



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 02:53 AM UTC

PDB ID : 1BPY / pdb_00001bpy
Title : HUMAN DNA POLYMERASE BETA COMPLEXED WITH GAPPED DNA AND DDCTP
Authors : Sawaya, M.R.; Pelletier, H.; Prasad, R.; Wilson, S.H.; Kraut, J.
Deposited on : 1997-04-15
Resolution : 2.20 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
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

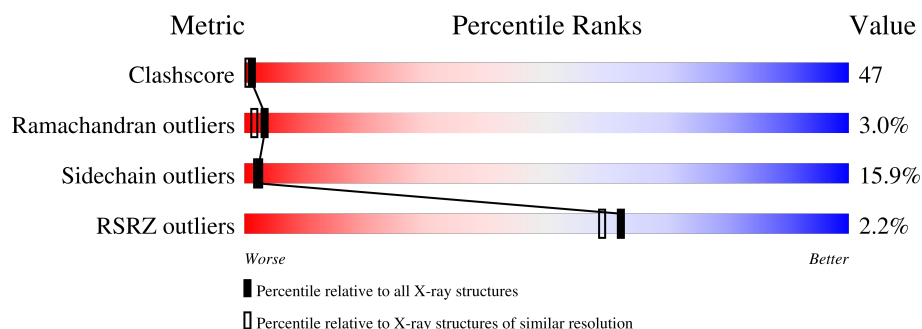
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	T	16	 25% 38% 38%
2	P	10	 10% 80% 10%
3	D	5	 20% 60% 20%
4	A	335	 2% 33% 47% 16% . .

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 3597 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*CP*GP*AP*CP*GP*GP*CP*GP*CP*AP*TP*CP*AP*GP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	T	16	Total	C	N	O	P	0	0	0
			323	153	63	92	15			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	P	10	Total	C	N	O	P	0	0	0
			202	97	38	58	9			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	5	Total	C	N	O	P	0	0	0
			106	49	20	32	5			

- Molecule 4 is a protein called PROTEIN (DNA POLYMERASE BETA).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	A	326	Total	C	N	O	S	0	0	0
			2614	1652	457	496	9			

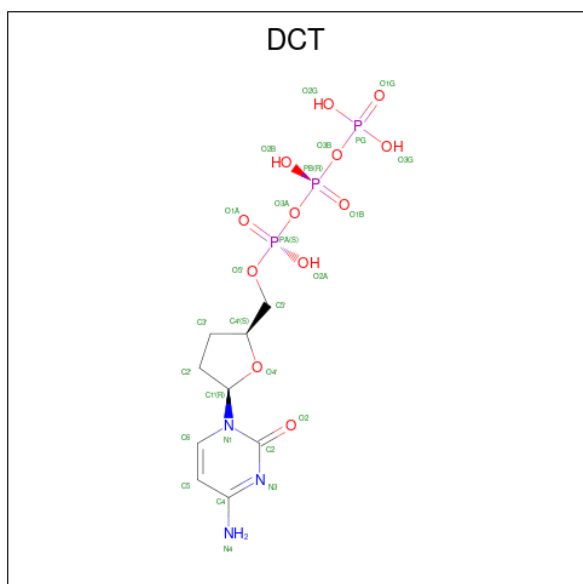
- Molecule 5 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

- Molecule 6 is SODIUM ION (CCD ID: NA) (formula: Na).

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

- Molecule 7 is 2',3'-DIDEOXYCYTIDINE 5'-TRIPHOSPHATE (CCD ID: DCT) (formula: $C_9H_{16}N_3O_{12}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	A	1	Total	C	N	O	P	0	0
			27	9	3	12	3		

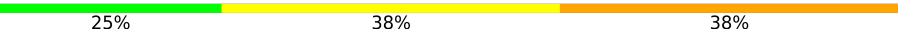
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	T	50	Total	O	0	0
			50	50		
8	P	21	Total	O	0	0
			21	21		
8	D	27	Total	O	0	0
			27	27		
8	A	223	Total	O	0	0
			223	223		

3 Residue-property plots [i](#)

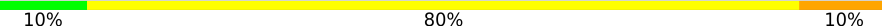
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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.

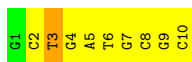
- Molecule 1: DNA (5'-D(*CP*CP*GP*AP*CP*GP*GP*CP*GP*CP*AP*TP*CP*AP*GP*C)-3')

Chain T: 

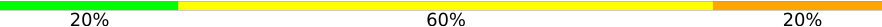


- Molecule 2: DNA (5'-D(*GP*CP*TP*GP*AP*TP*GP*CP*GP*DOC)-3')

Chain P: 

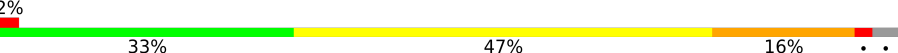


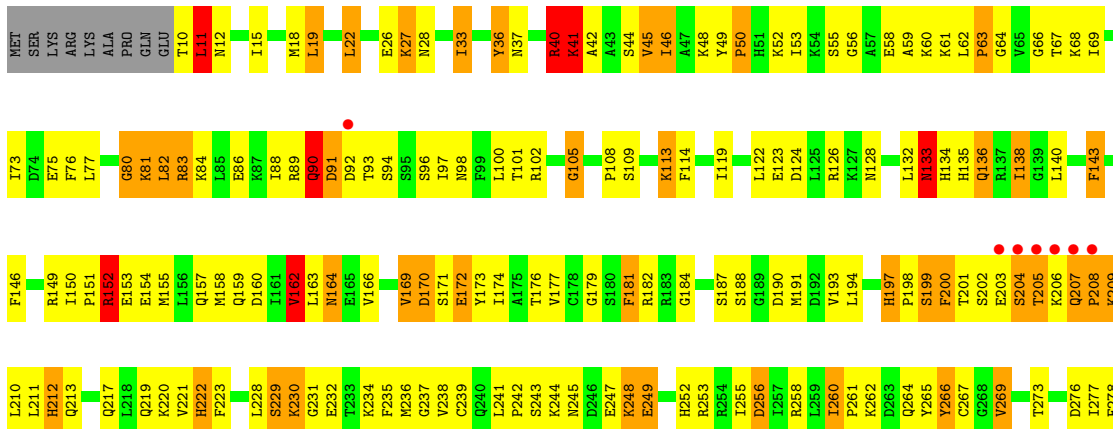
- Molecule 3: DNA (5'-D(*GP*TP*CP*GP*G)-3')

Chain D: 



- Molecule 4: PROTEIN (DNA POLYMERASE BETA)

Chain A: 





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	50.54Å 79.82Å 54.94Å 90.00° 107.11° 90.00°	Depositor
Resolution (Å)	20.00 – 2.20 20.00 – 2.20	Depositor EDS
% Data completeness (in resolution range)	95.4 (20.00-2.20) 91.6 (20.00-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.12 (at 2.19Å)	Xtriage
Refinement program	TNT 5D	Depositor
R, R_{free}	0.232 , 0.330 (Not available) , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	20.9	Xtriage
Anisotropy	0.849	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 99.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3597	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.05% 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

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

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	1.01	0/362	1.71	12/556 (2.2%)
2	P	1.00	0/206	1.60	1/317 (0.3%)
3	D	1.00	0/118	1.79	3/179 (1.7%)
4	A	1.50	7/2663 (0.3%)	2.13	95/3578 (2.7%)
All	All	1.42	7/3349 (0.2%)	2.04	111/4630 (2.4%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	230	LYS	N-CA	-6.33	1.38	1.46
4	A	262	LYS	CA-C	-6.04	1.44	1.52
4	A	27	LYS	CA-C	5.96	1.60	1.52
4	A	26	GLU	C-N	-5.41	1.26	1.33
4	A	212	HIS	ND1-CE1	5.21	1.37	1.32
4	A	174	ILE	CA-C	-5.08	1.46	1.52
4	A	320	PHE	CA-C	5.07	1.59	1.52

All (111) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	204	SER	N-CA-C	18.97	137.75	112.90
4	A	205	THR	N-CA-C	-13.39	93.97	111.92
4	A	248	LYS	N-CA-CB	9.75	125.77	109.87
4	A	297	THR	CB-CA-C	-9.59	92.37	111.17
4	A	207	GLN	N-CA-C	9.13	124.82	108.47
4	A	300	PRO	CB-CA-C	-8.84	99.45	111.21
3	D	2	DT	C5'-C4'-C3'	-8.63	101.95	114.90
4	A	41	LYS	N-CA-CB	8.48	122.58	110.12
1	T	14	DA	C4'-C3'-O3'	-8.38	97.44	110.00
4	A	333	ARG	N-CA-C	8.33	123.49	113.16

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	205	THR	N-CA-CB	8.15	125.26	112.46
4	A	318	ASP	CA-CB-CG	-8.09	104.51	112.60
4	A	212	HIS	CB-CG-CD2	-7.99	120.81	131.20
4	A	152	ARG	N-CA-CB	7.73	121.46	109.94
4	A	293	ILE	CB-CA-C	-7.61	100.25	110.98
4	A	80	GLY	N-CA-C	-7.53	104.92	115.32
4	A	222	HIS	CB-CG-CD2	-7.46	121.50	131.20
4	A	229	SER	N-CA-CB	7.21	122.53	111.56
4	A	164	ASN	CA-CB-CG	-7.16	105.44	112.60
4	A	205	THR	CA-C-N	7.08	131.50	121.02
4	A	205	THR	C-N-CA	7.08	131.50	121.02
4	A	105	GLY	N-CA-C	-7.06	105.47	114.37
4	A	252	HIS	CB-CG-CD2	-6.95	122.17	131.20
4	A	262	LYS	O-C-N	6.90	130.04	122.11
4	A	298	ILE	N-CA-CB	6.82	123.83	111.21
4	A	37	ASN	CA-CB-CG	-6.65	105.95	112.60
1	T	1	DC	P-O3'-C3'	-6.64	110.24	120.20
4	A	75	GLU	N-CA-C	-6.63	104.05	111.28
4	A	40	ARG	N-CA-CB	6.63	119.97	110.16
4	A	162	VAL	CB-CA-C	-6.62	103.49	111.97
4	A	204	SER	N-CA-CB	-6.60	100.43	110.33
4	A	200	PHE	CA-CB-CG	6.60	120.40	113.80
4	A	133	ASN	N-CA-CB	6.59	120.53	110.70
1	T	3	DG	N9-C1'-C2'	6.55	123.33	113.50
4	A	320	PHE	CA-CB-CG	6.45	120.25	113.80
4	A	208	PRO	N-CA-C	-6.44	106.11	114.03
4	A	269	VAL	CB-CA-C	-6.41	102.22	112.16
4	A	63	PRO	CB-CA-C	-6.39	105.59	111.40
4	A	323	ILE	N-CA-C	6.38	119.14	112.83
4	A	45	VAL	N-CA-CB	6.29	118.62	110.57
4	A	133	ASN	N-CA-C	-6.25	100.95	110.14
4	A	143	PHE	N-CA-C	6.24	124.09	110.80
4	A	297	THR	N-CA-C	6.22	117.68	108.60
4	A	146	PHE	CA-CB-CG	-6.22	107.58	113.80
4	A	199	SER	N-CA-CB	-6.17	100.07	110.49
1	T	2	DC	C4'-C3'-O3'	-6.14	100.80	110.00
4	A	81	LYS	N-CA-C	6.11	117.48	108.86
4	A	128	ASN	CA-CB-CG	6.10	118.70	112.60
4	A	36	TYR	N-CA-C	-6.06	104.59	111.07
4	A	302	GLY	CA-C-N	6.05	132.86	121.97
4	A	302	GLY	C-N-CA	6.05	132.86	121.97
1	T	2	DC	P-O5'-C5'	-6.04	110.94	120.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	92	ASP	N-CA-C	5.93	117.82	111.36
1	T	10	DC	N1-C1'-C2'	5.81	122.21	113.50
4	A	172	GLU	N-CA-C	5.77	118.52	111.82
4	A	280	LYS	CA-C-N	5.73	128.43	120.29
4	A	280	LYS	C-N-CA	5.73	128.43	120.29
4	A	213	GLN	CA-C-N	-5.69	112.00	120.82
4	A	213	GLN	C-N-CA	-5.69	112.00	120.82
4	A	140	LEU	N-CA-C	-5.67	105.01	111.07
4	A	169	VAL	N-CA-C	5.64	116.41	110.72
4	A	92	ASP	N-CA-CB	5.63	118.50	110.16
4	A	44	SER	N-CA-CB	5.55	118.37	110.16
4	A	321	ASP	CA-CB-CG	-5.54	107.06	112.60
4	A	152	ARG	N-CA-C	5.54	117.75	111.11
4	A	55	SER	CA-C-N	5.51	126.06	119.94
4	A	55	SER	C-N-CA	5.51	126.06	119.94
4	A	123	GLU	N-CA-CB	5.47	117.91	109.98
4	A	248	LYS	CA-CB-CG	-5.46	103.18	114.10
4	A	266	TYR	CA-C-N	-5.46	112.54	120.29
4	A	266	TYR	C-N-CA	-5.46	112.54	120.29
1	T	2	DC	O3'-P-O5'	-5.44	95.83	104.00
1	T	5	DC	C2'-C3'-O3'	-5.44	103.34	111.50
4	A	256	ASP	CA-CB-CG	5.39	117.99	112.60
1	T	2	DC	N1-C1'-C2'	5.38	121.56	113.50
1	T	12	DT	C4'-C3'-O3'	-5.36	101.97	110.00
4	A	301	LEU	CB-CA-C	-5.35	103.35	110.79
4	A	91	ASP	CA-CB-CG	5.33	117.93	112.60
4	A	27	LYS	N-CA-CB	-5.33	102.34	110.07
4	A	197	HIS	CA-C-N	5.33	126.50	119.84
4	A	197	HIS	C-N-CA	5.33	126.50	119.84
4	A	45	VAL	CB-CA-C	-5.31	105.06	112.02
4	A	213	GLN	CB-CA-C	-5.31	98.29	110.18
4	A	204	SER	CA-C-O	-5.30	113.40	119.60
4	A	61	LYS	CA-CB-CG	-5.27	103.56	114.10
4	A	58	GLU	CA-C-N	-5.25	111.76	120.68
4	A	58	GLU	C-N-CA	-5.25	111.76	120.68
4	A	260	ILE	N-CA-CB	-5.23	103.89	111.21
4	A	316	GLU	CA-C-N	5.22	127.58	120.54
4	A	316	GLU	C-N-CA	5.22	127.58	120.54
4	A	90	GLN	CB-CG-CD	5.22	121.47	112.60
4	A	297	THR	CA-C-N	-5.20	116.39	123.10
4	A	297	THR	C-N-CA	-5.20	116.39	123.10
1	T	3	DG	O4'-C1'-N9	-5.17	100.64	108.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	4	DG	P-O3'-C3'	5.16	127.94	120.20
4	A	258	ARG	CA-C-N	5.16	129.27	122.30
4	A	258	ARG	C-N-CA	5.16	129.27	122.30
2	P	3	DT	N1-C1'-C2'	5.16	121.24	113.50
1	T	6	DG	C4'-C3'-O3'	-5.15	102.28	110.00
4	A	108	PRO	N-CA-C	5.14	123.05	112.47
4	A	193	VAL	N-CA-C	5.12	115.22	107.75
4	A	261	PRO	N-CA-CB	5.11	107.79	103.19
4	A	222	HIS	CB-CG-ND1	5.10	130.35	122.70
4	A	209	LYS	N-CA-C	-5.10	105.64	111.14
4	A	181	PHE	O-C-N	5.08	127.31	122.03
3	D	2	DT	N1-C1'-C2'	5.05	121.08	113.50
4	A	287	LEU	CA-C-N	5.04	127.04	120.28
4	A	287	LEU	C-N-CA	5.04	127.04	120.28
4	A	50	PRO	N-CA-C	5.03	119.91	114.68
4	A	267	CYS	N-CA-C	-5.03	105.88	111.36
4	A	170	ASP	N-CA-CB	5.01	117.68	110.37

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	T	323	0	179	13	0
2	P	202	0	114	6	0
3	D	106	0	57	9	0
4	A	2614	0	2635	277	0
5	A	2	0	0	0	0
6	A	2	0	0	0	0
7	A	27	0	12	0	0
8	A	223	0	0	18	0
8	D	27	0	0	2	0
8	P	21	0	0	2	0
8	T	50	0	0	3	0
All	All	3597	0	2997	294	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 47.

All (294) 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)
4:A:285:HIS:HD2	4:A:323:ILE:HD12	1.17	1.07
4:A:230:LYS:HG3	4:A:235:PHE:HD1	1.13	1.06
4:A:316:GLU:H	4:A:317:LYS:NZ	1.53	1.06
4:A:155:MET:HE3	4:A:188:SER:HB2	1.09	1.05
4:A:207:GLN:HB3	4:A:209:LYS:HG2	1.41	1.03
4:A:83:ARG:HG2	4:A:83:ARG:HH11	1.24	1.02
4:A:230:LYS:HG3	4:A:235:PHE:CD1	2.03	0.92
4:A:228:LEU:HD22	4:A:237:GLY:HA2	1.51	0.92
4:A:285:HIS:CD2	4:A:323:ILE:HD12	2.04	0.90
4:A:33:ILE:HD12	4:A:33:ILE:H	1.37	0.90
4:A:18:MET:HE2	4:A:82:LEU:HB2	1.54	0.90
4:A:12:ASN:HD21	4:A:53:ILE:H	1.09	0.89
4:A:316:GLU:H	4:A:317:LYS:HZ3	1.20	0.87
4:A:113:LYS:HG2	4:A:114:PHE:N	1.91	0.86
4:A:315:SER:HB2	4:A:317:LYS:HE2	1.58	0.86
4:A:150:ILE:CG2	4:A:158:MET:HE1	2.06	0.85
1:T:2:DC:H2''	1:T:3:DG:C8	2.13	0.84
4:A:11:LEU:HD11	4:A:52:LYS:HB2	1.60	0.83
4:A:133:ASN:ND2	4:A:136:GLN:HG3	1.92	0.83
4:A:150:ILE:HD12	4:A:155:MET:HE2	1.62	0.82
4:A:194:LEU:HD11	4:A:260:ILE:HG13	1.61	0.81
3:D:2:DT:H5'	4:A:66:GLY:HA3	1.62	0.81
4:A:207:GLN:HB3	4:A:209:LYS:CG	2.10	0.81
2:P:5:DA:H2'	2:P:6:DT:H71	1.63	0.80
4:A:243:SER:OG	4:A:249:GLU:HA	1.82	0.80
4:A:197:HIS:ND1	4:A:199:SER:HB3	1.96	0.79
4:A:248:LYS:HG2	4:A:249:GLU:N	1.95	0.79
4:A:18:MET:CE	4:A:82:LEU:HB2	2.13	0.79
4:A:150:ILE:HD12	4:A:155:MET:CE	2.13	0.79
4:A:328:ARG:HH11	4:A:328:ARG:HB3	1.46	0.78
4:A:191:MET:HG2	4:A:255:ILE:HG13	1.65	0.78
4:A:18:MET:HE1	4:A:76:PHE:HB2	1.67	0.77
4:A:234:LYS:HE3	8:A:755:HOH:O	1.84	0.77
4:A:316:GLU:N	4:A:317:LYS:NZ	2.32	0.77
4:A:200:PHE:CD1	4:A:206:LYS:HB2	2.20	0.77
4:A:236:MET:HG3	4:A:256:ASP:OD1	1.85	0.77
4:A:150:ILE:HG21	4:A:158:MET:HE1	1.67	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:2:DT:C5'	4:A:66:GLY:HA3	2.14	0.76
4:A:69:ILE:O	4:A:73:ILE:HG13	1.85	0.75
4:A:41:LYS:HE2	4:A:64:GLY:CA	2.17	0.75
4:A:265:TYR:O	4:A:269:VAL:HG22	1.85	0.75
4:A:41:LYS:HE2	4:A:64:GLY:HA2	1.69	0.74
4:A:41:LYS:HD3	4:A:42:ALA:N	2.02	0.74
4:A:194:LEU:HD11	4:A:260:ILE:CG1	2.17	0.74
4:A:155:MET:CE	4:A:188:SER:HB2	2.04	0.74
4:A:316:GLU:H	4:A:317:LYS:HZ1	1.32	0.74
4:A:315:SER:CB	4:A:317:LYS:HE2	2.19	0.73
4:A:287:LEU:HD12	4:A:291:PHE:O	1.88	0.72
4:A:122:LEU:O	4:A:126:ARG:HG3	1.90	0.72
4:A:170:ASP:HB3	4:A:173:TYR:CD2	2.24	0.72
4:A:12:ASN:ND2	4:A:53:ILE:H	1.87	0.72
4:A:315:SER:HB2	4:A:317:LYS:CE	2.19	0.71
4:A:300:PRO:HD3	4:A:311:LEU:CD1	2.21	0.70
4:A:300:PRO:HD3	4:A:311:LEU:HD11	1.72	0.70
4:A:154:GLU:O	4:A:158:MET:HG3	1.90	0.69
4:A:12:ASN:HD21	4:A:53:ILE:N	1.86	0.69
4:A:33:ILE:H	4:A:33:ILE:CD1	1.94	0.69
4:A:207:GLN:CB	4:A:209:LYS:HG2	2.22	0.69
4:A:18:MET:HE2	4:A:82:LEU:CB	2.22	0.68
4:A:200:PHE:HA	4:A:204:SER:OG	1.94	0.68
4:A:228:LEU:HD22	4:A:237:GLY:CA	2.24	0.68
4:A:289:LYS:HZ3	4:A:324:GLN:HG3	1.59	0.67
4:A:138:ILE:N	4:A:138:ILE:HD13	2.08	0.67
4:A:209:LYS:HG3	4:A:210:LEU:N	2.10	0.67
4:A:301:LEU:HD23	8:A:520:HOH:O	1.95	0.66
4:A:328:ARG:HH11	4:A:328:ARG:CB	2.07	0.66
4:A:42:ALA:O	4:A:46:ILE:HG23	1.96	0.66
4:A:158:MET:O	4:A:162:VAL:HG22	1.96	0.66
4:A:113:LYS:HE3	4:A:114:PHE:CD1	2.31	0.66
4:A:162:VAL:HG23	4:A:163:LEU:H	1.61	0.66
4:A:328:ARG:HD2	8:A:578:HOH:O	1.95	0.66
4:A:160:ASP:OD1	4:A:164:ASN:ND2	2.30	0.65
3:D:3:DC:H5'	8:D:649:HOH:O	1.96	0.65
4:A:150:ILE:HG21	4:A:155:MET:HE2	1.79	0.65
4:A:133:ASN:H	4:A:136:GLN:HE21	1.43	0.65
4:A:303:VAL:HG23	4:A:304:THR:N	2.12	0.65
4:A:279:ASN:O	4:A:283:ARG:HG3	1.97	0.64
4:A:150:ILE:CB	4:A:155:MET:HE2	2.27	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:207:GLN:OE1	4:A:208:PRO:HD2	1.98	0.64
1:T:8:DC:H4'	4:A:295:GLU:OE2	1.98	0.64
4:A:11:LEU:N	8:A:712:HOH:O	2.30	0.64
4:A:289:LYS:NZ	4:A:324:GLN:HG3	2.12	0.64
4:A:328:ARG:O	4:A:333:ARG:NH1	2.29	0.64
4:A:133:ASN:HD21	4:A:135:HIS:HB3	1.62	0.64
3:D:1:DG:H5''	8:D:590:HOH:O	1.98	0.64
1:T:13:DC:H2''	1:T:14:DA:C8	2.34	0.63
4:A:113:LYS:HA	8:A:661:HOH:O	1.98	0.63
4:A:285:HIS:O	4:A:288:GLU:HB2	1.97	0.63
4:A:328:ARG:HH11	4:A:328:ARG:CG	2.10	0.63
4:A:41:LYS:HD3	4:A:42:ALA:H	1.63	0.63
4:A:278:PHE:CZ	4:A:333:ARG:HD2	2.33	0.63
4:A:40:ARG:NH1	4:A:276:ASP:OD2	2.31	0.63
4:A:109:SER:HB3	8:A:533:HOH:O	1.99	0.63
4:A:150:ILE:CG2	4:A:155:MET:HE2	2.29	0.63
4:A:36:TYR:CZ	4:A:40:ARG:HD2	2.34	0.62
4:A:113:LYS:HE3	4:A:114:PHE:CE1	2.34	0.62
4:A:212:HIS:N	8:A:752:HOH:O	2.30	0.62
4:A:303:VAL:HG23	4:A:304:THR:H	1.63	0.62
4:A:162:VAL:O	4:A:166:VAL:HG23	2.00	0.62
1:T:1:DC:H2''	1:T:2:DC:O5'	2.00	0.62
4:A:244:LYS:H	4:A:247:GLU:HG2	1.64	0.62
3:D:2:DT:H5'	4:A:66:GLY:CA	2.29	0.62
4:A:150:ILE:CD1	4:A:155:MET:HE2	2.29	0.62
4:A:200:PHE:HD1	4:A:206:LYS:HB2	1.65	0.62
4:A:172:GLU:OE1	4:A:198:PRO:HG2	2.00	0.62
4:A:301:LEU:HD11	4:A:305:GLY:C	2.25	0.62
4:A:299:ARG:HB3	4:A:308:GLY:O	2.00	0.61
4:A:155:MET:HE3	4:A:188:SER:CB	2.05	0.61
1:T:13:DC:H2''	1:T:14:DA:H8	1.66	0.61
4:A:248:LYS:HG2	4:A:249:GLU:H	1.64	0.60
4:A:244:LYS:N	4:A:247:GLU:HG2	2.16	0.60
4:A:11:LEU:HD11	4:A:52:LYS:CB	2.31	0.60
4:A:264:GLN:HB3	4:A:296:TYR:O	2.02	0.60
4:A:317:LYS:N	4:A:317:LYS:HD3	2.16	0.60
4:A:328:ARG:HB3	4:A:328:ARG:NH1	2.16	0.60
4:A:150:ILE:HB	4:A:155:MET:HE2	1.82	0.60
4:A:317:LYS:HD3	4:A:317:LYS:H	1.67	0.59
4:A:133:ASN:ND2	4:A:136:GLN:H	2.00	0.59
4:A:317:LYS:H	4:A:317:LYS:CD	2.15	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:133:ASN:ND2	4:A:135:HIS:HB3	2.18	0.58
4:A:228:LEU:N	4:A:236:MET:O	2.33	0.58
4:A:303:VAL:HG23	4:A:304:THR:HG23	1.86	0.58
3:D:1:DG:OP1	4:A:68:LYS:NZ	2.33	0.57
4:A:316:GLU:N	4:A:317:LYS:HZ1	1.95	0.57
4:A:133:ASN:N	4:A:136:GLN:HE21	2.02	0.57
3:D:1:DG:OP2	4:A:68:LYS:HD3	2.04	0.57
1:T:10:DC:OP1	4:A:231:GLY:HA3	2.05	0.57
4:A:162:VAL:HG23	4:A:163:LEU:N	2.18	0.57
4:A:132:LEU:HA	4:A:136:GLN:NE2	2.20	0.56
4:A:201:THR:H	4:A:204:SER:HB2	1.68	0.56
2:P:9:DG:H4'	4:A:236:MET:HE1	1.87	0.56
4:A:159:GLN:O	4:A:162:VAL:HG23	2.04	0.56
4:A:219:GLN:O	4:A:222:HIS:N	2.29	0.56
4:A:326:LYS:HG3	4:A:326:LYS:O	2.06	0.56
4:A:201:THR:H	4:A:204:SER:CB	2.19	0.56
4:A:113:LYS:NZ	8:A:696:HOH:O	2.38	0.55
4:A:207:GLN:CD	4:A:209:LYS:HD3	2.32	0.55
4:A:133:ASN:HD22	4:A:135:HIS:N	2.05	0.55
4:A:80:GLY:O	4:A:81:LYS:HD3	2.07	0.54
4:A:201:THR:O	4:A:204:SER:HB2	2.07	0.54
4:A:318:ASP:N	4:A:318:ASP:OD1	2.36	0.54
4:A:15:ILE:HG22	4:A:19:LEU:HD22	1.89	0.54
4:A:155:MET:HA	4:A:158:MET:HE3	1.89	0.54
4:A:154:GLU:C	4:A:158:MET:HE3	2.33	0.54
4:A:150:ILE:HD12	4:A:155:MET:HE1	1.90	0.53
4:A:150:ILE:HG23	4:A:158:MET:HE1	1.89	0.53
4:A:244:LYS:CB	4:A:247:GLU:HG2	2.38	0.53
4:A:197:HIS:HA	8:A:774:HOH:O	2.07	0.53
4:A:205:THR:HG22	4:A:205:THR:O	2.07	0.53
4:A:207:GLN:CG	4:A:209:LYS:HD3	2.37	0.53
4:A:217:GLN:HG3	4:A:217:GLN:O	2.06	0.53
4:A:83:ARG:HG2	4:A:83:ARG:NH1	2.01	0.53
4:A:323:ILE:C	4:A:324:GLN:HG2	2.33	0.53
4:A:154:GLU:HB3	4:A:158:MET:HE2	1.90	0.53
4:A:228:LEU:HB2	4:A:236:MET:O	2.09	0.53
4:A:18:MET:O	4:A:22:LEU:HD22	2.09	0.53
4:A:84:LYS:O	4:A:88:ILE:HG13	2.08	0.52
4:A:134:HIS:CE1	4:A:138:ILE:HG12	2.44	0.52
4:A:153:GLU:O	4:A:157:GLN:HG3	2.09	0.52
4:A:291:PHE:CB	4:A:298:ILE:HD11	2.39	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:320:PHE:HB3	4:A:325:TRP:O	2.09	0.52
4:A:83:ARG:HH11	4:A:83:ARG:CG	2.12	0.52
4:A:207:GLN:HB3	4:A:209:LYS:HE2	1.91	0.52
4:A:243:SER:N	8:A:558:HOH:O	2.43	0.52
4:A:203:GLU:OE1	4:A:203:GLU:O	2.28	0.52
4:A:48:LYS:HE3	4:A:334:SER:HB2	1.92	0.51
4:A:197:HIS:CE1	4:A:199:SER:HB3	2.45	0.51
4:A:291:PHE:CG	4:A:298:ILE:HD11	2.45	0.51
4:A:199:SER:O	4:A:204:SER:OG	2.28	0.51
4:A:287:LEU:HA	4:A:291:PHE:O	2.09	0.51
4:A:133:ASN:H	4:A:136:GLN:NE2	2.09	0.51
4:A:315:SER:CA	4:A:317:LYS:HZ1	2.24	0.50
4:A:329:GLU:HB3	4:A:330:PRO:HD2	1.92	0.50
4:A:11:LEU:HD12	4:A:52:LYS:HA	1.93	0.50
4:A:166:VAL:CG1	4:A:173:TYR:HB3	2.41	0.50
4:A:219:GLN:NE2	8:A:631:HOH:O	2.41	0.50
4:A:208:PRO:HA	8:A:577:HOH:O	2.11	0.50
2:P:2:DC:H4'	2:P:3:DT:OP1	2.12	0.49
4:A:207:GLN:HG3	4:A:209:LYS:NZ	2.26	0.49
4:A:278:PHE:CE1	4:A:333:ARG:HD2	2.47	0.49
1:T:11:DA:H5'	4:A:229:SER:HB3	1.95	0.49
4:A:201:THR:O	4:A:203:GLU:N	2.45	0.49
4:A:98:ASN:O	4:A:102:ARG:HG3	2.11	0.49
4:A:209:LYS:HG3	4:A:210:LEU:H	1.78	0.49
4:A:266:TYR:HA	4:A:269:VAL:CG2	2.43	0.49
4:A:291:PHE:HB3	4:A:298:ILE:HD11	1.94	0.49
3:D:2:DT:H5''	4:A:66:GLY:HA3	1.92	0.48
4:A:200:PHE:HD1	4:A:206:LYS:CB	2.25	0.48
4:A:300:PRO:HD3	4:A:311:LEU:HD13	1.94	0.48
1:T:2:DC:H2''	1:T:3:DG:H8	1.73	0.48
4:A:323:ILE:O	4:A:324:GLN:HG2	2.13	0.48
4:A:238:VAL:HA	4:A:253:ARG:O	2.13	0.48
4:A:40:ARG:HH12	4:A:276:ASP:CG	2.21	0.47
4:A:56:GLY:O	4:A:59:ALA:N	2.46	0.47
4:A:301:LEU:HD11	4:A:305:GLY:O	2.14	0.47
1:T:6:DG:C6	1:T:7:DG:C6	3.03	0.47
4:A:303:VAL:CG2	4:A:304:THR:H	2.28	0.47
4:A:90:GLN:O	4:A:90:GLN:HG2	2.01	0.47
4:A:45:VAL:HG23	4:A:46:ILE:N	2.30	0.47
4:A:207:GLN:HG3	4:A:209:LYS:HD3	1.95	0.47
4:A:300:PRO:HD2	4:A:309:GLU:O	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:86:GLU:OE1	4:A:89:ARG:NH1	2.47	0.47
4:A:160:ASP:O	4:A:164:ASN:ND2	2.44	0.47
4:A:244:LYS:H	4:A:247:GLU:CG	2.28	0.47
4:A:328:ARG:HH11	4:A:328:ARG:HG2	1.79	0.47
4:A:91:ASP:OD1	4:A:94:SER:OG	2.31	0.46
8:T:517:HOH:O	4:A:232:GLU:HG2	2.15	0.46
4:A:326:LYS:HG2	8:A:578:HOH:O	2.14	0.46
4:A:149:ARG:NH2	4:A:187:SER:O	2.49	0.46
4:A:182:ARG:C	4:A:184:GLY:H	2.23	0.46
4:A:211:LEU:HB3	8:A:752:HOH:O	2.15	0.46
4:A:207:GLN:CB	4:A:209:LYS:HE2	2.46	0.46
4:A:62:LEU:HB3	4:A:63:PRO:HD2	1.97	0.45
4:A:197:HIS:CE1	4:A:199:SER:CB	2.98	0.45
1:T:13:DC:C2'	1:T:14:DA:C8	2.99	0.45
4:A:105:GLY:CA	4:A:135:HIS:CD2	2.99	0.45
4:A:314:ASP:O	4:A:315:SER:HB3	2.16	0.45
4:A:330:PRO:HA	4:A:333:ARG:HG3	1.97	0.45
1:T:11:DA:N3	8:T:805:HOH:O	2.36	0.45
4:A:105:GLY:HA3	4:A:135:HIS:CD2	2.52	0.45
4:A:40:ARG:HB2	4:A:277:ILE:HD11	1.97	0.45
4:A:133:ASN:HD21	4:A:136:GLN:HG3	1.79	0.45
2:P:7:DG:H2''	2:P:8:DC:O5'	2.17	0.45
4:A:12:ASN:HB3	4:A:46:ILE:HD11	1.99	0.45
4:A:18:MET:CE	4:A:76:PHE:CD1	3.00	0.45
4:A:289:LYS:NZ	4:A:324:GLN:CG	2.79	0.45
4:A:60:LYS:HE3	4:A:67:THR:HB	1.99	0.45
4:A:93:THR:O	4:A:97:ILE:HG13	2.17	0.45
4:A:207:GLN:HB3	4:A:209:LYS:CD	2.46	0.44
4:A:177:VAL:CG1	4:A:181:PHE:CD2	3.00	0.44
4:A:303:VAL:CG2	4:A:304:THR:N	2.79	0.44
4:A:265:TYR:OH	4:A:269:VAL:HG11	2.17	0.44
4:A:265:TYR:CE1	4:A:269:VAL:HG13	2.53	0.44
4:A:91:ASP:HB3	4:A:94:SER:OG	2.16	0.44
4:A:209:LYS:CG	4:A:210:LEU:N	2.79	0.44
4:A:289:LYS:HD3	4:A:289:LYS:HA	1.59	0.44
4:A:334:SER:O	4:A:335:GLU:HB2	2.18	0.44
4:A:86:GLU:O	4:A:89:ARG:HB2	2.17	0.44
4:A:132:LEU:CA	4:A:136:GLN:HE21	2.31	0.44
4:A:18:MET:HE3	4:A:76:PHE:CD1	2.53	0.44
4:A:119:ILE:HG23	4:A:124:ASP:HB3	2.00	0.44
4:A:177:VAL:HG12	4:A:181:PHE:CD2	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:255:ILE:HG12	4:A:256:ASP:N	2.33	0.44
4:A:329:GLU:HB3	4:A:330:PRO:CD	2.48	0.44
4:A:328:ARG:CG	4:A:328:ARG:NH1	2.79	0.43
4:A:223:PHE:O	4:A:239:CYS:HA	2.18	0.43
4:A:157:GLN:NE2	8:A:516:HOH:O	2.29	0.43
4:A:122:LEU:HG	4:A:126:ARG:HD2	1.99	0.43
4:A:244:LYS:CA	4:A:247:GLU:HG2	2.48	0.43
1:T:7:DG:OP1	4:A:287:LEU:HD11	2.17	0.43
4:A:209:LYS:HE2	4:A:210:LEU:HG	2.00	0.43
4:A:28:ASN:HB3	4:A:101:THR:HG21	2.00	0.43
2:P:4:DG:H5'	8:P:677:HOH:O	2.18	0.43
8:T:598:HOH:O	4:A:280:LYS:HD2	2.18	0.43
4:A:155:MET:HE1	4:A:190:ASP:O	2.19	0.43
4:A:207:GLN:HB3	4:A:209:LYS:CE	2.49	0.43
4:A:309:GLU:HA	4:A:310:PRO:HD2	1.76	0.42
4:A:113:LYS:CE	4:A:114:PHE:CD1	3.02	0.42
4:A:133:ASN:N	4:A:136:GLN:NE2	2.67	0.42
4:A:49:TYR:HA	4:A:50:PRO:HD3	1.71	0.42
4:A:311:LEU:HD13	4:A:311:LEU:N	2.34	0.42
4:A:320:PHE:CD1	4:A:327:TYR:HD1	2.38	0.42
2:P:5:DA:H1'	8:P:731:HOH:O	2.19	0.42
4:A:201:THR:C	4:A:203:GLU:H	2.26	0.42
4:A:283:ARG:O	4:A:286:ALA:HB3	2.19	0.42
1:T:2:DC:H2''	1:T:3:DG:N7	2.33	0.42
4:A:40:ARG:HG3	8:A:683:HOH:O	2.19	0.42
4:A:159:GLN:O	4:A:163:LEU:HD23	2.20	0.42
4:A:245:ASN:HD22	4:A:245:ASN:HA	1.75	0.42
4:A:155:MET:CA	4:A:158:MET:HE3	2.50	0.42
4:A:289:LYS:HZ2	4:A:324:GLN:CG	2.33	0.41
4:A:152:ARG:NH2	4:A:184:GLY:C	2.78	0.41
4:A:194:LEU:HD11	4:A:260:ILE:HG12	2.00	0.41
4:A:293:ILE:CG2	4:A:294:ASN:N	2.79	0.41
3:D:1:DG:H2''	3:D:2:DT:O5'	2.20	0.41
4:A:11:LEU:CD1	4:A:52:LYS:HB2	2.41	0.41
4:A:151:PRO:HD2	4:A:154:GLU:HB2	2.02	0.41
4:A:201:THR:C	4:A:203:GLU:N	2.78	0.41
4:A:311:LEU:CD1	4:A:311:LEU:N	2.84	0.41
4:A:113:LYS:CE	4:A:114:PHE:CE1	3.03	0.41
4:A:135:HIS:CD2	4:A:135:HIS:C	2.99	0.41
4:A:241:LEU:HA	4:A:242:PRO:HD3	1.91	0.41
4:A:83:ARG:O	4:A:86:GLU:HB2	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:219:GLN:O	4:A:221:VAL:N	2.54	0.41
4:A:41:LYS:N	8:A:683:HOH:O	2.52	0.41
4:A:273:THR:HB	8:A:622:HOH:O	2.21	0.41
4:A:278:PHE:CD1	4:A:278:PHE:C	2.99	0.40
4:A:281:ASN:HD22	4:A:281:ASN:HA	1.66	0.40
4:A:150:ILE:HG12	4:A:253:ARG:HD2	2.02	0.40
4:A:173:TYR:OH	4:A:210:LEU:HA	2.21	0.40
4:A:22:LEU:HD13	4:A:22:LEU:HA	1.60	0.40
4:A:150:ILE:HG22	4:A:155:MET:HG2	2.03	0.40
4:A:207:GLN:CB	4:A:209:LYS:CE	3.00	0.40
4:A:207:GLN:CG	4:A:209:LYS:CD	3.00	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
4	A	236/335 (70%)	203 (86%)	26 (11%)	7 (3%)	3 1

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
4	A	11	LEU
4	A	179	GLY
4	A	202	SER
4	A	220	LYS
4	A	143	PHE
4	A	249	GLU
4	A	303	VAL

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
4	A	157/295 (53%)	132 (84%)	25 (16%)	2 2

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
4	A	10	THR
4	A	11	LEU
4	A	19	LEU
4	A	22	LEU
4	A	27	LYS
4	A	33	ILE
4	A	40	ARG
4	A	41	LYS
4	A	46	ILE
4	A	77	LEU
4	A	82	LEU
4	A	83	ARG
4	A	90	GLN
4	A	96	SER
4	A	100	LEU
4	A	113	LYS
4	A	133	ASN
4	A	136	GLN
4	A	138	ILE
4	A	152	ARG
4	A	162	VAL
4	A	169	VAL
4	A	171	SER
4	A	176	THR
4	A	298	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
4	A	12	ASN
4	A	28	ASN
4	A	90	GLN
4	A	133	ASN
4	A	136	GLN
4	A	157	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

1 non-standard protein/DNA/RNA residue is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	DOC	P	10	2,1	16,19,20	0.75	0	20,26,29	1.56	3 (15%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	DOC	P	10	2,1	-	0/7/18/19	0/2/2/2

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	P	10	DOC	C2'-C1'-N1	4.37	120.69	112.40
2	P	10	DOC	O4'-C1'-C2'	3.93	111.07	106.41
2	P	10	DOC	C2'-C3'-C4'	2.40	107.21	102.75

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 4 are monoatomic - leaving 1 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
7	DCT	A	338	5	27,28,28	1.86	6 (22%)	37,43,43	1.28	4 (10%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	DCT	A	338	5	-	5/22/31/31	0/2/2/2

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	A	338	DCT	PB-O3B	-4.70	1.54	1.59
7	A	338	DCT	C1'-N1	-3.95	1.38	1.48
7	A	338	DCT	O4'-C1'	-3.08	1.35	1.42
7	A	338	DCT	C2-N1	-2.98	1.33	1.40
7	A	338	DCT	PA-O2A	-2.71	1.42	1.55
7	A	338	DCT	PB-O3A	-2.25	1.57	1.59

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	A	338	DCT	O3G-PG-O3B	-2.85	95.07	104.64
7	A	338	DCT	C4'-O4'-C1'	2.74	112.40	109.81
7	A	338	DCT	O4'-C1'-N1	2.61	112.50	107.86
7	A	338	DCT	O2-C2-N3	-2.23	118.81	122.33

There are no chirality outliers.

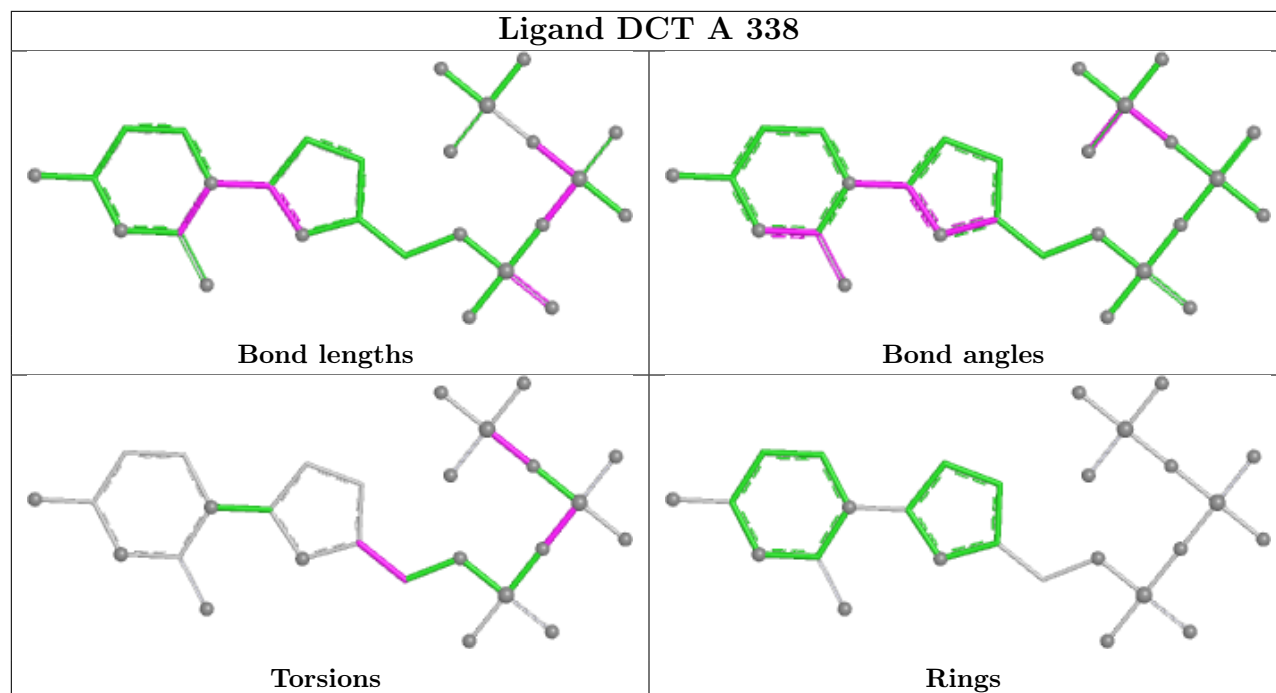
All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	A	338	DCT	O4'-C4'-C5'-O5'
7	A	338	DCT	PB-O3B-PG-O3G
7	A	338	DCT	C3'-C4'-C5'-O5'
7	A	338	DCT	PA-O3A-PB-O1B
7	A	338	DCT	PA-O3A-PB-O2B

There are no ring outliers.

No monomer is involved in short contacts.

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



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

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	16/16 (100%)	0.01	0 100 100	41, 49, 68, 69	0
2	P	9/10 (90%)	0.23	0 100 100	49, 53, 64, 66	0
3	D	5/5 (100%)	-0.25	0 100 100	38, 41, 44, 47	0
4	A	326/335 (97%)	0.55	8 (2%) 58 55	32, 48, 70, 97	0
All	All	356/366 (97%)	0.50	8 (2%) 62 59	32, 49, 70, 97	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	A	205	THR	3.9
4	A	204	SER	2.7
4	A	208	PRO	2.7
4	A	206	LYS	2.7
4	A	203	GLU	2.4
4	A	92	ASP	2.2
4	A	303	VAL	2.2
4	A	207	GLN	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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, 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	B-factors(Å ²)	Q<0.9
2	DOC	P	10	18/19	0.96	0.09	32,47,56,56	0

6.3 Carbohydrates [i](#)

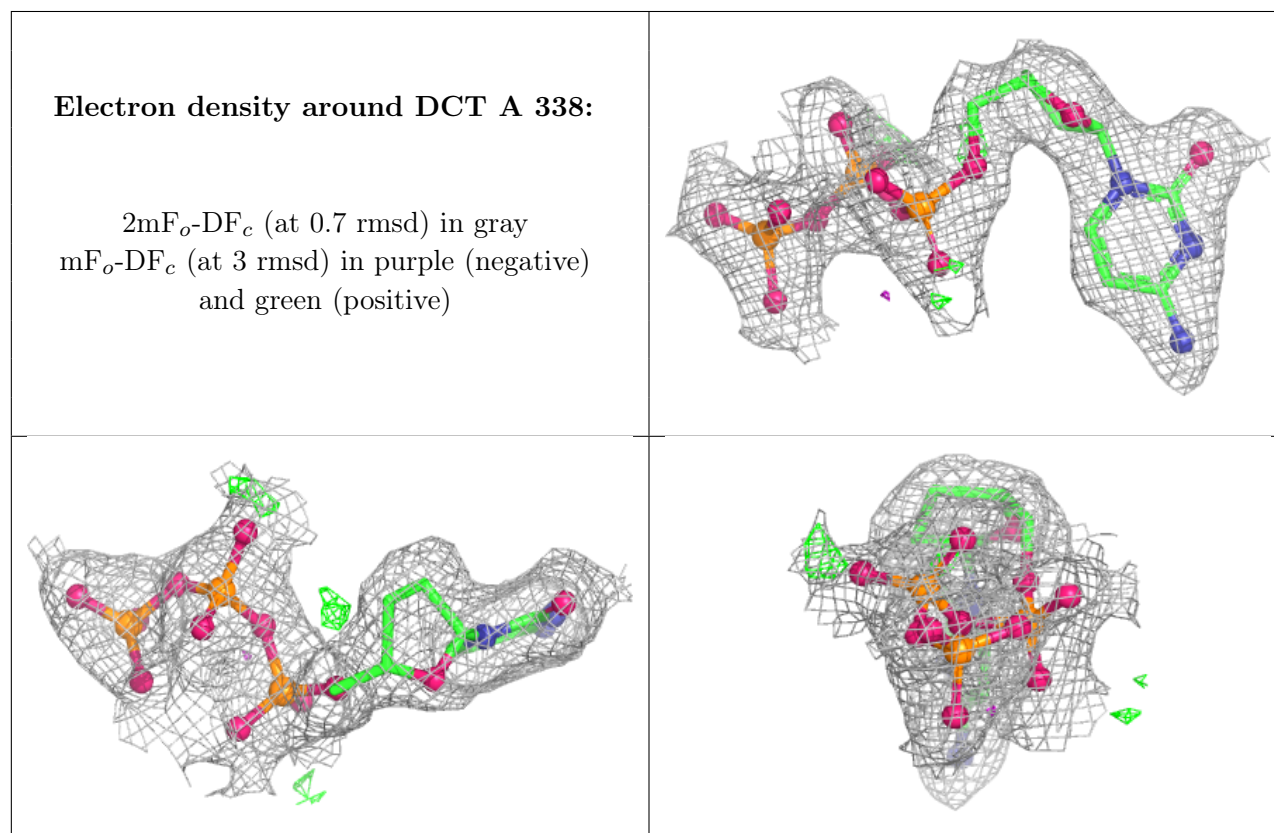
There are no oligosaccharides in this entry.

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, 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	B-factors($\text{\AA}^2$)	Q<0.9
5	MG	A	340	1/1	0.94	0.09	49,49,49,49	0
5	MG	A	339	1/1	0.96	0.07	45,45,45,45	0
7	DCT	A	338	27/27	0.96	0.08	27,41,48,57	0
6	NA	A	342	1/1	0.97	0.07	38,38,38,38	0
6	NA	A	341	1/1	0.98	0.05	29,29,29,29	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



6.5 Other polymers [i](#)

There are no such residues in this entry.