



# Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 10:47 pm BST

PDB ID : 2ZXU  
Title : Crystal structure of tRNA modification enzyme MiaA in the complex with tRNA(Phe) and DMASPP  
Authors : Sakai, J.; Yao, M.; Chimnaronk, S.; Tanaka, I.  
Deposited on : 2009-01-07  
Resolution : 2.75 Å(reported)

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

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
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

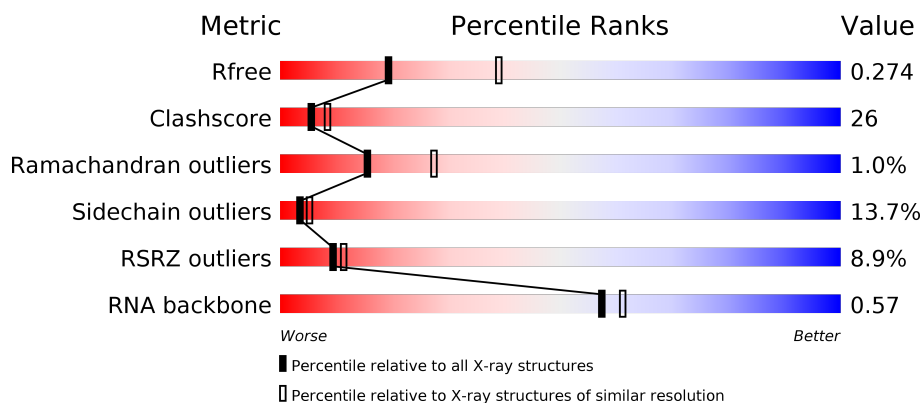
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.75 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	1235 (2.78-2.74)
Clashscore	141614	1277 (2.78-2.74)
Ramachandran outliers	138981	1257 (2.78-2.74)
Sidechain outliers	138945	1257 (2.78-2.74)
RSRZ outliers	127900	1207 (2.78-2.74)
RNA backbone	3102	1060 (3.02-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	316	<div> <div>2%</div> <div>66%</div> <div>23%</div> <div>8%</div> <div>.</div> </div>
1	B	316	<div> <div>9%</div> <div>48%</div> <div>39%</div> <div>9%</div> <div>..</div> </div>
2	C	76	<div> <div>3%</div> <div>49%</div> <div>21%</div> <div>21%</div> <div>7%</div> <div>.</div> </div>
2	D	76	<div> <div>41%</div> <div>29%</div> <div>41%</div> <div>18%</div> <div>.</div> <div>9%</div> </div>

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 8081 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.

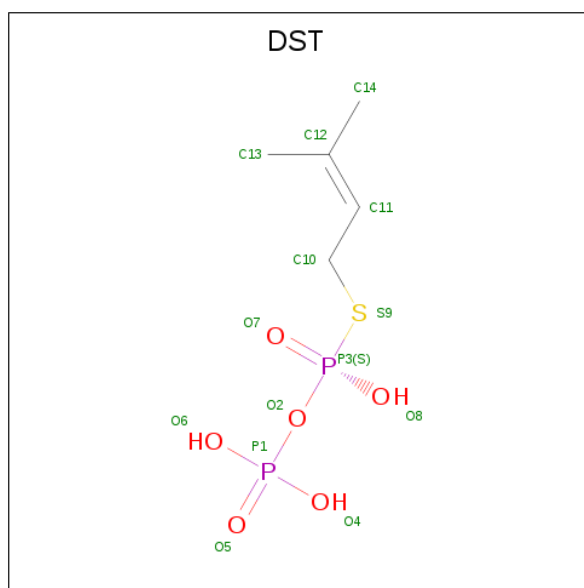
- Molecule 1 is a protein called tRNA delta(2)-isopentenylpyrophosphate transferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	306	Total	C	N	O	S	0	0	0
			2403	1523	428	443	9			
1	B	305	Total	C	N	O	S	0	0	0
			2397	1518	427	443	9			

- Molecule 2 is a RNA chain called tRNA(Phe).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	74	Total	C	N	O	P	0	0	0
			1577	704	282	518	73			
2	D	69	Total	C	N	O	P	0	0	0
			1469	656	261	484	68			

- Molecule 3 is DIMETHYLALLYL S-THIOLODIPHOSPHATE (three-letter code: DST) (formula:  $C_5H_{12}O_6P_2S$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	O	P	S	0	0
			14	5	6	2	1		
3	B	1	Total	C	O	P	S	0	0
			14	5	6	2	1		

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Mg	0	0
			1	1		
4	A	1	Total	Mg	0	0
			1	1		
4	D	1	Total	Mg	0	0
			1	1		
4	C	7	Total	Mg	0	0
			7	7		

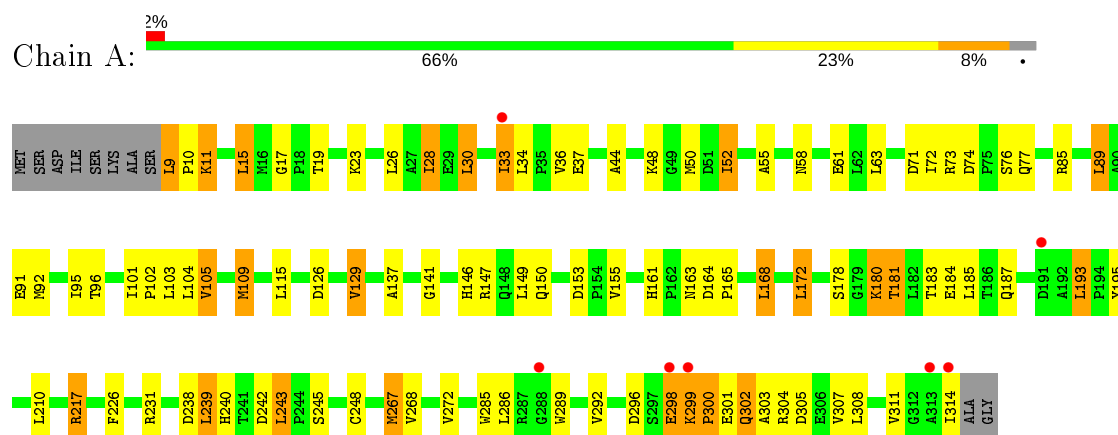
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	67	Total	O	0	0
			67	67		
5	B	42	Total	O	0	0
			42	42		
5	C	59	Total	O	0	0
			59	59		
5	D	29	Total	O	0	0
			29	29		

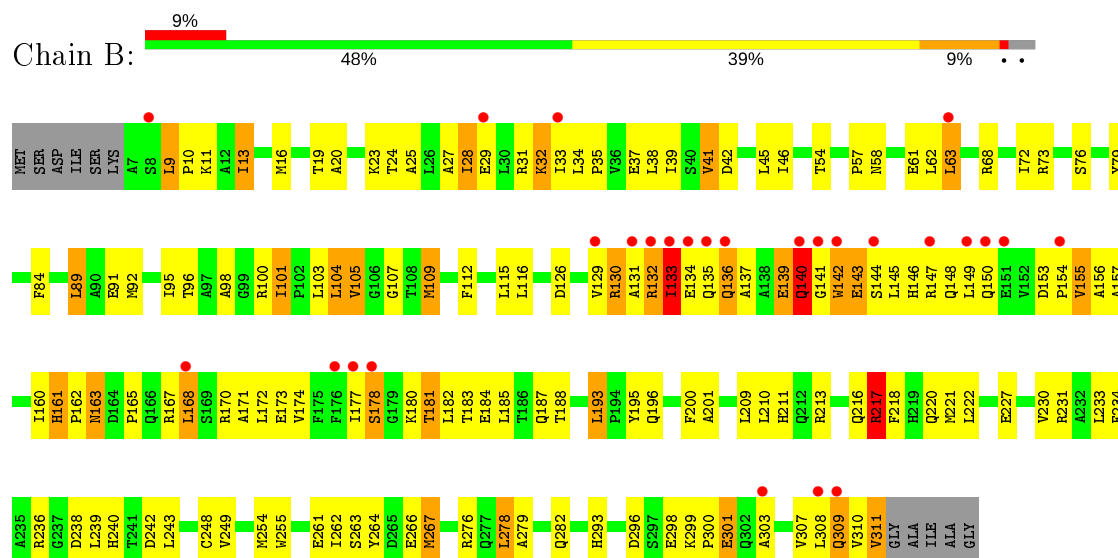
### 3 Residue-property plots [i](#)

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: tRNA delta(2)-isopentenylpyrophosphate transferase

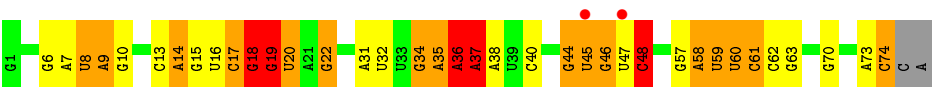


- Molecule 1: tRNA delta(2)-isopentenylpyrophosphate transferase

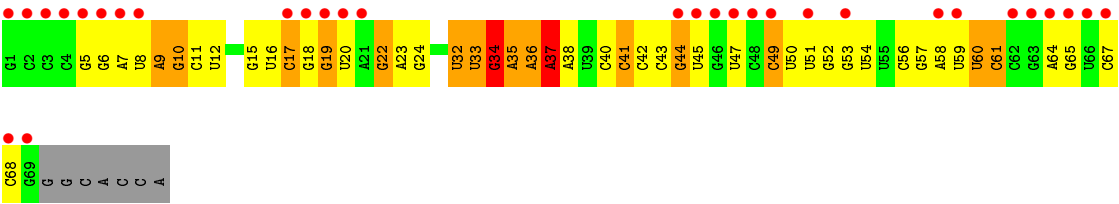


- Molecule 2: tRNA(Phe)





● Molecule 2: tRNA(Phe)



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	88.30 Å 90.00 Å 150.50 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.75 48.32 – 2.75	Depositor EDS
% Data completeness (in resolution range)	99.8 (20.00-2.75) 99.7 (48.32-2.75)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	0.05	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	5.70 (at 2.77 Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.237 , 0.281 0.231 , 0.274	Depositor DCC
$R_{free}$ test set	2223 reflections (6.99%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	39.0	Xtriage
Anisotropy	0.546	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 49.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.012 for k,h,-l	Xtriage
$F_o, F_c$ correlation	0.90	EDS
Total number of atoms	8081	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.04% of the height of the origin peak. No significant pseudotranslation is detected.*

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

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality

### 5.1 Standard geometry

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

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	0.39	0/2454	0.65	2/3329 (0.1%)
1	B	0.37	0/2448	0.66	1/3321 (0.0%)
2	C	0.51	0/1762	0.94	8/2746 (0.3%)
2	D	0.43	0/1641	0.86	3/2557 (0.1%)
All	All	0.42	0/8305	0.77	14/11953 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
2	C	0	2
2	D	0	3
All	All	0	5

There are no bond length outliers.

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	217	ARG	NE-CZ-NH2	-6.90	116.85	120.30
2	C	19	G	N9-C1'-C2'	6.78	122.81	114.00
2	D	34	G	N9-C1'-C2'	6.66	122.65	114.00
2	C	18	G	N9-C1'-C2'	6.53	122.49	114.00
2	C	34	G	N9-C1'-C2'	5.91	121.68	114.00
2	C	46	G	N9-C1'-C2'	5.91	121.68	114.00
2	D	35	A	N9-C1'-C2'	5.77	121.50	114.00
2	D	37	A	N9-C1'-C2'	5.50	121.16	114.00
2	C	36	A	N9-C1'-C2'	5.45	121.09	114.00
1	A	217	ARG	NE-CZ-NH2	5.30	122.95	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	217	ARG	NE-CZ-NH1	-5.24	117.68	120.30
2	C	48	C	N1-C1'-C2'	5.19	120.75	114.00
2	C	37	A	C4'-C3'-C2'	5.13	107.73	102.60
2	C	22	G	N9-C1'-C2'	5.02	120.53	114.00

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	C	19	G	Sidechain
2	C	8	U	Sidechain
2	D	22	G	Sidechain
2	D	34	G	Sidechain
2	D	37	A	Sidechain

## 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	2403	0	2407	102	0
1	B	2397	0	2398	193	0
2	C	1577	0	800	34	0
2	D	1469	0	745	47	0
3	A	14	0	10	3	0
3	B	14	0	10	5	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	7	0	0	0	0
4	D	1	0	0	0	0
5	A	67	0	0	3	0
5	B	42	0	0	3	0
5	C	59	0	0	1	0
5	D	29	0	0	0	0
All	All	8081	0	6370	368	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 26.

All (368) 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:B:129:VAL:O	1:B:133:ILE:HD11	1.60	1.01
1:A:101:ILE:HD11	1:A:314:ILE:HG21	1.43	0.99
1:B:136:GLN:HE22	1:B:145:LEU:HD11	1.24	0.99
1:B:136:GLN:HE21	1:B:136:GLN:C	1.68	0.97
1:A:52:ILE:HD11	1:A:226:PHE:HB2	1.48	0.94
2:C:13:C:H2'	2:C:14:A:H5''	1.49	0.94
1:B:23:LYS:HG2	3:B:317:DST:O6	1.72	0.89
1:A:33:ILE:O	1:A:33:ILE:HG13	1.73	0.87
1:B:37:GLU:HB3	1:B:95:ILE:HD13	1.57	0.87
1:A:193:LEU:HD13	1:A:195:TYR:HB2	1.57	0.85
1:B:9:LEU:HB2	1:B:10:PRO:HD2	1.59	0.85
2:C:13:C:C2'	2:C:14:A:H5''	2.06	0.84
1:B:161:HIS:HD2	1:B:163:ASN:OD1	1.60	0.83
1:A:307:VAL:O	1:A:311:VAL:HG12	1.79	0.82
1:B:147:ARG:HB3	1:B:147:ARG:HH11	1.45	0.82
2:C:18:G:O2'	2:C:57:G:N2	2.13	0.82
1:A:73:ARG:HH11	1:A:73:ARG:HG3	1.44	0.81
1:B:133:ILE:N	1:B:133:ILE:HD12	1.95	0.81
2:D:18:G:H2'	2:D:57:G:N2	1.96	0.81
1:B:181:THR:O	1:B:185:LEU:HD23	1.82	0.80
1:B:147:ARG:HB3	1:B:147:ARG:NH1	1.98	0.79
1:B:133:ILE:HD12	1:B:133:ILE:H	1.47	0.79
1:B:149:LEU:HD13	1:B:172:LEU:HD23	1.63	0.78
1:B:298:GLU:O	1:B:299:LYS:HD2	1.83	0.78
1:B:137:ALA:HA	1:B:141:GLY:HA3	1.68	0.76
1:B:136:GLN:NE2	1:B:145:LEU:HD11	2.00	0.76
1:A:311:VAL:HG22	1:A:314:ILE:HD11	1.68	0.76
1:B:27:ALA:HB1	1:B:38:LEU:HD11	1.67	0.76
1:B:41:VAL:HG11	1:B:112:PHE:CZ	2.20	0.76
1:A:161:HIS:HD2	1:A:163:ASN:H	1.34	0.75
1:A:183:THR:HG22	5:A:347:HOH:O	1.86	0.75
1:A:9:LEU:N	1:A:10:PRO:HD3	2.02	0.75
1:A:85:ARG:HG3	1:A:115:LEU:HA	1.69	0.74
1:A:155:VAL:HB	1:A:183:THR:HG21	1.70	0.73
2:C:9:A:H2	2:C:44:G:HO2'	1.36	0.73
1:B:147:ARG:HA	1:B:150:GLN:HG2	1.69	0.72
1:B:16:MET:HB3	1:B:109:MET:HE2	1.70	0.72
1:B:130:ARG:HA	1:B:133:ILE:HD11	1.71	0.71
1:A:299:LYS:C	1:A:301:GLU:H	1.93	0.71
1:A:92:MET:O	1:A:96:THR:HG23	1.91	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:248:CYS:HA	2:D:37:A:O2'	1.91	0.70
1:A:302:GLN:H	1:A:302:GLN:NE2	1.90	0.70
1:B:149:LEU:HB2	1:B:172:LEU:HD21	1.73	0.69
1:B:261:GLU:O	1:B:262:ILE:HD12	1.92	0.69
2:D:5:G:H2'	2:D:6:G:H8	1.56	0.69
1:B:132:ARG:O	1:B:136:GLN:HG3	1.93	0.69
1:B:193:LEU:HD13	1:B:195:TYR:HB2	1.74	0.69
2:D:18:G:H2'	2:D:57:G:H22	1.56	0.68
2:D:60:U:H5'	2:D:61:C:OP2	1.93	0.68
1:B:130:ARG:HA	1:B:133:ILE:CD1	2.23	0.68
1:B:58:ASN:OD1	1:B:61:GLU:HG3	1.93	0.68
1:B:299:LYS:N	1:B:300:PRO:HD3	2.07	0.68
1:B:126:ASP:CG	1:B:129:VAL:HG12	2.14	0.68
1:B:126:ASP:O	1:B:130:ARG:HB2	1.94	0.68
2:C:20:U:OP1	5:C:111:HOH:O	2.11	0.68
1:B:298:GLU:C	1:B:300:PRO:HD3	2.13	0.67
1:A:299:LYS:O	1:A:301:GLU:N	2.26	0.67
1:A:50:MET:HE1	1:A:245:SER:HB2	1.76	0.67
1:B:68:ARG:HD3	1:B:91:GLU:HG3	1.77	0.67
1:B:103:LEU:CD2	1:B:105:VAL:HG22	2.24	0.67
2:C:74:C:H6	2:C:74:C:H5'	1.60	0.66
1:A:181:THR:HG22	1:A:184:GLU:H	1.58	0.66
1:B:46:ILE:HG23	1:B:72:ILE:CG1	2.26	0.66
1:B:249:VAL:HG11	1:B:278:LEU:HD12	1.77	0.66
1:B:298:GLU:C	1:B:299:LYS:HD2	2.16	0.66
1:B:46:ILE:HG22	1:B:73:ARG:HB2	1.76	0.66
1:B:181:THR:HG22	1:B:184:GLU:H	1.59	0.66
1:B:32:LYS:HE2	1:B:32:LYS:N	2.10	0.66
1:B:155:VAL:HG12	1:B:183:THR:OG1	1.96	0.65
2:D:17:C:C5'	2:D:18:G:H5''	2.26	0.65
1:A:11:LYS:HD3	1:A:11:LYS:N	2.11	0.65
1:B:133:ILE:HA	1:B:136:GLN:OE1	1.96	0.65
1:B:142:TRP:HE3	1:B:168:LEU:HD13	1.61	0.65
1:B:147:ARG:CB	1:B:147:ARG:HH11	2.07	0.65
1:B:156:ALA:O	1:B:160:ILE:HG12	1.96	0.65
1:A:193:LEU:CD1	1:A:195:TYR:HB2	2.25	0.65
1:B:146:HIS:CD2	1:B:162:PRO:HG3	2.32	0.65
1:A:161:HIS:CD2	1:A:163:ASN:H	2.16	0.64
2:D:51:U:H2'	2:D:52:G:H8	1.61	0.64
1:B:126:ASP:HB3	1:B:129:VAL:CG1	2.27	0.64
2:C:18:G:HO2'	2:C:57:G:H22	1.44	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:139:GLU:OE2	1:B:140:GLN:HB3	1.97	0.64
1:B:137:ALA:HA	1:B:141:GLY:CA	2.26	0.64
2:D:17:C:H5'	2:D:18:G:H5''	1.80	0.64
1:B:161:HIS:CD2	1:B:163:ASN:OD1	2.46	0.63
2:D:43:C:H2'	2:D:44:G:H5''	1.81	0.63
2:D:5:G:H2'	2:D:6:G:C8	2.34	0.63
1:A:55:ALA:CB	1:A:217:ARG:NH1	2.62	0.63
1:B:296:ASP:HB3	1:B:299:LYS:HB2	1.81	0.62
2:D:6:G:H2'	2:D:7:A:O4'	1.99	0.62
1:B:216:GLN:O	1:B:220:GLN:HG3	1.99	0.61
1:B:240:HIS:CE1	1:B:242:ASP:HB2	2.35	0.61
1:B:136:GLN:HE21	1:B:137:ALA:N	1.97	0.61
2:D:33:U:HO2'	2:D:34:G:H21	1.47	0.61
1:A:267:MET:HE1	1:A:268:VAL:HA	1.81	0.61
2:D:32:U:O2'	2:D:36:A:N1	2.34	0.61
1:B:153:ASP:OD1	1:B:181:THR:HG23	2.01	0.60
2:C:73:A:H2'	2:C:74:C:C5'	2.31	0.60
2:D:51:U:H2'	2:D:52:G:C8	2.36	0.60
1:B:46:ILE:HG23	1:B:72:ILE:HG12	1.83	0.59
1:A:52:ILE:HD11	1:A:226:PHE:CB	2.27	0.59
1:B:134:GLU:O	1:B:137:ALA:HB3	2.02	0.59
2:D:37:A:H8	2:D:37:A:OP1	1.85	0.59
2:C:62:C:O2'	2:C:63:G:H5'	2.03	0.59
2:C:60:U:H3'	2:C:61:C:C5	2.38	0.58
1:A:89:LEU:HA	1:A:92:MET:HE3	1.85	0.58
1:B:181:THR:O	1:B:185:LEU:CD2	2.52	0.58
1:A:33:ILE:O	1:A:34:LEU:HD12	2.04	0.58
1:B:139:GLU:O	1:B:140:GLN:C	2.40	0.58
2:D:49:C:H6	2:D:49:C:O5'	1.87	0.58
1:B:19:THR:OG1	3:B:317:DST:H142	2.04	0.57
2:C:73:A:H2'	2:C:74:C:H5'	1.86	0.57
1:A:15:LEU:HB3	1:A:105:VAL:HG13	1.86	0.57
1:A:73:ARG:NH1	1:A:73:ARG:HG3	2.15	0.57
1:B:130:ARG:CA	1:B:133:ILE:HD11	2.35	0.57
1:B:79:TYR:OH	1:B:84:PHE:HA	2.05	0.57
1:B:173:GLU:O	1:B:177:ILE:HG22	2.05	0.56
1:B:155:VAL:HG11	1:B:183:THR:CG2	2.34	0.56
1:B:227:GLU:HG2	1:B:231:ARG:NH1	2.21	0.56
1:B:16:MET:HE1	1:B:109:MET:HG3	1.86	0.56
1:A:37:GLU:HB2	1:A:95:ILE:HD13	1.88	0.56
1:B:149:LEU:HB2	1:B:172:LEU:CD2	2.35	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:149:LEU:HD13	1:A:172:LEU:HD13	1.86	0.56
1:A:48:LYS:HG2	1:A:71:ASP:HB3	1.86	0.56
1:A:76:SER:HA	1:A:243:LEU:HD21	1.86	0.56
1:B:13:ILE:HG23	1:B:101:ILE:HD11	1.87	0.56
1:B:34:LEU:CD1	1:B:35:PRO:HD2	2.35	0.56
2:C:60:U:H3'	2:C:61:C:H5	1.71	0.56
2:D:6:G:O2'	2:D:7:A:H5'	2.06	0.56
1:A:301:GLU:C	1:A:303:ALA:H	2.09	0.55
1:B:218:PHE:HA	1:B:221:MET:HE2	1.88	0.55
2:D:33:U:O2'	2:D:34:G:N2	2.37	0.55
1:B:163:ASN:N	1:B:163:ASN:OD1	2.38	0.55
1:A:161:HIS:HD2	1:A:163:ASN:N	2.01	0.55
1:B:222:LEU:HD21	1:B:267:MET:HE1	1.89	0.55
1:B:109:MET:HE3	1:B:109:MET:HA	1.88	0.55
2:D:58:A:HO2'	2:D:60:U:H5	1.54	0.55
1:A:298:GLU:O	1:A:300:PRO:N	2.40	0.55
2:C:48:C:H2'	2:C:59:U:H1'	1.89	0.55
1:B:103:LEU:HD22	1:B:105:VAL:HG22	1.89	0.55
1:B:38:LEU:HD12	1:B:105:VAL:HG21	1.87	0.55
1:A:296:ASP:HB3	1:A:300:PRO:HB2	1.90	0.54
2:C:20:U:O5'	2:C:20:U:H6	1.90	0.54
1:B:38:LEU:HD13	1:B:103:LEU:HD22	1.89	0.54
1:B:129:VAL:HG22	1:B:129:VAL:O	2.07	0.54
1:B:167:ARG:HG3	1:B:167:ARG:HH21	1.72	0.54
1:B:309:GLN:HE21	1:B:309:GLN:HA	1.71	0.54
1:A:161:HIS:CD2	1:A:163:ASN:HB2	2.42	0.54
1:B:210:LEU:HD21	1:B:279:ALA:HB1	1.90	0.54
1:A:11:LYS:CD	1:A:11:LYS:N	2.70	0.53
1:A:126:ASP:HB3	1:A:129:VAL:HG13	1.90	0.53
1:B:146:HIS:NE2	1:B:157:ALA:O	2.42	0.53
1:B:68:ARG:HD3	1:B:91:GLU:CG	2.37	0.53
1:B:227:GLU:OE1	1:B:264:TYR:HE1	1.91	0.53
1:A:248:CYS:HA	2:C:37:A:O2'	2.08	0.53
1:B:126:ASP:CB	1:B:129:VAL:HG12	2.39	0.53
1:A:96:THR:HG21	1:A:195:TYR:HE1	1.73	0.53
1:A:231:ARG:HD2	1:B:264:TYR:CE2	2.43	0.53
1:A:268:VAL:O	1:A:272:VAL:HG23	2.09	0.53
1:B:27:ALA:HB1	1:B:38:LEU:CD1	2.36	0.53
1:B:161:HIS:CE1	2:D:41:C:H4'	2.44	0.53
1:B:13:ILE:CG1	1:B:103:LEU:HG	2.39	0.53
3:B:317:DST:H132	5:B:322:HOH:O	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:9:LEU:CD2	1:B:9:LEU:N	2.72	0.53
1:A:11:LYS:H	1:A:11:LYS:CD	2.22	0.53
1:A:74:ASP:O	1:A:77:GLN:HG2	2.08	0.52
1:B:34:LEU:HD12	1:B:35:PRO:HD2	1.91	0.52
1:B:92:MET:HE1	1:B:115:LEU:HD11	1.91	0.52
1:A:11:LYS:H	1:A:11:LYS:HD3	1.73	0.52
1:A:101:ILE:CD1	1:A:314:ILE:HG21	2.28	0.52
1:B:136:GLN:HE22	1:B:145:LEU:CD1	2.11	0.52
1:B:147:ARG:CA	1:B:150:GLN:HG2	2.39	0.52
2:C:60:U:O3'	2:C:61:C:H6	1.93	0.52
1:A:240:HIS:CE1	1:A:243:LEU:HD13	2.45	0.52
2:D:9:A:O2'	2:D:10:G:OP1	2.21	0.51
1:A:9:LEU:N	1:A:10:PRO:CD	2.73	0.51
1:B:141:GLY:C	1:B:143:GLU:H	2.14	0.51
1:B:145:LEU:HD22	1:B:172:LEU:CD1	2.41	0.51
3:B:317:DST:H133	5:B:348:HOH:O	2.10	0.51
1:B:89:LEU:HA	1:B:92:MET:HE3	1.93	0.51
1:B:131:ALA:O	1:B:134:GLU:HB3	2.10	0.51
1:B:147:ARG:C	1:B:149:LEU:N	2.63	0.51
1:B:45:LEU:HD12	1:B:54:THR:HB	1.93	0.51
1:A:153:ASP:OD2	1:A:181:THR:HG23	2.11	0.51
1:A:153:ASP:OD2	1:A:183:THR:HG23	2.11	0.51
1:A:289:TRP:HB3	1:A:292:VAL:HG21	1.92	0.50
1:A:103:LEU:HD22	1:A:105:VAL:HG22	1.92	0.50
2:C:9:A:H2	2:C:44:G:O2'	1.94	0.50
1:A:299:LYS:C	1:A:301:GLU:N	2.61	0.50
1:B:147:ARG:C	1:B:149:LEU:H	2.14	0.50
2:D:9:A:C2	2:D:44:G:H2'	2.46	0.50
1:B:130:ARG:CA	1:B:133:ILE:CD1	2.89	0.50
2:D:52:G:O2'	2:D:53:G:H5'	2.11	0.50
2:C:44:G:O2'	2:C:45:U:P	2.70	0.50
1:A:163:ASN:O	2:C:40:C:C5'	2.60	0.49
2:D:58:A:H1'	2:D:60:U:C5	2.46	0.49
1:A:180:LYS:HE3	1:A:180:LYS:HA	1.94	0.49
1:A:92:MET:HE1	1:A:115:LEU:HD11	1.94	0.49
1:A:55:ALA:HB2	1:A:217:ARG:NH1	2.26	0.49
1:B:218:PHE:HA	1:B:221:MET:CE	2.43	0.49
2:C:73:A:H2'	2:C:74:C:H5''	1.95	0.49
1:A:217:ARG:NH1	3:A:317:DST:S9	2.86	0.49
1:B:183:THR:O	1:B:187:GLN:HG2	2.13	0.49
1:B:19:THR:O	1:B:20:ALA:HB3	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:304:ARG:NH1	1:A:305:ASP:OD1	2.46	0.49
2:D:49:C:H2'	2:D:50:U:C6	2.48	0.49
1:A:153:ASP:OD1	1:A:183:THR:HG23	2.13	0.48
1:B:13:ILE:CG2	1:B:101:ILE:HD11	2.43	0.48
1:B:278:LEU:O	1:B:282:GLN:HG3	2.14	0.48
1:B:217:ARG:O	1:B:221:MET:HG3	2.13	0.48
1:B:299:LYS:N	1:B:300:PRO:CD	2.77	0.48
1:B:126:ASP:HB3	1:B:129:VAL:HG12	1.94	0.48
1:B:132:ARG:C	1:B:134:GLU:H	2.16	0.48
2:C:58:A:O2'	2:C:60:U:OP2	2.21	0.48
1:A:299:LYS:O	1:A:302:GLN:NE2	2.47	0.48
1:B:132:ARG:O	1:B:134:GLU:N	2.47	0.48
1:B:147:ARG:HA	1:B:150:GLN:CG	2.41	0.48
1:B:46:ILE:HG23	1:B:72:ILE:HG13	1.96	0.48
1:B:146:HIS:CE1	1:B:157:ALA:HB1	2.49	0.48
2:C:74:C:C6	2:C:74:C:H5'	2.47	0.48
1:B:150:GLN:O	1:B:154:PRO:HB3	2.14	0.48
1:B:149:LEU:HD21	1:B:171:ALA:CB	2.44	0.48
1:B:25:ALA:O	1:B:29:GLU:HG2	2.14	0.48
1:B:92:MET:CE	1:B:115:LEU:HD11	2.43	0.47
1:B:142:TRP:CE3	1:B:168:LEU:HD13	2.44	0.47
1:B:217:ARG:NH1	3:B:317:DST:S9	2.87	0.47
1:B:107:GLY:O	1:B:282:GLN:NE2	2.48	0.47
2:C:13:C:H2'	2:C:14:A:C5'	2.33	0.47
1:A:109:MET:HG2	1:A:285:TRP:HB3	1.97	0.47
1:A:33:ILE:C	1:A:34:LEU:HD12	2.35	0.47
1:B:236:ARG:HB3	1:B:238:ASP:OD2	2.13	0.47
1:B:155:VAL:HG12	1:B:183:THR:HG1	1.79	0.47
1:B:34:LEU:HD21	1:B:308:LEU:HD22	1.96	0.47
1:B:9:LEU:HB2	1:B:10:PRO:CD	2.37	0.47
2:D:36:A:H4'	2:D:37:A:O5'	2.14	0.47
1:B:301:GLU:H	1:B:301:GLU:CD	2.18	0.47
1:B:147:ARG:HG2	1:B:150:GLN:OE1	2.15	0.47
1:B:134:GLU:C	1:B:134:GLU:CD	2.74	0.47
1:B:136:GLN:NE2	1:B:137:ALA:N	2.61	0.47
2:C:73:A:C2'	2:C:74:C:H5''	2.44	0.47
1:A:146:HIS:HA	1:A:168:LEU:HD21	1.97	0.46
1:B:167:ARG:NH2	1:B:167:ARG:HG3	2.30	0.46
1:A:17:GLY:O	1:A:23:LYS:HE2	2.16	0.46
1:B:29:GLU:O	1:B:32:LYS:HB2	2.16	0.46
1:A:238:ASP:OD1	1:A:239:LEU:HD13	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:40:C:C2'	2:D:41:C:H5'	2.45	0.46
1:A:181:THR:HG23	1:A:183:THR:H	1.80	0.46
1:A:58:ASN:OD1	1:A:61:GLU:HG3	2.16	0.46
1:B:41:VAL:CG1	1:B:112:PHE:CZ	2.96	0.46
1:B:103:LEU:C	1:B:103:LEU:HD23	2.36	0.46
2:D:41:C:H2'	2:D:42:C:C6	2.51	0.46
1:A:9:LEU:CD2	1:A:9:LEU:N	2.79	0.46
1:B:16:MET:O	1:B:201:ALA:HA	2.16	0.46
2:D:17:C:H5''	2:D:18:G:H5''	1.98	0.46
1:A:302:GLN:H	1:A:302:GLN:CD	2.18	0.46
1:A:72:ILE:O	1:A:73:ARG:HG3	2.15	0.45
1:B:144:SER:O	1:B:148:GLN:N	2.43	0.45
1:B:234:PHE:HA	1:B:255:TRP:CH2	2.51	0.45
1:B:42:ASP:HB3	1:B:45:LEU:HB2	1.98	0.45
2:D:64:A:H2'	2:D:65:G:H8	1.81	0.45
1:B:230:VAL:HG21	1:B:254:MET:HG2	1.97	0.45
1:B:57:PRO:HA	1:B:61:GLU:OE1	2.16	0.45
2:D:50:U:O2'	2:D:51:U:H5'	2.17	0.45
3:A:317:DST:H132	5:A:343:HOH:O	2.16	0.45
1:A:30:LEU:HD12	1:A:30:LEU:HA	1.75	0.45
1:B:104:LEU:CD2	1:B:104:LEU:N	2.80	0.45
1:B:132:ARG:C	1:B:134:GLU:N	2.69	0.45
2:D:67:C:N4	2:D:68:C:H41	2.15	0.45
1:A:178:SER:C	1:A:180:LYS:H	2.20	0.45
1:A:34:LEU:HB3	1:A:36:VAL:HG13	1.98	0.45
1:B:39:ILE:HB	1:B:104:LEU:HD13	1.98	0.45
2:D:41:C:H2'	2:D:42:C:H6	1.82	0.45
1:A:240:HIS:CE1	1:A:242:ASP:HB2	2.52	0.45
1:A:183:THR:O	1:A:187:GLN:HG2	2.17	0.44
1:A:55:ALA:CB	1:A:217:ARG:HH12	2.30	0.44
1:A:311:VAL:O	1:A:314:ILE:HD11	2.17	0.44
1:B:139:GLU:O	1:B:141:GLY:N	2.51	0.44
1:B:16:MET:HE2	1:B:112:PHE:CD2	2.52	0.44
1:B:174:VAL:HG21	1:B:182:LEU:HD23	1.99	0.44
1:B:200:PHE:CD1	1:B:293:HIS:HB2	2.52	0.44
1:B:211:HIS:CD2	1:B:276:ARG:HD2	2.53	0.44
1:B:9:LEU:N	1:B:9:LEU:HD23	2.32	0.44
2:C:14:A:H2'	2:C:15:G:H5'	1.98	0.44
1:B:133:ILE:H	1:B:133:ILE:CD1	2.05	0.44
1:A:85:ARG:HD2	5:A:369:HOH:O	2.18	0.44
1:B:24:THR:O	1:B:28:ILE:HG23	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:15:G:H2'	2:D:16:U:C6	2.53	0.44
2:D:50:U:O5'	2:D:50:U:H6	2.00	0.44
2:D:53:G:O2'	2:D:54:U:H5'	2.18	0.44
1:A:301:GLU:C	1:A:303:ALA:N	2.71	0.44
1:B:143:GLU:CA	1:B:143:GLU:OE2	2.66	0.44
1:B:149:LEU:HD21	1:B:171:ALA:HB1	1.99	0.44
1:B:37:GLU:HG2	1:B:68:ARG:HE	1.82	0.44
1:A:147:ARG:HA	1:A:150:GLN:HE21	1.83	0.43
2:D:18:G:H4'	2:D:60:U:C2	2.53	0.43
1:B:155:VAL:HG11	1:B:183:THR:HG21	2.00	0.43
2:C:73:A:C2'	2:C:74:C:C5'	2.95	0.43
1:B:181:THR:HG23	1:B:183:THR:H	1.83	0.43
1:B:249:VAL:HG22	2:D:37:A:N3	2.33	0.43
1:B:104:LEU:HD22	1:B:104:LEU:N	2.34	0.43
1:B:309:GLN:CA	1:B:309:GLN:HE21	2.32	0.43
2:C:14:A:H2'	2:C:15:G:C5'	2.49	0.43
2:C:16:U:H5''	2:C:17:C:OP2	2.18	0.43
1:A:44:ALA:HB3	2:C:37:A:C6	2.54	0.43
1:B:170:ARG:HG2	2:D:34:G:OP1	2.19	0.43
1:B:96:THR:HA	1:B:100:ARG:O	2.19	0.43
2:C:35:A:O2'	2:C:36:A:OP1	2.28	0.43
1:B:38:LEU:N	1:B:38:LEU:CD2	2.82	0.43
1:A:19:THR:OG1	3:A:317:DST:S9	2.73	0.42
1:A:28:ILE:HD12	1:A:28:ILE:O	2.18	0.42
1:B:303:ALA:O	1:B:307:VAL:HG23	2.19	0.42
1:B:309:GLN:NE2	1:B:309:GLN:O	2.52	0.42
2:D:58:A:O2'	2:D:59:U:H3'	2.19	0.42
1:A:126:ASP:HB3	1:A:129:VAL:CG1	2.49	0.42
1:B:140:GLN:HG2	1:B:143:GLU:HG2	2.01	0.42
1:B:181:THR:O	1:B:184:GLU:HB2	2.20	0.42
2:D:23:A:H2'	2:D:24:G:C8	2.54	0.42
1:A:172:LEU:HA	1:A:172:LEU:HD12	1.91	0.42
1:B:141:GLY:O	1:B:143:GLU:N	2.52	0.42
1:B:89:LEU:HA	1:B:89:LEU:HD12	1.93	0.42
1:B:178:SER:C	1:B:180:LYS:H	2.22	0.42
1:A:301:GLU:H	1:A:301:GLU:HG3	1.69	0.42
1:B:200:PHE:CE2	1:B:307:VAL:HG13	2.55	0.42
1:B:213:ARG:O	1:B:217:ARG:HB2	2.20	0.42
1:A:165:PRO:HD2	2:C:40:C:H5'	2.01	0.42
1:A:9:LEU:HD23	1:A:9:LEU:N	2.35	0.42
1:B:100:ARG:HH11	1:B:100:ARG:HG2	1.85	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:16:MET:HG2	1:B:112:PHE:CE2	2.55	0.42
1:B:217:ARG:NH2	5:B:325:HOH:O	2.51	0.42
2:C:32:U:O2'	2:C:36:A:N1	2.51	0.42
1:A:240:HIS:NE2	1:A:243:LEU:HD13	2.35	0.42
2:D:40:C:H2'	2:D:41:C:H5'	2.02	0.42
1:A:163:ASN:O	2:C:40:C:H5''	2.20	0.41
1:B:155:VAL:HG11	1:B:183:THR:HG23	2.02	0.41
1:B:310:VAL:O	1:B:311:VAL:C	2.58	0.41
1:B:130:ARG:HA	1:B:133:ILE:HD13	2.02	0.41
2:D:11:C:H2'	2:D:12:U:C6	2.55	0.41
1:B:262:ILE:HG13	1:B:266:GLU:HG2	2.01	0.41
1:A:153:ASP:CG	1:A:183:THR:HG23	2.41	0.41
1:B:63:LEU:HA	1:B:63:LEU:HD22	1.88	0.41
1:B:76:SER:HA	1:B:243:LEU:HD13	2.01	0.41
2:D:7:A:N6	2:D:49:C:H42	2.19	0.41
1:B:145:LEU:HD22	1:B:172:LEU:HD13	2.02	0.41
1:B:68:ARG:CD	1:B:95:ILE:HD11	2.51	0.41
1:B:249:VAL:HG22	2:D:37:A:C2	2.56	0.41
2:D:19:G:N2	2:D:56:C:O2	2.53	0.41
1:B:116:LEU:HA	1:B:193:LEU:HB2	2.03	0.41
1:A:137:ALA:O	1:A:141:GLY:N	2.52	0.41
1:B:37:GLU:HB3	1:B:95:ILE:CD1	2.37	0.41
1:B:96:THR:C	1:B:98:ALA:N	2.74	0.41
1:A:11:LYS:CE	1:A:11:LYS:H	2.34	0.41
1:A:77:GLN:HE21	1:A:77:GLN:HB3	1.63	0.41
2:C:13:C:O2'	2:C:14:A:H5''	2.19	0.41
1:A:28:ILE:C	1:A:28:ILE:HD12	2.41	0.41
1:A:91:GLU:O	1:A:95:ILE:HG13	2.21	0.41
1:B:24:THR:HG21	1:B:57:PRO:HG3	2.02	0.41
1:A:101:ILE:HA	1:A:102:PRO:HD3	1.85	0.40
1:A:73:ARG:NH1	1:A:73:ARG:CG	2.81	0.40
1:B:262:ILE:HG23	1:B:263:SER:O	2.21	0.40
1:B:96:THR:C	1:B:98:ALA:H	2.23	0.40
1:A:164:ASP:O	1:A:168:LEU:HB2	2.21	0.40
1:B:145:LEU:HD22	1:B:172:LEU:HD11	2.03	0.40
1:B:311:VAL:HG22	1:B:311:VAL:O	2.22	0.40
1:A:153:ASP:OD1	1:A:183:THR:CG2	2.69	0.40
1:B:163:ASN:O	1:B:165:PRO:HD3	2.22	0.40
2:D:7:A:N6	2:D:49:C:N4	2.69	0.40
1:B:34:LEU:HD21	1:B:308:LEU:CD2	2.52	0.40
1:B:46:ILE:HD12	1:B:72:ILE:HD11	2.04	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	
1	A	304/316 (96%)	292 (96%)	9 (3%)	3 (1%)	15	27
1	B	303/316 (96%)	286 (94%)	14 (5%)	3 (1%)	15	27
All	All	607/632 (96%)	578 (95%)	23 (4%)	6 (1%)	15	27

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	299	LYS
1	A	298	GLU
1	B	140	GLN
1	B	142	TRP
1	B	133	ILE
1	A	300	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	
1	A	249/256 (97%)	222 (89%)	27 (11%)	6	10
1	B	249/256 (97%)	208 (84%)	41 (16%)	2	3
All	All	498/512 (97%)	430 (86%)	68 (14%)	3	5

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

Mol	Chain	Res	Type
1	A	9	LEU
1	A	11	LYS
1	A	15	LEU
1	A	26	LEU
1	A	28	ILE
1	A	30	LEU
1	A	33	ILE
1	A	52	ILE
1	A	63	LEU
1	A	89	LEU
1	A	104	LEU
1	A	105	VAL
1	A	109	MET
1	A	129	VAL
1	A	168	LEU
1	A	172	LEU
1	A	180	LYS
1	A	181	THR
1	A	185	LEU
1	A	193	LEU
1	A	210	LEU
1	A	239	LEU
1	A	243	LEU
1	A	267	MET
1	A	286	LEU
1	A	302	GLN
1	A	308	LEU
1	B	9	LEU
1	B	11	LYS
1	B	13	ILE
1	B	28	ILE
1	B	31	ARG
1	B	32	LYS
1	B	33	ILE
1	B	41	VAL
1	B	62	LEU
1	B	63	LEU
1	B	89	LEU
1	B	101	ILE
1	B	104	LEU
1	B	105	VAL
1	B	109	MET

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Mol	Chain	Res	Type
1	B	130	ARG
1	B	132	ARG
1	B	133	ILE
1	B	135	GLN
1	B	136	GLN
1	B	139	GLU
1	B	140	GLN
1	B	143	GLU
1	B	155	VAL
1	B	161	HIS
1	B	163	ASN
1	B	168	LEU
1	B	178	SER
1	B	181	THR
1	B	188	THR
1	B	193	LEU
1	B	196	GLN
1	B	209	LEU
1	B	217	ARG
1	B	233	LEU
1	B	239	LEU
1	B	267	MET
1	B	278	LEU
1	B	301	GLU
1	B	309	GLN
1	B	311	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	136	GLN
1	A	150	GLN
1	A	161	HIS
1	A	163	ASN
1	A	212	GLN
1	A	216	GLN
1	A	220	GLN
1	A	302	GLN
1	B	67	HIS
1	B	136	GLN
1	B	161	HIS
1	B	166	GLN

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Mol	Chain	Res	Type
1	B	211	HIS
1	B	216	GLN
1	B	220	GLN
1	B	277	GLN
1	B	302	GLN
1	B	309	GLN

### 5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	C	73/76 (96%)	23 (31%)	15 (20%)
2	D	68/76 (89%)	18 (26%)	9 (13%)
All	All	141/152 (92%)	41 (29%)	24 (17%)

All (41) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	C	6	G
2	C	7	A
2	C	8	U
2	C	10	G
2	C	14	A
2	C	17	C
2	C	18	G
2	C	19	G
2	C	20	U
2	C	31	A
2	C	34	G
2	C	35	A
2	C	36	A
2	C	37	A
2	C	38	A
2	C	45	U
2	C	46	G
2	C	47	U
2	C	48	C
2	C	59	U
2	C	61	C
2	C	70	G
2	C	74	C
2	D	8	U

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Mol	Chain	Res	Type
2	D	9	A
2	D	10	G
2	D	17	C
2	D	20	U
2	D	22	G
2	D	33	U
2	D	34	G
2	D	35	A
2	D	36	A
2	D	37	A
2	D	38	A
2	D	41	C
2	D	44	G
2	D	45	U
2	D	47	U
2	D	49	C
2	D	61	C

All (24) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	C	7	A
2	C	9	A
2	C	18	G
2	C	19	G
2	C	22	G
2	C	34	G
2	C	35	A
2	C	36	A
2	C	37	A
2	C	44	G
2	C	45	U
2	C	46	G
2	C	48	C
2	C	58	A
2	C	60	U
2	D	9	A
2	D	19	G
2	D	32	U
2	D	34	G
2	D	35	A
2	D	36	A

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Mol	Chain	Res	Type
2	D	37	A
2	D	44	G
2	D	60	U

## 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 12 ligands modelled in this entry, 10 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
3	DST	B	317	4	9,13,13	2.34	1 (11%)	11,19,19	1.16	1 (9%)
3	DST	A	317	4	9,13,13	2.27	1 (11%)	11,19,19	1.30	1 (9%)

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
3	DST	B	317	4	-	3/7/13/13	-
3	DST	A	317	4	-	3/7/13/13	-

All (2) bond length outliers are listed below:



Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	317	DST	C10-S9	-6.33	1.76	1.84
3	A	317	DST	C10-S9	-6.15	1.76	1.84

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	317	DST	O8-P3-O7	3.13	117.39	109.82
3	B	317	DST	O8-P3-O7	2.48	115.84	109.82

There are no chirality outliers.

All (6) torsion outliers are listed below:

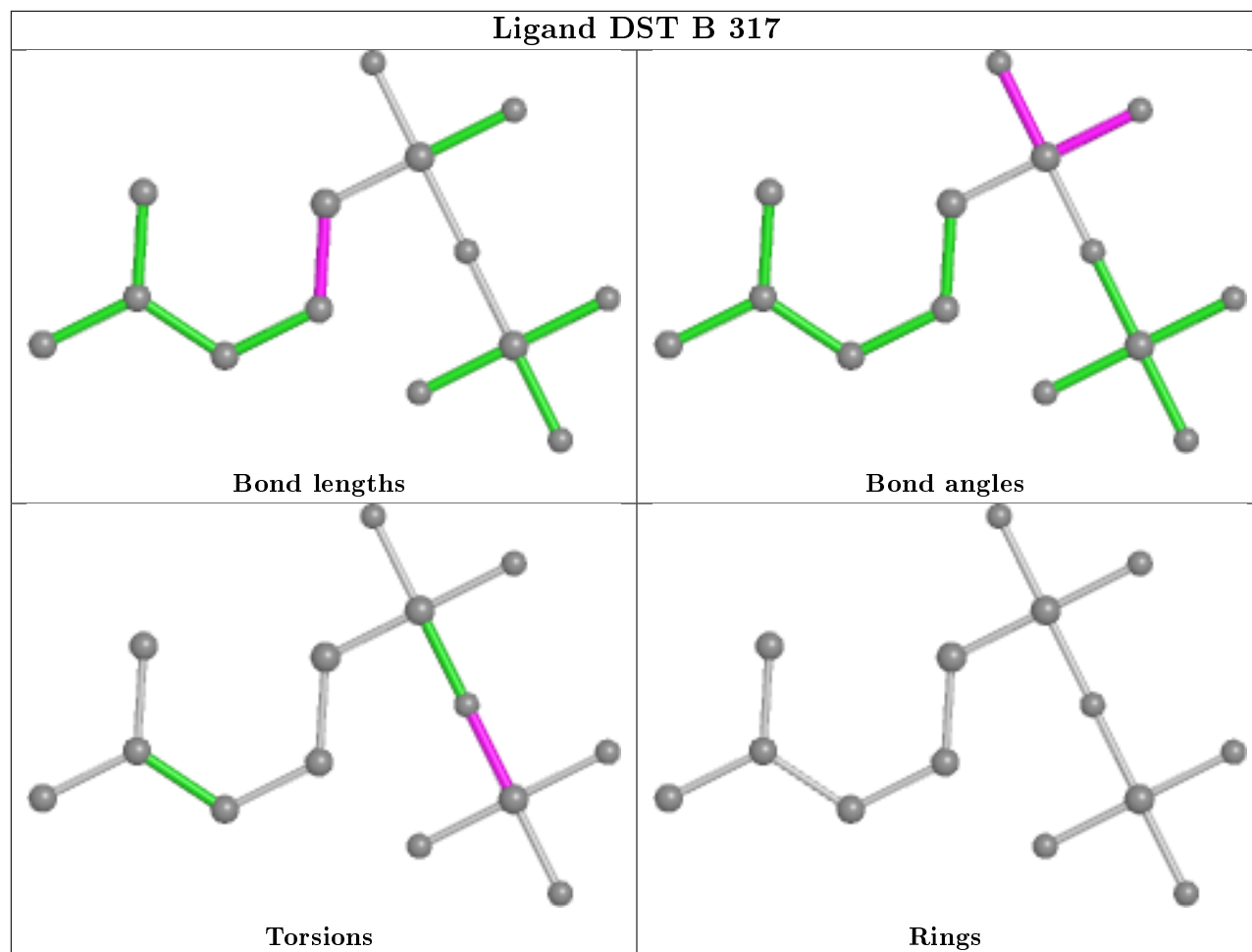
Mol	Chain	Res	Type	Atoms
3	A	317	DST	P3-O2-P1-O5
3	B	317	DST	P3-O2-P1-O5
3	A	317	DST	P3-O2-P1-O4
3	A	317	DST	P3-O2-P1-O6
3	B	317	DST	P3-O2-P1-O4
3	B	317	DST	P3-O2-P1-O6

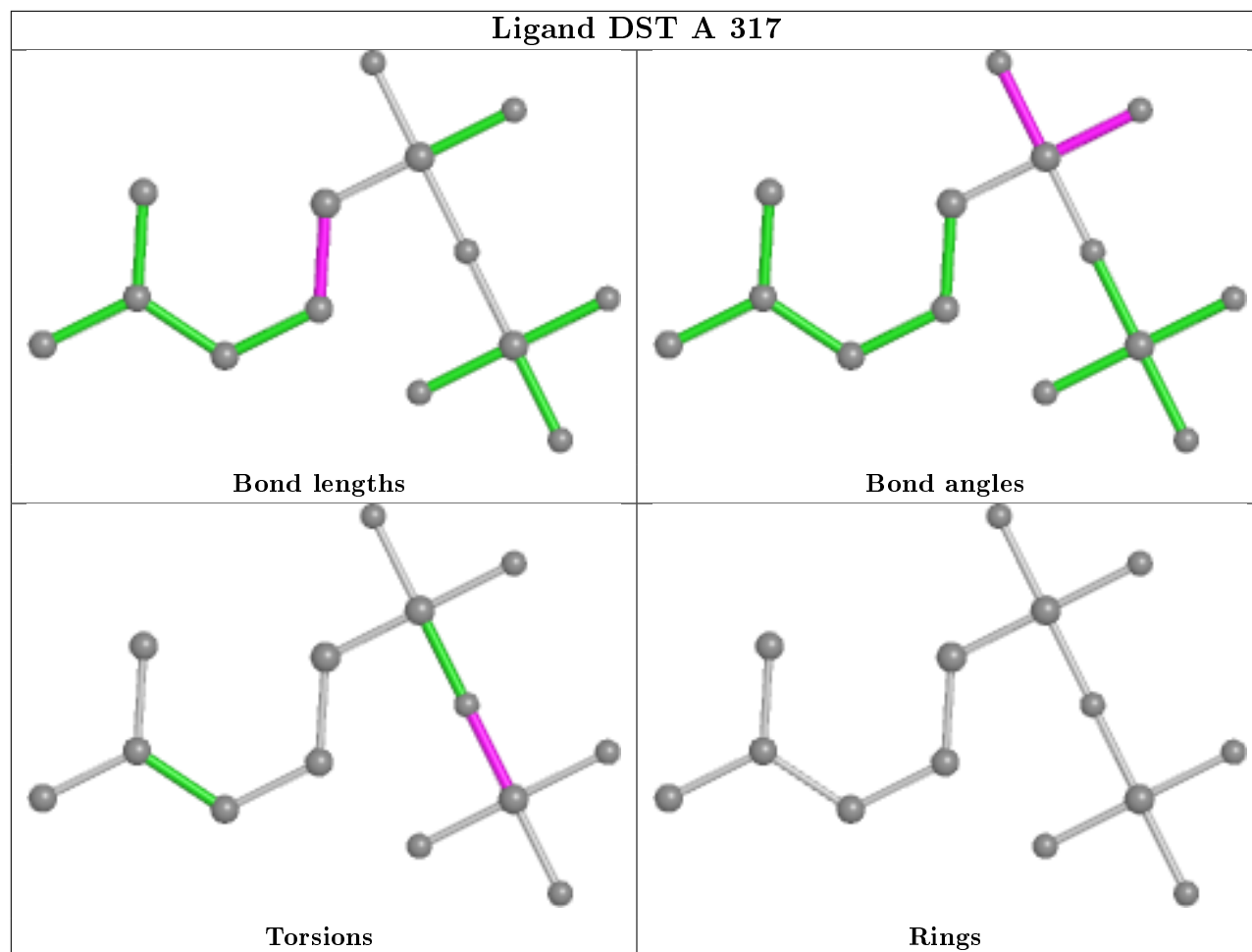
There are no ring outliers.

2 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	317	DST	5	0
3	A	317	DST	3	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, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	306/316 (96%)	0.29	7 (2%) 60 69	14, 29, 54, 73	0
1	B	305/316 (96%)	0.58	27 (8%) 9 11	19, 38, 77, 89	0
2	C	74/76 (97%)	0.24	2 (2%) 54 63	15, 31, 53, 72	0
2	D	69/76 (90%)	1.86	31 (44%) 0 0	25, 72, 90, 90	0
All	All	754/784 (96%)	0.55	67 (8%) 9 11	14, 34, 80, 90	0

All (67) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	3	C	6.3
1	B	140	GLN	6.2
1	B	150	GLN	6.1
1	B	154	PRO	5.7
2	D	4	C	5.5
2	D	64	A	5.5
2	D	47	U	5.4
2	D	2	C	5.2
2	D	65	G	5.1
1	A	33	ILE	5.1
2	D	20	U	5.1
2	D	63	G	5.0
1	B	136	GLN	4.8
2	D	1	G	4.5
2	D	45	U	4.5
2	C	47	U	4.5
1	B	134	GLU	4.2
1	A	313	ALA	4.0
1	A	288	GLY	4.0
1	B	135	GLN	4.0
2	D	66	U	3.9

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Mol	Chain	Res	Type	RSRZ
1	B	168	LEU	3.9
2	D	69	G	3.8
1	A	299	LYS	3.8
1	B	8	SER	3.7
1	B	131	ALA	3.7
1	B	308	LEU	3.6
2	D	17	C	3.6
2	D	67	C	3.6
1	B	177	ILE	3.5
2	D	48	C	3.4
2	D	18	G	3.3
1	A	314	ILE	3.3
1	B	149	LEU	3.2
2	D	49	C	3.2
1	B	132	ARG	3.2
1	A	191	ASP	3.0
1	B	129	VAL	3.0
2	D	68	C	2.9
2	D	7	A	2.7
1	B	303	ALA	2.7
1	A	298	GLU	2.7
1	B	147	ARG	2.6
2	D	62	C	2.6
1	B	309	GLN	2.6
2	D	44	G	2.5
1	B	63	LEU	2.5
2	D	51	U	2.5
2	D	59	U	2.5
1	B	142	TRP	2.4
2	D	58	A	2.4
2	D	21	A	2.4
1	B	141	GLY	2.3
1	B	178	SER	2.3
2	D	53	G	2.3
1	B	33	ILE	2.3
2	D	46	G	2.3
2	D	8	U	2.3
1	B	29	GLU	2.2
2	D	19	G	2.2
2	D	5	G	2.1
1	B	144	SER	2.1
1	B	151	GLU	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	176	PHE	2.0
2	C	45	U	2.0
1	B	133	ILE	2.0
2	D	6	G	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 carbohydrates in this entry.

## 6.4 Ligands [i](#)

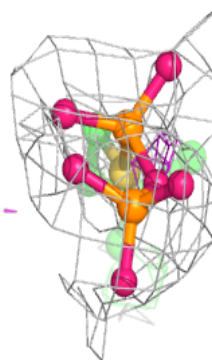
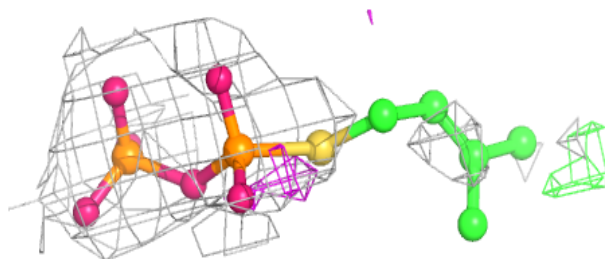
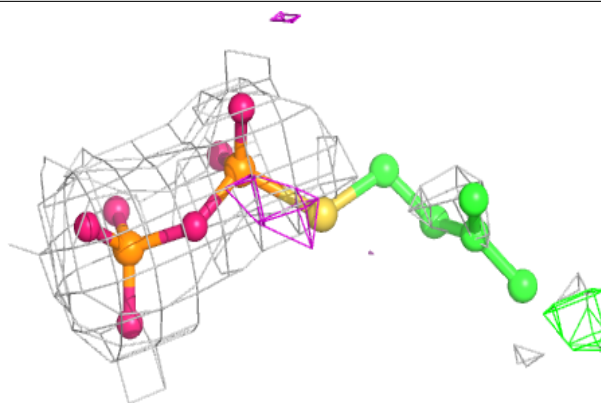
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
4	MG	B	318	1/1	0.49	0.28	63,63,63,63	0
4	MG	C	83	1/1	0.75	0.30	31,31,31,31	0
4	MG	A	318	1/1	0.81	0.15	52,52,52,52	0
4	MG	C	81	1/1	0.83	0.26	29,29,29,29	0
3	DST	B	317	14/14	0.85	0.36	73,81,83,83	0
4	MG	C	79	1/1	0.87	0.13	23,23,23,23	0
4	MG	C	82	1/1	0.89	0.22	36,36,36,36	0
4	MG	C	80	1/1	0.91	0.33	29,29,29,29	0
3	DST	A	317	14/14	0.91	0.23	44,51,56,56	0
4	MG	C	77	1/1	0.92	0.25	10,10,10,10	0
4	MG	C	78	1/1	0.92	0.37	16,16,16,16	0
4	MG	D	77	1/1	0.94	0.32	28,28,28,28	0

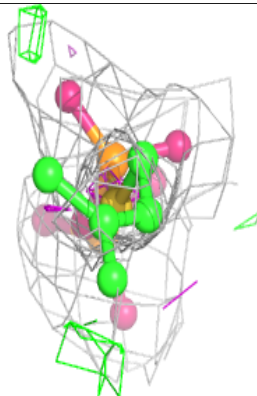
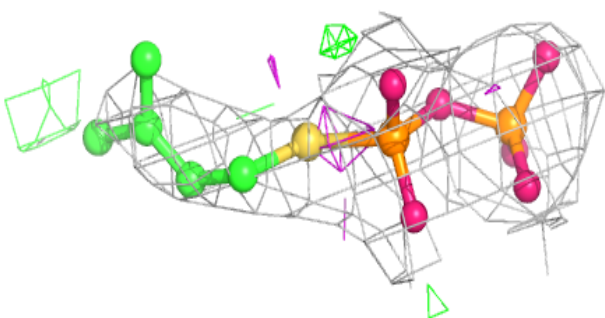
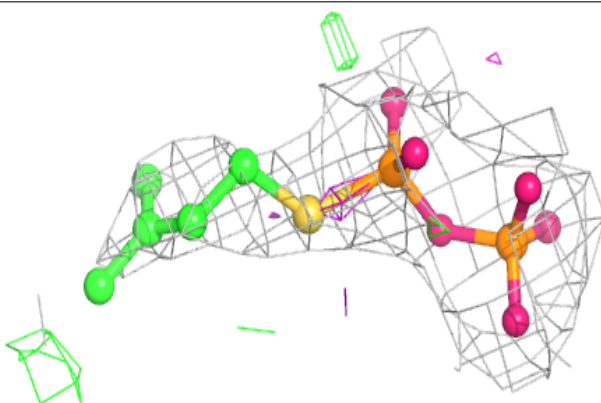
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

**Electron density around DST B 317:**

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

**Electron density around DST A 317:**

$2mF_o-DF_c$  (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.