



# wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 27, 2020 – 02:09 PM BST

PDB ID : 6UU7  
Title : E. coli sigma-S transcription initiation complex with a 6-nt RNA and an NTP ("Old" crystal soaked with UTP, CTP, ddGTP, and dinucleotide ApG for 30 minutes)  
Authors : Zuo, Y.; De, S.; Steitz, T.A.  
Deposited on : 2019-10-30  
Resolution : 4.40 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 : 2.13  
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) : 2.13

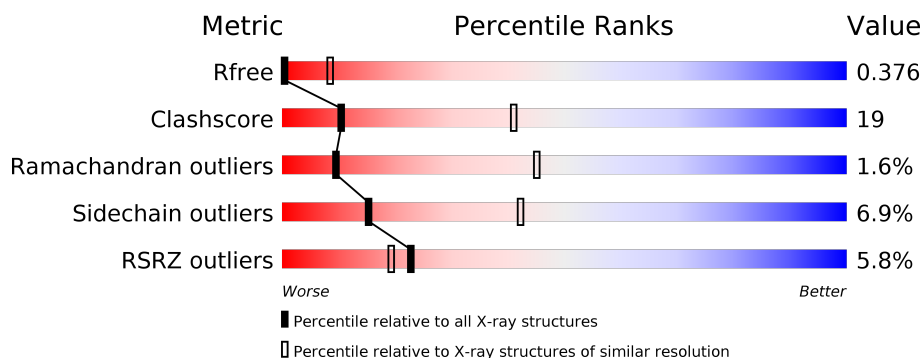
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 4.40 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.



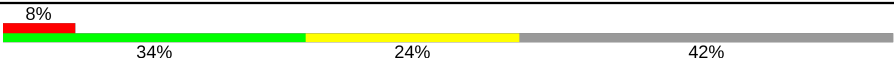

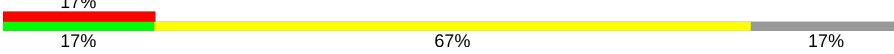
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	1043 (5.00-3.80)
Clashscore	141614	1111 (5.00-3.80)
Ramachandran outliers	138981	1059 (5.00-3.80)
Sidechain outliers	138945	1041 (5.00-3.80)
RSRZ outliers	127900	1095 (5.08-3.70)

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	AAA	242	<div> <div>8%</div> <div> <div></div> <div>60%</div> <div>31%</div> <div>5%</div> </div> </div>
1	BBB	242	<div> <div>5%</div> <div> <div></div> <div>61%</div> <div>31%</div> <div>6%</div> </div> </div>
2	CCC	1342	<div> <div>3%</div> <div> <div></div> <div>64%</div> <div>32%</div> <div></div> </div> </div>
3	DDD	1407	<div> <div>7%</div> <div> <div></div> <div>61%</div> <div>33%</div> <div></div> </div> </div>
4	EEE	90	<div> <div>3%</div> <div> <div></div> <div>59%</div> <div>27%</div> <div>12%</div> </div> </div>
5	FFF	336	<div> <div>5%</div> <div> <div></div> <div>54%</div> <div>24%</div> <div>20%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
6	111	50	
7	222	50	
8	333	6	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
11	DG3	DDD	1505	-	-	X	X
9	ZN	DDD	1502	-	-	X	-

## 2 Entry composition

There are 12 unique types of molecules in this entry. The entry contains 28934 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-directed RNA polymerase subunit alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	AAA	230	Total	C	N	O	S	0	0	0
			1787	1112	317	352	6			
1	BBB	228	Total	C	N	O	S	0	0	0
			1767	1100	312	349	6			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
AAA	-6	ALA	-	expression tag	UNP P0A7Z4
AAA	-5	HIS	-	expression tag	UNP P0A7Z4
AAA	-4	HIS	-	expression tag	UNP P0A7Z4
AAA	-3	HIS	-	expression tag	UNP P0A7Z4
AAA	-2	HIS	-	expression tag	UNP P0A7Z4
AAA	-1	HIS	-	expression tag	UNP P0A7Z4
AAA	0	HIS	-	expression tag	UNP P0A7Z4
BBB	-6	ALA	-	expression tag	UNP P0A7Z4
BBB	-5	HIS	-	expression tag	UNP P0A7Z4
BBB	-4	HIS	-	expression tag	UNP P0A7Z4
BBB	-3	HIS	-	expression tag	UNP P0A7Z4
BBB	-2	HIS	-	expression tag	UNP P0A7Z4
BBB	-1	HIS	-	expression tag	UNP P0A7Z4
BBB	0	HIS	-	expression tag	UNP P0A7Z4

- Molecule 2 is a protein called DNA-directed RNA polymerase subunit beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	CCC	1341	Total	C	N	O	S	0	0	0
			10576	6636	1842	2055	43			

- Molecule 3 is a protein called DNA-directed RNA polymerase subunit beta'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	DDD	1362	Total	C	N	O	S	0	0	0
			10568	6633	1887	1998	50			

- Molecule 4 is a protein called DNA-directed RNA polymerase subunit omega.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	EEE	79	Total	C	N	O	S	0	0	0
			627	382	118	126	1			

- Molecule 5 is a protein called RNA polymerase sigma factor RpoS.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	FFF	270	Total	C	N	O	S	0	0	0
			2205	1383	407	411	4			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
FFF	2	GLY	SER	conflict	UNP P13445
FFF	33	GLU	GLN	conflict	UNP P13445
FFF	329	LEU	-	expression tag	UNP P13445
FFF	330	GLU	-	expression tag	UNP P13445
FFF	331	HIS	-	expression tag	UNP P13445
FFF	332	HIS	-	expression tag	UNP P13445
FFF	333	HIS	-	expression tag	UNP P13445
FFF	334	HIS	-	expression tag	UNP P13445
FFF	335	HIS	-	expression tag	UNP P13445
FFF	336	HIS	-	expression tag	UNP P13445

- Molecule 6 is a DNA chain called Synthetic DNA 50-mer (promoter non-template strand).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	111	29	Total	C	N	O	P	0	0	0
			595	283	107	176	29			

- Molecule 7 is a DNA chain called Synthetic DNA 50-mer (promoter template strand).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	222	32	Total	C	N	O	P	0	0	0
			652	312	117	192	31			

- Molecule 8 is DNA/RNA hybrid called DNA/RNA (5'-R(\*AP\*GP\*UP\*CP\*U)-D(P\*(DDG

))-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	333	5	Total	C	N	O	P	0	0	0
			102	47	17	34	4			

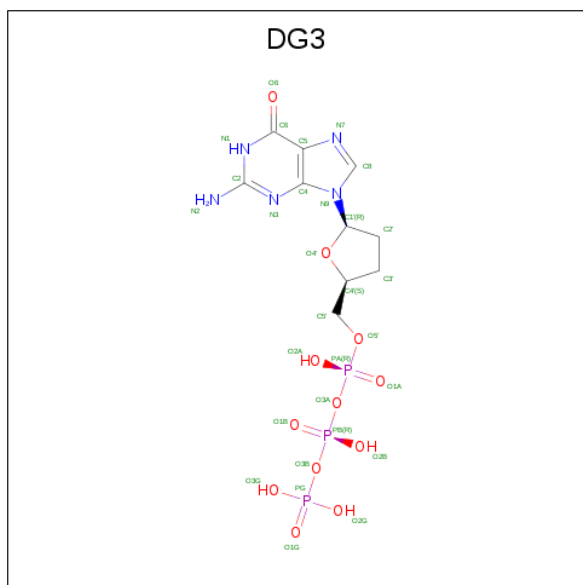
- Molecule 9 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	DDD	2	Total	Zn	0	0
			2	2		

- Molecule 10 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

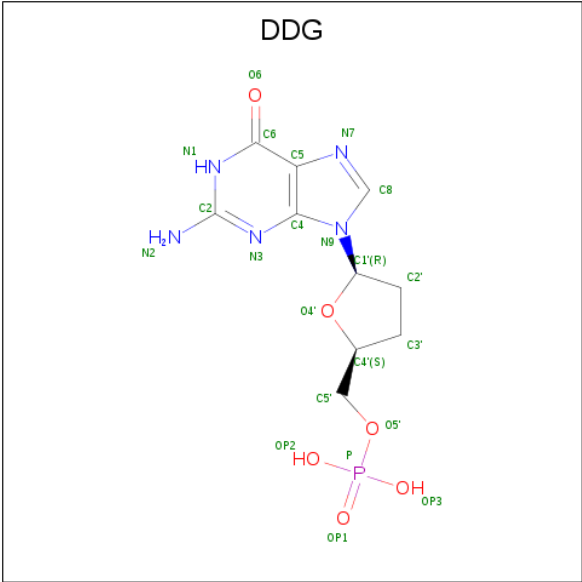
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	DDD	2	Total	Mg	0	0
			2	2		

- Molecule 11 is 2'-3'-DIDEOXYGUANOSINE-5'-TRIPHOSPHATE (three-letter code: DG3) (formula: C<sub>10</sub>H<sub>16</sub>N<sub>5</sub>O<sub>12</sub>P<sub>3</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
11	DDD	1	Total	C	N	O	P	0	0
			30	10	5	12	3		

- Molecule 12 is 2',3'-DIDEOXY-GUANOSINE-5'-MONOPHOSPHATE (three-letter code: DDG) (formula: C<sub>10</sub>H<sub>14</sub>N<sub>5</sub>O<sub>6</sub>P).

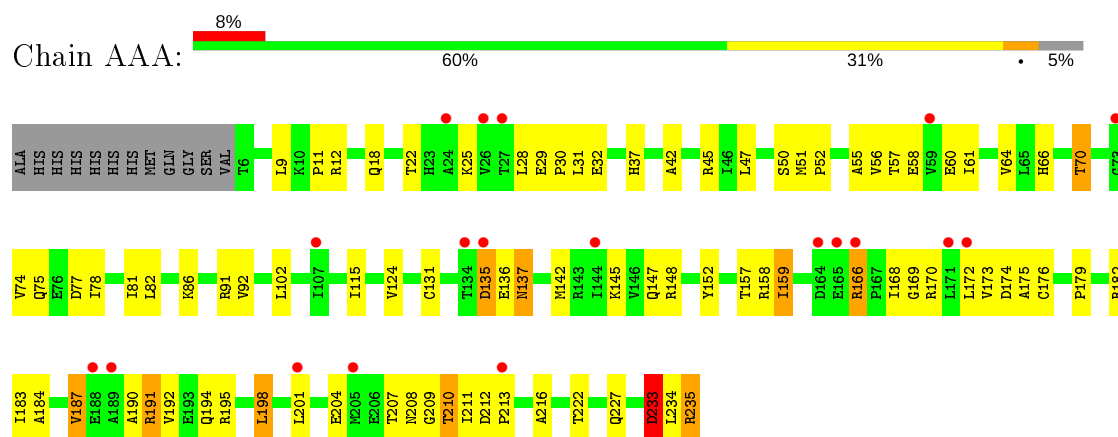


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
12	333	1	Total	C	N	O	P	0	0
			21	10	5	5	1		

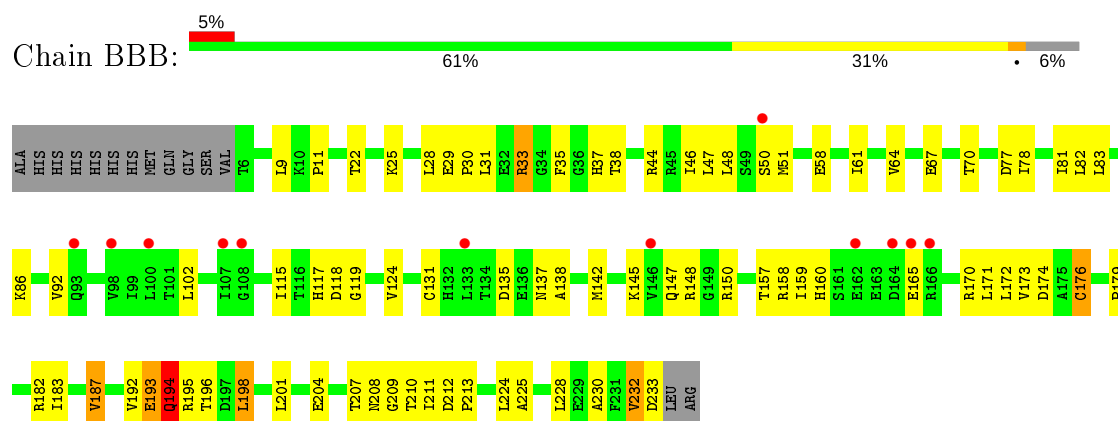
### 3 Residue-property plots [i](#)

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.

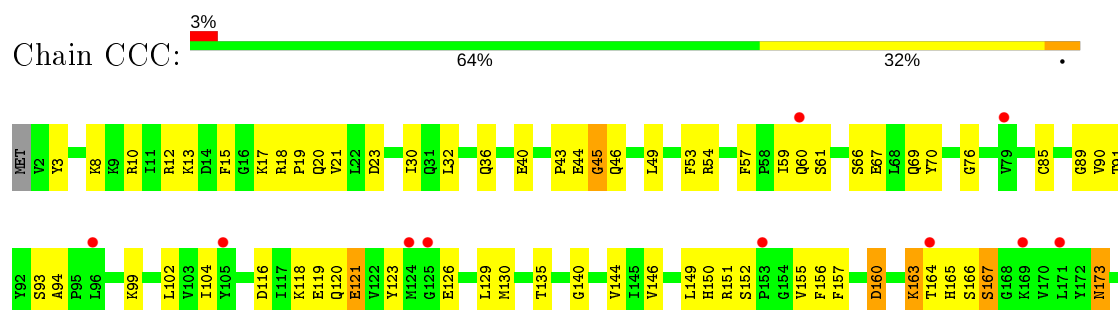
- Molecule 1: DNA-directed RNA polymerase subunit alpha



- Molecule 1: DNA-directed RNA polymerase subunit alpha



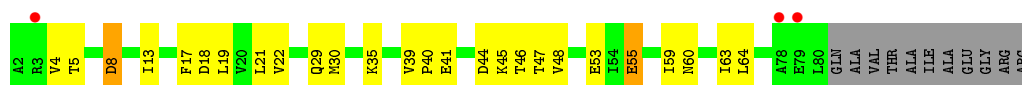
- Molecule 2: DNA-directed RNA polymerase subunit beta



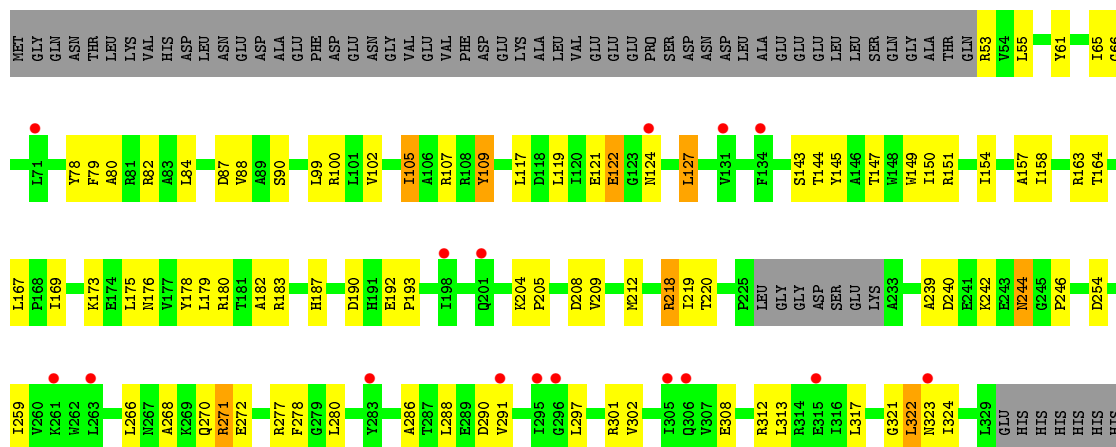




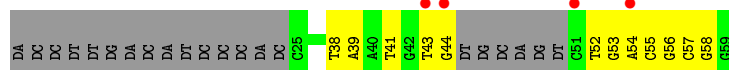




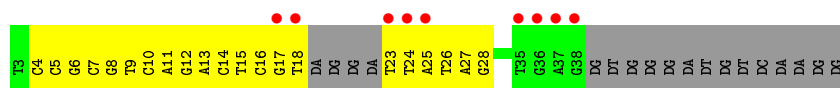
- Molecule 5: RNA polymerase sigma factor RpoS



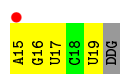
- Molecule 6: Synthetic DNA 50-mer (promoter non-template strand)



- Molecule 7: Synthetic DNA 50-mer (promoter template strand)



- Molecule 8: DNA/RNA (5'-R(\*AP\*GP\*UP\*CP\*U)-D(P\*(DDG))-3')



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	132.94Å 153.77Å 232.62Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.15 – 4.40 49.15 – 4.40	Depositor EDS
% Data completeness (in resolution range)	99.2 (49.15-4.40) 99.3 (49.15-4.40)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.07 (at 4.45Å)	Xtriage
Refinement program	REFMAC 5.8.0257	Depositor
R, $R_{free}$	0.344 , 0.390 0.332 , 0.376	Depositor DCC
$R_{free}$ test set	1464 reflections (4.77%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	179.5	Xtriage
Anisotropy	0.664	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.24 , 229.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.30$ , $\langle L^2 \rangle = 0.14$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.89	EDS
Total number of atoms	28934	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	304.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.32% 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: DG3, ZN, DDG, 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	AAA	0.63	0/1809	0.73	0/2450
1	BBB	0.64	0/1789	0.73	0/2425
2	CCC	0.62	0/10745	0.80	3/14499 (0.0%)
3	DDD	0.63	0/10729	0.78	0/14487
4	EEE	0.62	0/629	0.79	0/847
5	FFF	0.64	0/2233	0.63	0/3010
6	111	0.27	0/665	0.66	0/1022
7	222	0.26	0/729	0.62	0/1121
8	333	0.23	0/113	0.54	0/174
All	All	0.62	0/29441	0.76	3/40035 (0.0%)

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	CCC	0	1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	CCC	1088	ASP	CB-CA-C	-6.58	97.24	110.40
2	CCC	454	ARG	CB-CA-C	-5.48	99.44	110.40
2	CCC	1048	LYS	CB-CA-C	-5.36	99.69	110.40

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	CCC	1282	GLY	Peptide

## 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	AAA	1787	0	1813	82	0
1	BBB	1767	0	1789	74	0
2	CCC	10576	0	10591	422	0
3	DDD	10568	0	10781	455	0
4	EEE	627	0	634	17	0
5	FFF	2205	0	2249	116	0
6	111	595	0	329	30	0
7	222	652	0	364	39	0
8	333	102	0	54	5	0
9	DDD	2	0	0	2	0
10	DDD	2	0	0	0	0
11	DDD	30	0	12	11	0
12	333	21	0	12	3	0
All	All	28934	0	28628	1068	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 19.

The worst 5 of 1068 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:DDD:825:VAL:HG11	3:DDD:1242:ARG:NH1	1.51	1.24
3:DDD:572:THR:HG21	3:DDD:589:TYR:OH	1.35	1.19
5:FFF:266:LEU:HB3	5:FFF:312:ARG:HD2	1.29	1.13
5:FFF:266:LEU:HB3	5:FFF:312:ARG:CD	1.81	1.10
3:DDD:825:VAL:CG1	3:DDD:1242:ARG:HH12	1.70	1.04

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	AAA	228/242 (94%)	209 (92%)	14 (6%)	5 (2%)	6	38
1	BBB	226/242 (93%)	207 (92%)	13 (6%)	6 (3%)	5	34
2	CCC	1339/1342 (100%)	1239 (92%)	73 (6%)	27 (2%)	7	40
3	DDD	1360/1407 (97%)	1251 (92%)	92 (7%)	17 (1%)	12	48
4	EEE	77/90 (86%)	73 (95%)	4 (5%)	0	100	100
5	FFF	266/336 (79%)	247 (93%)	18 (7%)	1 (0%)	34	72
All	All	3496/3659 (96%)	3226 (92%)	214 (6%)	56 (2%)	9	45

5 of 56 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	BBB	117	HIS
1	BBB	193	GLU
1	BBB	194	GLN
2	CCC	46	GLN
2	CCC	247	ARG

### 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	AAA	198/208 (95%)	180 (91%)	18 (9%)	9	32
1	BBB	196/208 (94%)	185 (94%)	11 (6%)	21	48
2	CCC	1156/1157 (100%)	1065 (92%)	91 (8%)	12	38

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	DDD	1135/1168 (97%)	1061 (94%)	74 (6%)	17	44
4	EEE	67/74 (90%)	64 (96%)	3 (4%)	27	54
5	FFF	235/292 (80%)	225 (96%)	10 (4%)	29	55
All	All	2987/3107 (96%)	2780 (93%)	207 (7%)	15	42

5 of 207 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	CCC	817	LEU
2	CCC	1286	THR
3	DDD	1309	ILE
2	CCC	876	GLU
2	CCC	1135	GLN

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

### 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 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	$\# Z  > 2$
11	DG3	DDD	1505	10	25,32,32	1.10	1 (4%)	28,50,50	2.28	4 (14%)
12	DDG	333	101	-	17,23,24	1.38	2 (11%)	15,33,36	2.62	7 (46%)

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
11	DG3	DDD	1505	10	-	2/18/31/31	0/3/3/3
12	DDG	333	101	-	-	0/3/18/19	0/3/3/3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
12	333	101	DDG	C6-C5	4.60	1.49	1.41
11	DDD	1505	DG3	C6-N1	3.99	1.40	1.33
12	333	101	DDG	C5-C4	2.56	1.47	1.40

The worst 5 of 11 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
11	DDD	1505	DG3	C5-C6-N1	-8.74	111.47	123.43
11	DDD	1505	DG3	C6-N1-C2	5.87	125.25	115.93
12	333	101	DDG	C2-N3-C4	4.64	120.66	115.36
12	333	101	DDG	C5-C6-N1	-4.29	117.56	123.43
12	333	101	DDG	C6-N1-C2	3.63	121.70	115.93

There are no chirality outliers.

All (2) torsion outliers are listed below:

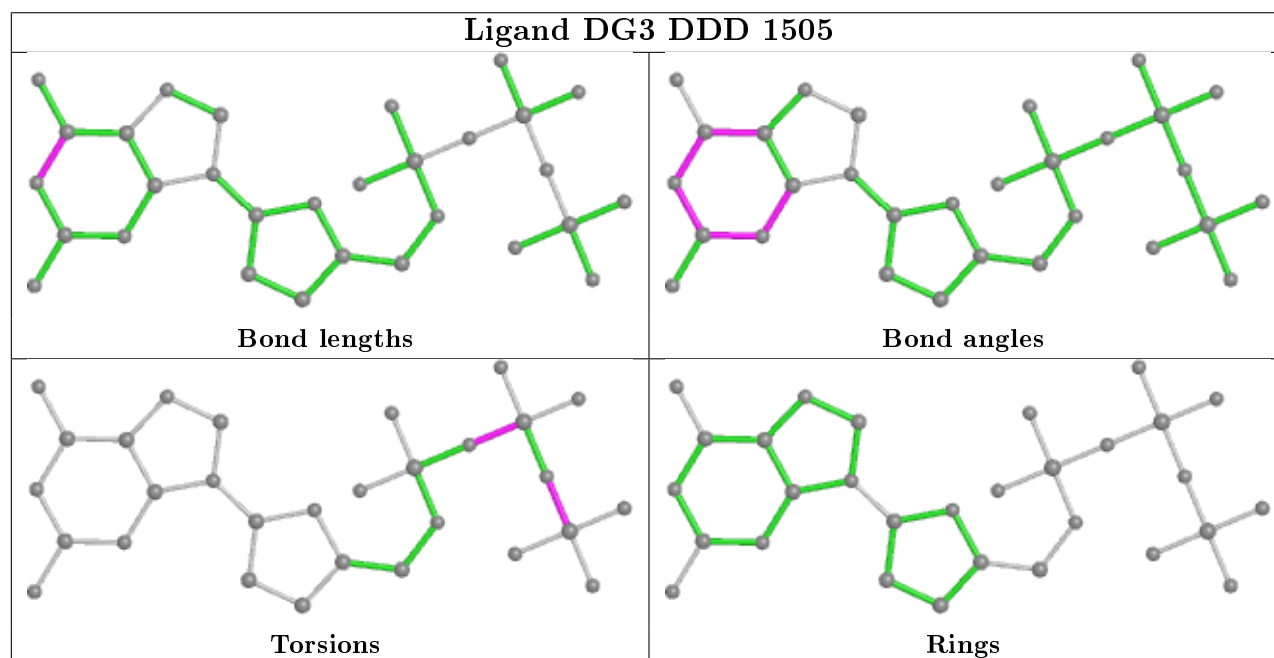
Mol	Chain	Res	Type	Atoms
11	DDD	1505	DG3	PB-O3B-PG-O3G
11	DDD	1505	DG3	PA-O3A-PB-O1B

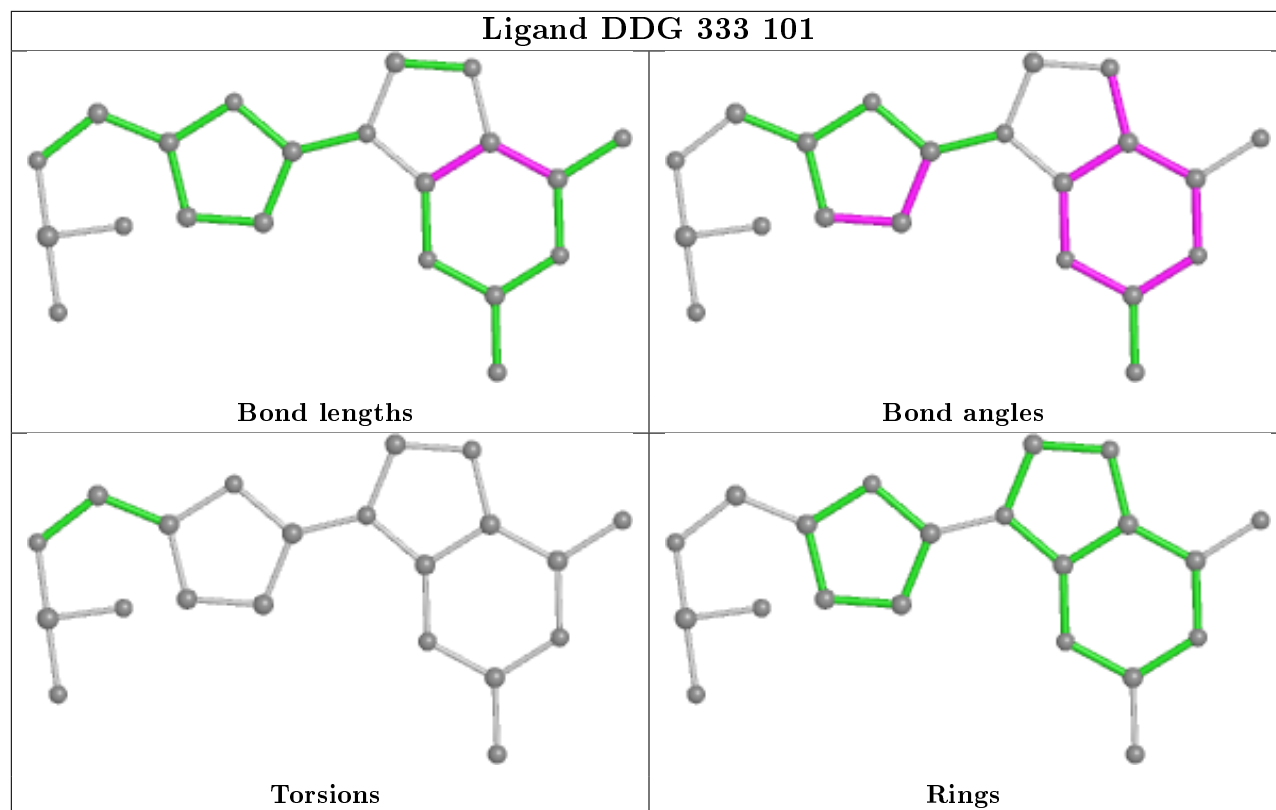
There are no ring outliers.

2 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
11	DDD	1505	DG3	11	0
12	333	101	DDG	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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å <sup>2</sup> )	Q<0.9
1	AAA	230/242 (95%)	0.36	19 (8%)	11 10	235, 332, 394, 420	0
1	BBB	228/242 (94%)	0.17	12 (5%)	26 23	219, 299, 405, 493	0
2	CCC	1341/1342 (99%)	0.00	43 (3%)	47 37	135, 265, 411, 503	0
3	DDD	1362/1407 (96%)	0.17	99 (7%)	15 12	138, 283, 458, 547	0
4	EEE	79/90 (87%)	0.07	3 (3%)	40 32	208, 315, 468, 542	0
5	FFF	270/336 (80%)	0.33	16 (5%)	22 19	263, 376, 482, 500	0
6	111	29/50 (58%)	0.42	4 (13%)	2 3	304, 372, 557, 600	0
7	222	32/50 (64%)	1.08	9 (28%)	0 1	256, 367, 577, 639	0
8	333	5/6 (83%)	0.99	1 (20%)	1 1	340, 354, 386, 388	0
All	All	3576/3765 (94%)	0.14	206 (5%)	23 19	135, 295, 457, 639	0

The worst 5 of 206 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	DDD	748	ALA	11.2
3	DDD	1059	LEU	6.9
7	222	36	DG	6.5
2	CCC	1004	ASP	6.4
2	CCC	1001	GLY	6.3

### 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](#)

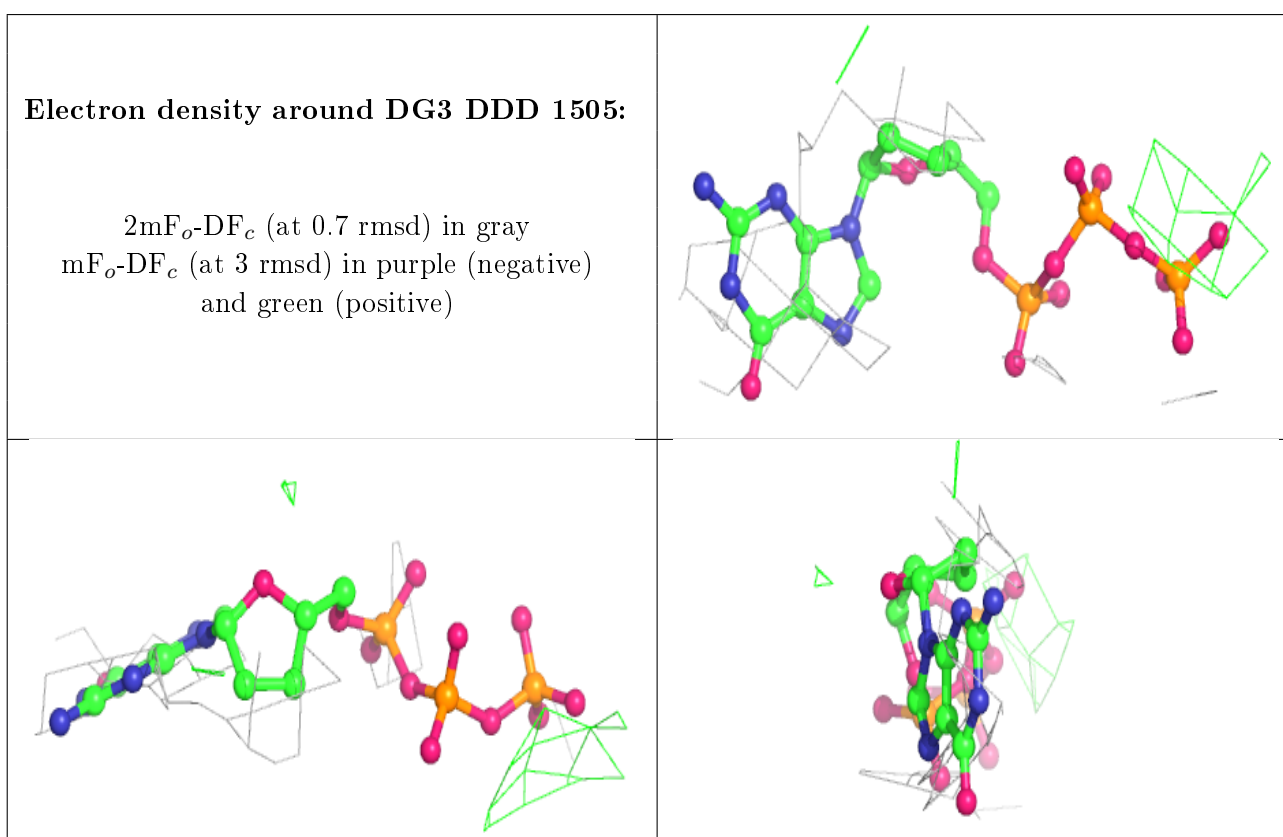
There are no monosaccharides in this entry.

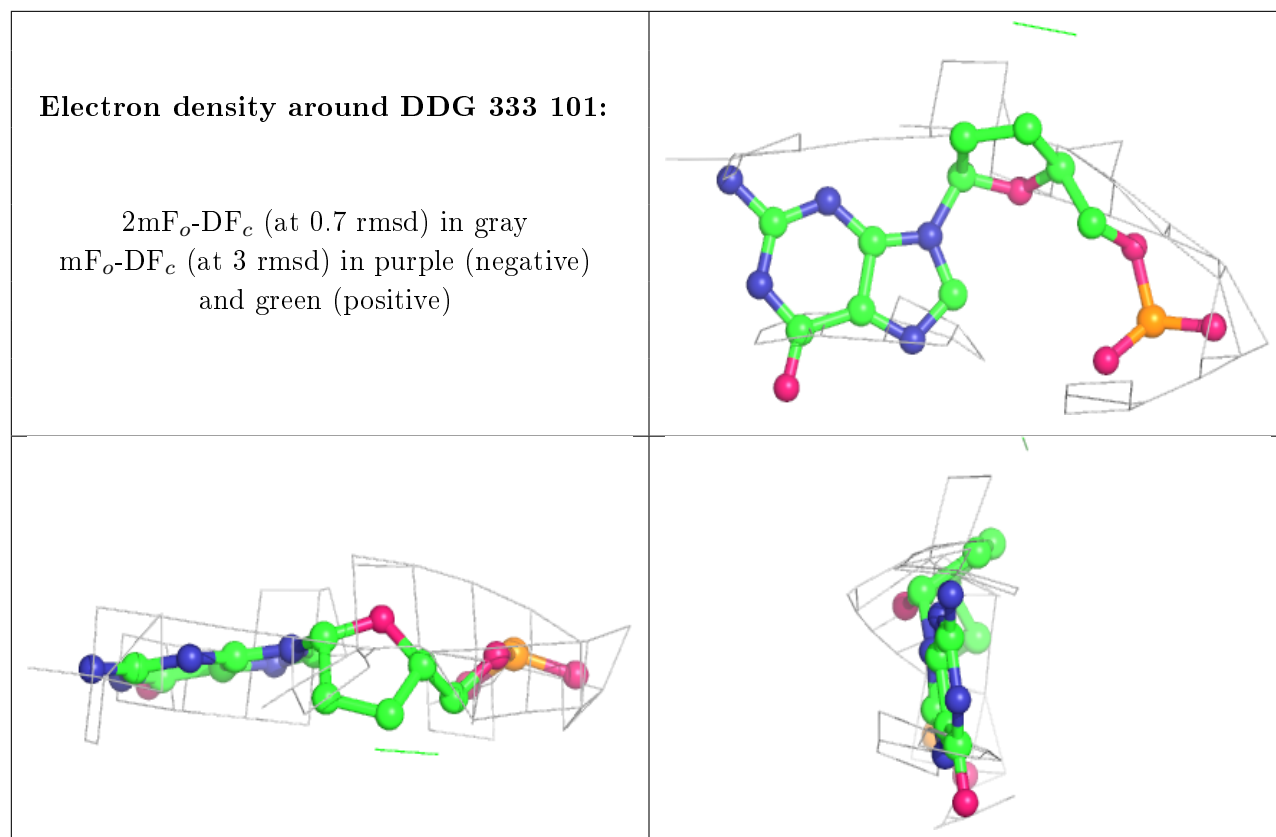
## 6.4 Ligands ⓘ

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
9	ZN	DDD	1501	1/1	0.73	0.05	406,406,406,406	0
11	DG3	DDD	1505	30/30	0.74	0.48	310,341,366,385	0
12	DDG	333	101	21/22	0.81	0.43	294,315,317,318	0
10	MG	DDD	1503	1/1	0.88	0.36	149,149,149,149	0
9	ZN	DDD	1502	1/1	0.93	0.13	348,348,348,348	0
10	MG	DDD	1504	1/1	0.99	0.31	113,113,113,113	0

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





## 6.5 Other polymers [i](#)

There are no such residues in this entry.