



# Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2020 – 11:46 am BST

PDB ID : 7ICN  
Title : DNA POLYMERASE BETA (E.C.2.7.7.7)/DNA COMPLEX, SOAKED IN  
THE PRESENCE OF NICL2  
Authors : Pelletier, H.; Sawaya, M.R.  
Deposited on : 1996-04-19  
Resolution : 2.80 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](mailto: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.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

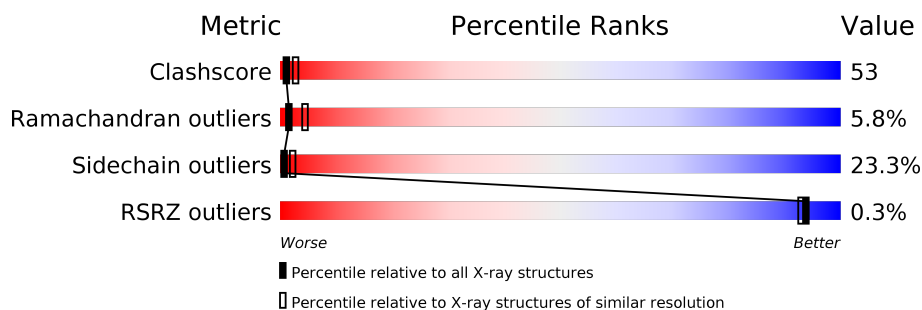
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*



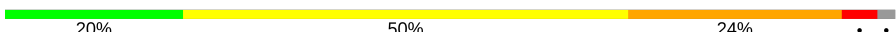
The reported resolution of this entry is 2.80 Å.

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	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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  $\geq 3$ , 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	T	7	
2	P	6	
3	A	335	

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 3018 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 SODIUM ION (three-letter code: NA) (formula: Na).

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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	T	14	Total	O	0	0
			14	14		
5	P	18	Total	O	0	0
			18	18		
5	A	113	Total	O	0	0
			113	113		

### 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 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.

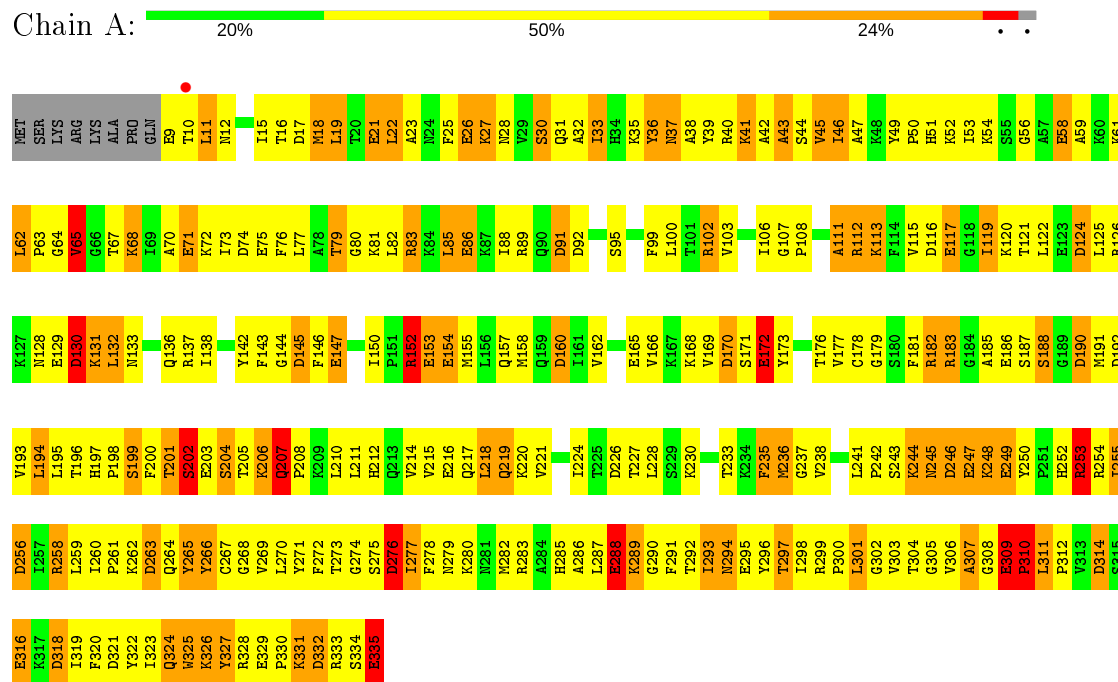
- Molecule 1: DNA (5'-D(\*CP\*AP\*TP\*CP\*TP\*GP\*T)-3')



- 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 2 <sub>1</sub> 2 <sub>1</sub> 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	178.57Å 57.75Å 48.51Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.80 11.93 – 2.79	Depositor EDS
% Data completeness (in resolution range)	89.0 (20.00-2.80) 74.0 (11.93-2.79)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	3.50	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.25 (at 2.78Å)	Xtriage
Refinement program	TNT 5-D	Depositor
R, $R_{free}$	0.162 , (Not available) 0.187 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	43.8	Xtriage
Anisotropy	0.216	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.12 , 74.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	3018	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.81% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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 ⓘ

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

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	2.25	1/135 (0.7%)	2.55	12/207 (5.8%)
2	P	2.97	10/141 (7.1%)	3.55	21/214 (9.8%)
3	A	1.29	26/2672 (1.0%)	1.79	60/3590 (1.7%)
All	All	1.47	37/2948 (1.3%)	1.97	93/4011 (2.3%)

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	5	0

All (37) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	P	5	DT	N1-C2	13.58	1.49	1.38
2	P	5	DT	C1'-N1	11.07	1.63	1.49
2	P	5	DT	C3'-O3'	-9.62	1.31	1.44
2	P	4	DA	C3'-O3'	-8.46	1.32	1.44
3	A	249	GLU	CD-OE2	7.67	1.34	1.25
3	A	21	GLU	CD-OE1	7.46	1.33	1.25
3	A	203	GLU	CD-OE1	7.42	1.33	1.25
3	A	71	GLU	CD-OE1	7.17	1.33	1.25
3	A	147	GLU	CD-OE2	7.16	1.33	1.25
3	A	75	GLU	CD-OE1	7.16	1.33	1.25
3	A	309	GLU	CD-OE2	7.16	1.33	1.25
2	P	5	DT	C4-C5	7.11	1.51	1.45
3	A	86	GLU	CD-OE1	7.04	1.33	1.25
2	P	3	DG	C5-C6	-6.93	1.35	1.42
3	A	26	GLU	CD-OE1	6.74	1.33	1.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	154	GLU	CD-OE2	6.70	1.33	1.25
3	A	329	GLU	CD-OE2	6.55	1.32	1.25
3	A	58	GLU	CD-OE1	6.50	1.32	1.25
3	A	153	GLU	CD-OE2	6.41	1.32	1.25
2	P	6	DG	N9-C4	-6.11	1.33	1.38
3	A	129	GLU	CD-OE1	6.07	1.32	1.25
3	A	316	GLU	CD-OE2	6.06	1.32	1.25
3	A	288	GLU	CD-OE2	5.99	1.32	1.25
3	A	172	GLU	CD-OE2	5.97	1.32	1.25
2	P	6	DG	N1-C2	5.90	1.42	1.37
3	A	165	GLU	CD-OE2	5.87	1.32	1.25
3	A	326	LYS	CE-NZ	-5.83	1.34	1.49
3	A	117	GLU	CD-OE2	5.81	1.32	1.25
3	A	186	GLU	CD-OE1	5.74	1.31	1.25
2	P	4	DA	O3'-P	-5.61	1.54	1.61
3	A	335	GLU	CD-OE2	5.56	1.31	1.25
1	T	7	DG	C5-C4	-5.51	1.34	1.38
2	P	1	DC	N1-C6	-5.43	1.33	1.37
3	A	9	GLU	CD-OE2	5.28	1.31	1.25
3	A	295	GLU	CD-OE1	5.17	1.31	1.25
3	A	247	GLU	CD-OE1	5.14	1.31	1.25
3	A	36	TYR	CB-CG	-5.13	1.44	1.51

All (93) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	P	5	DT	C6-N1-C1'	-18.41	92.78	120.40
2	P	5	DT	C2-N1-C1'	16.95	145.31	118.20
2	P	1	DC	P-O3'-C3'	13.60	136.02	119.70
3	A	192	ASP	N-CA-CB	12.37	132.86	110.60
1	T	2	DC	C2-N1-C1'	10.80	130.69	118.80
2	P	4	DA	O4'-C1'-N9	-10.54	100.62	108.00
1	T	2	DC	C6-N1-C1'	-10.50	108.20	120.80
2	P	2	DA	C4'-C3'-C2'	-10.24	93.88	103.10
2	P	2	DA	O4'-C4'-C3'	-10.21	99.88	106.00
3	A	126	ARG	NE-CZ-NH1	9.21	124.90	120.30
2	P	3	DG	C8-N9-C1'	9.19	138.94	127.00
3	A	102	ARG	NE-CZ-NH1	-9.17	115.71	120.30
3	A	263	ASP	CB-CG-OD2	-9.12	110.10	118.30
3	A	183	ARG	NE-CZ-NH1	9.08	124.84	120.30
2	P	5	DT	O4'-C1'-N1	8.77	114.14	108.00
3	A	126	ARG	NE-CZ-NH2	-8.40	116.10	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	170	ASP	CB-CG-OD1	8.32	125.78	118.30
3	A	190	ASP	CB-CG-OD2	-8.31	110.82	118.30
3	A	190	ASP	CB-CG-OD1	8.25	125.72	118.30
3	A	310	PRO	N-CA-CB	8.17	113.10	103.30
2	P	5	DT	N3-C2-O2	-8.14	117.42	122.30
1	T	4	DT	C6-N1-C1'	-8.06	108.31	120.40
3	A	256	ASP	CB-CG-OD1	7.99	125.49	118.30
2	P	4	DA	P-O3'-C3'	7.89	129.17	119.70
2	P	5	DT	C2-N3-C4	-7.80	122.52	127.20
3	A	314	ASP	CB-CG-OD1	-7.71	111.36	118.30
3	A	170	ASP	CB-CG-OD2	-7.62	111.44	118.30
3	A	92	ASP	CB-CG-OD2	-7.62	111.44	118.30
1	T	5	DC	P-O3'-C3'	7.56	128.78	119.70
1	T	4	DT	C2-N1-C1'	7.47	130.15	118.20
3	A	253	ARG	NE-CZ-NH1	7.37	123.98	120.30
2	P	3	DG	C4-N9-C1'	-7.36	116.93	126.50
1	T	6	DT	C6-N1-C1'	-7.33	109.40	120.40
3	A	263	ASP	CB-CG-OD1	7.15	124.73	118.30
1	T	6	DT	C6-N1-C2	7.14	124.87	121.30
3	A	276	ASP	CB-CG-OD1	7.09	124.68	118.30
3	A	116	ASP	CB-CG-OD2	-7.09	111.92	118.30
3	A	226	ASP	CB-CG-OD1	7.09	124.68	118.30
3	A	74	ASP	CB-CG-OD1	7.07	124.66	118.30
3	A	160	ASP	CB-CG-OD2	-6.90	112.09	118.30
2	P	5	DT	C6-C5-C7	-6.81	118.81	122.90
2	P	6	DG	C4-N9-C1'	-6.80	117.67	126.50
2	P	5	DT	C5-C6-N1	-6.79	119.63	123.70
3	A	145	ASP	CB-CG-OD2	-6.61	112.35	118.30
3	A	124	ASP	CB-CG-OD2	-6.55	112.40	118.30
1	T	7	DG	P-O3'-C3'	-6.51	111.89	119.70
1	T	6	DT	C1'-O4'-C4'	-6.46	103.64	110.10
3	A	102	ARG	CD-NE-CZ	-6.18	114.94	123.60
2	P	6	DG	C8-N9-C1'	6.07	134.89	127.00
3	A	91	ASP	CB-CG-OD1	-6.06	112.84	118.30
3	A	307	ALA	N-CA-CB	6.04	118.56	110.10
3	A	199	SER	N-CA-CB	-6.02	101.47	110.50
3	A	310	PRO	N-CA-C	5.96	127.61	112.10
3	A	266	TYR	CA-CB-CG	-5.96	102.07	113.40
3	A	246	ASP	CB-CG-OD1	-5.92	112.97	118.30
2	P	6	DG	P-O5'-C5'	-5.88	111.50	120.90
1	T	6	DT	C5-C6-N1	-5.85	120.19	123.70
3	A	266	TYR	CB-CG-CD1	-5.82	117.51	121.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	17	ASP	CB-CG-OD1	-5.80	113.08	118.30
3	A	130	ASP	CB-CG-OD2	-5.79	113.09	118.30
3	A	83	ARG	N-CA-CB	5.79	121.02	110.60
3	A	321	ASP	CB-CG-OD2	-5.75	113.12	118.30
2	P	3	DG	C8-N9-C4	-5.74	104.11	106.40
3	A	332	ASP	CB-CG-OD1	-5.71	113.16	118.30
3	A	89	ARG	NE-CZ-NH1	5.71	123.15	120.30
3	A	309	GLU	C-N-CD	-5.71	108.05	120.60
3	A	152	ARG	N-CA-CB	5.59	120.66	110.60
3	A	335	GLU	N-CA-CB	5.59	120.66	110.60
3	A	130	ASP	CB-CG-OD1	5.57	123.31	118.30
3	A	276	ASP	CB-CG-OD2	-5.55	113.30	118.30
2	P	5	DT	C4-C5-C6	5.52	121.31	118.00
3	A	70	ALA	N-CA-CB	-5.50	102.40	110.10
1	T	3	DA	C6-N1-C2	-5.42	115.35	118.60
3	A	256	ASP	CB-CG-OD2	-5.40	113.44	118.30
3	A	74	ASP	CB-CG-OD2	-5.37	113.47	118.30
3	A	91	ASP	CB-CG-OD2	5.32	123.09	118.30
2	P	5	DT	P-O5'-C5'	-5.31	112.40	120.90
3	A	17	ASP	CB-CG-OD2	5.30	123.07	118.30
3	A	307	ALA	CB-CA-C	5.28	118.01	110.10
3	A	11	LEU	CB-CA-C	5.26	120.19	110.20
2	P	3	DG	P-O3'-C3'	5.21	125.95	119.70
3	A	332	ASP	CB-CG-OD2	5.20	122.98	118.30
3	A	65	VAL	N-CA-CB	5.16	122.85	111.50
3	A	43	ALA	N-CA-CB	5.14	117.30	110.10
3	A	52	LYS	N-CA-CB	5.14	119.85	110.60
3	A	253	ARG	CD-NE-CZ	5.13	130.78	123.60
3	A	111	ALA	CB-CA-C	5.12	117.78	110.10
3	A	246	ASP	CB-CG-OD2	5.11	122.90	118.30
3	A	314	ASP	CB-CG-OD2	5.05	122.84	118.30
1	T	6	DT	O4'-C4'-C3'	-5.05	102.48	104.50
3	A	196	THR	N-CA-CB	5.03	119.86	110.30
3	A	318	ASP	CB-CG-OD1	5.01	122.81	118.30
3	A	89	ARG	NE-CZ-NH2	-5.00	117.80	120.30

All (5) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	A	65	VAL	CA
3	A	192	ASP	CA
3	A	246	ASP	CA

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Mol	Chain	Res	Type	Atom
3	A	292	THR	CA
3	A	310	PRO	CA

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	T	122	0	69	9	0
2	P	126	0	68	12	0
3	A	2623	0	2641	277	1
4	A	2	0	0	0	0
5	A	113	0	0	9	0
5	P	18	0	0	2	0
5	T	14	0	0	3	0
All	All	3018	0	2778	295	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 53.

All (295) 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:23:ALA:HB2	3:A:39:TYR:HB2	1.24	1.17
2:P:1:DC:H2"	2:P:2:DA:H5"	1.26	1.10
3:A:330:PRO:HA	3:A:333:ARG:HG3	1.37	1.05
3:A:201:THR:HG23	3:A:204:SER:HB3	1.47	0.93
3:A:152:ARG:HA	3:A:155:MET:HB2	1.50	0.93
3:A:218:LEU:HB2	3:A:224:ILE:HD12	1.50	0.92
3:A:49:TYR:CD1	3:A:50:PRO:HD2	2.05	0.92
3:A:277:ILE:HG13	3:A:335:GLU:HB3	1.52	0.92
3:A:197:HIS:ND1	3:A:199:SER:HB3	1.87	0.90
3:A:270:LEU:HD22	3:A:319:ILE:HD13	1.52	0.88
3:A:119:ILE:HG23	3:A:124:ASP:HB3	1.56	0.88
3:A:23:ALA:HB2	3:A:39:TYR:CB	2.04	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:31:GLN:NE2	3:A:112:ARG:HH12	1.74	0.85
3:A:194:LEU:CD1	3:A:258:ARG:HD3	2.07	0.85
3:A:293:ILE:CD1	3:A:298:ILE:HG13	2.07	0.84
3:A:261:PRO:HG2	3:A:264:GLN:HG3	1.59	0.84
2:P:1:DC:C2'	2:P:2:DA:H5''	2.06	0.84
3:A:201:THR:CG2	3:A:204:SER:HB3	2.08	0.83
3:A:293:ILE:HD12	3:A:298:ILE:HG13	1.62	0.81
3:A:12:ASN:HB3	3:A:46:ILE:HD12	1.61	0.81
3:A:182:ARG:HG2	3:A:273:THR:HG23	1.62	0.81
3:A:277:ILE:HG13	3:A:335:GLU:CB	2.10	0.81
3:A:157:GLN:HE22	3:A:244:LYS:NZ	1.78	0.80
3:A:197:HIS:ND1	3:A:198:PRO:HD2	1.96	0.80
3:A:154:GLU:O	3:A:158:MET:HG3	1.83	0.79
3:A:119:ILE:CG2	3:A:124:ASP:HB3	2.13	0.78
3:A:194:LEU:HD11	3:A:258:ARG:HD3	1.64	0.78
3:A:197:HIS:CG	3:A:198:PRO:HD2	2.19	0.78
3:A:207:GLN:O	3:A:210:LEU:HB2	1.82	0.78
3:A:12:ASN:HB3	3:A:46:ILE:CD1	2.14	0.76
2:P:5:DT:H1'	5:P:565:HOH:O	1.85	0.76
3:A:261:PRO:HG2	3:A:264:GLN:CG	2.16	0.76
3:A:179:GLY:O	3:A:182:ARG:HB3	1.86	0.76
1:T:6:DT:H5''	5:T:547:HOH:O	1.87	0.75
3:A:270:LEU:CD2	3:A:319:ILE:HD13	2.17	0.74
3:A:18:MET:CE	3:A:76:PHE:HB2	2.17	0.74
3:A:15:ILE:HG21	3:A:46:ILE:HD13	1.70	0.74
3:A:68:LYS:O	3:A:71:GLU:HB3	1.88	0.74
3:A:244:LYS:HB2	3:A:245:ASN:HD22	1.52	0.74
3:A:289:LYS:HD2	3:A:324:GLN:OE1	1.88	0.74
3:A:327:TYR:HE1	3:A:333:ARG:HH21	1.35	0.73
3:A:23:ALA:CB	3:A:39:TYR:HB2	2.12	0.73
2:P:3:DG:H1'	5:P:511:HOH:O	1.85	0.73
3:A:132:LEU:HD23	3:A:132:LEU:N	2.03	0.73
3:A:68:LYS:HB2	3:A:68:LYS:NZ	2.03	0.72
3:A:217:GLN:O	3:A:221:VAL:HG22	1.89	0.72
2:P:5:DT:H5''	3:A:107:GLY:H	1.54	0.72
3:A:41:LYS:HD3	3:A:42:ALA:N	2.05	0.72
3:A:111:ALA:O	3:A:115:VAL:HG23	1.90	0.71
3:A:292:THR:O	3:A:298:ILE:HA	1.91	0.71
3:A:162:VAL:O	3:A:166:VAL:HG23	1.91	0.71
3:A:41:LYS:HD3	3:A:42:ALA:H	1.56	0.70
3:A:278:PHE:HB2	3:A:333:ARG:O	1.91	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:271:TYR:HB2	5:A:592:HOH:O	1.90	0.70
3:A:331:LYS:HD2	3:A:332:ASP:N	2.06	0.70
3:A:285:HIS:CE1	3:A:289:LYS:HE3	2.27	0.70
3:A:18:MET:HE1	3:A:76:PHE:HB2	1.74	0.70
3:A:49:TYR:CG	3:A:50:PRO:HD2	2.28	0.69
3:A:182:ARG:HG2	3:A:273:THR:CG2	2.23	0.69
3:A:276:ASP:O	3:A:280:LYS:HG3	1.92	0.69
3:A:201:THR:HA	3:A:261:PRO:CB	2.22	0.69
2:P:5:DT:H5''	3:A:107:GLY:N	2.08	0.69
3:A:330:PRO:CA	3:A:333:ARG:HG3	2.20	0.68
3:A:41:LYS:O	3:A:45:VAL:HG13	1.91	0.68
2:P:2:DA:C8	2:P:2:DA:H5'	2.27	0.68
3:A:201:THR:HA	3:A:261:PRO:HB3	1.75	0.68
3:A:218:LEU:CB	3:A:224:ILE:HD12	2.21	0.68
3:A:245:ASN:H	3:A:245:ASN:HD22	1.41	0.68
3:A:150:ILE:HG21	3:A:158:MET:CE	2.25	0.67
3:A:266:TYR:HA	3:A:269:VAL:HB	1.77	0.67
3:A:301:LEU:HD12	3:A:302:GLY:H	1.60	0.67
3:A:178:CYS:SG	3:A:194:LEU:HD22	2.35	0.67
3:A:59:ALA:O	3:A:62:LEU:HB2	1.95	0.66
3:A:299:ARG:HB3	3:A:300:PRO:HD2	1.77	0.66
3:A:211:LEU:HB2	3:A:259:LEU:HD22	1.77	0.66
3:A:306:VAL:HG23	3:A:307:ALA:N	2.11	0.66
3:A:245:ASN:N	3:A:245:ASN:HD22	1.93	0.66
3:A:182:ARG:NH1	3:A:273:THR:OG1	2.28	0.65
3:A:172:GLU:HG2	3:A:198:PRO:HG3	1.78	0.65
3:A:253:ARG:HG3	3:A:253:ARG:HH11	1.61	0.65
3:A:152:ARG:NH2	3:A:181:PHE:O	2.29	0.65
3:A:82:LEU:HB3	3:A:85:LEU:HB2	1.79	0.65
3:A:19:LEU:HD23	3:A:43:ALA:HB2	1.78	0.65
3:A:327:TYR:HD1	3:A:328:ARG:N	1.95	0.65
3:A:103:VAL:HG22	3:A:143:PHE:CE1	2.32	0.64
3:A:31:GLN:HE21	3:A:112:ARG:HH12	1.45	0.64
3:A:157:GLN:HE22	3:A:244:LYS:HZ1	1.45	0.64
3:A:297:THR:HB	5:A:626:HOH:O	1.97	0.64
3:A:288:GLU:HA	3:A:288:GLU:OE1	1.89	0.64
3:A:202:SER:HB2	3:A:263:ASP:OD1	1.98	0.64
3:A:81:LYS:NZ	3:A:86:GLU:OE1	2.28	0.64
3:A:194:LEU:HD12	3:A:258:ARG:HD3	1.80	0.63
3:A:235:PHE:CZ	3:A:237:GLY:HA3	2.34	0.63
3:A:319:ILE:HG22	3:A:320:PHE:N	2.12	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:15:ILE:HD11	3:A:77:LEU:HD11	1.80	0.63
1:T:3:DA:H2"	5:T:573:HOH:O	1.99	0.63
3:A:308:GLY:O	3:A:309:GLU:HB2	1.99	0.62
3:A:262:LYS:O	3:A:262:LYS:HG3	2.00	0.61
3:A:133:ASN:HD21	3:A:136:GLN:H	1.48	0.61
3:A:288:GLU:O	3:A:290:GLY:N	2.32	0.61
3:A:150:ILE:HG21	3:A:158:MET:HE1	1.83	0.61
3:A:31:GLN:N	5:A:641:HOH:O	2.24	0.61
3:A:172:GLU:HG2	3:A:198:PRO:CG	2.31	0.60
3:A:215:VAL:O	3:A:219:GLN:HG2	2.01	0.60
3:A:177:VAL:HG22	3:A:193:VAL:HG22	1.84	0.59
3:A:286:ALA:HA	3:A:323:ILE:HG21	1.85	0.59
1:T:4:DT:H2"	1:T:5:DC:O5'	2.02	0.59
3:A:279:ASN:O	3:A:283:ARG:N	2.30	0.59
3:A:68:LYS:HB2	3:A:68:LYS:HZ1	1.65	0.59
3:A:68:LYS:O	3:A:72:LYS:HE3	2.03	0.58
3:A:33:ILE:O	3:A:37:ASN:HB2	2.03	0.58
3:A:172:GLU:CG	3:A:198:PRO:HG3	2.34	0.58
3:A:319:ILE:O	3:A:322:TYR:HB2	2.04	0.58
3:A:157:GLN:HE22	3:A:244:LYS:HZ3	1.49	0.57
3:A:208:PRO:O	3:A:212:HIS:HD2	1.87	0.57
3:A:12:ASN:HD21	3:A:53:ILE:H	1.51	0.57
3:A:298:ILE:HG23	3:A:311:LEU:HB2	1.85	0.57
3:A:207:GLN:OE1	3:A:210:LEU:HG	2.03	0.57
3:A:298:ILE:HG22	5:A:578:HOH:O	2.03	0.57
3:A:212:HIS:HB3	5:A:541:HOH:O	2.04	0.57
3:A:248:LYS:HG2	3:A:248:LYS:O	2.04	0.57
3:A:44:SER:O	3:A:47:ALA:HB3	2.04	0.57
3:A:254:ARG:HH11	3:A:255:ILE:H	1.52	0.57
3:A:150:ILE:O	3:A:188:SER:N	2.30	0.56
2:P:5:DT:H2"	2:P:6:DG:O5'	2.06	0.56
1:T:6:DT:H2"	1:T:7:DG:C8	2.40	0.56
3:A:255:ILE:HG12	3:A:256:ASP:N	2.20	0.56
3:A:166:VAL:O	3:A:169:VAL:HG12	2.06	0.55
3:A:18:MET:HG2	3:A:22:LEU:HD23	1.88	0.55
3:A:18:MET:O	3:A:21:GLU:HB2	2.06	0.55
3:A:197:HIS:CE1	3:A:198:PRO:HD2	2.42	0.55
3:A:309:GLU:OE1	3:A:309:GLU:HA	2.07	0.55
3:A:99:PHE:HD2	3:A:100:LEU:HD12	1.72	0.55
3:A:43:ALA:O	3:A:47:ALA:HB2	2.08	0.55
3:A:62:LEU:HD12	3:A:63:PRO:HD2	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:49:TYR:CZ	3:A:51:HIS:HB2	2.43	0.54
3:A:176:THR:HG22	3:A:178:CYS:SG	2.48	0.54
3:A:132:LEU:HB3	3:A:136:GLN:HB3	1.89	0.54
3:A:218:LEU:HB2	3:A:224:ILE:CD1	2.32	0.53
3:A:270:LEU:HD22	3:A:319:ILE:CD1	2.32	0.53
3:A:35:LYS:O	3:A:38:ALA:HB3	2.08	0.53
3:A:41:LYS:NZ	3:A:64:GLY:O	2.39	0.53
3:A:150:ILE:HD12	3:A:155:MET:HE2	1.89	0.53
3:A:63:PRO:C	3:A:65:VAL:H	2.10	0.53
3:A:15:ILE:CG2	3:A:46:ILE:HD13	2.38	0.53
3:A:253:ARG:HH11	3:A:253:ARG:CG	2.22	0.53
3:A:327:TYR:CD1	3:A:328:ARG:N	2.76	0.53
3:A:200:PHE:O	3:A:262:LYS:N	2.38	0.53
3:A:227:THR:HG21	3:A:230:LYS:HB2	1.91	0.53
3:A:157:GLN:NE2	3:A:244:LYS:NZ	2.52	0.53
3:A:12:ASN:CB	3:A:46:ILE:HD12	2.36	0.52
3:A:266:TYR:O	3:A:270:LEU:HB2	2.08	0.52
3:A:27:LYS:HG3	3:A:28:ASN:N	2.25	0.52
3:A:244:LYS:CB	3:A:245:ASN:HD22	2.19	0.52
3:A:326:LYS:O	3:A:326:LYS:HG3	2.10	0.52
3:A:31:GLN:HE21	3:A:112:ARG:NH1	2.08	0.51
3:A:150:ILE:HG21	3:A:158:MET:HE3	1.92	0.51
3:A:303:VAL:C	3:A:305:GLY:H	2.13	0.51
3:A:157:GLN:NE2	3:A:244:LYS:HZ1	2.07	0.50
3:A:205:THR:O	3:A:206:LYS:O	2.29	0.50
3:A:150:ILE:CG2	3:A:158:MET:HE1	2.40	0.50
3:A:76:PHE:O	3:A:79:THR:O	2.29	0.50
3:A:292:THR:O	3:A:293:ILE:HD13	2.11	0.50
3:A:128:ASN:HB3	3:A:131:LYS:HD2	1.92	0.50
3:A:145:ASP:CG	3:A:252:HIS:H	2.14	0.50
3:A:292:THR:C	3:A:293:ILE:HD13	2.32	0.50
3:A:286:ALA:O	3:A:291:PHE:HB2	2.12	0.50
3:A:293:ILE:HD13	3:A:298:ILE:HG13	1.92	0.50
3:A:150:ILE:HD12	3:A:155:MET:CE	2.42	0.50
3:A:288:GLU:C	3:A:290:GLY:H	2.14	0.49
3:A:113:LYS:O	3:A:117:GLU:HG3	2.11	0.49
3:A:170:ASP:HB3	3:A:173:TYR:CD2	2.47	0.49
3:A:260:ILE:HG22	3:A:261:PRO:HD2	1.94	0.49
3:A:206:LYS:O	3:A:207:GLN:HB2	2.10	0.49
3:A:28:ASN:HD22	3:A:28:ASN:N	2.10	0.49
1:T:2:DC:N4	5:T:655:HOH:O	2.46	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:18:MET:HG2	3:A:22:LEU:CD2	2.42	0.49
3:A:245:ASN:N	3:A:245:ASN:ND2	2.61	0.49
3:A:241:LEU:HB2	3:A:250:TYR:CD2	2.47	0.49
3:A:267:CYS:SG	3:A:297:THR:HA	2.53	0.49
3:A:310:PRO:O	3:A:312:PRO:HD3	2.12	0.49
3:A:300:PRO:HD3	3:A:311:LEU:HD13	1.95	0.49
3:A:312:PRO:O	3:A:322:TYR:OH	2.28	0.49
3:A:282:MET:HE3	3:A:323:ILE:HD11	1.94	0.48
3:A:18:MET:HE3	3:A:76:PHE:CG	2.48	0.48
3:A:202:SER:HB2	3:A:263:ASP:CG	2.34	0.48
3:A:36:TYR:CD2	3:A:37:ASN:N	2.81	0.48
3:A:282:MET:CE	3:A:323:ILE:HD11	2.44	0.48
3:A:254:ARG:HH11	3:A:255:ILE:N	2.12	0.47
3:A:121:THR:O	3:A:124:ASP:HB2	2.13	0.47
3:A:298:ILE:HD13	3:A:322:TYR:CD2	2.49	0.47
3:A:125:LEU:HD23	3:A:125:LEU:HA	1.74	0.47
3:A:217:GLN:HA	3:A:217:GLN:NE2	2.29	0.47
3:A:331:LYS:CD	3:A:332:ASP:N	2.76	0.47
3:A:12:ASN:HB3	3:A:46:ILE:HD11	1.94	0.47
3:A:272:PHE:HD1	5:A:587:HOH:O	1.98	0.47
3:A:133:ASN:O	3:A:137:ARG:HG3	2.15	0.47
3:A:182:ARG:NH2	3:A:316:GLU:OE1	2.47	0.47
3:A:243:SER:HB3	3:A:249:GLU:HA	1.97	0.47
3:A:81:LYS:HZ3	3:A:86:GLU:HB3	1.79	0.47
3:A:301:LEU:HD12	3:A:302:GLY:N	2.26	0.47
3:A:142:TYR:O	3:A:146:PHE:HB2	2.15	0.46
3:A:254:ARG:NH1	3:A:255:ILE:H	2.14	0.46
3:A:58:GLU:O	3:A:61:LYS:HB2	2.15	0.46
3:A:253:ARG:NH1	5:A:503:HOH:O	2.48	0.46
1:T:5:DC:H2"	1:T:6:DT:O5'	2.15	0.46
3:A:217:GLN:HE22	3:A:220:LYS:HD3	1.81	0.46
3:A:49:TYR:HA	3:A:50:PRO:HD3	1.38	0.46
3:A:243:SER:CB	3:A:249:GLU:HA	2.46	0.46
3:A:277:ILE:CG1	3:A:335:GLU:HA	2.46	0.46
3:A:106:ILE:HG22	3:A:107:GLY:O	2.17	0.45
3:A:298:ILE:CD1	3:A:322:TYR:HD2	2.29	0.45
3:A:16:THR:HG23	3:A:46:ILE:CG1	2.45	0.45
3:A:298:ILE:HD13	3:A:322:TYR:HD2	1.81	0.45
3:A:142:TYR:CE2	3:A:238:VAL:HG11	2.52	0.45
1:T:2:DC:H2"	1:T:3:DA:C8	2.51	0.45
3:A:285:HIS:NE2	3:A:289:LYS:CG	2.80	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:283:ARG:O	3:A:286:ALA:HB3	2.16	0.45
3:A:138:ILE:HG22	3:A:142:TYR:CD2	2.50	0.45
3:A:219:GLN:HB3	3:A:219:GLN:HE21	1.49	0.45
3:A:244:LYS:CB	3:A:245:ASN:ND2	2.80	0.45
3:A:316:GLU:O	3:A:320:PHE:HD2	1.99	0.45
3:A:331:LYS:HD3	3:A:332:ASP:HB3	1.98	0.45
2:P:4:DA:C8	2:P:5:DT:H72	2.52	0.45
3:A:157:GLN:O	3:A:160:ASP:HB3	2.17	0.45
3:A:274:GLY:HA3	3:A:275:SER:HA	1.76	0.45
3:A:26:GLU:HA	3:A:30:SER:OG	2.17	0.45
3:A:254:ARG:NH1	3:A:255:ILE:N	2.65	0.45
3:A:241:LEU:HA	3:A:242:PRO:HD3	1.68	0.45
3:A:325:TRP:HD1	5:A:583:HOH:O	1.99	0.44
3:A:100:LEU:HD11	3:A:120:LYS:O	2.17	0.44
3:A:133:ASN:ND2	3:A:136:GLN:H	2.13	0.44
3:A:197:HIS:CE1	3:A:199:SER:HB3	2.50	0.44
3:A:200:PHE:HB2	5:A:625:HOH:O	2.15	0.44
3:A:299:ARG:HB3	3:A:300:PRO:CD	2.45	0.44
3:A:49:TYR:CE2	3:A:51:HIS:HB2	2.52	0.44
3:A:285:HIS:NE2	3:A:289:LYS:HG3	2.32	0.44
3:A:56:GLY:O	3:A:59:ALA:N	2.51	0.44
1:T:6:DT:C2'	1:T:7:DG:C8	3.01	0.44
3:A:73:ILE:HG22	3:A:77:LEU:HD22	1.99	0.44
3:A:289:LYS:HD3	3:A:289:LYS:HA	1.66	0.44
3:A:18:MET:HE3	3:A:76:PHE:HB2	1.96	0.44
3:A:103:VAL:HG22	3:A:143:PHE:HE1	1.80	0.44
3:A:132:LEU:N	3:A:132:LEU:CD2	2.76	0.44
3:A:155:MET:HA	3:A:158:MET:HE2	2.00	0.44
3:A:23:ALA:O	3:A:36:TYR:HD1	2.00	0.44
3:A:150:ILE:O	3:A:187:SER:HA	2.18	0.43
3:A:311:LEU:HA	3:A:311:LEU:HD12	1.64	0.43
3:A:330:PRO:HA	3:A:333:ARG:CG	2.27	0.43
3:A:303:VAL:O	3:A:305:GLY:N	2.51	0.43
3:A:108:PRO:O	3:A:112:ARG:HG2	2.18	0.43
3:A:194:LEU:HA	3:A:194:LEU:HD12	1.66	0.43
3:A:265:TYR:O	3:A:268:GLY:N	2.51	0.43
3:A:216:GLU:HA	3:A:219:GLN:HG3	2.01	0.43
3:A:228:LEU:HB2	3:A:236:MET:O	2.19	0.43
3:A:255:ILE:HG23	3:A:255:ILE:O	2.18	0.43
3:A:296:TYR:O	3:A:297:THR:HG22	2.19	0.43
3:A:132:LEU:HB2	3:A:137:ARG:HG3	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:59:ALA:O	3:A:62:LEU:N	2.28	0.43
2:P:2:DA:C8	2:P:2:DA:C5'	3.00	0.42
3:A:182:ARG:HG2	3:A:182:ARG:HH11	1.83	0.42
3:A:197:HIS:CG	3:A:198:PRO:CD	2.99	0.42
3:A:244:LYS:HB2	3:A:245:ASN:ND2	2.28	0.42
3:A:291:PHE:CD1	3:A:300:PRO:HA	2.54	0.42
3:A:144:GLY:O	3:A:147:GLU:N	2.51	0.42
3:A:241:LEU:O	3:A:250:TYR:HD2	2.02	0.42
3:A:88:ILE:HD13	3:A:88:ILE:HG21	1.84	0.42
3:A:236:MET:HG2	3:A:254:ARG:HH22	1.84	0.42
3:A:195:LEU:HA	3:A:195:LEU:HD12	1.75	0.42
1:T:4:DT:H2''	1:T:5:DC:C5'	2.49	0.42
3:A:18:MET:HE3	3:A:76:PHE:CB	2.50	0.41
3:A:277:ILE:HG13	3:A:335:GLU:HA	2.02	0.41
3:A:81:LYS:NZ	3:A:86:GLU:HB3	2.35	0.41
3:A:15:ILE:CG2	3:A:46:ILE:CD1	2.98	0.41
3:A:172:GLU:O	3:A:198:PRO:HD3	2.21	0.41
3:A:201:THR:HG22	3:A:204:SER:HB3	1.97	0.41
2:P:5:DT:P	3:A:107:GLY:HA3	2.60	0.41
3:A:119:ILE:HG22	3:A:124:ASP:HB3	1.99	0.41
3:A:120:LYS:N	3:A:124:ASP:OD2	2.28	0.41
3:A:280:LYS:O	3:A:283:ARG:N	2.54	0.41
3:A:177:VAL:O	3:A:182:ARG:HB2	2.20	0.41
3:A:191:MET:HG2	3:A:255:ILE:HG13	2.02	0.41
3:A:294:ASN:HD22	3:A:294:ASN:N	2.18	0.41
3:A:286:ALA:HA	3:A:323:ILE:CG2	2.49	0.41
3:A:268:GLY:O	3:A:271:TYR:HB3	2.20	0.41
3:A:128:ASN:C	3:A:130:ASP:N	2.74	0.40
3:A:128:ASN:O	3:A:131:LYS:HG3	2.20	0.40
3:A:228:LEU:HD23	3:A:228:LEU:HA	1.57	0.40
2:P:2:DA:H5'	2:P:2:DA:H2'	1.27	0.40
3:A:183:ARG:HD3	3:A:273:THR:O	2.22	0.40
3:A:208:PRO:O	3:A:211:LEU:N	2.51	0.40
3:A:15:ILE:O	3:A:19:LEU:HD22	2.21	0.40
3:A:16:THR:CG2	3:A:46:ILE:CG1	3.00	0.40
3:A:41:LYS:CD	3:A:42:ALA:N	2.80	0.40
3:A:102:ARG:HH11	3:A:102:ARG:HD3	1.48	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)
3:A:83:ARG:NH1	3:A:117:GLU:CB[3_558]	2.10	0.10

## 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%)	276 (85%)	30 (9%)	19 (6%)	<b>1</b> <b>4</b>

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	32	ALA
3	A	206	LYS
3	A	207	GLN
3	A	244	LYS
3	A	246	ASP
3	A	289	LYS
3	A	309	GLU
3	A	334	SER
3	A	247	GLU
3	A	304	THR
3	A	10	THR
3	A	91	ASP
3	A	185	ALA
3	A	202	SER
3	A	265	TYR
3	A	310	PRO
3	A	318	ASP
3	A	324	GLN
3	A	80	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	288 / 295 (98%)	221 (77%)	67 (23%)	<b>1</b> <b>2</b>

All (67) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	11	LEU
3	A	18	MET
3	A	19	LEU
3	A	22	LEU
3	A	25	PHE
3	A	27	LYS
3	A	30	SER
3	A	33	ILE
3	A	37	ASN
3	A	40	ARG
3	A	41	LYS
3	A	45	VAL
3	A	46	ILE
3	A	54	LYS
3	A	62	LEU
3	A	65	VAL
3	A	67	THR
3	A	68	LYS
3	A	79	THR
3	A	85	LEU
3	A	95	SER
3	A	112	ARG
3	A	113	LYS
3	A	119	ILE
3	A	122	LEU
3	A	130	ASP
3	A	131	LYS
3	A	132	LEU
3	A	152	ARG
3	A	153	GLU

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Mol	Chain	Res	Type
3	A	168	LYS
3	A	171	SER
3	A	172	GLU
3	A	182	ARG
3	A	188	SER
3	A	190	ASP
3	A	194	LEU
3	A	201	THR
3	A	202	SER
3	A	204	SER
3	A	207	GLN
3	A	214	VAL
3	A	218	LEU
3	A	219	GLN
3	A	233	THR
3	A	235	PHE
3	A	236	MET
3	A	245	ASN
3	A	248	LYS
3	A	253	ARG
3	A	255	ILE
3	A	258	ARG
3	A	276	ASP
3	A	277	ILE
3	A	287	LEU
3	A	288	GLU
3	A	293	ILE
3	A	294	ASN
3	A	297	THR
3	A	301	LEU
3	A	309	GLU
3	A	311	LEU
3	A	314	ASP
3	A	325	TRP
3	A	327	TYR
3	A	331	LYS
3	A	335	GLU

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

*Continued on next page...*

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Mol	Chain	Res	Type
3	A	28	ASN
3	A	31	GLN
3	A	37	ASN
3	A	90	GLN
3	A	134	HIS
3	A	136	GLN
3	A	157	GLN
3	A	159	GLN
3	A	212	HIS
3	A	213	GLN
3	A	217	GLN
3	A	219	GLN
3	A	245	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 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	T	7/7 (100%)	-0.56	0	100 100	19, 30, 61, 96	0
2	P	6/6 (100%)	-0.92	0	100 100	13, 23, 29, 38	0
3	A	324/335 (96%)	-0.63	1 (0%)	94 93	4, 34, 82, 100	0
All	All	337/348 (96%)	-0.64	1 (0%)	94 93	4, 34, 82, 100	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	A	10	THR	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. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
4	NA	A	342	1/1	0.92	0.19	20,20,20,20	0
4	NA	A	341	1/1	0.92	0.20	23,23,23,23	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.