



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:39 AM GMT

PDB ID : 2I05
Title : Escherichia Coli Replication Terminator Protein (Tus) Complexed With TerA DNA
Authors : Oakley, A.J.; Mulcair, M.D.; Schaeffer, P.M.; Dixon, N.E.
Deposited on : 2006-08-10
Resolution : 2.60 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
<http://wwpdb.org/validation/2016/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
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

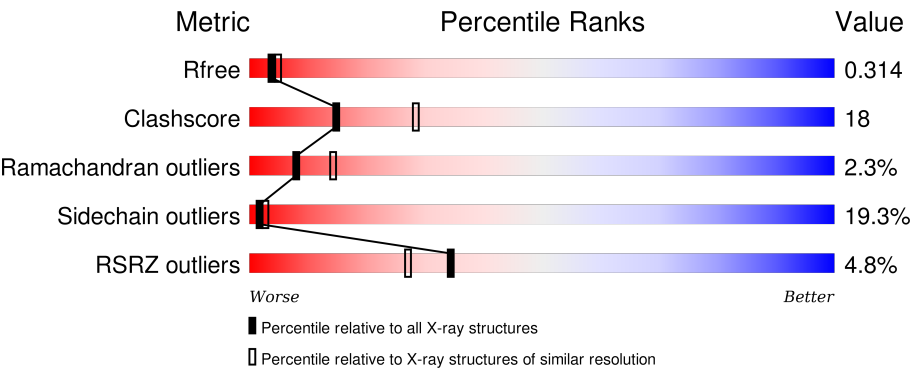
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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}	91344	2328 (2.60-2.60)
Clashscore	102246	2679 (2.60-2.60)
Ramachandran outliers	100387	2635 (2.60-2.60)
Sidechain outliers	100360	2635 (2.60-2.60)
RSRZ outliers	91569	2334 (2.60-2.60)

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. 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	B	16	
2	C	16	
3	A	309	

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 3115 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 5'-D(*T*TP*AP*GP*TP*TP*AP*CP*AP*AP*CP*AP*TP*AP*CP*T)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	15	Total	C	N	O	P	0	0	0
			302	147	54	87	14			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	15	Total	C	N	O	P	0	0	0
			306	149	52	91	14			

- Molecule 3 is a protein called DNA replication terminus site-binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	305	Total	C	N	O	S	0	0	0
			2495	1584	469	439	3			

- Molecule 4 is IODIDE ION (three-letter code: IOD) (formula: I).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	I	0	0
			2	2		

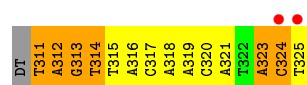
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	8	Total	O	0	0
			8	8		
5	C	2	Total	O	0	0
			2	2		

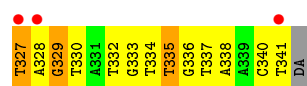
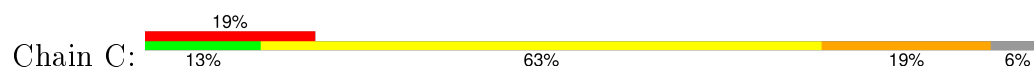
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

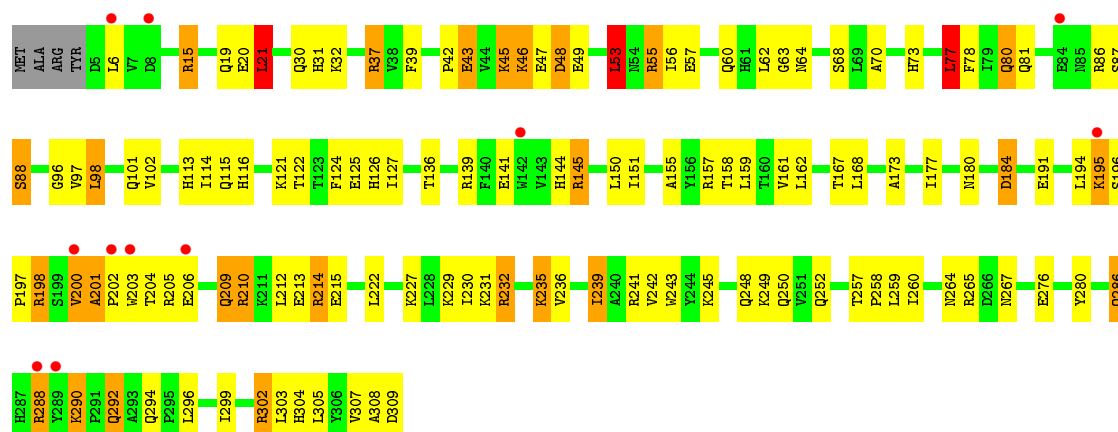
- Molecule 1: 5'-D(*T*TP*AP*GP*TP*TP*AP*CP*AP*AP*CP*AP*TP*AP*CP*T)-3'



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



- Molecule 3: DNA replication terminus site-binding protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	63.37Å 63.37Å 252.65Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.60 36.55 – 2.59	Depositor EDS
% Data completeness (in resolution range)	82.5 (50.00-2.60) 81.7 (36.55-2.59)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.71 (at 2.58Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.226 , 0.312 0.224 , 0.314	Depositor DCC
R_{free} test set	690 reflections (5.24%)	DCC
Wilson B-factor (Å ²)	56.2	Xtriage
Anisotropy	0.506	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 41.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 13904 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3115	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.01% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: IOD

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	B	1.63	2/338 (0.6%)	2.66	32/519 (6.2%)
2	C	1.54	3/342 (0.9%)	2.46	28/527 (5.3%)
3	A	0.88	2/2553 (0.1%)	0.93	3/3465 (0.1%)
All	All	1.07	7/3233 (0.2%)	1.48	63/4511 (1.4%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	141	GLU	CG-CD	6.87	1.62	1.51
2	C	334	DT	C3'-O3'	-6.27	1.35	1.44
2	C	327	DT	C1'-N1	5.49	1.56	1.49
3	A	141	GLU	CB-CG	5.48	1.62	1.52
2	C	335	DT	C2-N3	-5.37	1.33	1.37
1	B	318	DA	C3'-O3'	5.15	1.50	1.44
1	B	316	DA	P-O5'	5.01	1.64	1.59

All (63) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	318	DA	O4'-C1'-N9	17.04	119.93	108.00
1	B	315	DT	O4'-C4'-C3'	-15.67	96.60	106.00
2	C	338	DA	O4'-C1'-N9	13.09	117.17	108.00
2	C	332	DT	O4'-C1'-N1	-11.06	100.25	108.00
1	B	324	DC	O4'-C1'-N1	10.75	115.52	108.00
1	B	317	DC	C1'-O4'-C4'	-10.55	99.55	110.10
2	C	336	DG	O4'-C4'-C3'	-10.30	99.82	106.00
1	B	316	DA	O4'-C1'-N9	10.16	115.11	108.00
2	C	340	DC	O4'-C1'-N1	10.12	115.08	108.00
1	B	312	DA	O4'-C1'-N9	-9.03	101.68	108.00
1	B	316	DA	OP1-P-OP2	7.89	131.44	119.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	332	DT	P-O3'-C3'	7.84	129.11	119.70
2	C	329	DG	O4'-C4'-C3'	-7.67	101.39	106.00
1	B	314	DT	C6-C5-C7	-7.35	118.49	122.90
1	B	323	DA	P-O3'-C3'	7.31	128.47	119.70
1	B	314	DT	N1-C2-N3	7.29	118.98	114.60
2	C	340	DC	P-O3'-C3'	-7.25	111.00	119.70
1	B	317	DC	P-O5'-C5'	-7.01	109.68	120.90
2	C	335	DT	O5'-P-OP1	-6.80	99.58	105.70
2	C	336	DG	C4'-C3'-C2'	-6.58	97.18	103.10
3	A	21	LEU	CA-CB-CG	6.49	130.23	115.30
2	C	335	DT	C5'-C4'-O4'	-6.45	97.05	109.30
2	C	336	DG	C5-C6-O6	6.44	132.47	128.60
1	B	321	DA	O4'-C1'-N9	6.34	112.44	108.00
2	C	338	DA	C1'-O4'-C4'	-6.32	103.78	110.10
1	B	317	DC	O4'-C4'-C3'	-6.29	101.98	104.50
1	B	315	DT	P-O3'-C3'	6.23	127.17	119.70
1	B	316	DA	C8-N9-C4	-6.15	103.34	105.80
1	B	318	DA	P-O3'-C3'	6.11	127.03	119.70
2	C	341	DT	C4-C5-C7	6.07	122.64	119.00
2	C	336	DG	C1'-O4'-C4'	-6.00	104.10	110.10
2	C	330	DT	C5-C4-O4	-5.99	120.70	124.90
2	C	337	DT	N3-C2-O2	-5.97	118.72	122.30
1	B	313	DG	C5-C6-N1	5.92	114.46	111.50
2	C	327	DT	P-O3'-C3'	5.84	126.71	119.70
1	B	314	DT	P-O3'-C3'	-5.75	112.80	119.70
2	C	338	DA	C3'-C2'-C1'	-5.73	95.63	102.50
1	B	320	DC	O4'-C4'-C3'	-5.68	102.23	104.50
1	B	321	DA	O4'-C4'-C3'	-5.62	102.25	104.50
1	B	318	DA	C2-N3-C4	5.62	113.41	110.60
2	C	332	DT	N1-C1'-C2'	5.59	123.22	112.60
1	B	324	DC	P-O3'-C3'	5.58	126.39	119.70
1	B	311	DT	OP2-P-O3'	5.54	117.38	105.20
1	B	316	DA	O5'-P-OP1	-5.50	100.75	105.70
2	C	335	DT	N3-C2-O2	-5.49	119.01	122.30
1	B	321	DA	N1-C6-N6	-5.48	115.31	118.60
2	C	327	DT	C6-N1-C2	-5.48	118.56	121.30
1	B	317	DC	O4'-C1'-N1	5.45	111.81	108.00
2	C	333	DG	N1-C6-O6	-5.40	116.66	119.90
2	C	333	DG	O4'-C1'-N9	5.35	111.75	108.00
1	B	319	DA	O4'-C1'-C2'	5.35	110.18	105.90
1	B	311	DT	O3'-P-O5'	-5.32	93.89	104.00
2	C	336	DG	P-O5'-C5'	-5.21	112.57	120.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	332	DT	C1'-O4'-C4'	-5.20	104.90	110.10
2	C	327	DT	N3-C4-O4	5.17	123.00	119.90
2	C	341	DT	C6-C5-C7	-5.17	119.80	122.90
2	C	332	DT	C4'-C3'-O3'	-5.15	96.83	109.70
1	B	323	DA	OP1-P-OP2	5.12	127.27	119.60
1	B	323	DA	P-O5'-C5'	-5.11	112.72	120.90
3	A	53	LEU	CA-CB-CG	5.08	126.99	115.30
1	B	313	DG	C2-N3-C4	5.04	114.42	111.90
1	B	316	DA	P-O3'-C3'	5.00	125.70	119.70
3	A	77	LEU	CA-CB-CG	5.00	126.81	115.30

There are no chirality outliers.

There are no planarity outliers.

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	B	302	0	172	8	0
2	C	306	0	174	4	0
3	A	2495	0	2533	99	0
4	A	2	0	0	1	0
5	A	8	0	0	1	0
5	C	2	0	0	0	0
All	All	3115	0	2879	106	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 18.

All (106) 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)
3:A:198:ARG:HG3	3:A:198:ARG:HH11	1.11	1.04
1:B:324:DC:H2''	1:B:325:DT:H5'	1.45	0.99
3:A:195:LYS:HB3	3:A:195:LYS:NZ	1.79	0.94
3:A:55:ARG:HH11	3:A:55:ARG:CG	1.83	0.92

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:55:ARG:HH11	3:A:55:ARG:HG2	1.37	0.90
3:A:73:HIS:NE2	3:A:257:THR:HG22	1.87	0.89
3:A:215:GLU:OE1	3:A:302:ARG:NH1	2.08	0.86
3:A:78:PHE:H	3:A:81:GLN:NE2	1.75	0.84
3:A:121:LYS:HE2	3:A:155:ALA:O	1.79	0.82
3:A:198:ARG:HG3	3:A:198:ARG:NH1	1.91	0.78
3:A:195:LYS:HZ1	3:A:195:LYS:HB3	1.50	0.77
3:A:159:LEU:HD22	3:A:260:ILE:HD11	1.66	0.77
3:A:198:ARG:HH11	3:A:198:ARG:CG	1.94	0.76
3:A:210:ARG:HG3	3:A:210:ARG:HH11	1.52	0.74
3:A:292:GLN:HE21	3:A:292:GLN:H	1.38	0.71
3:A:55:ARG:HD2	3:A:56:ILE:N	2.06	0.71
3:A:286:GLN:HA	3:A:286:GLN:HE21	1.55	0.71
3:A:241:ARG:HD2	3:A:250:GLN:HE22	1.56	0.70
3:A:195:LYS:O	3:A:197:PRO:HD3	1.95	0.67
3:A:184:ASP:N	3:A:184:ASP:OD2	2.28	0.67
3:A:173:ALA:HB2	3:A:239:ILE:HD13	1.76	0.66
3:A:55:ARG:NH1	3:A:55:ARG:HG2	2.07	0.66
3:A:145:ARG:HH11	3:A:145:ARG:HB2	1.60	0.66
1:B:311:DT:O2	3:A:288:ARG:HD2	1.96	0.65
3:A:31:HIS:HD2	3:A:101:GLN:O	1.79	0.65
3:A:30:GLN:OE1	4:A:1002:IOD:I	2.85	0.65
3:A:77:LEU:H	3:A:81:GLN:HE22	1.47	0.63
3:A:78:PHE:H	3:A:81:GLN:HE21	1.47	0.62
3:A:102:VAL:O	3:A:264:ASN:HA	2.00	0.60
3:A:78:PHE:N	3:A:81:GLN:NE2	2.49	0.59
3:A:264:ASN:HD22	3:A:267:ASN:CG	2.05	0.59
3:A:46:LYS:HD3	3:A:87:SER:HA	1.84	0.59
3:A:286:GLN:HA	3:A:286:GLN:NE2	2.17	0.59
3:A:200:VAL:HG12	3:A:201:ALA:N	2.18	0.59
3:A:180:ASN:HD21	3:A:229:LYS:HZ3	1.51	0.58
3:A:136:THR:HA	3:A:139:ARG:HD2	1.84	0.58
3:A:201:ALA:HB3	3:A:202:PRO:CD	2.34	0.58
3:A:144:HIS:HE1	3:A:150:LEU:O	1.88	0.57
3:A:159:LEU:HD22	3:A:260:ILE:CD1	2.36	0.56
3:A:55:ARG:HD2	3:A:56:ILE:H	1.70	0.55
3:A:161:VAL:HA	3:A:260:ILE:O	2.07	0.55
3:A:210:ARG:HG3	3:A:210:ARG:NH1	2.20	0.54
3:A:55:ARG:HD3	3:A:280:TYR:O	2.07	0.54
3:A:78:PHE:HB2	3:A:81:GLN:HE21	1.73	0.53
3:A:210:ARG:HH22	3:A:214:ARG:HG3	1.73	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:327:DT:H2''	2:C:328:DA:N7	2.23	0.53
3:A:47:GLU:HG3	3:A:48:ASP:OD1	2.08	0.53
3:A:43:GLU:HG2	3:A:45:LYS:HD3	1.91	0.53
3:A:195:LYS:CB	3:A:195:LYS:NZ	2.65	0.52
3:A:15:ARG:N	3:A:15:ARG:HD2	2.24	0.52
3:A:202:PRO:HD2	3:A:203:TRP:CE3	2.45	0.52
3:A:39:PHE:CE1	3:A:60:GLN:HG3	2.45	0.52
3:A:45:LYS:H	3:A:45:LYS:HE3	1.75	0.51
3:A:46:LYS:CD	3:A:87:SER:HA	2.40	0.51
3:A:70:ALA:O	3:A:73:HIS:HB3	2.10	0.51
3:A:144:HIS:CE1	3:A:150:LEU:O	2.63	0.51
3:A:201:ALA:CB	3:A:202:PRO:CD	2.89	0.50
3:A:243:TRP:CE2	3:A:250:GLN:HB3	2.45	0.50
3:A:209:GLN:O	3:A:213:GLU:HG3	2.12	0.50
3:A:197:PRO:HB2	3:A:205:ARG:HD2	1.93	0.50
1:B:313:DG:H2''	1:B:314:DT:OP2	2.12	0.50
1:B:314:DT:H4'	3:A:290:LYS:HD2	1.93	0.50
3:A:124:PHE:HA	3:A:127:ILE:HD12	1.95	0.49
2:C:327:DT:H2''	2:C:328:DA:C8	2.48	0.49
3:A:60:GLN:OE1	3:A:62:LEU:HD11	2.12	0.48
3:A:206:GLU:HA	3:A:209:GLN:HE22	1.78	0.48
3:A:78:PHE:N	3:A:81:GLN:HE21	2.08	0.48
3:A:162:LEU:HD11	3:A:259:LEU:CD1	2.43	0.48
3:A:168:LEU:HD13	3:A:242:VAL:HG22	1.95	0.48
3:A:230:ILE:HA	3:A:304:HIS:O	2.15	0.47
3:A:167:THR:C	3:A:168:LEU:HD22	2.35	0.46
3:A:37:ARG:HH11	3:A:60:GLN:NE2	2.13	0.46
3:A:43:GLU:HB2	3:A:86:ARG:NH2	2.31	0.46
1:B:323:DA:P	3:A:88:SER:HG	2.39	0.46
3:A:55:ARG:NH1	3:A:55:ARG:CG	2.55	0.45
2:C:335:DT:OP1	3:A:157:ARG:NH1	2.49	0.45
3:A:42:PRO:HD3	3:A:57:GLU:O	2.16	0.45
3:A:97:VAL:HG22	3:A:98:LEU:N	2.32	0.44
3:A:198:ARG:CD	3:A:200:VAL:HG22	2.46	0.44
3:A:53:LEU:HD22	3:A:235:LYS:HE2	2.00	0.44
3:A:162:LEU:HD11	3:A:259:LEU:HD13	1.98	0.44
2:C:328:DA:H2''	2:C:329:DG:OP2	2.18	0.44
1:B:324:DC:H2''	1:B:325:DT:C5'	2.31	0.44
3:A:198:ARG:HD3	3:A:200:VAL:HG22	2.00	0.43
3:A:236:VAL:O	3:A:236:VAL:HG13	2.18	0.43
3:A:55:ARG:HH11	3:A:55:ARG:HG3	1.75	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:150:LEU:HG	3:A:151:ILE:N	2.34	0.43
3:A:232:ARG:NH1	5:A:1003:HOH:O	2.52	0.43
3:A:55:ARG:CD	3:A:280:TYR:O	2.66	0.43
3:A:126:HIS:O	3:A:127:ILE:C	2.57	0.42
3:A:37:ARG:NH1	3:A:60:GLN:NE2	2.68	0.42
3:A:201:ALA:CB	3:A:202:PRO:HD3	2.50	0.42
3:A:114:ILE:C	3:A:116:HIS:N	2.73	0.42
3:A:201:ALA:HB3	3:A:202:PRO:HD3	2.02	0.41
3:A:46:LYS:HG3	3:A:49:GLU:OE1	2.19	0.41
3:A:229:LYS:NZ	3:A:294:GLN:HE22	2.18	0.41
3:A:210:ARG:O	3:A:213:GLU:HB2	2.20	0.41
1:B:312:DA:OP2	3:A:241:ARG:NE	2.46	0.41
3:A:296:LEU:HD23	3:A:308:ALA:HA	2.03	0.41
3:A:194:LEU:HG	3:A:212:LEU:HD23	2.03	0.41
3:A:37:ARG:HH12	3:A:276:GLU:HB3	1.86	0.40
3:A:20:GLU:O	3:A:21:LEU:C	2.58	0.40
3:A:96:GLY:O	3:A:258:PRO:HA	2.21	0.40
3:A:37:ARG:NH1	3:A:276:GLU:HB3	2.36	0.40
3:A:161:VAL:HG13	3:A:161:VAL:O	2.21	0.40
1:B:311:DT:H2"	1:B:312:DA:C8	2.56	0.40

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
3	A	303/309 (98%)	276 (91%)	20 (7%)	7 (2%)	8 14

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	200	VAL

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Mol	Chain	Res	Type
3	A	201	ALA
3	A	80	GLN
3	A	196	SER
3	A	248	GLN
3	A	204	THR
3	A	63	GLY

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
3	A	270/273 (99%)	218 (81%)	52 (19%)	2 3

All (52) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	6	LEU
3	A	15	ARG
3	A	19	GLN
3	A	21	LEU
3	A	32	LYS
3	A	37	ARG
3	A	43	GLU
3	A	45	LYS
3	A	46	LYS
3	A	48	ASP
3	A	53	LEU
3	A	55	ARG
3	A	64	ASN
3	A	68	SER
3	A	77	LEU
3	A	80	GLN
3	A	88	SER
3	A	98	LEU
3	A	113	HIS
3	A	115	GLN

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Mol	Chain	Res	Type
3	A	122	THR
3	A	125	GLU
3	A	145	ARG
3	A	158	THR
3	A	177	ILE
3	A	184	ASP
3	A	191	GLU
3	A	195	LYS
3	A	198	ARG
3	A	209	GLN
3	A	210	ARG
3	A	214	ARG
3	A	222	LEU
3	A	227	LYS
3	A	231	LYS
3	A	232	ARG
3	A	235	LYS
3	A	239	ILE
3	A	245	LYS
3	A	249	LYS
3	A	252	GLN
3	A	265	ARG
3	A	286	GLN
3	A	288	ARG
3	A	290	LYS
3	A	292	GLN
3	A	299	ILE
3	A	302	ARG
3	A	303	LEU
3	A	305	LEU
3	A	307	VAL
3	A	309	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
3	A	31	HIS
3	A	51	ASN
3	A	64	ASN
3	A	80	GLN
3	A	81	GLN
3	A	82	GLN

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Mol	Chain	Res	Type
3	A	144	HIS
3	A	176	HIS
3	A	180	ASN
3	A	209	GLN
3	A	237	GLN
3	A	250	GLN
3	A	264	ASN
3	A	286	GLN
3	A	292	GLN
3	A	294	GLN

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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

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	B	15/16 (93%)	0.21	2 (13%) 4 3	41, 51, 104, 106	0
2	C	15/16 (93%)	0.42	3 (20%) 1 1	41, 54, 108, 110	0
3	A	305/309 (98%)	0.14	11 (3%) 46 38	33, 49, 72, 86	0
All	All	335/341 (98%)	0.16	16 (4%) 34 27	33, 49, 74, 110	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	C	327	DT	6.0
1	B	324	DC	5.1
3	A	288	ARG	4.8
1	B	325	DT	4.2
2	C	328	DA	4.1
3	A	289	TYR	3.9
3	A	84	GLU	3.5
3	A	206	GLU	3.4
3	A	195	LYS	3.2
3	A	203	TRP	3.2
3	A	202	PRO	3.0
3	A	6	LEU	2.8
3	A	8	ASP	2.5
3	A	200	VAL	2.3
3	A	142	TRP	2.2
2	C	341	DT	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 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
4	IOD	A	1001	1/1	0.99	0.11	-1.45	38,38,38,38	1
4	IOD	A	1002	1/1	0.94	0.09	-5.62	79,79,79,79	1

6.5 Other polymers [i](#)

There are no such residues in this entry.