



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 22, 2023 – 08:11 AM EDT

PDB ID : 2R8J
Title : Structure of the Eukaryotic DNA Polymerase eta in complex with 1,2-d(GpG)-cisplatin containing DNA
Authors : Carell, T.; Alt, A.; Lammens, K.
Deposited on : 2007-09-11
Resolution : 3.10 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.35
buster-report : 1.1.7 (2018)
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) : 2.35

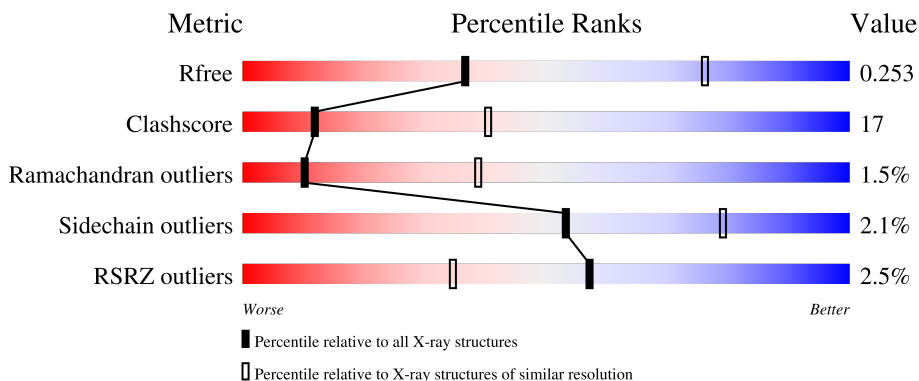
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.10 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



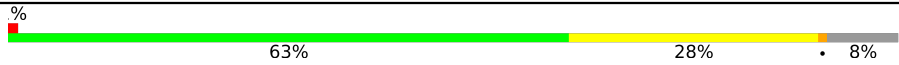
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1094 (3.10-3.10)
Clashscore	141614	1184 (3.10-3.10)
Ramachandran outliers	138981	1141 (3.10-3.10)
Sidechain outliers	138945	1141 (3.10-3.10)
RSRZ outliers	127900	1067 (3.10-3.10)

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	P	9	
1	Q	9	
2	T	11	
2	U	11	
3	A	554	

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Mol	Chain	Length	Quality of chain
3	B	554	 <p>A horizontal bar chart representing the quality of chain. The bar is divided into three segments: a green segment on the left labeled '63%', a yellow segment in the middle labeled '28%', and a red segment on the right labeled '8%'. A small red square is positioned at the start of the bar, and a small black dot is at the end of the red segment. A '%' symbol is located above the bar.</p>

2 Entry composition i

There are 7 unique types of molecules in this entry. The entry contains 9108 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(*DGP*DTP*DGP*DGP*DAP*DTP*DGP*DAP*DG)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
1	Q	9	188	90	39	51	8	0	0	0
1	P	9	188	90	39	51	8	0	0	0

- Molecule 2 is a DNA chain called DNA (5'-D(P*DGP*DGP*DCP*DTP*DCP*DAP*DTP*DCP*DCP*DAP*DC)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	U	11	221	105	39	66	11	0	0	0
2	T	11	221	105	39	66	11	0	0	0

- Molecule 3 is a protein called DNA polymerase eta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	A	511	4059	2585	684	766	24	112	0	0
3	B	509	4044	2577	679	764	24	93	0	0

There are 46 discrepancies between the modelled and reference sequences:

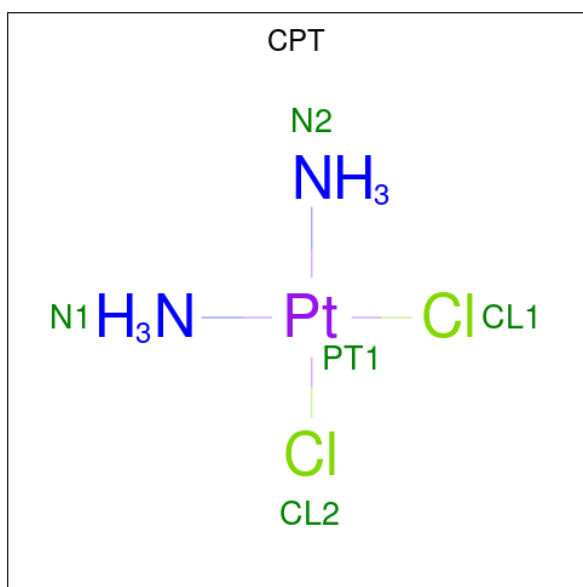
Chain	Residue	Modelled	Actual	Comment	Reference
A	-22	MET	-	expression tag	UNP Q04049
A	-21	ALA	-	expression tag	UNP Q04049
A	-20	SER	-	expression tag	UNP Q04049
A	-19	TRP	-	expression tag	UNP Q04049
A	-18	SER	-	expression tag	UNP Q04049
A	-17	HIS	-	expression tag	UNP Q04049

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-16	PRO	-	expression tag	UNP Q04049
A	-15	GLN	-	expression tag	UNP Q04049
A	-14	PHE	-	expression tag	UNP Q04049
A	-13	GLU	-	expression tag	UNP Q04049
A	-12	LYS	-	expression tag	UNP Q04049
A	-11	GLY	-	expression tag	UNP Q04049
A	-10	ALA	-	expression tag	UNP Q04049
A	-9	SER	-	expression tag	UNP Q04049
A	-8	THR	-	expression tag	UNP Q04049
A	-7	SER	-	expression tag	UNP Q04049
A	-6	LEU	-	expression tag	UNP Q04049
A	-5	TYR	-	expression tag	UNP Q04049
A	-4	LYS	-	expression tag	UNP Q04049
A	-3	LYS	-	expression tag	UNP Q04049
A	-2	ALA	-	expression tag	UNP Q04049
A	-1	GLY	-	expression tag	UNP Q04049
A	0	ARG	-	expression tag	UNP Q04049
B	-22	MET	-	expression tag	UNP Q04049
B	-21	ALA	-	expression tag	UNP Q04049
B	-20	SER	-	expression tag	UNP Q04049
B	-19	TRP	-	expression tag	UNP Q04049
B	-18	SER	-	expression tag	UNP Q04049
B	-17	HIS	-	expression tag	UNP Q04049
B	-16	PRO	-	expression tag	UNP Q04049
B	-15	GLN	-	expression tag	UNP Q04049
B	-14	PHE	-	expression tag	UNP Q04049
B	-13	GLU	-	expression tag	UNP Q04049
B	-12	LYS	-	expression tag	UNP Q04049
B	-11	GLY	-	expression tag	UNP Q04049
B	-10	ALA	-	expression tag	UNP Q04049
B	-9	SER	-	expression tag	UNP Q04049
B	-8	THR	-	expression tag	UNP Q04049
B	-7	SER	-	expression tag	UNP Q04049
B	-6	LEU	-	expression tag	UNP Q04049
B	-5	TYR	-	expression tag	UNP Q04049
B	-4	LYS	-	expression tag	UNP Q04049
B	-3	LYS	-	expression tag	UNP Q04049
B	-2	ALA	-	expression tag	UNP Q04049
B	-1	GLY	-	expression tag	UNP Q04049
B	0	ARG	-	expression tag	UNP Q04049

- Molecule 4 is Cisplatin (three-letter code: CPT) (formula: Cl₂H₆N₂Pt).

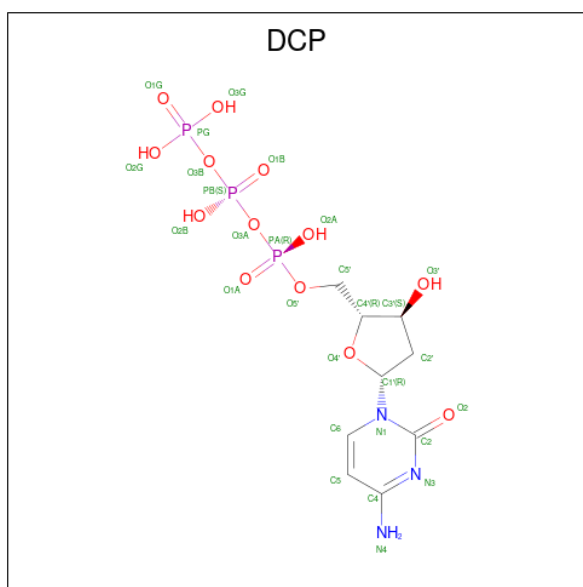


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	U	1	Total N Pt 3 2 1	0	0
4	T	1	Total N Pt 3 2 1	0	0

- Molecule 5 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	2	Total Ca 2 2	0	0
5	B	2	Total Ca 2 2	0	0

- Molecule 6 is 2'-DEOXYCYTIDINE-5'-TRIPHOSPHATE (three-letter code: DCP) (formula: C₉H₁₆N₃O₁₃P₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
6	A	1	Total	C	N	O	P	0	0
			28	9	3	13	3		
6	B	1	Total	C	N	O	P	0	0
			28	9	3	13	3		

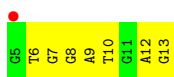
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	Q	4	Total	O	0	0
			4	4		
7	U	3	Total	O	0	0
			3	3		
7	P	2	Total	O	0	0
			2	2		
7	T	2	Total	O	0	0
			2	2		
7	A	49	Total	O	0	0
			49	49		
7	B	61	Total	O	0	0
			61	61		

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.

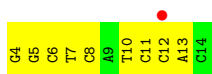
- Molecule 1: DNA (5'-D(*DGP*DTP*DGP*DGP*DAP*DTP*DGP*DAP*DG)-3')



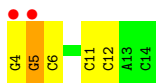
- Molecule 1: DNA (5'-D(*DGP*DTP*DGP*DGP*DAP*DTP*DGP*DAP*DG)-3')



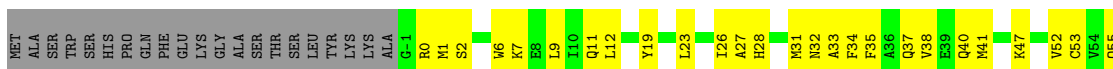
- Molecule 2: DNA (5'-D(P*DGP*DGP*DCP*DTP*DCP*DAP*DTP*DCP*DCP*DAP*DC)-3')

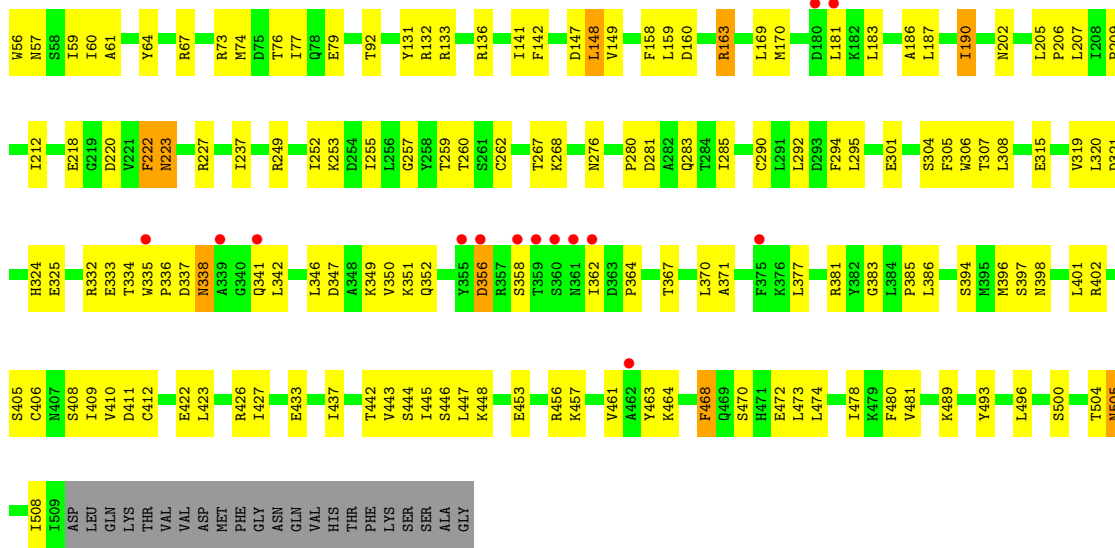


- Molecule 2: DNA (5'-D(P*DGP*DGP*DCP*DTP*DCP*DAP*DTP*DCP*DCP*DAP*DC)-3')

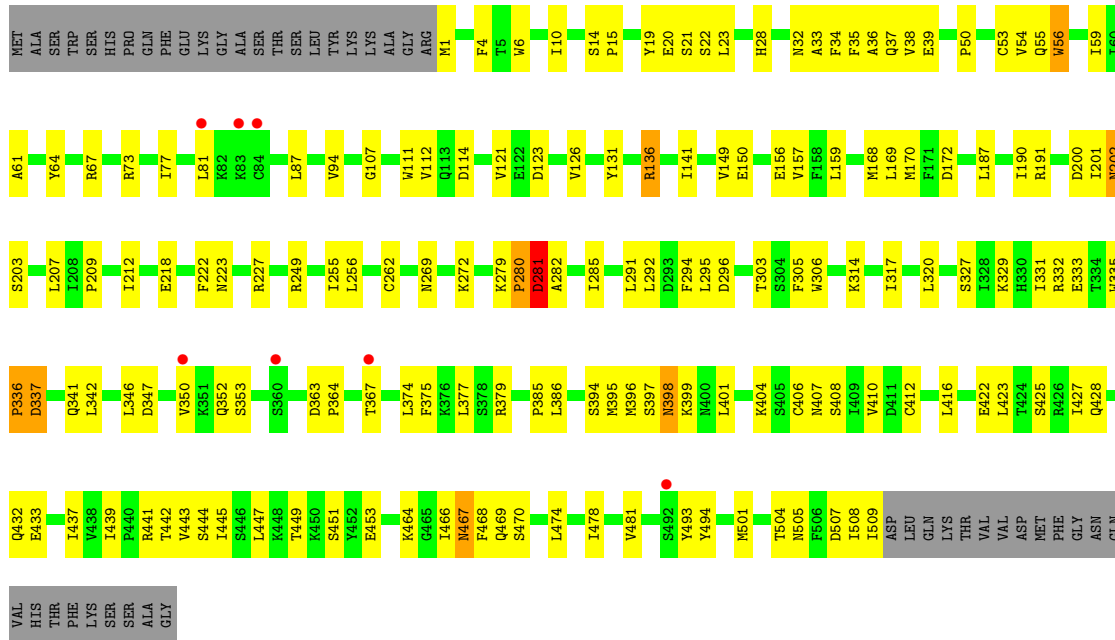


- Molecule 3: DNA polymerase eta





• Molecule 3: DNA polymerase eta



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	104.09Å 104.09Å 293.05Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 3.10 46.55 – 3.10	Depositor EDS
% Data completeness (in resolution range)	99.6 (20.00-3.10) 99.6 (46.55-3.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.15	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.31 (at 3.12Å)	Xtrriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.208 , 0.257 0.199 , 0.253	Depositor DCC
R_{free} test set	2412 reflections (7.99%)	wwPDB-VP
Wilson B-factor (Å ²)	51.5	Xtrriage
Anisotropy	0.048	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 58.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	9108	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.41% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: CA, CPT, DCP

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	P	0.24	0/212	0.68	0/327
1	Q	0.25	0/212	0.63	0/327
2	T	1.22	3/246 (1.2%)	2.21	10/376 (2.7%)
2	U	1.39	6/246 (2.4%)	1.44	7/376 (1.9%)
3	A	0.29	0/4137	0.50	0/5578
3	B	0.28	0/4122	0.51	0/5559
All	All	0.41	9/9175 (0.1%)	0.68	17/12543 (0.1%)

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	U	4	DG	C8-N7	-9.81	1.25	1.30
2	T	5	DG	N3-C4	7.02	1.40	1.35
2	U	4	DG	N3-C4	6.24	1.39	1.35
2	T	5	DG	C3'-O3'	6.05	1.51	1.44
2	U	4	DG	C2-N3	5.81	1.37	1.32
2	T	5	DG	C2-N3	5.69	1.37	1.32
2	U	5	DG	N3-C4	5.52	1.39	1.35
2	U	5	DG	C2-N3	5.20	1.36	1.32
2	U	5	DG	C8-N7	-5.02	1.27	1.30

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	T	5	DG	O5'-P-OP1	-23.40	82.62	110.70
2	T	5	DG	O5'-P-OP2	-19.72	87.04	110.70
2	T	4	DG	O3'-P-O5'	-12.62	80.02	104.00
2	T	4	DG	O4'-C1'-N9	10.87	115.61	108.00
2	U	4	DG	C4'-C3'-C2'	-8.61	95.35	103.10
2	U	4	DG	N7-C8-N9	7.48	116.84	113.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	U	4	DG	O4'-C1'-N9	7.37	113.16	108.00
2	T	5	DG	N3-C2-N2	7.35	125.05	119.90
2	U	4	DG	C8-N9-C4	-7.04	103.58	106.40
2	U	5	DG	N3-C2-N2	6.58	124.50	119.90
2	T	4	DG	C8-N9-C4	-5.99	104.00	106.40
2	T	4	DG	OP1-P-O3'	5.80	117.96	105.20
2	T	4	DG	C4'-C3'-C2'	-5.64	98.02	103.10
2	T	5	DG	O4'-C1'-C2'	-5.46	101.53	105.90
2	U	5	DG	OP2-P-O3'	5.35	116.97	105.20
2	U	5	DG	O4'-C1'-C2'	-5.18	101.75	105.90
2	T	4	DG	C4'-C3'-O3'	5.05	122.39	112.30

There are no chirality outliers.

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	P	188	0	101	6	0
1	Q	188	0	101	12	0
2	T	221	0	124	3	0
2	U	221	0	124	5	0
3	A	4059	0	4105	152	0
3	B	4044	0	4089	128	0
4	T	3	0	0	1	0
4	U	3	0	0	1	0
5	A	2	0	0	0	0
5	B	2	0	0	0	0
6	A	28	0	12	1	0
6	B	28	0	12	1	0
7	A	49	0	0	4	0
7	B	61	0	0	4	0
7	P	2	0	0	0	0
7	Q	4	0	0	0	0
7	T	2	0	0	0	0
7	U	3	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	9108	0	8668	297	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 17.

All (297) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:P:12:DA:H2''	1:P:13:DG:H5'	1.38	1.01
3:A:202:ASN:ND2	3:B:437:ILE:HD11	1.81	0.95
3:B:449:THR:HG22	3:B:453:GLU:H	1.32	0.94
3:B:136:ARG:NH2	3:B:433:GLU:HB2	1.84	0.93
3:B:469:GLN:HG2	3:B:470:SER:H	1.32	0.92
3:A:31:MET:HA	3:A:260:THR:HG22	1.57	0.86
3:A:437:ILE:HD11	3:B:202:ASN:HD22	1.41	0.86
1:Q:7:DG:H2''	1:Q:8:DG:H5''	1.59	0.84
3:A:202:ASN:HD22	3:B:437:ILE:HD11	1.41	0.81
3:A:394:SER:HB3	3:A:504:THR:HG22	1.62	0.79
1:Q:6:DT:H2''	1:Q:7:DG:N7	1.98	0.78
3:B:437:ILE:HB	3:B:509:ILE:HB	1.67	0.77
3:A:442:THR:HG21	3:A:505:ASN:HD22	1.49	0.77
3:A:136:ARG:HH22	3:A:433:GLU:HG3	1.49	0.76
3:A:223:ASN:HD21	3:A:227:ARG:H	1.35	0.75
3:A:338:ASN:ND2	3:A:341:GLN:HG2	2.01	0.75
3:A:338:ASN:HD22	3:A:341:GLN:HG2	1.52	0.73
3:B:439:ILE:HG21	3:B:467:ASN:HA	1.70	0.73
3:A:446:SER:HB3	3:A:456:ARG:HD3	1.69	0.73
3:B:209:PRO:O	3:B:212:ILE:HG22	1.89	0.72
3:B:136:ARG:HH22	3:B:433:GLU:HB2	1.54	0.71
3:B:347:ASP:O	3:B:350:VAL:HG22	1.89	0.71
1:P:9:DA:H2''	1:P:10:DT:C5'	2.20	0.71
3:B:38:VAL:HG21	3:B:131:TYR:CD2	2.26	0.71
3:B:305:PHE:CE1	3:B:377:LEU:HB3	2.27	0.70
3:B:412:CYS:O	3:B:416:LEU:HD23	1.91	0.70
3:A:53:CYS:HB3	3:A:61:ALA:HB3	1.75	0.69
3:B:449:THR:HG22	3:B:453:GLU:N	2.06	0.69
3:A:223:ASN:ND2	3:A:227:ARG:H	1.90	0.68
3:A:292:LEU:HD22	3:A:332:ARG:CZ	2.24	0.68
3:B:223:ASN:HD21	3:B:227:ARG:H	1.41	0.67
3:A:442:THR:HG21	3:A:505:ASN:ND2	2.08	0.67
3:B:333:GLU:O	3:B:336:PRO:HD3	1.96	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:449:THR:CG2	3:B:453:GLU:H	2.06	0.65
3:A:209:PRO:O	3:A:212:ILE:HG22	1.96	0.65
3:B:303:THR:HG21	3:B:317:ILE:HD11	1.78	0.64
3:B:64:TYR:CD2	3:B:280:PRO:HG3	2.33	0.64
3:A:396:MET:HG2	3:A:397:SER:N	2.12	0.63
3:A:437:ILE:HD11	3:B:202:ASN:ND2	2.13	0.63
1:Q:8:DG:H2''	1:Q:9:DA:H5'	1.80	0.62
3:A:169:LEU:HG	3:A:170:MET:HE2	1.82	0.62
3:A:132:ARG:HH11	3:A:132:ARG:HG3	1.65	0.61
3:A:333:GLU:O	3:A:336:PRO:HD3	2.00	0.61
1:P:9:DA:H2''	1:P:10:DT:H5''	1.81	0.61
3:A:442:THR:HG22	3:A:505:ASN:HB2	1.82	0.61
3:B:428:GLN:O	3:B:432:GLN:HG2	2.00	0.60
3:A:149:VAL:HG22	3:A:159:LEU:CD2	2.32	0.60
3:A:40:GLN:HE21	3:A:41:MET:CE	2.14	0.60
3:A:38:VAL:HG21	3:A:131:TYR:CD2	2.35	0.60
3:B:337:ASP:HB2	3:B:341:GLN:OE1	2.02	0.60
3:A:406:CYS:HA	3:A:411:ASP:HB3	1.83	0.60
1:Q:7:DG:C2'	1:Q:8:DG:H5''	2.31	0.60
3:B:394:SER:CB	3:B:504:THR:HG22	2.32	0.59
1:P:9:DA:H2''	1:P:10:DT:H5'	1.83	0.59
3:B:447:LEU:HD12	3:B:447:LEU:O	2.02	0.59
3:A:412:CYS:HB3	3:A:481:VAL:HG21	1.84	0.59
3:B:94:VAL:HG12	3:B:126:VAL:HA	1.84	0.59
3:B:156:GLU:OE2	3:B:272:LYS:HE2	2.02	0.59
3:A:442:THR:CG2	3:A:505:ASN:HD22	2.16	0.59
3:B:469:GLN:HG2	3:B:470:SER:N	2.12	0.58
1:Q:9:DA:H2''	1:Q:10:DT:O5'	2.03	0.58
2:U:6:DC:N4	4:U:45:CPT:N2	2.49	0.58
3:A:398:ASN:HB2	3:A:500:SER:HB3	1.85	0.58
3:B:474:LEU:O	3:B:478:ILE:HG13	2.03	0.58
3:A:437:ILE:HD13	3:B:201:ILE:HD11	1.85	0.58
3:A:320:LEU:HD21	3:A:346:LEU:HD21	1.86	0.57
3:B:223:ASN:ND2	3:B:227:ARG:H	2.02	0.57
3:A:349:LYS:HA	3:A:352:GLN:HE21	1.70	0.57
3:B:364:PRO:HA	3:B:367:THR:HG23	1.85	0.57
3:B:469:GLN:CG	3:B:470:SER:H	2.12	0.57
3:B:329:LYS:O	3:B:333:GLU:HG2	2.05	0.56
3:B:439:ILE:CG2	3:B:467:ASN:HA	2.35	0.56
3:B:136:ARG:HH21	3:B:433:GLU:HB2	1.69	0.56
3:A:190:ILE:HD13	3:A:190:ILE:O	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Q:6:DT:H2''	1:Q:7:DG:C8	2.40	0.56
3:B:38:VAL:HG21	3:B:131:TYR:CE2	2.40	0.56
3:A:409:ILE:HG23	3:A:410:VAL:N	2.20	0.56
3:B:269:ASN:HB2	7:B:3025:HOH:O	2.06	0.56
3:A:401:LEU:HD13	3:A:406:CYS:HB3	1.88	0.56
3:A:52:VAL:HG21	3:A:59:ILE:HD13	1.87	0.55
3:A:64:TYR:CD2	3:A:280:PRO:HD3	2.42	0.55
3:B:77:ILE:O	3:B:81:LEU:HG	2.06	0.55
3:A:398:ASN:CB	3:A:500:SER:HB3	2.36	0.55
3:A:307:THR:O	3:A:308:LEU:HD12	2.07	0.55
3:A:443:VAL:HG23	3:A:461:VAL:HB	1.87	0.55
3:A:183:LEU:HD22	3:A:187:LEU:HD12	1.88	0.55
3:B:53:CYS:HB3	3:B:61:ALA:HB3	1.89	0.55
3:A:0:ARG:HA	3:A:207:LEU:HD11	1.90	0.54
3:B:112:VAL:HG12	3:B:114:ASP:H	1.71	0.54
3:A:267:THR:HG23	3:A:383:GLY:HA2	1.90	0.54
3:A:262:CYS:HB3	3:A:283:GLN:HG3	1.89	0.54
3:A:301:GLU:H	3:A:304:SER:CB	2.21	0.54
3:B:292:LEU:HD22	3:B:332:ARG:CZ	2.38	0.54
3:B:303:THR:CG2	3:B:317:ILE:HD11	2.38	0.54
3:A:133:ARG:HD3	7:A:2015:HOH:O	2.07	0.54
3:A:463:TYR:HA	3:A:472:GLU:OE1	2.07	0.54
3:B:121:VAL:HG12	7:B:3012:HOH:O	2.08	0.54
3:A:337:ASP:HB2	3:A:341:GLN:HG3	1.90	0.54
3:B:449:THR:OG1	3:B:493:TYR:HA	2.08	0.54
3:A:396:MET:HE3	3:A:398:ASN:HD22	1.73	0.54
3:B:1:MET:HE2	3:B:1:MET:HA	1.90	0.53
3:A:315:GLU:O	3:A:319:VAL:HG23	2.09	0.53
3:A:349:LYS:HA	3:A:352:GLN:NE2	2.24	0.53
3:B:21:SER:O	3:B:22:SER:HB2	2.08	0.53
3:B:394:SER:HB2	3:B:504:THR:HG22	1.90	0.53
1:Q:12:DA:H2''	1:Q:13:DG:OP2	2.07	0.53
3:A:38:VAL:HG21	3:A:131:TYR:CE2	2.43	0.53
1:P:9:DA:C2'	1:P:10:DT:H5''	2.39	0.53
3:B:379:ARG:HD3	7:B:3047:HOH:O	2.09	0.53
3:B:50:PRO:O	3:B:87:LEU:HD12	2.09	0.52
3:A:26:ILE:HG21	3:A:268:LYS:HG2	1.90	0.52
3:B:412:CYS:HB3	3:B:481:VAL:HG21	1.92	0.52
3:B:222:PHE:CE2	3:B:294:PHE:HA	2.45	0.52
3:A:28:HIS:HD2	3:A:158:PHE:CZ	2.27	0.52
3:A:342:LEU:O	3:A:342:LEU:HD23	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:474:LEU:O	3:A:478:ILE:HG13	2.09	0.52
3:A:12:LEU:O	3:A:163:ARG:HD3	2.11	0.51
3:A:142:PHE:CE2	3:A:252:ILE:HD11	2.45	0.51
3:A:478:ILE:O	3:A:481:VAL:HG12	2.10	0.51
3:B:493:TYR:HD2	3:B:494:TYR:CE1	2.28	0.51
3:B:6:TRP:O	3:B:10:ILE:HG12	2.10	0.51
3:A:443:VAL:HG12	3:A:444:SER:H	1.74	0.51
3:B:190:ILE:HG23	3:B:191:ARG:N	2.25	0.51
3:A:32:ASN:ND2	3:A:281:ASP:H	2.10	0.50
3:B:4:PHE:CD2	3:B:23:LEU:HG	2.46	0.50
3:B:291:LEU:O	3:B:295:LEU:HD23	2.10	0.50
3:A:40:GLN:HE21	3:A:41:MET:HE2	1.75	0.49
2:T:5:DG:H1'	2:T:6:DC:H5'	1.94	0.49
3:A:305:PHE:CE1	3:A:377:LEU:HB3	2.47	0.49
3:B:141:ILE:HD13	3:B:255:ILE:HD12	1.93	0.49
3:B:320:LEU:HD11	3:B:374:LEU:HD13	1.94	0.49
3:A:301:GLU:H	3:A:304:SER:HB3	1.77	0.49
3:B:35:PHE:O	3:B:38:VAL:HG22	2.13	0.49
3:B:38:VAL:HG21	3:B:131:TYR:HD2	1.72	0.49
3:B:442:THR:CB	3:B:505:ASN:HD22	2.25	0.49
3:A:136:ARG:NH2	3:A:433:GLU:HA	2.27	0.49
3:B:249:ARG:HD2	3:B:262:CYS:HB2	1.95	0.49
3:A:32:ASN:O	3:A:33:ALA:C	2.52	0.49
3:B:508:ILE:HG13	3:B:508:ILE:O	2.12	0.49
2:T:11:DC:H2''	2:T:12:DC:C5	2.48	0.49
3:A:40:GLN:HE21	3:A:41:MET:HE3	1.78	0.49
3:A:32:ASN:HD21	3:A:281:ASP:H	1.60	0.48
3:A:183:LEU:HD22	3:A:187:LEU:CD1	2.43	0.48
3:A:132:ARG:HG3	3:A:132:ARG:NH1	2.26	0.48
3:A:447:LEU:C	3:A:447:LEU:HD12	2.32	0.48
3:B:54:VAL:HG21	3:B:111:TRP:CZ3	2.48	0.48
3:B:169:LEU:HG	3:B:170:MET:HE2	1.93	0.48
3:A:443:VAL:HG12	3:A:444:SER:N	2.28	0.48
3:B:255:ILE:O	3:B:256:LEU:HD23	2.13	0.48
3:A:337:ASP:HB2	3:A:341:GLN:CG	2.44	0.48
3:B:77:ILE:HG21	3:B:111:TRP:HB3	1.95	0.48
3:A:342:LEU:HD22	3:A:371:ALA:HA	1.95	0.48
3:B:296:ASP:OD2	3:B:332:ARG:NH2	2.45	0.48
3:A:19:TYR:CZ	3:A:385:PRO:HB3	2.48	0.48
3:A:170:MET:HE3	3:A:190:ILE:HD12	1.95	0.48
3:B:168:MET:HA	3:B:172:ASP:HB2	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:U:12:DC:H2''	2:U:13:DA:N7	2.29	0.47
3:A:73:ARG:NH1	6:A:2001:DCP:H2'1	2.29	0.47
3:B:404:LYS:HB3	3:B:407:ASN:HD22	1.79	0.47
3:B:423:LEU:O	3:B:427:ILE:HG13	2.14	0.47
3:A:448:LYS:HA	3:A:453:GLU:O	2.15	0.47
3:B:21:SER:O	3:B:22:SER:CB	2.61	0.47
3:B:478:ILE:HA	3:B:481:VAL:HG12	1.96	0.47
3:A:92:THR:HB	7:A:2019:HOH:O	2.13	0.47
3:B:200:ASP:HB3	3:B:203:SER:OG	2.15	0.47
3:B:449:THR:HG23	3:B:451:SER:H	1.79	0.47
3:A:60:ILE:HG22	3:A:74:MET:CE	2.45	0.47
3:A:324:HIS:O	3:A:325:GLU:HG3	2.14	0.47
3:B:422:GLU:O	3:B:425:SER:HB3	2.15	0.47
1:Q:12:DA:H4'	3:A:306:TRP:CH2	2.50	0.47
3:B:21:SER:C	3:B:23:LEU:H	2.16	0.47
3:A:47:LYS:NZ	3:A:259:THR:HG21	2.30	0.46
3:B:33:ALA:HB1	3:B:36:ALA:HB3	1.96	0.46
3:A:141:ILE:HD13	3:A:255:ILE:HD12	1.97	0.46
3:B:445:ILE:HD11	3:B:501:MET:HE3	1.97	0.46
3:A:9:LEU:HA	3:A:23:LEU:HD12	1.98	0.46
3:A:457:LYS:HD3	3:A:480:PHE:CG	2.50	0.46
3:A:1:MET:H	3:A:207:LEU:HD21	1.81	0.46
3:A:347:ASP:O	3:A:351:LYS:HG3	2.16	0.46
3:B:56:TRP:CE2	3:B:121:VAL:HG23	2.50	0.46
3:B:394:SER:HB3	3:B:504:THR:HG22	1.97	0.46
3:A:160:ASP:HB2	3:A:386:LEU:HD22	1.98	0.46
3:B:306:TRP:HA	3:B:306:TRP:CE3	2.51	0.46
3:A:32:ASN:O	3:A:37:GLN:HG3	2.15	0.45
3:B:474:LEU:HG	3:B:478:ILE:HD11	1.97	0.45
2:U:7:DT:H2''	2:U:8:DC:H5'	1.97	0.45
3:B:187:LEU:HD21	3:B:212:ILE:HD13	1.98	0.45
3:A:57:ASN:O	3:A:77:ILE:HG13	2.16	0.45
3:A:307:THR:C	3:A:308:LEU:HD12	2.37	0.45
2:T:6:DC:N4	4:T:45:CPT:N2	2.65	0.45
3:A:64:TYR:HA	3:A:67:ARG:HG3	1.98	0.45
3:A:73:ARG:O	3:A:74:MET:HB2	2.17	0.45
3:A:169:LEU:HG	3:A:170:MET:CE	2.47	0.45
3:B:327:SER:O	3:B:331:ILE:HG13	2.16	0.45
3:A:227:ARG:HD2	3:A:290:CYS:SG	2.56	0.45
3:B:32:ASN:ND2	3:B:280:PRO:O	2.49	0.45
3:A:142:PHE:CZ	3:A:252:ILE:HD11	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:150:GLU:OE1	3:B:386:LEU:HA	2.17	0.45
3:A:347:ASP:OD1	3:A:367:THR:HG21	2.16	0.45
3:A:346:LEU:HD13	3:A:370:LEU:HD23	1.99	0.44
3:A:457:LYS:HD3	3:A:480:PHE:HA	1.99	0.44
3:B:281:ASP:HB3	3:B:282:ALA:H	1.56	0.44
3:A:169:LEU:HD22	3:A:237:ILE:HD11	1.99	0.44
3:A:406:CYS:SG	3:A:496:LEU:HD12	2.58	0.44
3:B:32:ASN:O	3:B:37:GLN:HG3	2.18	0.44
3:B:314:LYS:HB3	3:B:314:LYS:NZ	2.33	0.44
1:Q:6:DT:H5'	3:A:456:ARG:NH1	2.33	0.44
3:A:346:LEU:O	3:A:350:VAL:HG13	2.18	0.44
3:B:443:VAL:HG12	3:B:444:SER:N	2.32	0.44
3:B:466:ILE:O	3:B:467:ASN:HB3	2.18	0.44
3:A:218:GLU:O	3:A:285:ILE:HD11	2.18	0.44
3:B:396:MET:HG2	3:B:397:SER:N	2.33	0.44
3:A:334:THR:HB	3:A:335:TRP:CE3	2.53	0.43
3:B:28:HIS:CE1	3:B:272:LYS:HD2	2.53	0.43
3:B:169:LEU:HD12	3:B:187:LEU:HD13	2.00	0.43
1:P:12:DA:C2'	1:P:13:DG:H5'	2.28	0.43
3:A:6:TRP:HA	3:A:205:LEU:HD21	2.00	0.43
3:A:26:ILE:HG22	3:A:27:ALA:N	2.33	0.43
3:A:260:THR:O	3:A:281:ASP:HA	2.18	0.43
3:A:402:ARG:HB2	7:A:2017:HOH:O	2.17	0.43
3:A:364:PRO:O	3:A:367:THR:HG22	2.18	0.43
3:B:67:ARG:NH2	7:B:3036:HOH:O	2.52	0.43
1:Q:6:DT:H2''	1:Q:7:DG:C5	2.52	0.43
1:Q:12:DA:H5''	3:A:306:TRP:CZ3	2.53	0.43
3:A:60:ILE:HG22	3:A:74:MET:HE3	2.00	0.43
3:A:362:ILE:O	3:A:364:PRO:HD3	2.18	0.43
3:A:396:MET:CE	3:A:398:ASN:HD22	2.30	0.43
3:A:27:ALA:O	3:A:158:PHE:HA	2.19	0.43
3:A:401:LEU:HD13	3:A:406:CYS:CB	2.48	0.43
3:B:107:GLY:O	3:B:112:VAL:HG21	2.19	0.43
3:A:2:SER:HB3	3:A:205:LEU:O	2.18	0.43
3:A:181:LEU:HD11	3:A:186:ALA:HB2	2.01	0.43
3:A:423:LEU:O	3:A:427:ILE:HG13	2.18	0.43
3:A:468:PHE:C	3:A:470:SER:H	2.21	0.43
3:B:59:ILE:O	3:B:73:ARG:HG3	2.18	0.43
3:B:33:ALA:HA	6:B:3001:DCP:O1B	2.19	0.43
3:B:408:SER:OG	3:B:410:VAL:HG22	2.18	0.43
3:A:73:ARG:HG3	3:A:74:MET:N	2.34	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:76:THR:OG1	3:A:79:GLU:HG3	2.18	0.43
3:A:202:ASN:HD21	3:B:437:ILE:HD11	1.73	0.43
3:A:409:ILE:CG2	3:A:410:VAL:N	2.81	0.43
3:A:253:LYS:O	3:A:257:GLY:HA2	2.18	0.43
3:B:218:GLU:O	3:B:285:ILE:HD11	2.19	0.43
3:A:52:VAL:CG2	3:A:59:ILE:HG23	2.50	0.42
3:A:445:ILE:O	3:A:456:ARG:HG3	2.18	0.42
3:B:64:TYR:OH	3:B:279:LYS:HD3	2.19	0.42
3:A:147:ASP:O	3:A:148:LEU:HB2	2.18	0.42
3:A:422:GLU:O	3:A:426:ARG:HG3	2.19	0.42
3:A:35:PHE:O	3:A:38:VAL:HG22	2.20	0.42
3:A:222:PHE:CE2	3:A:294:PHE:HA	2.55	0.42
3:A:223:ASN:HD21	3:A:227:ARG:N	2.11	0.42
3:A:281:ASP:CG	3:A:281:ASP:O	2.57	0.42
3:A:489:LYS:N	7:A:2022:HOH:O	2.52	0.42
1:Q:7:DG:H2''	1:Q:8:DG:C5'	2.40	0.42
3:A:56:TRP:HA	3:A:56:TRP:CE3	2.55	0.42
3:A:381:ARG:HG2	3:A:381:ARG:HH11	1.85	0.42
3:B:64:TYR:HA	3:B:67:ARG:HG3	2.01	0.42
3:B:306:TRP:HA	3:B:306:TRP:HE3	1.84	0.42
3:B:375:PHE:CE1	3:B:379:ARG:NH2	2.87	0.42
3:B:398:ASN:CG	3:B:399:LYS:N	2.73	0.42
3:A:35:PHE:HB2	3:A:61:ALA:HB2	2.01	0.42
3:A:55:GLN:O	3:A:56:TRP:HB2	2.20	0.42
2:U:10:DT:H2'	2:U:11:DC:C6	2.55	0.42
3:A:7:LYS:O	3:A:11:GLN:HG3	2.19	0.42
3:B:38:VAL:HG23	3:B:39:GLU:N	2.35	0.42
3:B:352:GLN:HG3	3:B:353:SER:H	1.84	0.42
3:B:1:MET:N	3:B:207:LEU:HG	2.35	0.41
3:B:441:ARG:HG2	3:B:441:ARG:HH11	1.85	0.41
3:A:461:VAL:O	3:A:461:VAL:HG13	2.20	0.41
3:B:149:VAL:HG22	3:B:159:LEU:CD2	2.51	0.41
3:B:395:MET:HE1	3:B:427:ILE:HG12	2.02	0.41
3:A:338:ASN:HD22	3:A:338:ASN:N	2.17	0.41
3:B:449:THR:HG23	3:B:451:SER:N	2.35	0.41
3:A:209:PRO:HG2	3:A:212:ILE:HB	2.02	0.41
3:B:346:LEU:O	3:B:350:VAL:HG13	2.20	0.41
3:B:350:VAL:HG11	3:B:364:PRO:HB3	2.01	0.41
3:B:14:SER:HA	3:B:15:PRO:HD3	1.91	0.41
3:B:401:LEU:HD13	3:B:406:CYS:HB3	2.03	0.41
2:U:6:DC:H6	2:U:6:DC:H2'	1.70	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:55:GLN:HG3	3:B:126:VAL:HG21	2.03	0.41
3:B:281:ASP:O	3:B:282:ALA:HB3	2.20	0.41
3:B:363:ASP:HA	3:B:364:PRO:HD2	1.97	0.41
3:A:447:LEU:HD21	3:A:493:TYR:OH	2.20	0.41
3:A:356:ASP:C	3:A:358:SER:H	2.25	0.40
3:A:408:SER:OG	3:A:410:VAL:HG22	2.22	0.40
3:A:410:VAL:HG23	3:A:411:ASP:N	2.35	0.40
3:B:305:PHE:HE1	3:B:377:LEU:HB3	1.82	0.40
3:A:223:ASN:CG	3:A:227:ARG:H	2.25	0.40
3:A:1:MET:CE	3:A:206:PRO:HA	2.51	0.40
3:A:163:ARG:HD3	3:A:163:ARG:N	2.36	0.40
3:B:19:TYR:CZ	3:B:385:PRO:HB3	2.56	0.40
3:B:169:LEU:HD12	3:B:187:LEU:CD1	2.51	0.40
3:B:335:TRP:CE3	3:B:342:LEU:HD13	2.56	0.40
3:A:169:LEU:HD12	3:A:187:LEU:HD13	2.02	0.40
3:A:249:ARG:NH1	3:A:283:GLN:HB2	2.36	0.40
3:B:32:ASN:O	3:B:33:ALA:C	2.58	0.40
3:B:150:GLU:O	3:B:157:VAL:HA	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	A	509/554 (92%)	451 (89%)	51 (10%)	7 (1%)	11	40
3	B	507/554 (92%)	440 (87%)	59 (12%)	8 (2%)	9	37
All	All	1016/1108 (92%)	891 (88%)	110 (11%)	15 (2%)	10	39

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	222	PHE
3	A	405	SER
3	B	337	ASP
3	B	468	PHE
3	A	148	LEU
3	A	468	PHE
3	A	508	ILE
3	B	34	PHE
3	B	281	ASP
3	B	467	ASN
3	A	34	PHE
3	A	505	ASN
3	B	56	TRP
3	B	336	PRO
3	B	280	PRO

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.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
3	A	457/493 (93%)	446 (98%)	11 (2%)	49 76
3	B	456/493 (92%)	448 (98%)	8 (2%)	59 82
All	All	913/986 (93%)	894 (98%)	19 (2%)	53 79

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	163	ARG
3	A	190	ILE
3	A	220	ASP
3	A	223	ASN
3	A	276	ASN
3	A	295	LEU
3	A	321	ASP
3	A	338	ASN
3	A	356	ASP

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Mol	Chain	Res	Type
3	A	464	LYS
3	A	473	LEU
3	B	20	GLU
3	B	123	ASP
3	B	136	ARG
3	B	202	ASN
3	B	281	ASP
3	B	398	ASN
3	B	464	LYS
3	B	507	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
3	A	11	GLN
3	A	28	HIS
3	A	32	ASN
3	A	40	GLN
3	A	57	ASN
3	A	202	ASN
3	A	223	ASN
3	A	243	GLN
3	A	283	GLN
3	A	288	ASN
3	A	338	ASN
3	A	352	GLN
3	A	398	ASN
3	A	505	ASN
3	B	32	ASN
3	B	40	GLN
3	B	57	ASN
3	B	202	ASN
3	B	223	ASN
3	B	288	ASN
3	B	326	ASN
3	B	330	HIS
3	B	361	ASN
3	B	407	ASN
3	B	505	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 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
4	CPT	T	45	2	0,2,4	-	-	-	-	-
6	DCP	B	3001	5	25,29,29	1.93	4 (16%)	37,45,45	1.56	9 (24%)
6	DCP	A	2001	5	25,29,29	1.88	4 (16%)	37,45,45	1.57	10 (27%)
4	CPT	U	45	2	0,2,4	-	-	-	-	-

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
6	DCP	B	3001	5	-	8/22/34/34	0/2/2/2
6	DCP	A	2001	5	-	5/22/34/34	0/2/2/2

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	3001	DCP	C6-C5	7.88	1.53	1.35
6	A	2001	DCP	C6-C5	7.64	1.52	1.35
6	A	2001	DCP	C4-N3	3.32	1.41	1.34
6	B	3001	DCP	C4-N3	3.14	1.40	1.34
6	B	3001	DCP	PG-O1G	2.15	1.57	1.50
6	B	3001	DCP	PA-O5'	2.07	1.67	1.59
6	A	2001	DCP	PG-O1G	2.05	1.57	1.50
6	A	2001	DCP	PA-O5'	2.04	1.67	1.59

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	B	3001	DCP	O2-C2-N3	-4.53	114.95	122.33
6	A	2001	DCP	O2-C2-N3	-4.33	115.28	122.33
6	A	2001	DCP	C5'-C4'-C3'	-2.90	97.73	114.74
6	A	2001	DCP	O4'-C4'-C5'	2.76	118.45	109.37
6	B	3001	DCP	C5'-C4'-C3'	-2.73	98.68	114.74
6	B	3001	DCP	PB-O3B-PG	-2.66	123.69	132.83
6	B	3001	DCP	O4'-C4'-C5'	2.65	118.08	109.37
6	A	2001	DCP	C2'-C3'-C4'	2.59	108.16	102.76
6	A	2001	DCP	C1'-N1-C2	2.41	121.97	117.74
6	A	2001	DCP	O5'-C5'-C4'	2.30	116.92	108.99
6	B	3001	DCP	N1-C2-N3	2.27	122.94	118.81
6	A	2001	DCP	C6-C5-C4	-2.25	113.87	117.50
6	B	3001	DCP	C6-N1-C2	-2.22	116.64	120.49
6	A	2001	DCP	C6-N1-C2	-2.20	116.68	120.49
6	B	3001	DCP	O4'-C1'-N1	2.13	111.67	107.86
6	B	3001	DCP	C6-C5-C4	-2.12	114.07	117.50
6	A	2001	DCP	N1-C2-N3	2.11	122.64	118.81
6	A	2001	DCP	PB-O3B-PG	-2.02	125.91	132.83
6	B	3001	DCP	C2'-C3'-C4'	2.01	106.95	102.76

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	2001	DCP	O4'-C4'-C5'-O5'
6	A	2001	DCP	C5'-O5'-PA-O1A
6	B	3001	DCP	O4'-C4'-C5'-O5'
6	B	3001	DCP	C5'-O5'-PA-O3A
6	A	2001	DCP	C3'-C4'-C5'-O5'
6	B	3001	DCP	C3'-C4'-C5'-O5'

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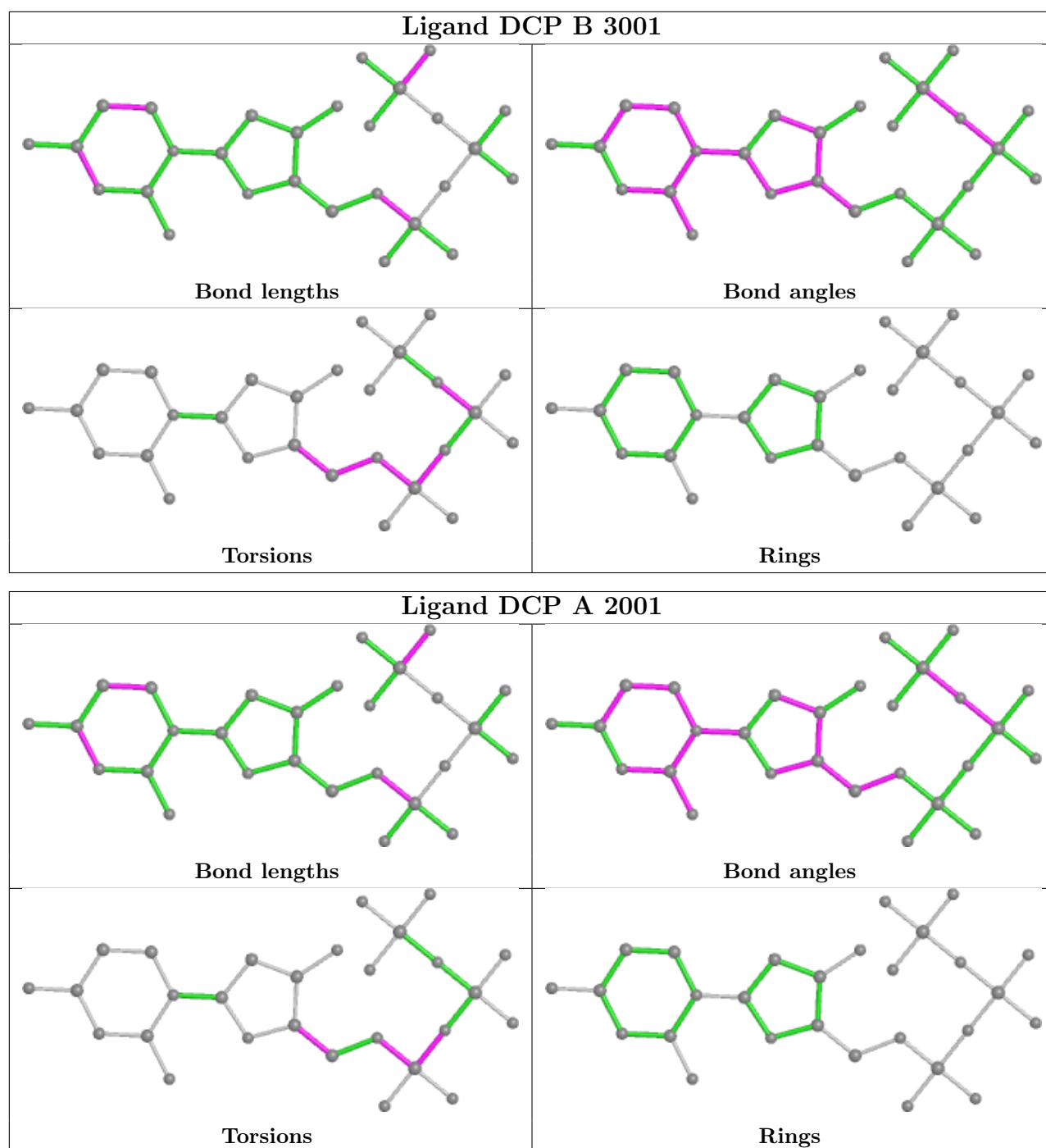
Mol	Chain	Res	Type	Atoms
6	A	2001	DCP	PB-O3A-PA-O1A
6	B	3001	DCP	PB-O3A-PA-O1A
6	B	3001	DCP	PG-O3B-PB-O2B
6	B	3001	DCP	C5'-O5'-PA-O1A
6	B	3001	DCP	C5'-O5'-PA-O2A
6	B	3001	DCP	C4'-C5'-O5'-PA
6	A	2001	DCP	PB-O3A-PA-O2A

There are no ring outliers.

4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	T	45	CPT	1	0
6	B	3001	DCP	1	0
6	A	2001	DCP	1	0
4	U	45	CPT	1	0

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

6.1 Protein, DNA and RNA chains

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	P	9/9 (100%)	0.40	1 (11%) 5 2	43, 50, 56, 60	0
1	Q	9/9 (100%)	0.64	1 (11%) 5 2	37, 51, 60, 66	0
2	T	11/11 (100%)	1.22	2 (18%) 1 0	43, 55, 78, 82	0
2	U	11/11 (100%)	0.44	1 (9%) 9 3	35, 48, 63, 66	0
3	A	508/554 (91%)	-0.16	14 (2%) 53 30	15, 37, 85, 111	23 (4%)
3	B	509/554 (91%)	-0.21	7 (1%) 75 56	9, 37, 70, 92	31 (6%)
All	All	1057/1148 (92%)	-0.15	26 (2%) 57 34	9, 38, 79, 111	54 (5%)

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	A	361	ASN	4.7
3	A	360	SER	4.2
2	T	5	DG	4.1
3	A	180	ASP	4.0
2	T	4	DG	3.7
3	A	355	TYR	3.6
3	B	367	THR	3.4
3	A	356	ASP	3.2
3	B	81	LEU	3.0
1	Q	5	DG	2.8
3	A	359	THR	2.7
3	A	181	LEU	2.6
3	B	84	CYS	2.6
3	A	339	ALA	2.6
1	P	5	DG	2.4
3	B	492	SER	2.3
3	B	83	LYS	2.3
3	A	358	SER	2.3
3	A	335	TRP	2.2

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Mol	Chain	Res	Type	RSRZ
3	A	462	ALA	2.2
2	U	12	DC	2.2
3	A	375	PHE	2.1
3	B	350	VAL	2.1
3	B	360	SER	2.1
3	A	362	ILE	2.1
3	A	341	GLN	2.0

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 monosaccharides in this entry.

6.4 Ligands [i](#)

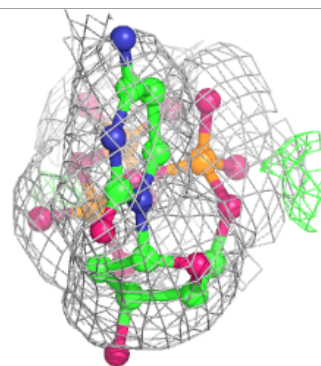
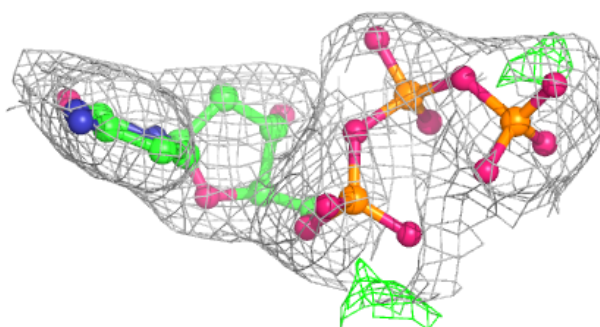
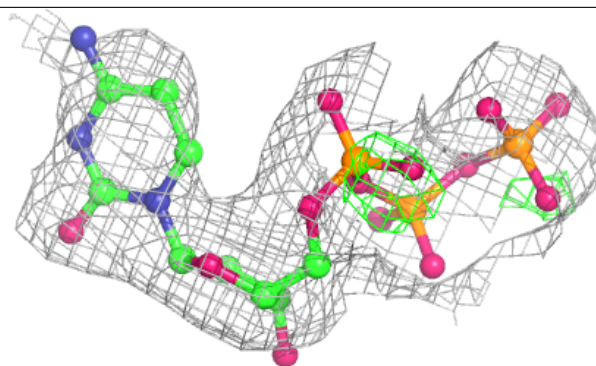
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
5	CA	A	1501	1/1	0.89	0.19	91,91,91,91	0
5	CA	B	1701	1/1	0.90	0.17	39,39,39,39	0
6	DCP	A	2001	28/28	0.96	0.22	29,34,38,40	0
6	DCP	B	3001	28/28	0.97	0.26	32,42,48,50	0
5	CA	A	1502	1/1	0.98	0.15	21,21,21,21	0
5	CA	B	1702	1/1	0.98	0.06	21,21,21,21	0
4	CPT	T	45	3/5	0.99	0.17	50,50,55,57	0
4	CPT	U	45	3/5	1.00	0.14	38,38,39,42	0

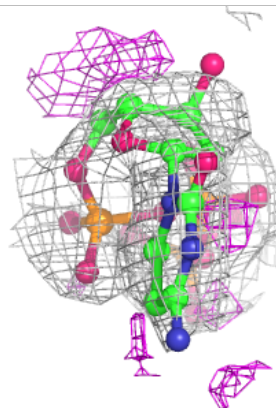
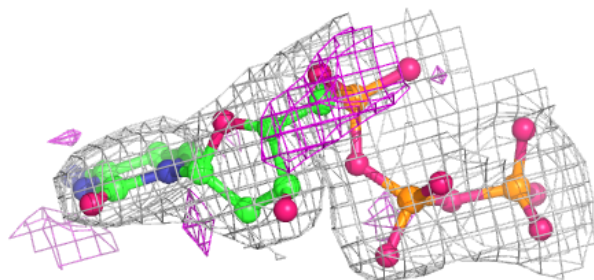
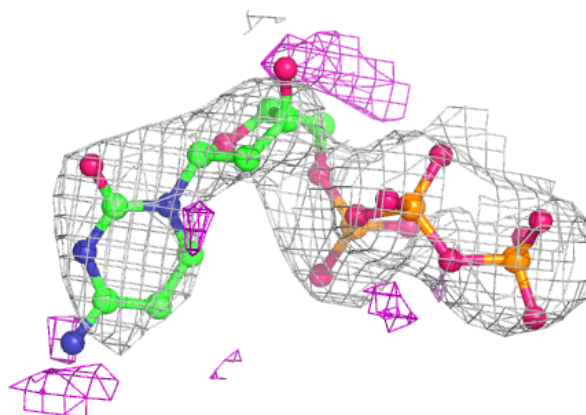
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around DCP A 2001:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around DCP B 3001:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.