



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 2, 2016 – 12:00 AM GMT

PDB ID : 7ICR
Title : DNA POLYMERASE BETA (E.C.2.7.7.7)/DNA COMPLEX, SOAKED IN
THE PRESENCE OF ZNCL2
Authors : Pelletier, H.; Sawaya, M.R.
Deposited on : 1996-04-19
Resolution : 3.00 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

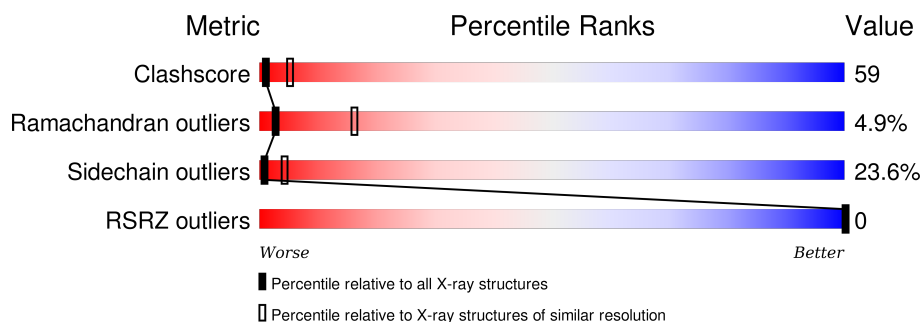
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION



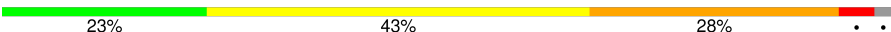
The reported resolution of this entry is 3.00 Å.

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	102246	1912 (3.00-3.00)
Ramachandran outliers	100387	1853 (3.00-3.00)
Sidechain outliers	100360	1856 (3.00-3.00)
RSRZ outliers	91569	1592 (3.00-3.00)

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	T	7	
2	P	6	
3	A	335	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	NA	A	341	-	-	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 3029 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(*CP*AP*TP*CP*TP*GP*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	T	7	Total	C	N	O	P	0	0	0
			122	58	20	38	6			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	P	6	Total	C	N	O	P	0	0	0
			126	59	25	36	6			

- Molecule 3 is a protein called PROTEIN (DNA POLYMERASE BETA (E.C.2.7.7.7)).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	327	Total	C	N	O	S	26	0	0
			2623	1657	458	499	9			

- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 5 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	2	Total	Na	0	0
			2	2		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	121	Total	O	0	0
			121	121		

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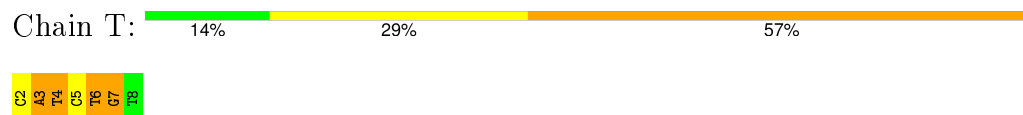
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	P	22	Total 22	O 22	0	0
6	T	11	Total 11	O 11	0	0

3 Residue-property plots

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.

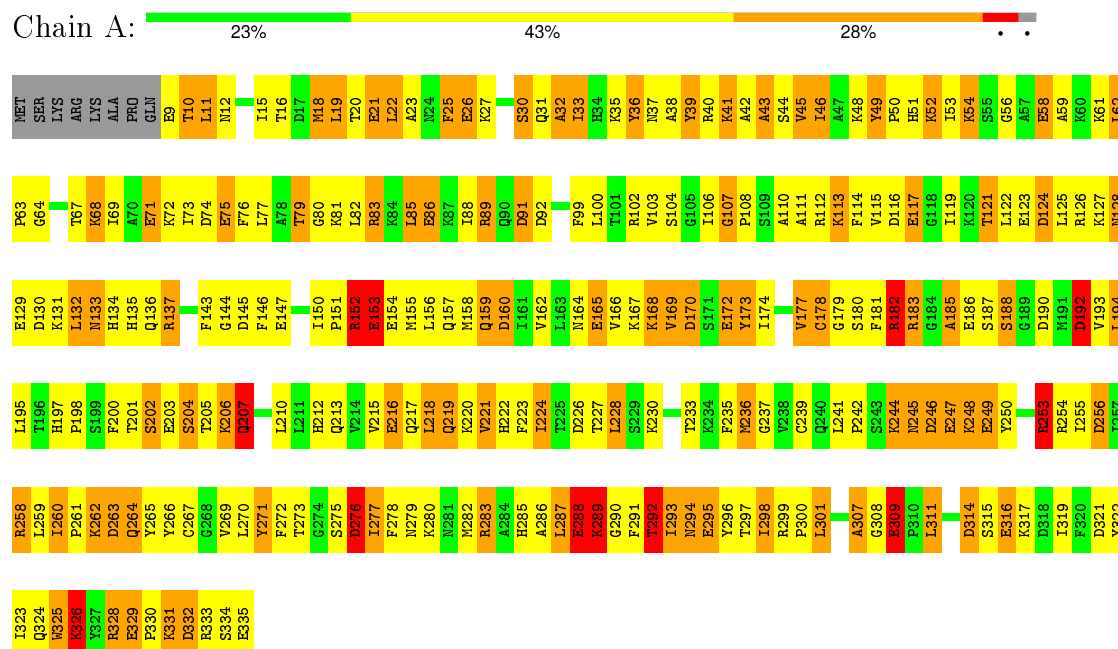
- Molecule 1: DNA (5'-D(*CP*AP*TP*CP*TP*GP*T)-3')



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



- Molecule 3: PROTEIN (DNA POLYMERASE BETA (E.C.2.7.7.7))



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	179.28 Å 57.74 Å 48.48 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 3.00 11.88 – 2.70	Depositor EDS
% Data completeness (in resolution range)	92.0 (20.00-3.00) 88.3 (11.88-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.52 (at 2.70 Å)	Xtriage
Refinement program	TNT V. 5-D	Depositor
R, R_{free}	0.160 , (Not available) 0.154 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	37.3	Xtriage
Anisotropy	0.059	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.16 , 131.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Outliers	1 of 12632 reflections (0.008%)	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3029	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.17% 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: NA, ZN

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	T	2.07	7/135 (5.2%)	2.81	14/207 (6.8%)
2	P	2.17	5/141 (3.5%)	3.50	22/214 (10.3%)
3	A	1.23	23/2672 (0.9%)	1.75	63/3590 (1.8%)
All	All	1.34	35/2948 (1.2%)	1.95	99/4011 (2.5%)

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
3	A	1	0

All (35) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	58	GLU	CD-OE1	7.68	1.34	1.25
2	P	5	DT	N1-C2	7.63	1.44	1.38
2	P	5	DT	C4-C5	7.47	1.51	1.45
3	A	129	GLU	CD-OE1	7.27	1.33	1.25
3	A	249	GLU	CD-OE2	7.08	1.33	1.25
3	A	309	GLU	CD-OE2	7.05	1.33	1.25
3	A	203	GLU	CD-OE1	6.90	1.33	1.25
3	A	75	GLU	CD-OE1	6.81	1.33	1.25
3	A	86	GLU	CD-OE1	6.80	1.33	1.25
2	P	5	DT	C1'-N1	6.79	1.58	1.49
1	T	4	DT	C1'-N1	6.42	1.57	1.49
3	A	117	GLU	CD-OE2	6.39	1.32	1.25
3	A	21	GLU	CD-OE1	6.37	1.32	1.25
3	A	329	GLU	CD-OE2	6.30	1.32	1.25
3	A	216	GLU	CD-OE2	6.16	1.32	1.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	T	7	DG	N3-C4	-6.10	1.31	1.35
3	A	165	GLU	CD-OE2	5.96	1.32	1.25
3	A	123	GLU	CD-OE1	5.75	1.31	1.25
3	A	71	GLU	CD-OE1	5.70	1.31	1.25
3	A	295	GLU	CD-OE1	5.64	1.31	1.25
1	T	7	DG	C1'-N9	-5.56	1.39	1.47
1	T	6	DT	C3'-O3'	-5.54	1.36	1.44
3	A	316	GLU	CD-OE1	5.47	1.31	1.25
3	A	288	GLU	CD-OE2	5.44	1.31	1.25
2	P	2	DA	C3'-O3'	5.43	1.51	1.44
3	A	153	GLU	CD-OE2	5.40	1.31	1.25
3	A	172	GLU	CD-OE2	5.34	1.31	1.25
2	P	4	DA	C1'-N9	-5.27	1.39	1.47
1	T	7	DG	C3'-O3'	-5.25	1.37	1.44
3	A	186	GLU	CD-OE1	5.25	1.31	1.25
3	A	247	GLU	CD-OE1	5.23	1.31	1.25
3	A	26	GLU	CD-OE1	5.22	1.31	1.25
3	A	9	GLU	CD-OE2	5.21	1.31	1.25
1	T	3	DA	N9-C4	5.15	1.41	1.37
1	T	7	DG	N9-C4	-5.12	1.33	1.38

All (99) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	P	6	DG	C8-N9-C1'	19.89	152.86	127.00
2	P	6	DG	C4-N9-C1'	-19.28	101.44	126.50
1	T	7	DG	C4-N9-C1'	-14.46	107.70	126.50
2	P	5	DT	C6-N1-C1'	-13.56	100.05	120.40
1	T	7	DG	C8-N9-C1'	13.53	144.59	127.00
2	P	5	DT	C2-N1-C1'	12.57	138.32	118.20
3	A	263	ASP	CB-CG-OD2	-11.00	108.40	118.30
2	P	4	DA	O4'-C1'-N9	-10.67	100.53	108.00
1	T	4	DT	P-O3'-C3'	9.92	131.60	119.70
1	T	4	DT	C6-N1-C1'	-9.53	106.11	120.40
3	A	192	ASP	CB-CG-OD2	-9.37	109.87	118.30
3	A	192	ASP	N-CA-CB	9.36	127.45	110.60
1	T	4	DT	C2-N1-C1'	9.07	132.72	118.20
3	A	263	ASP	CB-CG-OD1	8.96	126.37	118.30
3	A	190	ASP	CB-CG-OD2	-8.66	110.50	118.30
3	A	124	ASP	CB-CG-OD2	-8.62	110.54	118.30
1	T	6	DT	C6-N1-C1'	-8.44	107.73	120.40
2	P	2	DA	C4'-C3'-C2'	-8.41	95.53	103.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	182	ARG	NE-CZ-NH1	8.33	124.47	120.30
3	A	126	ARG	NE-CZ-NH1	8.29	124.44	120.30
3	A	271	TYR	CA-CB-CG	-8.23	97.76	113.40
3	A	253	ARG	NE-CZ-NH1	8.16	124.38	120.30
3	A	226	ASP	CB-CG-OD2	-7.90	111.19	118.30
3	A	145	ASP	CB-CG-OD2	-7.89	111.19	118.30
1	T	4	DT	O4'-C1'-N1	7.76	113.44	108.00
2	P	2	DA	O4'-C4'-C3'	-7.72	101.37	106.00
2	P	5	DT	N3-C4-O4	-7.60	115.34	119.90
1	T	2	DC	C2-N1-C1'	7.59	127.15	118.80
3	A	256	ASP	CB-CG-OD1	7.57	125.11	118.30
1	T	6	DT	C2-N1-C1'	7.51	130.21	118.20
3	A	126	ARG	NE-CZ-NH2	-7.43	116.58	120.30
2	P	6	DG	P-O5'-C5'	-7.43	109.02	120.90
2	P	2	DA	P-O5'-C5'	-7.32	109.19	120.90
3	A	328	ARG	NE-CZ-NH2	-7.20	116.70	120.30
2	P	5	DT	N3-C2-O2	-7.17	118.00	122.30
3	A	192	ASP	CB-CA-C	-7.12	96.17	110.40
3	A	74	ASP	CB-CG-OD1	7.05	124.65	118.30
3	A	256	ASP	CB-CG-OD2	-7.04	111.96	118.30
3	A	246	ASP	CB-CG-OD1	-7.03	111.97	118.30
2	P	3	DG	P-O3'-C3'	7.00	128.11	119.70
2	P	5	DT	C6-C5-C7	-6.89	118.77	122.90
3	A	92	ASP	CB-CG-OD2	-6.88	112.10	118.30
3	A	170	ASP	CB-CG-OD1	6.84	124.45	118.30
3	A	116	ASP	CB-CG-OD2	-6.78	112.20	118.30
3	A	226	ASP	CB-CG-OD1	6.66	124.29	118.30
3	A	124	ASP	CB-CG-OD1	6.58	124.22	118.30
3	A	102	ARG	NE-CZ-NH1	-6.52	117.04	120.30
1	T	3	DA	C4-N9-C1'	6.39	137.80	126.30
3	A	91	ASP	CB-CG-OD1	-6.31	112.62	118.30
2	P	1	DC	C6-N1-C1'	-6.27	113.27	120.80
3	A	292	THR	N-CA-CB	6.18	122.04	110.30
1	T	3	DA	C8-N9-C1'	-6.18	116.58	127.70
3	A	152	ARG	N-CA-CB	6.17	121.70	110.60
2	P	1	DC	C2-N1-C1'	6.14	125.56	118.80
3	A	74	ASP	CB-CG-OD2	-6.12	112.79	118.30
3	A	332	ASP	CB-CG-OD1	-6.10	112.81	118.30
3	A	192	ASP	CB-CG-OD1	6.09	123.78	118.30
1	T	2	DC	C6-N1-C1'	-6.08	113.50	120.80
3	A	326	LYS	CD-CE-NZ	6.06	125.64	111.70
2	P	4	DA	N1-C6-N6	6.05	122.23	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	130	ASP	CB-CG-OD1	6.02	123.72	118.30
2	P	1	DC	P-O5'-C5'	-6.01	111.29	120.90
3	A	314	ASP	CB-CG-OD1	-5.93	112.97	118.30
3	A	130	ASP	CB-CG-OD2	-5.82	113.06	118.30
3	A	183	ARG	NE-CZ-NH1	5.82	123.21	120.30
3	A	246	ASP	CB-CG-OD2	5.80	123.52	118.30
2	P	4	DA	C8-N9-C4	5.79	108.11	105.80
3	A	20	THR	N-CA-CB	5.78	121.29	110.30
3	A	107	GLY	N-CA-C	-5.78	98.66	113.10
3	A	43	ALA	CB-CA-C	5.76	118.74	110.10
3	A	224	ILE	CA-CB-CG1	-5.73	100.11	111.00
2	P	5	DT	C4-C5-C7	5.70	122.42	119.00
3	A	332	ASP	CB-CG-OD2	5.70	123.43	118.30
3	A	121	THR	CA-CB-CG2	-5.64	104.51	112.40
1	T	7	DG	P-O5'-C5'	5.55	129.79	120.90
3	A	116	ASP	CB-CG-OD1	5.53	123.28	118.30
2	P	1	DC	P-O3'-C3'	5.52	126.32	119.70
3	A	137	ARG	NE-CZ-NH2	-5.50	117.55	120.30
1	T	7	DG	P-O3'-C3'	-5.47	113.13	119.70
3	A	83	ARG	NE-CZ-NH1	5.47	123.03	120.30
3	A	321	ASP	CB-CG-OD2	-5.45	113.40	118.30
3	A	271	TYR	CB-CG-CD1	-5.41	117.75	121.00
3	A	222	HIS	CA-CB-CG	-5.38	104.45	113.60
3	A	49	TYR	C-N-CD	-5.38	108.77	120.60
3	A	283	ARG	NE-CZ-NH2	-5.37	117.61	120.30
3	A	99	PHE	CA-CB-CG	-5.35	101.06	113.90
3	A	307	ALA	N-CA-CB	5.35	117.59	110.10
3	A	221	VAL	CA-CB-CG1	-5.31	102.93	110.90
3	A	160	ASP	CB-CG-OD2	-5.23	113.60	118.30
3	A	170	ASP	CB-CG-OD2	-5.22	113.60	118.30
3	A	89	ARG	NE-CZ-NH1	5.20	122.90	120.30
3	A	276	ASP	CB-CG-OD2	-5.17	113.65	118.30
3	A	276	ASP	CB-CG-OD1	5.15	122.94	118.30
2	P	4	DA	N9-C4-C5	-5.13	103.75	105.80
3	A	190	ASP	CB-CG-OD1	5.12	122.90	118.30
3	A	177	VAL	CB-CA-C	5.09	121.08	111.40
3	A	102	ARG	CD-NE-CZ	-5.06	116.52	123.60
3	A	39	TYR	CB-CG-CD2	-5.06	117.97	121.00
2	P	5	DT	C1'-O4'-C4'	-5.01	105.09	110.10

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	A	292	THR	CA

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	T	122	0	69	4	0
2	P	126	0	68	13	0
3	A	2623	0	2641	314	2
4	A	2	0	0	0	0
5	A	2	0	0	0	0
6	A	121	0	0	21	0
6	P	22	0	0	5	0
6	T	11	0	0	1	0
All	All	3029	0	2778	330	2

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 59.

All (330) 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:151:PRO:HG2	3:A:154:GLU:HG3	1.35	1.07
3:A:23:ALA:HB2	3:A:39:TYR:HB2	1.30	1.07
3:A:15:ILE:HD11	3:A:77:LEU:HD11	1.38	1.06
3:A:119:ILE:HG23	3:A:124:ASP:HB3	1.37	1.04
2:P:1:DC:H2''	2:P:2:DA:H5''	1.38	1.01
3:A:293:ILE:HG23	3:A:298:ILE:HG13	1.49	0.91
3:A:182:ARG:HG2	3:A:182:ARG:HH11	1.36	0.90
3:A:172:GLU:HG2	3:A:198:PRO:HG2	1.51	0.90
3:A:330:PRO:HA	3:A:333:ARG:HG3	1.56	0.88
3:A:201:THR:HG22	3:A:204:SER:HB3	1.56	0.88
3:A:31:GLN:NE2	3:A:112:ARG:HH12	1.72	0.86
3:A:330:PRO:HA	3:A:333:ARG:CG	2.06	0.86
2:P:5:DT:H2''	2:P:6:DG:H5'	1.58	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:177:VAL:HG22	3:A:193:VAL:HG22	1.57	0.84
2:P:1:DC:C2'	2:P:2:DA:H5''	2.08	0.83
3:A:201:THR:CG2	3:A:204:SER:HB3	2.08	0.83
3:A:25:PHE:CE2	3:A:88:ILE:HG12	2.15	0.82
3:A:254:ARG:HH11	3:A:255:ILE:H	1.26	0.82
3:A:254:ARG:NH1	3:A:255:ILE:H	1.78	0.82
3:A:41:LYS:HE2	3:A:64:GLY:HA2	1.62	0.80
3:A:12:ASN:HB3	3:A:46:ILE:CD1	2.11	0.80
3:A:68:LYS:O	3:A:72:LYS:HE3	1.82	0.80
3:A:73:ILE:HG22	3:A:77:LEU:HD13	1.64	0.80
3:A:68:LYS:HB2	3:A:68:LYS:NZ	1.95	0.80
3:A:15:ILE:O	3:A:19:LEU:HD22	1.82	0.79
3:A:286:ALA:CB	3:A:293:ILE:HD11	2.12	0.79
3:A:31:GLN:NE2	3:A:112:ARG:NH1	2.29	0.79
3:A:155:MET:CE	3:A:188:SER:HB2	2.13	0.78
3:A:103:VAL:HG22	3:A:143:PHE:CE1	2.18	0.78
3:A:32:ALA:HB1	3:A:35:LYS:HB2	1.64	0.78
3:A:218:LEU:CB	3:A:224:ILE:HD12	2.13	0.78
3:A:207:GLN:O	3:A:210:LEU:HB2	1.83	0.77
3:A:23:ALA:CB	3:A:39:TYR:HB2	2.15	0.77
3:A:162:VAL:O	3:A:166:VAL:HG23	1.84	0.77
3:A:212:HIS:HB3	6:A:541:HOH:O	1.84	0.76
3:A:119:ILE:CG2	3:A:124:ASP:HB3	2.14	0.76
3:A:166:VAL:O	3:A:169:VAL:HG12	1.86	0.76
3:A:155:MET:HE3	3:A:188:SER:HB2	1.68	0.76
3:A:52:LYS:HB2	6:A:554:HOH:O	1.85	0.76
3:A:178:CYS:SG	3:A:194:LEU:HD22	2.26	0.76
3:A:31:GLN:HE21	3:A:112:ARG:NH1	1.83	0.75
3:A:286:ALA:HB1	3:A:293:ILE:HD11	1.67	0.75
3:A:271:TYR:HB2	6:A:592:HOH:O	1.85	0.75
3:A:41:LYS:HD3	3:A:42:ALA:N	2.02	0.75
3:A:68:LYS:O	3:A:71:GLU:HB3	1.87	0.74
3:A:289:LYS:HD2	3:A:324:GLN:OE1	1.87	0.74
3:A:235:PHE:CZ	3:A:237:GLY:HA3	2.22	0.74
3:A:23:ALA:HB2	3:A:39:TYR:CB	2.14	0.74
3:A:103:VAL:HB	3:A:106:ILE:HD12	1.70	0.74
3:A:331:LYS:HD2	3:A:332:ASP:N	2.01	0.74
3:A:151:PRO:HG2	3:A:154:GLU:CG	2.15	0.73
3:A:279:ASN:O	3:A:283:ARG:HG3	1.89	0.73
3:A:194:LEU:CD1	3:A:258:ARG:HD3	2.19	0.72
3:A:201:THR:HA	3:A:261:PRO:HB3	1.70	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:227:THR:HG23	3:A:235:PHE:CE1	2.24	0.72
3:A:278:PHE:HB2	3:A:333:ARG:O	1.89	0.71
3:A:244:LYS:HB3	3:A:245:ASN:ND2	2.05	0.71
3:A:245:ASN:N	3:A:245:ASN:HD22	1.87	0.71
3:A:264:GLN:HB2	3:A:296:TYR:O	1.91	0.71
3:A:114:PHE:HB3	3:A:119:ILE:HB	1.71	0.71
3:A:132:LEU:HD23	3:A:132:LEU:N	2.05	0.71
2:P:5:DT:H2''	2:P:6:DG:C5'	2.19	0.70
6:P:568:HOH:O	3:A:110:ALA:HB2	1.90	0.70
3:A:218:LEU:HB2	3:A:224:ILE:HD12	1.73	0.70
3:A:278:PHE:CE1	3:A:328:ARG:HD2	2.26	0.70
3:A:254:ARG:HH11	3:A:255:ILE:N	1.89	0.70
3:A:155:MET:HA	3:A:158:MET:HE3	1.74	0.69
3:A:108:PRO:O	3:A:112:ARG:HG3	1.92	0.69
3:A:319:ILE:O	3:A:322:TYR:HB2	1.93	0.69
3:A:41:LYS:O	3:A:45:VAL:HG13	1.93	0.68
3:A:262:LYS:HG3	3:A:262:LYS:O	1.93	0.68
3:A:254:ARG:NH1	3:A:255:ILE:N	2.41	0.68
3:A:82:LEU:HB3	3:A:85:LEU:HB2	1.75	0.68
3:A:218:LEU:HB3	3:A:224:ILE:HD12	1.75	0.68
3:A:227:THR:HG23	3:A:235:PHE:HE1	1.57	0.68
3:A:154:GLU:O	3:A:158:MET:HG3	1.93	0.67
3:A:31:GLN:HE21	3:A:112:ARG:HH12	1.41	0.67
3:A:12:ASN:HB3	3:A:46:ILE:HD12	1.75	0.67
3:A:150:ILE:HG21	3:A:158:MET:HE1	1.77	0.67
3:A:293:ILE:HG23	3:A:298:ILE:CG1	2.23	0.67
3:A:152:ARG:NH2	3:A:181:PHE:O	2.29	0.66
3:A:31:GLN:N	6:A:641:HOH:O	2.27	0.66
3:A:202:SER:HB2	3:A:263:ASP:CG	2.16	0.66
3:A:293:ILE:CG2	3:A:298:ILE:HG13	2.24	0.66
3:A:215:VAL:O	3:A:219:GLN:HG2	1.96	0.66
3:A:286:ALA:HA	3:A:323:ILE:HG21	1.79	0.65
3:A:26:GLU:OE1	3:A:32:ALA:HB3	1.97	0.65
3:A:277:ILE:HG13	3:A:335:GLU:HB3	1.79	0.65
3:A:253:ARG:NH1	6:A:503:HOH:O	2.29	0.64
3:A:49:TYR:CG	3:A:50:PRO:HD2	2.31	0.64
3:A:202:SER:HB2	3:A:263:ASP:OD1	1.98	0.64
3:A:12:ASN:HD21	3:A:53:ILE:H	1.46	0.64
3:A:21:GLU:OE2	3:A:89:ARG:NH2	2.30	0.64
3:A:18:MET:HE1	3:A:76:PHE:HB2	1.79	0.64
3:A:242:PRO:HG2	6:A:589:HOH:O	1.97	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:56:GLY:O	3:A:59:ALA:HB3	1.97	0.64
3:A:326:LYS:O	3:A:326:LYS:HG3	1.97	0.64
3:A:42:ALA:O	3:A:46:ILE:HG23	1.97	0.64
3:A:124:ASP:O	3:A:128:ASN:ND2	2.29	0.63
3:A:287:LEU:O	3:A:287:LEU:HD12	1.97	0.63
3:A:330:PRO:HA	3:A:333:ARG:HG2	1.80	0.63
3:A:41:LYS:HE2	3:A:64:GLY:CA	2.29	0.63
2:P:2:DA:C8	2:P:2:DA:H5'	2.34	0.63
3:A:286:ALA:HA	3:A:323:ILE:CG2	2.28	0.62
3:A:12:ASN:HB3	3:A:46:ILE:HD11	1.81	0.62
3:A:267:CYS:SG	3:A:297:THR:HA	2.39	0.62
3:A:286:ALA:CB	3:A:323:ILE:HG21	2.30	0.62
3:A:293:ILE:CD1	3:A:298:ILE:HG13	2.30	0.62
3:A:197:HIS:CE1	3:A:198:PRO:HD2	2.35	0.62
3:A:25:PHE:CD2	3:A:88:ILE:HG12	2.34	0.62
3:A:172:GLU:CG	3:A:198:PRO:HG2	2.26	0.62
1:T:4:DT:H2''	1:T:5:DC:O5'	2.00	0.62
3:A:18:MET:HE3	3:A:76:PHE:CG	2.35	0.62
3:A:103:VAL:HG22	3:A:143:PHE:HE1	1.62	0.62
3:A:113:LYS:O	3:A:117:GLU:HG3	2.00	0.62
3:A:308:GLY:O	3:A:309:GLU:HB2	1.99	0.62
3:A:273:THR:HB	6:A:584:HOH:O	2.00	0.61
3:A:228:LEU:HB2	3:A:236:MET:O	2.00	0.61
3:A:194:LEU:HD12	3:A:258:ARG:HD3	1.82	0.61
3:A:202:SER:HB2	3:A:263:ASP:OD2	2.00	0.61
3:A:293:ILE:HD13	3:A:298:ILE:HG13	1.81	0.61
3:A:133:ASN:HD22	3:A:134:HIS:N	1.98	0.61
3:A:294:ASN:HB2	3:A:295:GLU:OE2	2.01	0.61
3:A:285:HIS:CE1	3:A:289:LYS:HE3	2.35	0.61
3:A:253:ARG:HH11	3:A:253:ARG:CG	2.14	0.61
3:A:288:GLU:HA	3:A:288:GLU:OE1	1.97	0.60
3:A:197:HIS:ND1	3:A:198:PRO:HD2	2.16	0.60
3:A:293:ILE:HD13	3:A:298:ILE:CG1	2.30	0.60
3:A:244:LYS:HB3	3:A:245:ASN:HD22	1.67	0.59
3:A:216:GLU:O	3:A:220:LYS:N	2.32	0.59
3:A:150:ILE:HD12	3:A:155:MET:HE1	1.85	0.59
3:A:195:LEU:O	3:A:260:ILE:N	2.37	0.57
3:A:41:LYS:HD3	3:A:42:ALA:H	1.68	0.57
3:A:23:ALA:O	3:A:36:TYR:HD2	1.87	0.57
3:A:298:ILE:O	3:A:311:LEU:HB2	2.04	0.57
3:A:245:ASN:N	3:A:245:ASN:ND2	2.53	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:291:PHE:HD2	3:A:323:ILE:HG22	1.70	0.56
3:A:182:ARG:NH1	3:A:182:ARG:HG2	2.11	0.56
3:A:172:GLU:HB3	3:A:198:PRO:CD	2.35	0.56
3:A:194:LEU:HD11	3:A:258:ARG:HD3	1.88	0.56
3:A:68:LYS:HB2	3:A:68:LYS:HZ1	1.67	0.56
3:A:299:ARG:HB3	3:A:300:PRO:HD2	1.87	0.56
3:A:276:ASP:O	3:A:280:LYS:HG3	2.06	0.56
3:A:248:LYS:HG2	3:A:248:LYS:O	2.05	0.56
3:A:210:LEU:HB3	3:A:259:LEU:HD21	1.87	0.56
3:A:18:MET:HG2	3:A:22:LEU:CD2	2.36	0.55
3:A:49:TYR:CD2	3:A:50:PRO:HD2	2.41	0.55
3:A:125:LEU:O	3:A:128:ASN:N	2.37	0.55
3:A:58:GLU:O	3:A:61:LYS:HB2	2.06	0.55
3:A:48:LYS:HG3	6:A:607:HOH:O	2.07	0.55
3:A:11:LEU:O	3:A:12:ASN:ND2	2.40	0.55
3:A:133:ASN:ND2	3:A:135:HIS:H	2.05	0.55
3:A:206:LYS:O	3:A:207:GLN:HB2	2.05	0.55
3:A:52:LYS:HG2	3:A:52:LYS:O	2.06	0.55
3:A:165:GLU:HA	3:A:168:LYS:HG3	1.88	0.54
3:A:152:ARG:HA	3:A:155:MET:HB2	1.88	0.54
3:A:19:LEU:O	3:A:39:TYR:HB3	2.08	0.54
3:A:217:GLN:O	3:A:221:VAL:HG22	2.08	0.54
2:P:4:DA:H2'	2:P:5:DT:H71	1.89	0.54
3:A:300:PRO:HD3	3:A:311:LEU:HD13	1.91	0.53
3:A:329:GLU:O	3:A:333:ARG:HG2	2.07	0.53
2:P:5:DT:H1'	6:P:565:HOH:O	2.08	0.53
3:A:30:SER:HB3	6:A:561:HOH:O	2.07	0.53
3:A:155:MET:HE2	3:A:188:SER:HB2	1.89	0.53
3:A:309:GLU:OE1	3:A:309:GLU:HA	2.08	0.53
3:A:275:SER:HB2	3:A:277:ILE:HD13	1.91	0.53
3:A:286:ALA:CA	3:A:323:ILE:HG21	2.39	0.53
3:A:295:GLU:HA	6:A:592:HOH:O	2.09	0.52
3:A:12:ASN:CB	3:A:46:ILE:HD12	2.40	0.52
3:A:69:ILE:O	3:A:73:ILE:HG13	2.09	0.52
3:A:286:ALA:HB2	3:A:293:ILE:HD11	1.90	0.52
3:A:166:VAL:HG12	3:A:173:TYR:HB3	1.91	0.52
3:A:18:MET:CE	3:A:76:PHE:HB2	2.40	0.52
3:A:73:ILE:HG22	3:A:77:LEU:CD1	2.39	0.51
3:A:106:ILE:HG22	3:A:107:GLY:O	2.10	0.51
3:A:182:ARG:NH1	3:A:273:THR:OG1	2.44	0.51
3:A:260:ILE:HG22	3:A:261:PRO:HD2	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:266:TYR:HA	3:A:269:VAL:HB	1.92	0.51
3:A:255:ILE:HG12	3:A:256:ASP:N	2.26	0.51
3:A:286:ALA:O	3:A:291:PHE:HB2	2.10	0.51
3:A:106:ILE:HG23	3:A:110:ALA:HB3	1.94	0.51
3:A:26:GLU:HA	3:A:30:SER:OG	2.10	0.51
3:A:172:GLU:HG2	3:A:198:PRO:CG	2.32	0.50
3:A:111:ALA:O	3:A:115:VAL:HG23	2.11	0.50
3:A:278:PHE:CZ	3:A:328:ARG:HD2	2.46	0.50
3:A:133:ASN:O	3:A:137:ARG:HG3	2.10	0.50
2:P:3:DG:H1'	6:P:511:HOH:O	2.10	0.50
3:A:151:PRO:HB2	3:A:153:GLU:HG2	1.94	0.50
3:A:38:ALA:O	3:A:41:LYS:HD2	2.11	0.50
3:A:277:ILE:HG13	3:A:335:GLU:CB	2.42	0.49
3:A:150:ILE:HG21	3:A:158:MET:CE	2.42	0.49
3:A:291:PHE:CD2	3:A:323:ILE:HG22	2.46	0.49
3:A:223:PHE:O	3:A:239:CYS:HB2	2.12	0.49
3:A:121:THR:O	3:A:124:ASP:HB2	2.12	0.49
3:A:174:ILE:O	3:A:195:LEU:HD12	2.12	0.49
3:A:119:ILE:HG21	3:A:125:LEU:HD23	1.94	0.49
3:A:134:HIS:HE1	3:A:227:THR:O	1.95	0.49
3:A:59:ALA:O	3:A:62:LEU:N	2.30	0.49
3:A:182:ARG:HG2	3:A:273:THR:HG23	1.94	0.49
3:A:253:ARG:HG3	3:A:253:ARG:HH11	1.78	0.49
3:A:205:THR:O	3:A:205:THR:HG23	2.12	0.49
3:A:244:LYS:C	3:A:245:ASN:HD22	2.16	0.48
3:A:200:PHE:HA	6:A:625:HOH:O	2.13	0.48
3:A:159:GLN:HG2	3:A:159:GLN:O	2.13	0.48
3:A:197:HIS:CG	3:A:198:PRO:HD2	2.49	0.48
3:A:166:VAL:HG12	3:A:173:TYR:CB	2.43	0.48
3:A:11:LEU:HD23	3:A:11:LEU:H	1.78	0.48
3:A:32:ALA:HB1	3:A:35:LYS:CB	2.41	0.48
3:A:201:THR:HA	3:A:261:PRO:CB	2.42	0.48
3:A:317:LYS:HG3	6:A:501:HOH:O	2.13	0.48
3:A:297:THR:HB	6:A:626:HOH:O	2.14	0.47
3:A:288:GLU:C	3:A:290:GLY:H	2.18	0.47
3:A:27:LYS:HB2	3:A:36:TYR:CG	2.50	0.47
3:A:291:PHE:CD1	3:A:300:PRO:HA	2.50	0.47
3:A:114:PHE:CD1	3:A:119:ILE:HD12	2.50	0.47
3:A:16:THR:HG23	3:A:46:ILE:HG13	1.96	0.47
3:A:172:GLU:O	3:A:198:PRO:HD3	2.14	0.47
2:P:5:DT:P	3:A:107:GLY:HA3	2.55	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:275:SER:CB	3:A:277:ILE:HD13	2.45	0.47
3:A:164:ASN:O	3:A:167:LYS:HB3	2.15	0.47
3:A:183:ARG:HD3	3:A:273:THR:O	2.15	0.47
3:A:114:PHE:O	3:A:119:ILE:N	2.45	0.47
3:A:12:ASN:HD21	3:A:52:LYS:HA	1.80	0.47
3:A:41:LYS:NZ	3:A:64:GLY:O	2.38	0.47
3:A:26:GLU:OE1	3:A:26:GLU:HA	2.13	0.47
2:P:2:DA:H5'	2:P:2:DA:H2'	1.53	0.47
3:A:182:ARG:NH1	3:A:182:ARG:CG	2.76	0.47
3:A:241:LEU:O	3:A:250:TYR:HD2	1.97	0.47
3:A:218:LEU:HB3	3:A:224:ILE:CD1	2.44	0.46
3:A:267:CYS:O	3:A:271:TYR:HB2	2.16	0.46
3:A:85:LEU:O	3:A:89:ARG:HG3	2.15	0.46
3:A:218:LEU:CD1	3:A:218:LEU:N	2.78	0.46
3:A:125:LEU:HD23	3:A:125:LEU:HA	1.34	0.46
3:A:19:LEU:CB	3:A:43:ALA:HB2	2.46	0.46
3:A:137:ARG:HB3	6:A:622:HOH:O	2.16	0.46
3:A:75:GLU:O	3:A:79:THR:HG23	2.16	0.46
3:A:288:GLU:O	3:A:290:GLY:N	2.49	0.46
3:A:299:ARG:HB3	3:A:300:PRO:CD	2.46	0.46
3:A:41:LYS:O	3:A:44:SER:HB3	2.16	0.46
3:A:280:LYS:O	3:A:283:ARG:N	2.48	0.46
3:A:85:LEU:HA	3:A:85:LEU:HD12	1.65	0.46
3:A:144:GLY:O	3:A:147:GLU:N	2.40	0.46
3:A:201:THR:HG23	3:A:204:SER:HB3	1.91	0.45
3:A:197:HIS:CG	3:A:198:PRO:CD	2.99	0.45
3:A:201:THR:HG22	3:A:204:SER:CB	2.35	0.45
3:A:157:GLN:O	3:A:160:ASP:HB3	2.16	0.45
3:A:300:PRO:HD3	3:A:309:GLU:O	2.16	0.45
3:A:194:LEU:HA	3:A:194:LEU:HD12	1.62	0.45
3:A:296:TYR:HB2	3:A:297:THR:HG23	1.99	0.45
3:A:18:MET:CE	3:A:76:PHE:CG	2.99	0.45
3:A:213:GLN:HB2	3:A:213:GLN:HE21	1.41	0.45
3:A:18:MET:CE	3:A:76:PHE:CD1	2.99	0.45
3:A:292:THR:O	3:A:298:ILE:HA	2.17	0.45
3:A:31:GLN:O	3:A:33:ILE:N	2.50	0.45
3:A:253:ARG:NH1	3:A:253:ARG:CG	2.79	0.45
3:A:217:GLN:O	3:A:220:LYS:HB3	2.17	0.45
3:A:295:GLU:H	3:A:295:GLU:CD	2.21	0.44
3:A:156:LEU:HD23	3:A:181:PHE:CE1	2.52	0.44
3:A:179:GLY:O	3:A:182:ARG:HB3	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:333:ARG:NH1	6:A:584:HOH:O	2.49	0.44
3:A:287:LEU:HD13	3:A:287:LEU:HA	1.55	0.44
3:A:180:SER:HB2	3:A:185:ALA:CB	2.47	0.44
3:A:228:LEU:HA	3:A:228:LEU:HD12	1.65	0.44
3:A:254:ARG:NH1	3:A:254:ARG:HB3	2.33	0.44
3:A:169:VAL:CG1	3:A:170:ASP:N	2.80	0.44
3:A:301:LEU:HD12	3:A:301:LEU:HA	1.55	0.44
3:A:81:LYS:NZ	3:A:86:GLU:OE1	2.47	0.44
3:A:316:GLU:CD	3:A:333:ARG:HH22	2.20	0.44
3:A:127:LYS:HB2	3:A:128:ASN:ND2	2.33	0.44
3:A:26:GLU:HG3	3:A:35:LYS:HB3	2.00	0.44
3:A:19:LEU:HA	3:A:19:LEU:HD12	1.57	0.43
1:T:3:DA:H1'	6:T:617:HOH:O	2.17	0.43
3:A:227:THR:CG2	3:A:230:LYS:HB2	2.48	0.43
3:A:156:LEU:CD2	3:A:181:PHE:HE1	2.31	0.43
3:A:270:LEU:HD21	3:A:282:MET:CE	2.48	0.43
3:A:193:VAL:N	6:A:664:HOH:O	2.51	0.43
3:A:210:LEU:HA	3:A:210:LEU:HD23	1.57	0.43
3:A:54:LYS:HD2	3:A:54:LYS:HA	1.85	0.43
3:A:317:LYS:HD2	3:A:317:LYS:HA	1.50	0.43
3:A:133:ASN:ND2	3:A:135:HIS:N	2.67	0.43
3:A:62:LEU:HA	3:A:63:PRO:HD2	1.70	0.43
3:A:172:GLU:HB3	3:A:198:PRO:CG	2.48	0.43
3:A:258:ARG:O	3:A:258:ARG:HG2	2.19	0.43
3:A:289:LYS:HD3	3:A:289:LYS:HA	1.63	0.43
3:A:12:ASN:HA	6:A:642:HOH:O	2.19	0.43
3:A:194:LEU:HD23	3:A:269:VAL:HA	2.01	0.43
3:A:322:TYR:C	3:A:324:GLN:H	2.21	0.43
3:A:132:LEU:HB3	3:A:136:GLN:HE21	1.84	0.43
3:A:22:LEU:HD13	3:A:85:LEU:CD1	2.48	0.43
3:A:270:LEU:HD21	3:A:282:MET:HE1	2.01	0.43
3:A:19:LEU:HB3	3:A:43:ALA:HB2	2.00	0.42
3:A:25:PHE:C	3:A:25:PHE:CD1	2.92	0.42
3:A:169:VAL:HG13	3:A:170:ASP:N	2.33	0.42
3:A:22:LEU:HD11	3:A:85:LEU:HD13	2.01	0.42
3:A:325:TRP:HA	6:A:583:HOH:O	2.20	0.42
3:A:132:LEU:HB3	3:A:136:GLN:HB3	2.01	0.42
3:A:272:PHE:HD1	6:A:587:HOH:O	2.01	0.42
1:T:6:DT:H1'	1:T:7:DG:H5'	2.02	0.42
3:A:255:ILE:HG23	3:A:255:ILE:O	2.19	0.42
3:A:146:PHE:HA	3:A:146:PHE:HD1	1.33	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:P:1:DC:H5	6:P:572:HOH:O	2.03	0.42
3:A:210:LEU:HD12	6:A:625:HOH:O	2.18	0.42
3:A:331:LYS:CD	3:A:332:ASP:N	2.79	0.42
3:A:265:TYR:O	3:A:269:VAL:N	2.45	0.42
3:A:172:GLU:CG	3:A:198:PRO:CG	2.96	0.41
2:P:3:DG:N3	6:P:511:HOH:O	2.37	0.41
3:A:49:TYR:HA	3:A:50:PRO:HD3	1.24	0.41
3:A:294:ASN:ND2	3:A:297:THR:H	2.17	0.41
3:A:49:TYR:CZ	3:A:51:HIS:HB2	2.55	0.41
3:A:286:ALA:O	3:A:291:PHE:N	2.53	0.41
3:A:16:THR:CG2	3:A:46:ILE:HG13	2.51	0.41
3:A:18:MET:O	3:A:22:LEU:HD22	2.20	0.41
3:A:156:LEU:HD23	3:A:181:PHE:HE1	1.85	0.41
2:P:5:DT:H2''	2:P:6:DG:C8	2.56	0.41
3:A:128:ASN:ND2	3:A:128:ASN:N	2.69	0.41
3:A:236:MET:HG2	3:A:254:ARG:NH2	2.36	0.41
3:A:230:LYS:HB3	3:A:230:LYS:HE3	1.53	0.41
3:A:292:THR:O	3:A:298:ILE:HG12	2.20	0.41
3:A:299:ARG:CB	3:A:300:PRO:CD	2.99	0.41
3:A:315:SER:C	3:A:317:LYS:N	2.74	0.41
3:A:40:ARG:HB2	3:A:40:ARG:HE	1.73	0.41
3:A:155:MET:O	3:A:158:MET:HB2	2.21	0.41
3:A:174:ILE:HG21	3:A:174:ILE:HD13	1.78	0.41
3:A:16:THR:HG23	3:A:46:ILE:CG1	2.51	0.40
3:A:11:LEU:CD2	3:A:11:LEU:H	2.35	0.40
1:T:5:DC:H2''	1:T:6:DT:O5'	2.20	0.40
3:A:192:ASP:HB3	6:A:664:HOH:O	2.21	0.40
3:A:18:MET:O	3:A:21:GLU:HB2	2.21	0.40
3:A:104:SER:OG	3:A:135:HIS:NE2	2.40	0.40
3:A:49:TYR:CE1	3:A:51:HIS:HB2	2.56	0.40

All (2) 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)
3:A:83:ARG:NH1	3:A:117:GLU:CB[3_558]	1.73	0.47
3:A:83:ARG:NH1	3:A:117:GLU:CA[3_558]	2.15	0.05

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	325/335 (97%)	269 (83%)	40 (12%)	16 (5%)	3	16

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	10	THR
3	A	32	ALA
3	A	206	LYS
3	A	207	GLN
3	A	244	LYS
3	A	246	ASP
3	A	247	GLU
3	A	309	GLU
3	A	80	GLY
3	A	185	ALA
3	A	289	LYS
3	A	307	ALA
3	A	334	SER
3	A	178	CYS
3	A	264	GLN
3	A	91	ASP

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.

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
3	A	288/295 (98%)	220 (76%)	68 (24%)	1 4

All (68) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	10	THR
3	A	11	LEU
3	A	18	MET
3	A	19	LEU
3	A	22	LEU
3	A	25	PHE
3	A	30	SER
3	A	33	ILE
3	A	36	TYR
3	A	37	ASN
3	A	41	LYS
3	A	45	VAL
3	A	46	ILE
3	A	52	LYS
3	A	54	LYS
3	A	62	LEU
3	A	67	THR
3	A	68	LYS
3	A	79	THR
3	A	85	LEU
3	A	100	LEU
3	A	113	LYS
3	A	122	LEU
3	A	128	ASN
3	A	131	LYS
3	A	132	LEU
3	A	133	ASN
3	A	152	ARG
3	A	153	GLU
3	A	159	GLN
3	A	168	LYS
3	A	169	VAL
3	A	173	TYR
3	A	182	ARG
3	A	187	SER

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Mol	Chain	Res	Type
3	A	188	SER
3	A	192	ASP
3	A	194	LEU
3	A	202	SER
3	A	204	SER
3	A	207	GLN
3	A	218	LEU
3	A	219	GLN
3	A	228	LEU
3	A	233	THR
3	A	236	MET
3	A	245	ASN
3	A	248	LYS
3	A	249	GLU
3	A	253	ARG
3	A	258	ARG
3	A	260	ILE
3	A	262	LYS
3	A	276	ASP
3	A	277	ILE
3	A	287	LEU
3	A	288	GLU
3	A	289	LYS
3	A	292	THR
3	A	293	ILE
3	A	294	ASN
3	A	298	ILE
3	A	301	LEU
3	A	311	LEU
3	A	314	ASP
3	A	325	TRP
3	A	326	LYS
3	A	331	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
3	A	12	ASN
3	A	28	ASN
3	A	31	GLN
3	A	37	ASN
3	A	133	ASN

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Mol	Chain	Res	Type
3	A	134	HIS
3	A	136	GLN
3	A	157	GLN
3	A	213	GLN
3	A	217	GLN
3	A	219	GLN
3	A	245	ASN
3	A	264	GLN
3	A	279	ASN
3	A	294	ASN

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 4 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 [i](#)

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	T	7/7 (100%)	-1.28	0 100 100	15, 26, 59, 100	0
2	P	6/6 (100%)	-1.45	0 100 100	13, 23, 31, 36	0
3	A	324/335 (96%)	-1.22	0 100 100	3, 34, 86, 100	0
All	All	337/348 (96%)	-1.23	0 100 100	3, 33, 86, 100	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 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(Å ²)	Q<0.9
5	NA	A	341	1/1	0.95	0.12	5.74	18,18,18,18	0
5	NA	A	342	1/1	0.99	0.09	-0.28	24,24,24,24	0
4	ZN	A	343	1/1	0.98	0.04	-	60,60,60,60	0
4	ZN	A	340	1/1	0.97	0.20	-	87,87,87,87	0

6.5 Other polymers [i](#)

There are no such residues in this entry.