



Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 02:59 am BST

PDB ID : 4F4Y
Title : Y-family DNA polymerase chimera Dbh-Dpo4-Dbh
Authors : Pata, J.D.; Wilson, R.C.
Deposited on : 2012-05-11
Resolution : 2.34 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 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 : 2.11
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) : 2.11

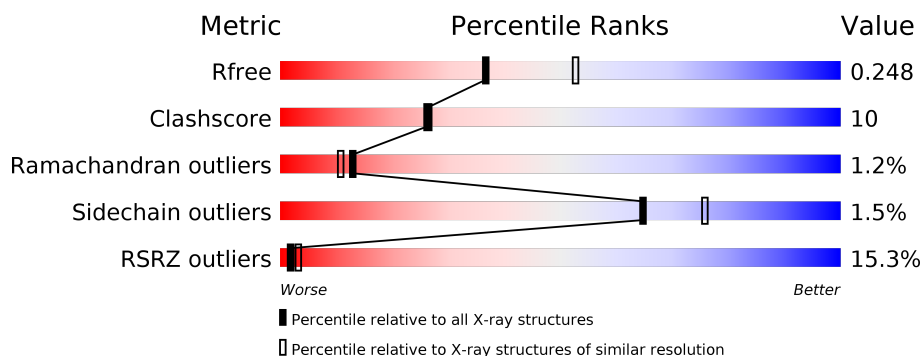
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.34 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	2096 (2.36-2.32)
Clashscore	141614	2193 (2.36-2.32)
Ramachandran outliers	138981	2159 (2.36-2.32)
Sidechain outliers	138945	2160 (2.36-2.32)
RSRZ outliers	127900	2067 (2.36-2.32)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria 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	A	362	<div> <div>13%</div> <div> <div>74%</div> <div>20%</div> <div>• 5%</div> </div> </div>
1	B	362	<div> <div>19%</div> <div> <div>68%</div> <div>25%</div> <div>• 5%</div> </div> </div>
2	C	16	<div> <div>6%</div> <div> <div>63%</div> <div>25%</div> <div>13%</div> </div> </div>
2	P	16	<div> <div>56%</div> <div>25%</div> <div>19%</div> </div>
3	D	19	<div> <div>5%</div> <div> <div>63%</div> <div>21%</div> <div>11%</div> <div>5%</div> </div> </div>
3	T	19	<div> <div>63%</div> <div>26%</div> <div>5%</div> <div>5%</div> </div>

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 7148 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 protein called DNA polymerase IV.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	343	Total	C	N	O	S	0	1	0
			2736	1758	476	494	8			
1	B	343	Total	C	N	O	S	0	0	0
			2730	1753	475	494	8			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	355	GLY	-	EXPRESSION TAG	UNP Q4JB80
A	356	GLY	-	EXPRESSION TAG	UNP Q4JB80
A	357	HIS	-	EXPRESSION TAG	UNP Q4JB80
A	358	HIS	-	EXPRESSION TAG	UNP Q4JB80
A	359	HIS	-	EXPRESSION TAG	UNP Q4JB80
A	360	HIS	-	EXPRESSION TAG	UNP Q4JB80
A	361	HIS	-	EXPRESSION TAG	UNP Q4JB80
A	362	HIS	-	EXPRESSION TAG	UNP Q4JB80
B	355	GLY	-	EXPRESSION TAG	UNP Q4JB80
B	356	GLY	-	EXPRESSION TAG	UNP Q4JB80
B	357	HIS	-	EXPRESSION TAG	UNP Q4JB80
B	358	HIS	-	EXPRESSION TAG	UNP Q4JB80
B	359	HIS	-	EXPRESSION TAG	UNP Q4JB80
B	360	HIS	-	EXPRESSION TAG	UNP Q4JB80
B	361	HIS	-	EXPRESSION TAG	UNP Q4JB80
B	362	HIS	-	EXPRESSION TAG	UNP Q4JB80

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	16	Total	C	N	O	P	0	0	0
			325	154	62	94	15			

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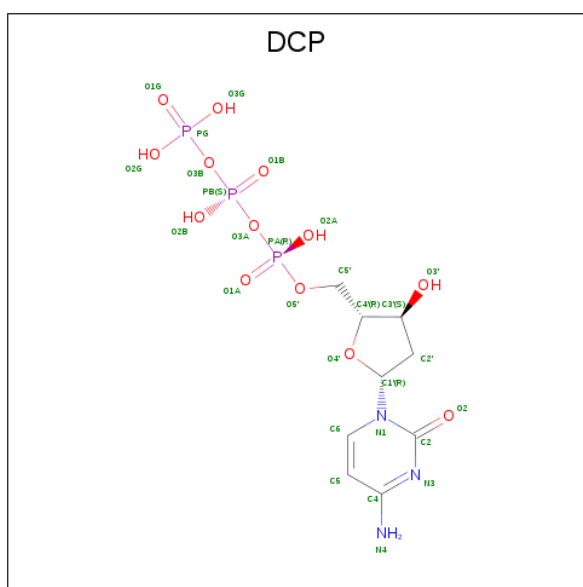
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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	P	16	Total	C	N	O	P	0	0	0
			325	154	62	94	15			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	18	Total	C	N	O	P	0	0	0
			364	173	64	109	18			
3	T	18	Total	C	N	O	P	0	0	0
			364	173	64	109	18			

- Molecule 4 is 2'-DEOXYCYTIDINE-5'-TRIPHOSPHATE (three-letter code: DCP) (formula: C₉H₁₆N₃O₁₃P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			24	9	3	10	2		
4	C	1	Total	C	N	O	P	0	0
			24	9	3	10	2		

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

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

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

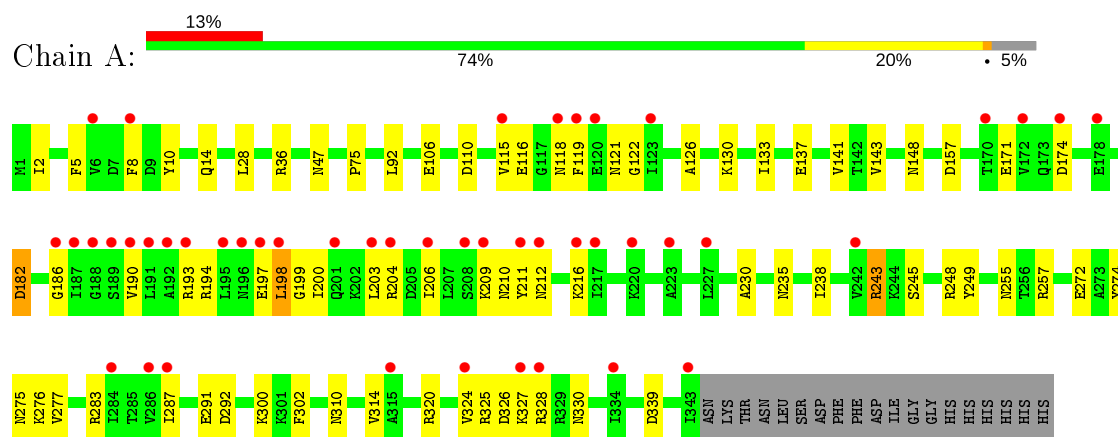
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	114	Total	O	0	0
			114	114		
6	B	77	Total	O	0	0
			77	77		
6	C	18	Total	O	0	0
			18	18		
6	D	13	Total	O	0	0
			13	13		
6	P	17	Total	O	0	0
			17	17		
6	T	13	Total	O	0	0
			13	13		

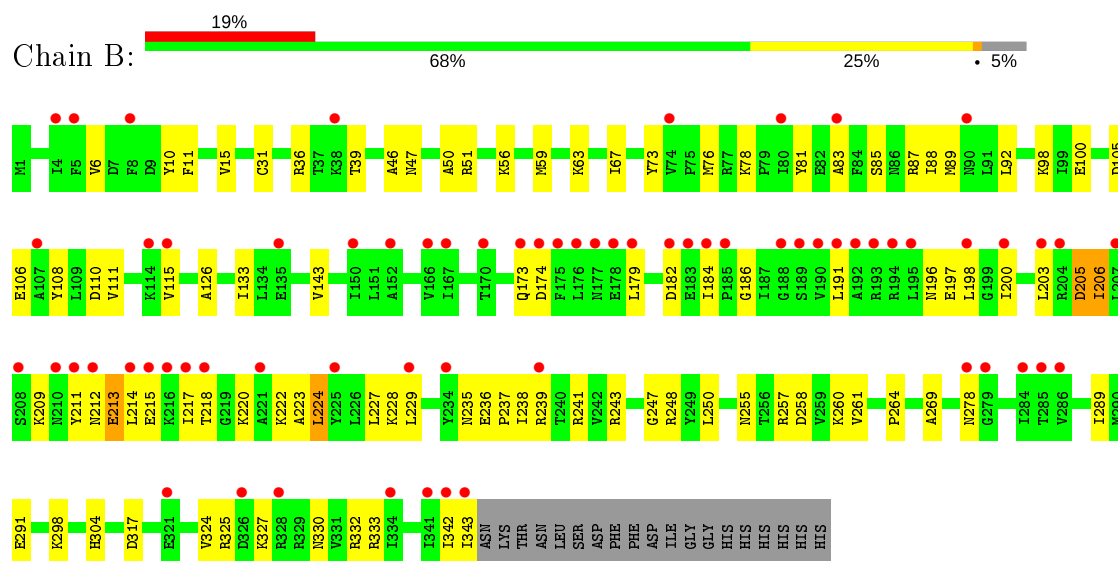
3 Residue-property plots [i](#)

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 polymerase IV

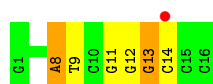


• Molecule 1: DNA polymerase IV

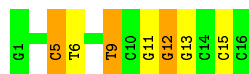


• Molecule 2: DNA (5'-D(*GP*GP*CP*AP*CP*TP*GP*AP*TP*CP*GP*GP*GP*CP*CP*C)-3')

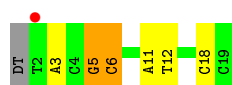




- Molecule 2: DNA (5'-D(*GP*GP*CP*AP*CP*TP*GP*AP*TP*CP*GP*GP*GP*CP*CP*C)-3')



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



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



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	53.08 Å 103.39 Å 112.99 Å 90.00° 101.76° 90.00°	Depositor
Resolution (Å)	29.25 – 2.34 29.25 – 2.34	Depositor EDS
% Data completeness (in resolution range)	94.1 (29.25-2.34) 94.2 (29.25-2.34)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.71 (at 2.34 Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.3_928)	Depositor
R, R_{free}	0.222 , 0.259 0.212 , 0.248	Depositor DCC
R_{free} test set	1893 reflections (3.98%)	wwPDB-VP
Wilson B-factor (Å ²)	31.8	Xtriage
Anisotropy	0.602	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 56.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.027 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	7148	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 25.50 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 3.1373e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: CA, 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	A	0.32	0/2777	0.51	0/3741
1	B	0.31	0/2768	0.52	1/3730 (0.0%)
2	C	0.55	0/364	1.29	4/560 (0.7%)
2	P	0.63	0/364	1.50	6/560 (1.1%)
3	D	0.55	0/406	1.37	2/623 (0.3%)
3	T	0.60	0/406	1.38	4/623 (0.6%)
All	All	0.39	0/7085	0.81	17/9837 (0.2%)

There are no bond length outliers.

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	T	5	DG	O4'-C1'-N9	-13.59	98.49	108.00
3	D	5	DG	O4'-C1'-N9	-11.58	99.89	108.00
3	D	6	DC	O4'-C1'-N1	-8.68	101.92	108.00
2	P	12	DG	O4'-C1'-N9	8.13	113.69	108.00
2	P	13	DG	O4'-C1'-N9	-7.04	103.07	108.00
2	P	9	DT	O4'-C1'-N1	-6.92	103.16	108.00
3	T	6	DC	O4'-C1'-N1	-6.03	103.78	108.00
2	C	14	DC	O4'-C1'-N1	5.95	112.17	108.00
3	T	9	DT	N3-C4-O4	5.87	123.42	119.90
2	C	8	DA	O4'-C1'-N9	-5.84	103.91	108.00
2	P	12	DG	C1'-O4'-C4'	-5.71	104.39	110.10
2	C	13	DG	O4'-C1'-N9	-5.69	104.02	108.00
2	P	5	DC	C1'-O4'-C4'	-5.66	104.44	110.10
3	T	9	DT	C5-C4-O4	-5.64	120.95	124.90
2	C	13	DG	C3'-C2'-C1'	-5.34	96.09	102.50
2	P	13	DG	C3'-C2'-C1'	-5.29	96.15	102.50
1	B	174	ASP	CB-CG-OD2	5.18	122.96	118.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2736	0	2917	54	0
1	B	2730	0	2904	66	0
2	C	325	0	180	7	0
2	P	325	0	180	6	0
3	D	364	0	203	9	0
3	T	364	0	203	3	1
4	A	24	0	12	0	0
4	C	24	0	12	1	0
5	A	2	0	0	0	0
5	B	2	0	0	0	0
6	A	114	0	0	13	0
6	B	77	0	0	8	1
6	C	18	0	0	3	0
6	D	13	0	0	6	0
6	P	17	0	0	2	0
6	T	13	0	0	0	0
All	All	7148	0	6611	137	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 10.

All (137) 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:A:198:LEU:O	6:A:588:HOH:O	1.96	0.82
1:B:325:ARG:NH2	2:C:9:DT:OP1	2.13	0.81
2:C:9:DT:O4	6:C:216:HOH:O	1.99	0.80
1:B:182:ASP:O	6:B:560:HOH:O	2.00	0.79
1:A:36:ARG:NH2	1:A:255:ASN:OD1	2.15	0.75
1:B:317:ASP:OD1	6:B:506:HOH:O	2.05	0.74
1:A:137:GLU:OE2	6:A:613:HOH:O	2.06	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:204:ARG:NH1	1:A:230:ALA:O	2.24	0.71
1:A:272:GLU:OE2	6:A:520:HOH:O	2.09	0.70
1:A:14:GLN:OE1	6:A:610:HOH:O	2.10	0.67
3:D:6:DC:N4	6:D:101:HOH:O	1.99	0.67
1:B:243:ARG:NH2	1:B:278:ASN:O	2.29	0.66
1:B:186:GLY:O	1:B:222:LYS:NZ	2.29	0.65
4:C:101:DCP:O3B	6:C:210:HOH:O	2.13	0.65
3:D:11:DA:O5'	6:D:108:HOH:O	2.15	0.64
1:B:235:ASN:ND2	6:B:567:HOH:O	2.31	0.62
1:B:289:ILE:HB	1:B:333:ARG:HB2	1.81	0.62
1:A:235:ASN:ND2	6:A:592:HOH:O	2.22	0.61
1:B:36:ARG:NH2	1:B:255:ASN:OD1	2.18	0.61
2:C:13:DG:O6	6:C:205:HOH:O	2.16	0.61
1:B:214:LEU:HD22	1:B:217:ILE:HD12	1.83	0.60
1:A:339:ASP:HB2	6:A:502:HOH:O	2.01	0.60
1:A:116:GLU:OE1	1:B:98:LYS:NZ	2.24	0.59
1:A:272:GLU:HA	1:A:275:ASN:HB3	1.82	0.59
2:C:12:DG:H2''	2:C:13:DG:C8	2.37	0.59
1:A:243:ARG:NH2	3:T:8:DC:H5''	2.18	0.58
1:B:332:ARG:NH2	3:D:5:DG:OP1	2.35	0.58
1:A:75:PRO:O	6:A:509:HOH:O	2.18	0.57
2:C:11:DG:H2''	2:C:12:DG:O5'	2.05	0.57
1:B:126:ALA:HB1	1:B:143:VAL:HG11	1.86	0.56
1:A:182:ASP:OD1	1:A:182:ASP:N	2.33	0.56
1:A:126:ALA:HB1	1:A:143:VAL:HG11	1.89	0.55
1:A:5:PHE:HZ	1:A:106:GLU:OE1	1.90	0.55
3:D:18:DC:N4	6:D:110:HOH:O	2.34	0.55
1:B:173:GLN:NE2	6:B:544:HOH:O	2.36	0.54
1:A:325:ARG:HH22	2:P:9:DT:P	2.30	0.54
1:B:78:LYS:HA	1:B:81:TYR:CD1	2.42	0.54
1:A:325:ARG:NH2	2:P:9:DT:OP1	2.40	0.54
1:B:100:GLU:HB3	1:B:108:TYR:HB2	1.89	0.53
1:B:250:LEU:HD13	1:B:269:ALA:HB2	1.90	0.53
1:B:236:GLU:HG2	1:B:237:PRO:HD2	1.89	0.53
1:B:211:TYR:HE1	1:B:227:LEU:HD22	1.74	0.52
1:B:257:ARG:NH2	1:B:327:LYS:HA	2.25	0.52
1:A:110:ASP:HB2	1:A:238:ILE:HG13	1.91	0.52
2:P:11:DG:O3'	6:P:116:HOH:O	2.19	0.52
1:B:218:THR:HB	1:B:222:LYS:HB2	1.91	0.51
1:A:115:VAL:HG11	1:A:122:GLY:HA2	1.92	0.51
1:A:243:ARG:HD3	1:A:245:SER:H	1.76	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:P:5:DC:H2'	2:P:6:DT:H71	1.93	0.51
1:B:247:GLY:O	1:B:248:ARG:NH1	2.34	0.51
1:B:342:ILE:HA	6:B:569:HOH:O	2.10	0.51
1:B:200:ILE:HD11	1:B:209:LYS:HG3	1.92	0.51
2:C:12:DG:H2"	2:C:13:DG:H8	1.76	0.51
1:B:63:LYS:HE2	3:D:3:DA:N7	2.26	0.50
1:B:76:MET:O	6:B:551:HOH:O	2.19	0.50
1:B:6:VAL:HG22	1:B:143:VAL:HG22	1.93	0.50
1:B:237:PRO:HG3	1:B:239:ARG:HH21	1.76	0.50
1:B:258:ASP:OD2	1:B:260:LYS:HD3	2.12	0.50
1:B:11:PHE:O	1:B:15:VAL:HG23	2.12	0.49
1:B:184:ILE:HD11	1:B:203:LEU:HD13	1.93	0.49
1:A:174:ASP:HB2	6:A:539:HOH:O	2.13	0.49
1:B:92:LEU:HD21	1:B:133:ILE:HD11	1.94	0.49
1:A:283:ARG:HH21	1:A:339:ASP:CG	2.16	0.48
2:P:12:DG:H5'	6:P:116:HOH:O	2.13	0.48
3:D:18:DC:OP2	6:D:113:HOH:O	2.20	0.48
1:B:224:LEU:O	1:B:228:LYS:HG2	2.13	0.48
1:A:200:ILE:HD11	1:A:209:LYS:HG2	1.96	0.48
1:A:212:ASN:O	1:A:216:LYS:HG2	2.13	0.48
3:T:9:DT:H2"	3:T:10:DG:O5'	2.13	0.48
1:A:8:PHE:HA	1:A:141:VAL:HG12	1.95	0.48
1:A:193:ARG:O	1:A:197:GLU:HB2	2.13	0.48
1:A:243:ARG:NH1	1:A:245:SER:O	2.46	0.47
1:B:215:GLU:OE2	1:B:223:ALA:HB3	2.13	0.47
1:A:249:TYR:HE2	1:A:287:ILE:HD12	1.79	0.47
1:B:203:LEU:HD21	1:B:229:LEU:HD13	1.97	0.47
1:A:118:ASN:ND2	1:A:121:ASN:OD1	2.48	0.47
1:B:304:HIS:NE2	6:B:530:HOH:O	2.36	0.47
3:D:11:DA:H2'	6:D:108:HOH:O	2.14	0.46
1:B:59:MET:HE1	1:B:67:ILE:HD11	1.96	0.46
1:A:310:ASN:O	1:A:314:VAL:HG23	2.16	0.46
1:B:205:ASP:N	1:B:205:ASP:OD1	2.47	0.45
1:A:326:ASP:O	1:A:328:ARG:N	2.48	0.45
3:D:11:DA:P	6:D:108:HOH:O	2.75	0.45
1:A:92:LEU:HD21	1:A:133:ILE:HD11	1.98	0.44
1:B:206:ILE:HD11	1:B:227:LEU:HA	1.98	0.44
1:A:276:LYS:HA	1:A:276:LYS:HD2	1.81	0.44
1:B:11:PHE:HB2	1:B:105:ASP:OD1	2.17	0.44
1:B:179:LEU:HD23	1:B:203:LEU:HD22	1.99	0.44
3:D:11:DA:H2"	3:D:12:DT:H5"	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:292:ASP:N	1:A:292:ASP:OD2	2.42	0.44
1:B:111:VAL:O	1:B:115:VAL:HB	2.18	0.44
1:B:110:ASP:HB2	1:B:238:ILE:HD12	1.99	0.44
1:B:291:GLU:CD	1:B:330:ASN:HB2	2.37	0.44
1:A:190:VAL:O	1:A:194:ARG:HG2	2.17	0.44
1:A:197:GLU:O	1:A:199:GLY:N	2.45	0.44
1:A:248:ARG:HD2	6:A:520:HOH:O	2.18	0.44
1:B:248:ARG:HA	1:B:248:ARG:HD3	1.86	0.44
1:B:191:LEU:HD21	1:B:218:THR:HG22	2.00	0.43
1:B:56:LYS:HE3	1:B:56:LYS:HB3	1.79	0.43
1:A:257:ARG:HG3	1:A:330:ASN:OD1	2.18	0.43
1:A:203:LEU:HD12	1:A:206:ILE:HD12	2.01	0.43
1:A:272:GLU:N	1:A:272:GLU:OE1	2.47	0.43
1:A:28:LEU:HB3	1:A:47:ASN:ND2	2.32	0.43
1:B:46:ALA:HB1	1:B:50:ALA:HB3	2.01	0.43
1:B:224:LEU:HG	1:B:228:LYS:NZ	2.33	0.43
1:A:248:ARG:HG3	1:A:272:GLU:HG2	2.01	0.43
1:A:171:GLU:OE2	6:A:596:HOH:O	2.21	0.43
1:B:100:GLU:HG3	1:B:241:ARG:HG2	2.00	0.43
1:B:211:TYR:C	1:B:213:GLU:H	2.21	0.43
1:A:291:GLU:HG2	6:A:549:HOH:O	2.17	0.43
1:B:298:LYS:HD2	2:C:8:DA:OP1	2.19	0.43
1:A:186:GLY:HA2	2:P:15:DC:H1'	2.00	0.42
1:A:274:TYR:O	1:A:277:VAL:HG12	2.19	0.42
1:B:83:ALA:O	1:B:87:ARG:HG3	2.19	0.42
1:A:212:ASN:OD1	1:A:216:LYS:HE3	2.19	0.42
1:B:243:ARG:HG3	1:B:243:ARG:O	2.18	0.42
1:B:343:ILE:HG12	6:B:569:HOH:O	2.20	0.42
1:A:148:ASN:HB2	6:A:593:HOH:O	2.19	0.42
1:B:47:ASN:O	1:B:51:ARG:HG3	2.20	0.42
1:B:196:ASN:C	1:B:198:LEU:H	2.23	0.42
1:B:215:GLU:CD	1:B:220:LYS:HA	2.40	0.42
3:T:9:DT:H2'	3:T:10:DG:C8	2.55	0.42
1:A:300:LYS:HE3	1:A:302:PHE:CE2	2.55	0.42
1:B:191:LEU:HD11	1:B:218:THR:HG22	2.02	0.42
1:A:320:ARG:O	1:A:324:VAL:HG23	2.20	0.41
1:B:261:VAL:O	1:B:264:PRO:HD2	2.20	0.41
1:B:220:LYS:HB3	1:B:220:LYS:HE3	1.88	0.41
1:B:198:LEU:HD23	1:B:198:LEU:HA	1.88	0.41
1:A:157:ASP:OD2	6:A:594:HOH:O	2.22	0.41
1:B:332:ARG:HG2	1:B:333:ARG:HG3	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:130:LYS:NZ	1:A:141:VAL:O	2.53	0.41
1:A:248:ARG:HD3	1:A:248:ARG:HA	1.86	0.41
1:B:88:ILE:O	1:B:92:LEU:HG	2.21	0.41
1:B:85:SER:O	1:B:89:MET:HG2	2.21	0.41
1:A:2:ILE:HD13	1:A:119:PHE:HA	2.03	0.40
1:B:31:CYS:SG	1:B:73:TYR:HD2	2.44	0.40
1:B:76:MET:HG3	1:B:81:TYR:HE1	1.86	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:T:2:DT:O4	6:B:515:HOH:O[1_554]	2.12	0.08

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	
1	A	342/362 (94%)	315 (92%)	22 (6%)	5 (2%)	10	7
1	B	341/362 (94%)	318 (93%)	20 (6%)	3 (1%)	17	17
All	All	683/724 (94%)	633 (93%)	42 (6%)	8 (1%)	13	11

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	10	TYR
1	A	211	TYR
1	A	210	ASN
1	A	198	LEU
1	A	327	LYS
1	B	10	TYR

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Mol	Chain	Res	Type
1	B	197	GLU
1	B	212	ASN

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	
1	A	299/315 (95%)	297 (99%)	2 (1%)	84	90
1	B	298/315 (95%)	291 (98%)	7 (2%)	50	61
All	All	597/630 (95%)	588 (98%)	9 (2%)	65	76

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	182	ASP
1	A	243	ARG
1	B	39	THR
1	B	106	GLU
1	B	205	ASP
1	B	206	ILE
1	B	213	GLU
1	B	224	LEU
1	B	324	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 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	DCP	C	101	5	21,25,29	1.58	4 (19%)	28,38,45	1.40	3 (10%)
4	DCP	A	401	5	21,25,29	1.67	5 (23%)	28,38,45	1.51	6 (21%)

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
4	DCP	C	101	5	-	3/13/28/34	0/2/2/2
4	DCP	A	401	5	-	1/13/28/34	0/2/2/2

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	401	DCP	C3'-C4'	-3.77	1.42	1.53
4	C	101	DCP	C3'-C4'	-3.35	1.43	1.53
4	C	101	DCP	C4-N4	3.29	1.44	1.35
4	A	401	DCP	C4-N4	3.27	1.44	1.35
4	A	401	DCP	O3'-C3'	-2.41	1.38	1.43
4	C	101	DCP	C2-N3	-2.28	1.33	1.38
4	A	401	DCP	C2'-C1'	-2.25	1.46	1.52
4	C	101	DCP	C2'-C1'	-2.24	1.46	1.52
4	A	401	DCP	C2-N3	-2.20	1.33	1.38

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	101	DCP	O2B-PB-O1B	-3.52	96.92	110.68
4	A	401	DCP	O2B-PB-O1B	-3.36	97.51	110.68
4	A	401	DCP	C2-N3-C4	3.30	119.69	116.34
4	C	101	DCP	C2-N3-C4	3.07	119.45	116.34
4	A	401	DCP	O2A-PA-O1A	-2.73	98.75	112.24
4	A	401	DCP	C2'-C1'-N1	-2.65	108.17	114.27
4	C	101	DCP	O2A-PA-O1A	-2.59	99.42	112.24
4	A	401	DCP	PA-O3A-PB	-2.32	124.86	132.83
4	A	401	DCP	O5'-C5'-C4'	2.01	115.92	108.99

There are no chirality outliers.

All (4) torsion outliers are listed below:

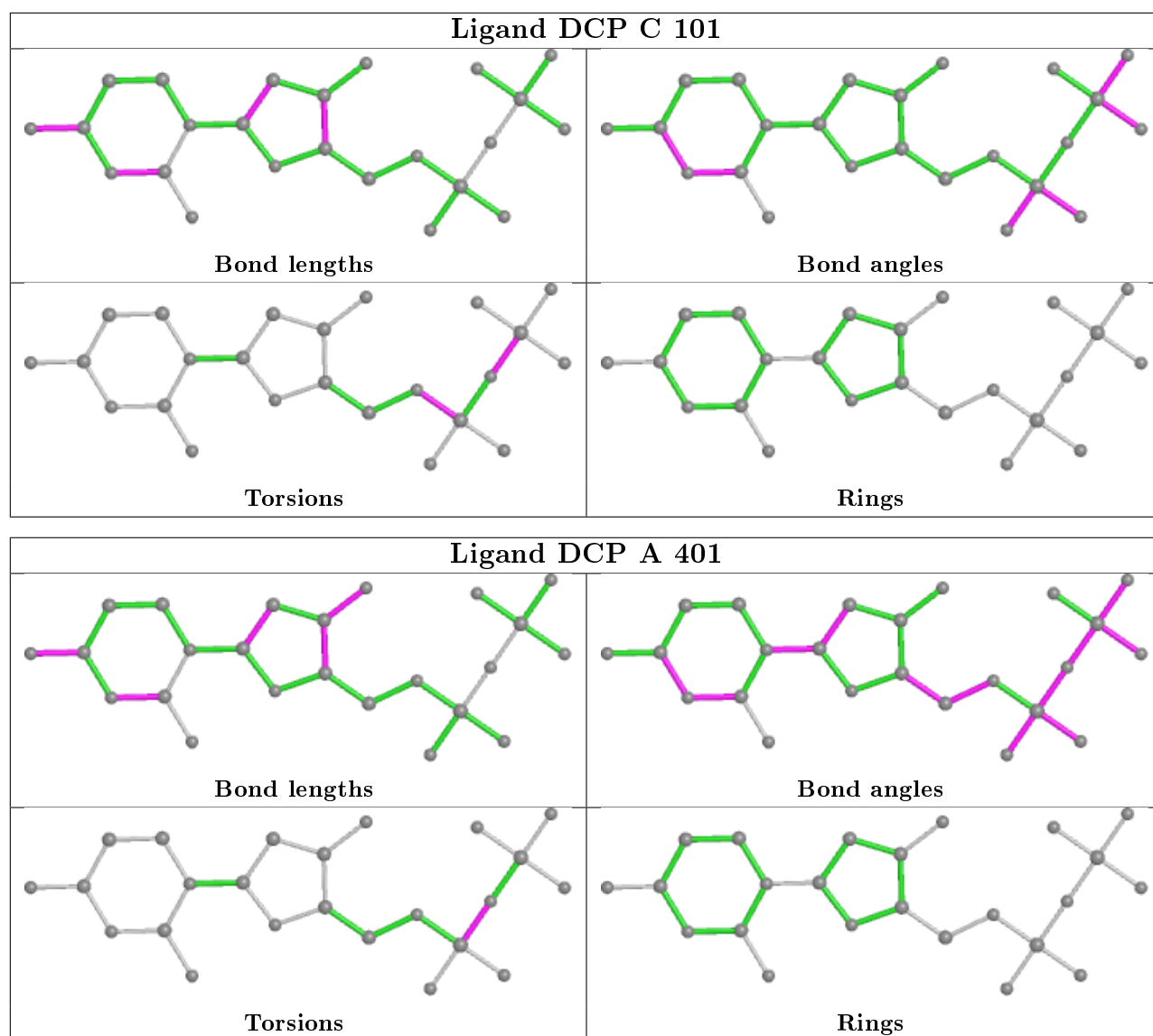
Mol	Chain	Res	Type	Atoms
4	A	401	DCP	PB-O3A-PA-O1A
4	C	101	DCP	C5'-O5'-PA-O1A
4	C	101	DCP	PA-O3A-PB-O1B
4	C	101	DCP	PA-O3A-PB-O3B

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	101	DCP	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	A	343/362 (94%)	0.93	46 (13%) 3 5	19, 38, 89, 112	0
1	B	343/362 (94%)	1.24	67 (19%) 1 2	25, 43, 91, 118	0
2	C	16/16 (100%)	0.69	1 (6%) 20 28	46, 59, 70, 71	0
2	P	16/16 (100%)	0.43	0 100 100	37, 48, 57, 72	0
3	D	18/19 (94%)	0.81	1 (5%) 24 33	27, 59, 72, 75	0
3	T	18/19 (94%)	0.49	0 100 100	32, 50, 64, 79	0
All	All	754/794 (94%)	1.04	115 (15%) 2 3	19, 42, 90, 118	0

All (115) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	211	TYR	9.6
1	B	217	ILE	8.3
1	B	343	ILE	7.3
1	B	215	GLU	6.5
1	B	191	LEU	6.5
1	B	214	LEU	6.4
1	B	174	ASP	6.1
1	A	343	ILE	6.1
1	B	192	ALA	6.0
1	B	170	THR	5.4
1	B	189	SER	5.4
1	B	204	ARG	5.2
1	A	212	ASN	4.8
1	A	324	VAL	4.5
1	A	328	ARG	4.5
1	A	190	VAL	4.5
1	A	196	ASN	4.4
1	B	218	THR	4.2
1	B	173	GLN	4.1

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Mol	Chain	Res	Type	RSRZ
1	B	212	ASN	4.0
1	B	286	VAL	4.0
1	A	189	SER	3.9
1	B	200	ILE	3.8
1	A	242	VAL	3.8
1	A	192	ALA	3.7
1	B	80	ILE	3.7
1	A	197	GLU	3.6
1	B	194	ARG	3.6
1	B	74	VAL	3.6
1	B	178	GLU	3.6
1	B	195	LEU	3.6
1	A	188	GLY	3.6
1	A	172	VAL	3.6
1	B	207	LEU	3.5
1	A	174	ASP	3.5
1	B	177	ASN	3.4
1	A	8	PHE	3.4
1	B	135	GLU	3.4
1	B	208	SER	3.4
1	B	234	TYR	3.4
1	B	328	ARG	3.3
1	B	239	ARG	3.3
1	A	120	GLU	3.3
1	A	198	LEU	3.2
1	B	166	VAL	3.2
1	B	342	ILE	3.1
1	B	190	VAL	3.1
1	A	170	THR	3.1
1	A	191	LEU	3.1
1	A	178	GLU	3.0
1	A	209	LYS	3.0
1	A	211	TYR	2.9
1	A	193	ARG	2.9
1	A	208	SER	2.9
1	B	284	ILE	2.9
1	A	217	ILE	2.8
1	A	284	ILE	2.8
1	B	198	LEU	2.8
1	B	203	LEU	2.8
1	B	285	THR	2.8
1	A	286	VAL	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	195	LEU	2.8
1	A	223	ALA	2.7
1	B	185	PRO	2.7
1	A	201	GLN	2.7
1	B	152	ALA	2.7
1	B	278	ASN	2.7
1	A	334	ILE	2.7
1	B	5	PHE	2.7
1	A	187	ILE	2.7
1	A	6	VAL	2.7
1	B	115	VAL	2.7
1	B	90	ASN	2.6
1	B	184	ILE	2.6
1	B	114	LYS	2.6
1	A	287	ILE	2.6
1	B	183	GLU	2.6
1	B	38	LYS	2.6
1	B	279	GLY	2.5
1	B	83	ALA	2.5
1	B	182	ASP	2.5
1	B	341	ILE	2.5
1	B	193	ARG	2.5
1	A	327	LYS	2.5
1	B	334	ILE	2.4
1	B	229	LEU	2.4
1	A	220	LYS	2.4
1	B	326	ASP	2.3
1	A	123	ILE	2.3
1	A	206	ILE	2.3
3	D	2	DT	2.3
1	A	203	LEU	2.3
1	B	175	PHE	2.3
1	B	210	ASN	2.3
1	B	150	ILE	2.3
1	A	227	LEU	2.2
1	A	186	GLY	2.2
1	A	204	ARG	2.2
1	B	167	ILE	2.2
1	B	221	ALA	2.2
1	A	216	LYS	2.2
1	B	8	PHE	2.1
1	B	321	GLU	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	115	VAL	2.1
1	A	315	ALA	2.1
1	B	179	LEU	2.1
2	C	14	DC	2.1
1	B	188	GLY	2.1
1	B	107	ALA	2.1
1	B	4	ILE	2.1
1	A	119	PHE	2.1
1	B	225	TYR	2.1
1	B	216	LYS	2.0
1	A	118	ASN	2.0
1	B	176	LEU	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

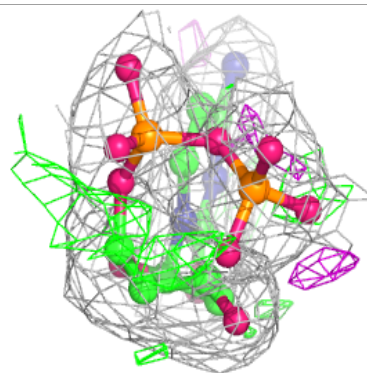
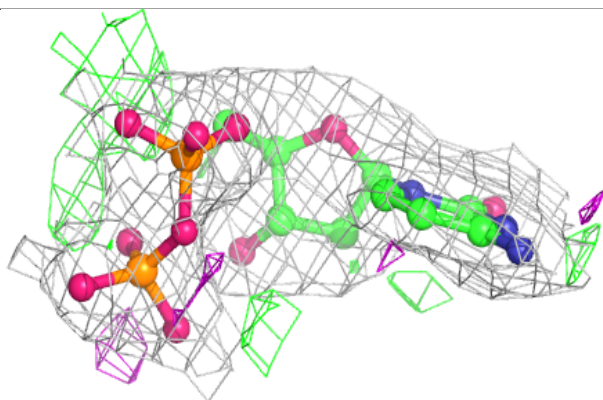
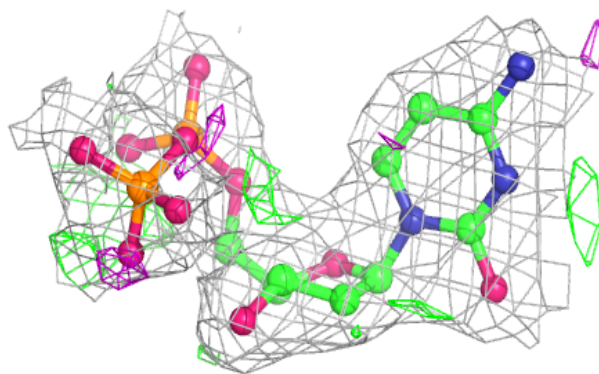
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	CA	A	403	1/1	0.89	0.29	36,36,36,36	0
5	CA	B	402	1/1	0.92	0.29	35,35,35,35	0
4	DCP	A	401	24/28	0.95	0.18	7,21,34,143	0
4	DCP	C	101	24/28	0.95	0.16	22,32,45,53	0
5	CA	B	401	1/1	0.96	0.29	47,47,47,47	0
5	CA	A	402	1/1	0.97	0.30	30,30,30,30	0

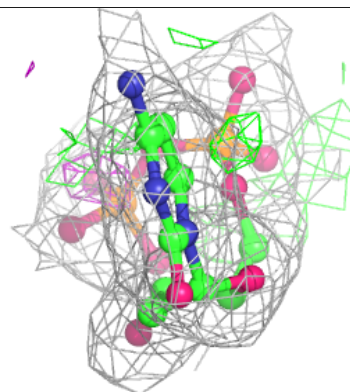
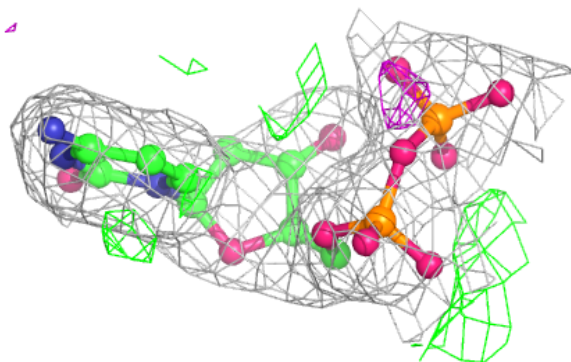
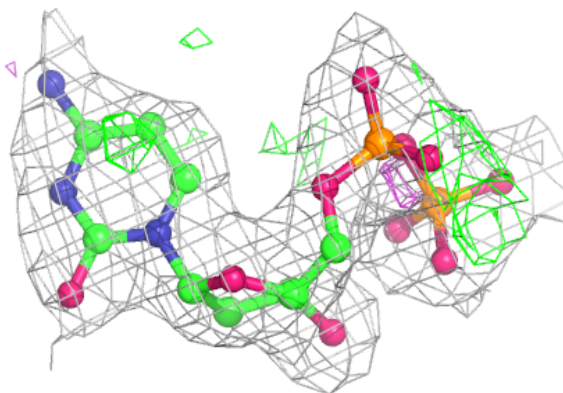
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around DCP A 401:

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

**Electron density around DCP C 101:**

$2mF_o-DF_c$ (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.