SER	CA-C-N	6.15	129.34	120.79
2	E	50	SER	C-N-CA	6.15	129.34	120.79
2	E	80	SER	O-C-N	6.07	130.14	123.04
2	E	30	MET	CG-SD-CE	-6.02	87.66	100.90
2	E	74	GLN	OE1-CD-NE2	-6.01	116.59	122.60
2	E	67	ASN	N-CA-CB	-5.99	101.78	110.47
2	E	50	SER	O-C-N	5.99	130.24	122.87
1	C	302	DT	C6-N1-C1'	5.98	128.32	119.35
2	E	36	THR	N-CA-CB	-5.98	101.57	111.49
2	E	67	ASN	CA-CB-CG	5.95	118.55	112.60
2	E	39	ARG	CA-CB-CG	5.93	125.96	114.10
2	E	39	ARG	CG-CD-NE	5.92	125.03	112.00
2	E	11	GLN	N-CA-CB	-5.88	100.62	110.39
2	E	71	ASN	N-CA-C	5.88	119.22	112.57
2	E	55	GLN	OE1-CD-NE2	-5.85	116.75	122.60
2	E	35	LEU	N-CA-CB	-5.80	102.05	110.70
2	E	24	ASN	CA-CB-CG	-5.76	106.84	112.60
2	E	45	THR	CB-CA-C	5.75	120.66	109.33
2	E	64	ALA	CA-C-O	5.72	127.80	121.56
2	E	89	SER	N-CA-CB	5.72	120.15	110.49
2	E	27	ASN	OD1-CG-ND2	-5.65	116.95	122.60
1	C	303	DT	C5'-C4'-C3'	-5.63	106.45	114.90
2	E	23	SER	CA-C-O	5.55	125.85	119.35
2	E	98	LYS	CB-CG-CD	5.52	124.00	111.30
2	E	93	PRO	CA-C-N	5.49	131.27	122.32
2	E	93	PRO	C-N-CA	5.49	131.27	122.32
2	E	71	ASN	CB-CG-ND2	5.48	124.62	116.40
2	E	94	ASN	CA-CB-CG	5.46	118.06	112.60
2	E	119	HIS	CA-CB-CG	5.45	119.25	113.80
2	E	25	TYR	CA-C-O	-5.45	112.72	120.51
2	E	82	THR	CA-C-O	-5.42	114.96	120.71

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	300	DT	O3'-P-O5'	-5.41	95.89	104.00
2	E	93	PRO	N-CA-C	5.41	126.16	112.10
2	E	48	HIS	CA-CB-CG	5.39	119.19	113.80
2	E	85	ARG	CA-CB-CG	-5.39	103.32	114.10
2	E	69	GLN	OE1-CD-NE2	-5.31	117.29	122.60
2	E	90	SER	N-CA-C	-5.26	102.40	110.30
2	E	120	PHE	N-CA-C	-5.18	100.73	108.96
2	E	46	PHE	CB-CA-C	-5.17	101.87	110.14
1	C	301	DT	C6-N1-C1'	-5.14	111.64	119.35
2	E	40	CYS	N-CA-C	-5.13	97.53	107.44
2	E	7	LYS	N-CA-C	-5.13	105.38	110.97
2	E	106	ILE	CA-CB-CG1	5.13	119.11	110.40
2	E	76	TYR	N-CA-C	-5.12	105.13	111.33
2	E	118	VAL	CB-CA-C	-5.12	106.56	111.06
2	E	73	TYR	CA-C-N	5.11	128.10	120.95
2	E	73	TYR	C-N-CA	5.11	128.10	120.95
2	E	85	ARG	CG-CD-NE	5.09	123.19	112.00
2	E	71	ASN	OD1-CG-ND2	-5.08	117.52	122.60
2	E	86	GLU	CA-CB-CG	5.06	124.22	114.10
2	E	37	LYS	O-C-N	-5.06	116.92	122.38
2	E	24	ASN	CB-CG-ND2	5.05	123.98	116.40
2	E	67	ASN	O-C-N	-5.05	115.88	122.39
2	E	98	LYS	N-CA-CB	-5.03	101.72	110.17
2	E	107	ILE	N-CA-C	-5.03	101.08	108.11
2	E	51	LEU	O-C-N	5.01	127.45	122.08
2	E	86	GLU	CB-CA-C	-5.01	103.10	110.06
2	E	35	LEU	CA-C-N	5.01	130.50	122.14
2	E	35	LEU	C-N-CA	5.01	130.50	122.14

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	302	DT	Sidechain
2	E	25	TYR	Sidechain
2	E	33	ARG	Sidechain
2	E	92	TYR	Sidechain

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	C	81	0	49	43	0
2	E	950	0	907	71	1
All	All	1031	0	956	86	1

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 44.

All (86) 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:C:300:DT:C7	2:E:39:ARG:HD3	1.26	1.56
1:C:300:DT:C7	2:E:39:ARG:CD	2.07	1.33
1:C:300:DT:H72	2:E:39:ARG:CD	1.71	1.16
2:E:54:VAL:O	2:E:57:VAL:HG23	1.50	1.11
1:C:300:DT:OP1	1:C:300:DT:H73	1.52	1.09
1:C:303:DT:H73	2:E:71:ASN:HD21	0.96	1.07
1:C:303:DT:H73	2:E:71:ASN:ND2	1.70	1.06
1:C:303:DT:H71	2:E:111:GLU:OE1	1.57	1.05
1:C:300:DT:H73	2:E:39:ARG:HD3	1.38	1.03
1:C:300:DT:H71	2:E:39:ARG:CD	1.96	0.96
1:C:300:DT:H72	2:E:39:ARG:HD3	0.99	0.95
1:C:303:DT:C7	2:E:71:ASN:HD21	1.83	0.90
1:C:300:DT:C2	1:C:302:DT:OP2	2.25	0.90
2:E:109:ALA:HB3	2:E:119:HIS:HB3	1.55	0.89
2:E:34:ASN:HD22	2:E:37:LYS:HE2	1.39	0.85
1:C:301:DT:H4'	1:C:302:DT:H5''	1.61	0.81
1:C:303:DT:C7	2:E:71:ASN:ND2	2.42	0.80
1:C:301:DT:H5''	1:C:302:DT:OP1	1.84	0.77
2:E:38:ASP:HB2	2:E:39:ARG:NH2	2.00	0.76
1:C:303:DT:C7	2:E:111:GLU:OE1	2.33	0.76
2:E:79:MET:HE2	2:E:106:ILE:HD11	1.69	0.75
1:C:300:DT:H71	2:E:39:ARG:HD2	1.69	0.74
1:C:300:DT:C7	1:C:300:DT:OP1	2.32	0.74
2:E:34:ASN:ND2	2:E:37:LYS:HE2	2.08	0.68
2:E:61:LYS:HB3	2:E:74:GLN:HB3	1.76	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:27:ASN:OD1	2:E:96:ALA:HA	1.94	0.66
2:E:65:CYS:HB2	2:E:69:GLN:H	1.60	0.66
1:C:300:DT:C7	2:E:39:ARG:HD2	2.16	0.66
2:E:106:ILE:C	2:E:107:ILE:HD12	2.20	0.66
2:E:92:TYR:CD1	2:E:93:PRO:HA	2.31	0.65
2:E:53:ASP:O	2:E:56:ALA:HB3	1.97	0.64
2:E:107:ILE:CD1	2:E:107:ILE:N	2.62	0.63
1:C:300:DT:OP1	2:E:39:ARG:HB3	1.99	0.62
1:C:301:DT:C5'	1:C:302:DT:OP1	2.46	0.62
1:C:302:DT:O2	2:E:12:HIS:CE1	2.54	0.61
1:C:302:DT:H4'	1:C:303:DT:OP2	2.02	0.59
2:E:86:GLU:HA	2:E:97:TYR:CD1	2.38	0.58
1:C:302:DT:C4'	1:C:303:DT:OP2	2.51	0.57
1:C:300:DT:H2''	1:C:301:DT:OP1	2.04	0.56
1:C:302:DT:H3	2:E:45:THR:HG23	1.70	0.56
1:C:300:DT:O2	1:C:302:DT:OP2	2.24	0.56
2:E:107:ILE:HD12	2:E:107:ILE:N	2.21	0.56
1:C:301:DT:C4'	1:C:302:DT:OP1	2.54	0.55
2:E:43:VAL:HA	2:E:84:CYS:O	2.06	0.55
2:E:54:VAL:O	2:E:57:VAL:CG2	2.40	0.55
2:E:61:LYS:O	2:E:73:TYR:HA	2.08	0.54
1:C:301:DT:OP1	1:C:301:DT:H3'	2.08	0.53
2:E:38:ASP:HB2	2:E:39:ARG:CZ	2.39	0.52
2:E:14:ASP:OD2	2:E:29:MET:HE3	2.10	0.51
2:E:65:CYS:SG	2:E:69:GLN:HB3	2.51	0.51
2:E:80:SER:HA	2:E:102:ALA:O	2.11	0.51
2:E:81:ILE:HD11	2:E:104:LYS:HG3	1.91	0.51
2:E:2:GLU:OE1	2:E:10:ARG:NH1	2.44	0.50
1:C:301:DT:H4'	1:C:302:DT:C5'	2.35	0.50
1:C:302:DT:H2''	2:E:120:PHE:O	2.12	0.50
2:E:57:VAL:O	2:E:73:TYR:HB3	2.11	0.49
2:E:31:LYS:HA	2:E:36:THR:OG1	2.12	0.49
1:C:302:DT:O2	2:E:12:HIS:HE1	1.95	0.49
2:E:29:MET:HA	2:E:32:SER:HB2	1.96	0.48
1:C:300:DT:OP1	2:E:39:ARG:CB	2.63	0.47
1:C:303:DT:H5'	2:E:119:HIS:HA	1.95	0.47
2:E:25:TYR:CE2	2:E:82:THR:HG21	2.50	0.46
2:E:34:ASN:ND2	2:E:37:LYS:CE	2.79	0.46
2:E:13:MET:HB3	2:E:13:MET:HE3	1.75	0.46
2:E:58:CYS:HB3	2:E:115:TYR:CD2	2.52	0.45
1:C:301:DT:P	2:E:42:PRO:HG2	2.57	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:38:ASP:CB	2:E:39:ARG:NH2	2.76	0.45
2:E:33:ARG:C	2:E:35:LEU:HD12	2.42	0.44
1:C:300:DT:OP1	2:E:39:ARG:HA	2.18	0.44
1:C:300:DT:OP1	2:E:39:ARG:HD3	2.18	0.44
2:E:24:ASN:O	2:E:27:ASN:N	2.50	0.44
1:C:301:DT:H4'	1:C:302:DT:OP1	2.18	0.43
2:E:15:SER:OG	2:E:50:SER:HA	2.18	0.43
1:C:300:DT:C2'	1:C:301:DT:OP1	2.67	0.43
1:C:303:DT:H5''	2:E:118:VAL:O	2.18	0.42
2:E:40:CYS:SG	2:E:92:TYR:HB2	2.59	0.42
2:E:58:CYS:HB3	2:E:115:TYR:HD2	1.83	0.42
2:E:101:GLN:NE2	2:E:102:ALA:N	2.67	0.42
2:E:48:HIS:HB2	2:E:80:SER:OG	2.18	0.42
2:E:74:GLN:OE1	2:E:105:HIS:HB3	2.19	0.42
1:C:300:DT:H72	2:E:39:ARG:NE	2.29	0.42
1:C:303:DT:H5'	2:E:119:HIS:CB	2.50	0.41
1:C:302:DT:H3	2:E:45:THR:CG2	2.34	0.41
2:E:57:VAL:CG1	2:E:75:SER:HA	2.50	0.41
2:E:78:THR:HA	2:E:104:LYS:O	2.22	0.40
2:E:9:GLU:O	2:E:13:MET:HB2	2.20	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:66:LYS:O	2:E:101:GLN:NE2[4_457]	2.17	0.03

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
2	E	122/124 (98%)	108 (88%)	12 (10%)	2 (2%)	7 14

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	E	89	SER
2	E	93	PRO

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
2	E	109/109 (100%)	81 (74%)	28 (26%)	0 1

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	E	2	GLU
2	E	3	THR
2	E	7	LYS
2	E	16	SER
2	E	23	SER
2	E	31	LYS
2	E	33	ARG
2	E	37	LYS
2	E	39	ARG
2	E	40	CYS
2	E	43	VAL
2	E	45	THR
2	E	55	GLN
2	E	57	VAL
2	E	59	SER
2	E	61	LYS
2	E	70	THR
2	E	77	SER
2	E	78	THR
2	E	79	MET
2	E	81	ILE
2	E	86	GLU
2	E	91	LYS

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Mol	Chain	Res	Type
2	E	94	ASN
2	E	100	THR
2	E	107	ILE
2	E	113	ASN
2	E	124	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
2	E	11	GLN
2	E	24	ASN
2	E	34	ASN
2	E	55	GLN
2	E	71	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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 oligosaccharides 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 [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	C	4/4 (100%)	-1.35	0 100 100	26, 39, 40, 41	0
2	E	124/124 (100%)	-1.71	0 100 100	1, 5, 15, 28	0
All	All	128/128 (100%)	-1.70	0 100 100	1, 5, 19, 41	0

There are no RSRZ outliers to report.

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.