



wwPDB X-ray Structure Validation Summary Report ⓘ

Oct 15, 2023 – 05:42 AM EDT

PDB ID : 7M0D
Title : Pre-catalytic quaternary complex of DNA Polymerase Lambda with bound complementary DSB substrate and incoming dUMPNPP
Authors : Kaminski, A.M.; Bebenek, K.; Pedersen, L.C.; Kunkel, T.A.
Deposited on : 2021-03-10
Resolution : 1.80 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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtrriage (Phenix) : 1.13
EDS : 2.36
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) : 2.36

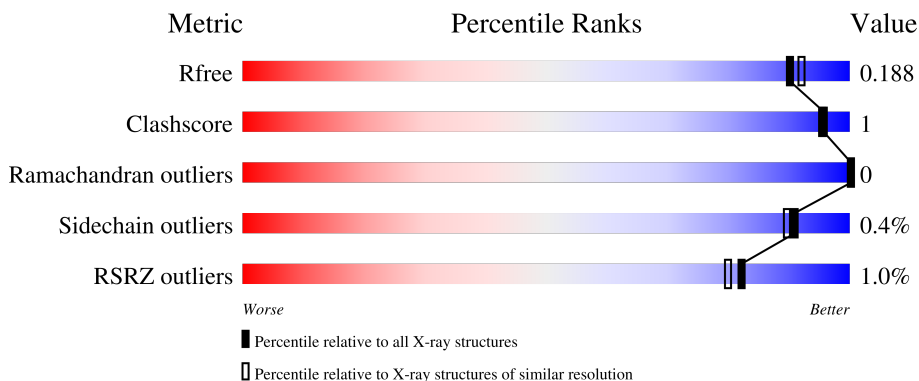
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

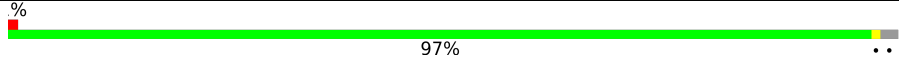
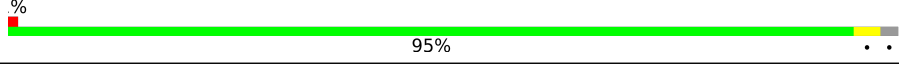
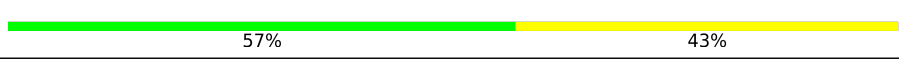
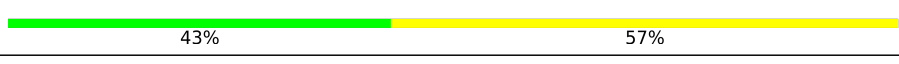

The reported resolution of this entry is 1.80 Å.

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	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

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

Mol	Chain	Length	Quality of chain
1	A	346	 97%
1	B	346	 95%
2	G	7	 57% 43%
2	K	7	 43% 57%
3	F	6	 83% 17%

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Mol	Chain	Length	Quality of chain
3	J	6	 83% 17%
4	H	4	 75% 25%
4	L	4	 75% 25%
5	E	4	 100%
5	I	4	 75% 25%

2 Entry composition [i](#)

There are 15 unique types of molecules in this entry. The entry contains 7458 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 lambda.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	339	2664	1672	486	493	13	0	3	0
1	B	340	2691	1690	493	495	13	0	6	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	230	GLY	-	expression tag	UNP Q9UGP5
A	231	SER	-	expression tag	UNP Q9UGP5
A	232	ALA	-	expression tag	UNP Q9UGP5
A	233	ALA	-	expression tag	UNP Q9UGP5
B	230	GLY	-	expression tag	UNP Q9UGP5
B	231	SER	-	expression tag	UNP Q9UGP5
B	232	ALA	-	expression tag	UNP Q9UGP5
B	233	ALA	-	expression tag	UNP Q9UGP5

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	G	7	141	67	29	39	6	0	0	0
2	K	7	141	67	29	39	6	0	0	0

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
3	F	6	120	58	23	34	5	0	0	0
3	J	6	120	58	23	34	5	0	0	0

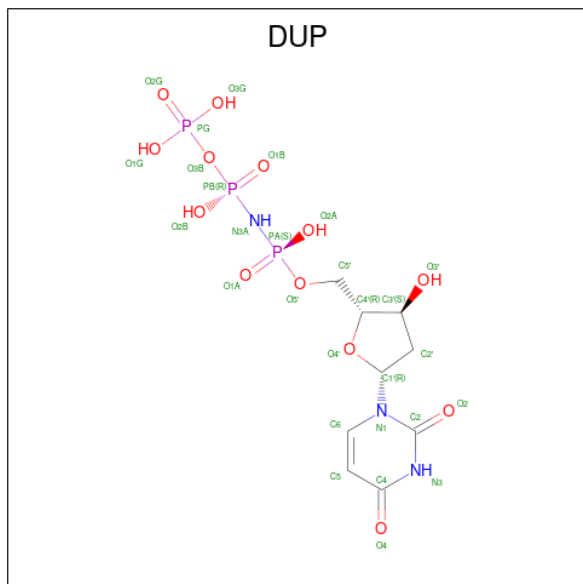
- Molecule 4 is a DNA chain called DNA (5'-D(P*GP*CP*CP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	H	4	Total	C	N	O	P	0	0	0
			83	38	16	25	4			
4	L	4	Total	C	N	O	P	0	0	0
			83	38	16	25	4			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	I	4	Total	C	N	O	P	0	0	0
			79	39	15	22	3			
5	E	4	Total	C	N	O	P	0	0	0
			79	39	15	22	3			

- Molecule 6 is 2'-DEOXYURIDINE 5'-ALPHA,BETA-IMIDO-TRIPHOSPHATE (three-letter code: DUP) (formula: C₉H₁₆N₃O₁₃P₃) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 7 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	2	Total Na 2 2	0	0
7	B	2	Total Na 2 2	0	0
7	K	1	Total Na 1 1	0	0

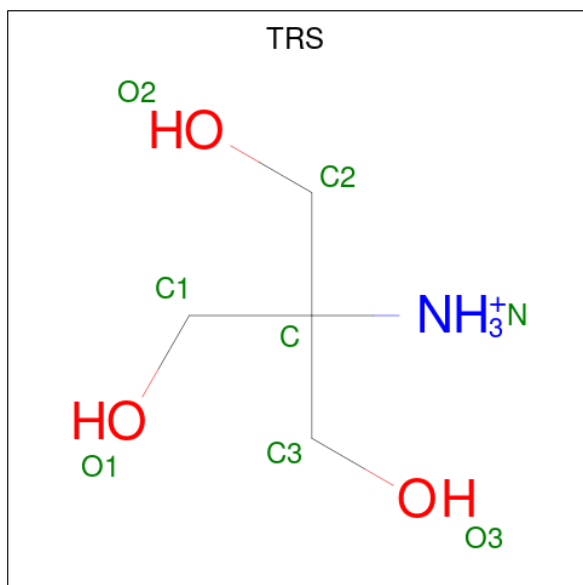
- Molecule 8 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	1	Total Mg 1 1	0	0
8	B	1	Total Mg 1 1	0	0

- Molecule 9 is POTASSIUM ION (three-letter code: K) (formula: K).

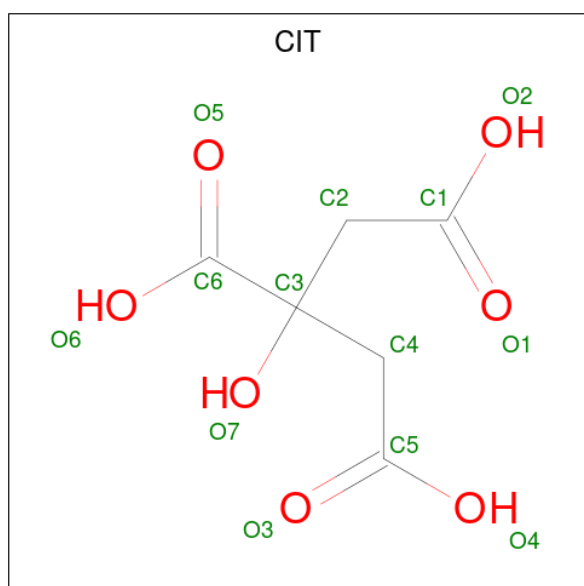
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	A	3	Total K 3 3	0	0
9	B	3	Total K 3 3	0	0

- Molecule 10 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (three-letter code: TRS) (formula: C₄H₁₂NO₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
10	A	1	Total	C	N	O	0	0
			8	4	1	3		
10	A	1	Total	C	N	O	0	1
			8	4	1	3		
10	B	1	Total	C	N	O	0	0
			8	4	1	3		
10	J	1	Total	C	N	O	0	1
			8	4	1	3		

- Molecule 11 is CITRIC ACID (three-letter code: CIT) (formula: C₆H₈O₇).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
11	A	1	Total	C	O	0	0
			13	6	7		
11	B	1	Total	C	O	0	1
			13	6	7		

- Molecule 12 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).

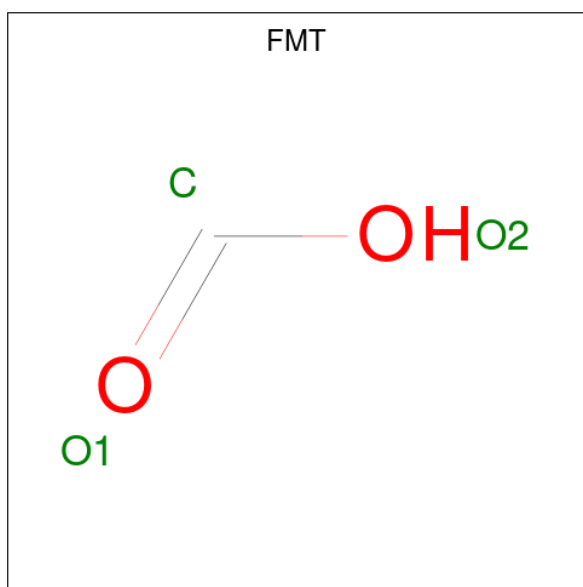


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
12	A	1	Total C O 4 2 2	0	0
12	A	1	Total C O 4 2 2	0	0
12	A	1	Total C O 4 2 2	0	0
12	A	1	Total C O 4 2 2	0	0
12	F	1	Total C O 4 2 2	0	0
12	B	1	Total C O 4 2 2	0	0
12	B	1	Total C O 4 2 2	0	0
12	J	1	Total C O 4 2 2	0	0

- Molecule 13 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
13	B	1	Total Cl 1 1	0	0

- Molecule 14 is FORMIC ACID (three-letter code: FMT) (formula: CH₂O₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
14	B	1	Total C O 3 1 2	0	0
14	B	1	Total C O 3 1 2	0	0

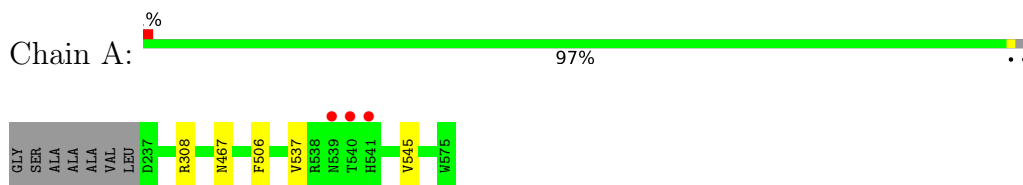
- Molecule 15 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
15	A	459	Total O 466 466	0	11
15	G	18	Total O 18 18	0	0
15	F	23	Total O 24 24	0	2
15	H	17	Total O 17 17	0	0
15	B	463	Total O 469 469	0	9
15	K	19	Total O 19 19	0	0
15	J	20	Total O 20 20	0	1
15	L	17	Total O 18 18	0	1
15	I	19	Total O 19 19	0	0
15	E	21	Total O 21 21	0	0

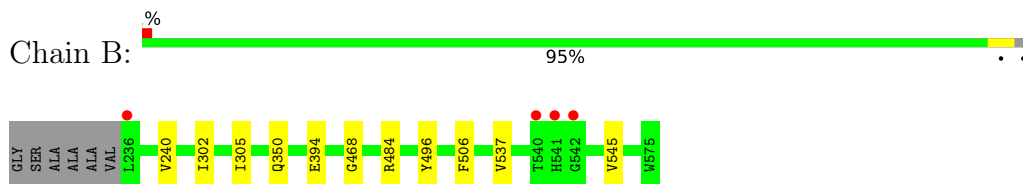
3 Residue-property plots [i](#)

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 polymerase lambda



- Molecule 1: DNA polymerase lambda



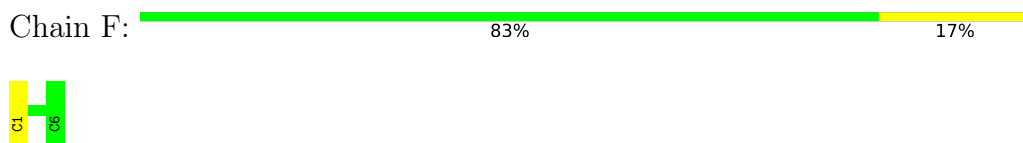
- Molecule 2: DNA (5'-D(*CP*GP*GP*CP*AP*GP*C)-3')




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



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




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

Chain J:  83% 17%




- Molecule 4: DNA (5'-D(P*GP*CP*CP*G)-3')

Chain H:  75% 25%



- Molecule 4: DNA (5'-D(P*GP*CP*CP*G)-3')

Chain L:  75% 25%



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

Chain I:  75% 25%



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

Chain E:  100%

There are no outlier residues recorded for this chain.

4 Data and refinement statistics i

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	95.44Å 151.99Å 86.47Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.78 – 1.80 41.78 – 1.80	Depositor EDS
% Data completeness (in resolution range)	98.8 (41.78-1.80) 94.1 (41.78-1.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.32 (at 1.81Å)	Xtrriage
Refinement program	PHENIX 1.19_4092	Depositor
R, R_{free}	0.164 , 0.189 0.163 , 0.188	Depositor DCC
R_{free} test set	4945 reflections (4.28%)	wwPDB-VP
Wilson B-factor (Å ²)	16.4	Xtrriage
Anisotropy	0.431	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 48.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7458	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 56.11 % 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 2.8699e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: EDO, TRS, NA, MG, CL, K, DUP, FMT, CIT

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	0.46	0/2720	0.62	0/3675
1	B	0.46	0/2747	0.61	0/3711
2	G	0.93	0/158	1.14	1/242 (0.4%)
2	K	0.94	0/158	1.09	2/242 (0.8%)
3	F	1.05	0/134	1.04	0/205
3	J	0.97	0/134	1.07	0/205
4	H	1.45	1/92 (1.1%)	0.90	0/138
4	L	1.45	1/92 (1.1%)	0.89	0/138
5	E	0.95	0/88	1.02	0/134
5	I	1.21	0/88	1.20	1/134 (0.7%)
All	All	0.60	2/6411 (0.0%)	0.71	4/8824 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	H	1	DG	OP3-P	-10.93	1.48	1.61
4	L	1	DG	OP3-P	-10.21	1.49	1.61

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	K	7	DC	O5'-P-OP1	-6.51	99.84	105.70
2	G	7	DC	O5'-P-OP2	-6.06	100.25	105.70
5	I	3	DT	O4'-C1'-N1	-5.48	104.17	108.00
2	K	5	DA	O4'-C1'-N9	-5.05	104.47	108.00

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	2664	0	2628	4	0
1	B	2691	0	2653	9	0
2	G	141	0	79	1	0
2	K	141	0	79	1	0
3	F	120	0	69	1	0
3	J	120	0	69	1	0
4	H	83	0	45	0	0
4	L	83	0	45	0	0
5	E	79	0	47	0	0
5	I	79	0	47	0	0
6	A	28	0	12	0	0
6	B	28	0	12	0	0
7	A	2	0	0	0	0
7	B	2	0	0	0	0
7	K	1	0	0	0	0
8	A	1	0	0	0	0
8	B	1	0	0	0	0
9	A	3	0	0	0	0
9	B	3	0	0	0	0
10	A	16	0	24	0	0
10	B	8	0	12	0	0
10	J	8	0	12	0	0
11	A	13	0	5	0	0
11	B	13	0	5	1	0
12	A	16	0	24	0	0
12	B	8	0	12	1	0
12	F	4	0	4	0	0
12	J	4	0	4	0	0
13	B	1	0	0	0	0
14	B	6	0	2	0	0
15	A	466	0	0	0	0
15	B	469	0	0	2	0
15	E	21	0	0	0	0
15	F	24	0	0	0	0
15	G	18	0	0	0	0
15	H	17	0	0	1	0
15	I	19	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
15	J	20	0	0	0	0
15	K	19	0	0	0	0
15	L	18	0	0	0	0
All	All	7458	0	5889	14	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 1.

The worst 5 of 14 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:545:VAL:O	1:B:545:VAL:HG22	2.05	0.56
1:B:350[A]:GLN:HG3	15:B:815:HOH:O	2.13	0.47
1:B:394:GLU:OE1	1:B:484:ARG:NH2	2.48	0.47
1:B:302:ILE:HB	1:B:305:ILE:HD12	1.96	0.46
3:F:1:DC:H5''	1:B:537:VAL:HG22	1.97	0.46

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	340/346 (98%)	337 (99%)	3 (1%)	0	100	100
1	B	344/346 (99%)	340 (99%)	4 (1%)	0	100	100
All	All	684/692 (99%)	677 (99%)	7 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	280/288 (97%)	279 (100%)	1 (0%)	91	89
1	B	281/288 (98%)	280 (100%)	1 (0%)	91	89
All	All	561/576 (97%)	559 (100%)	2 (0%)	91	89

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	506	PHE
1	B	506	PHE

Sometimes 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 32 ligands modelled in this entry, 14 are monoatomic - leaving 18 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	CIT	B	710[A]	-	12,12,12	1.03	0	17,17,17	1.62	2 (11%)
12	EDO	A	712	-	3,3,3	0.51	0	2,2,2	0.28	0
12	EDO	B	712	-	3,3,3	0.45	0	2,2,2	0.34	0
10	TRS	A	708	-	7,7,7	0.41	0	9,9,9	0.44	0
12	EDO	A	714	-	3,3,3	0.57	0	2,2,2	0.52	0
10	TRS	A	709[A]	-	7,7,7	0.42	0	9,9,9	1.09	1 (11%)
12	EDO	A	713	-	3,3,3	0.53	0	2,2,2	0.42	0
12	EDO	F	101	7,9	3,3,3	0.57	0	2,2,2	0.14	0
11	CIT	A	710	-	12,12,12	1.01	0	17,17,17	1.56	2 (11%)
14	FMT	B	713	-	2,2,2	0.63	0	1,1,1	0.49	0
12	EDO	J	102	7,9	3,3,3	0.51	0	2,2,2	0.17	0
12	EDO	B	711	-	3,3,3	0.61	0	2,2,2	0.11	0
14	FMT	B	714	-	2,2,2	0.62	0	1,1,1	0.37	0
12	EDO	A	711	-	3,3,3	0.53	0	2,2,2	0.28	0
10	TRS	B	709	-	7,7,7	0.35	0	9,9,9	0.75	0
6	DUP	B	701	7,8	25,29,29	1.31	4 (16%)	30,45,45	1.48	4 (13%)
10	TRS	J	101[A]	-	7,7,7	0.42	0	9,9,9	1.24	1 (11%)
6	DUP	A	701	7,8	25,29,29	1.52	6 (24%)	30,45,45	1.40	5 (16%)

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	CIT	B	710[A]	-	-	2/16/16/16	-
12	EDO	A	712	-	-	0/1/1/1	-
12	EDO	B	712	-	-	1/1/1/1	-
10	TRS	A	708	-	-	0/9/9/9	-
12	EDO	A	714	-	-	0/1/1/1	-
10	TRS	A	709[A]	-	-	6/9/9/9	-
12	EDO	A	713	-	-	0/1/1/1	-
12	EDO	F	101	7,9	-	0/1/1/1	-
11	CIT	A	710	-	-	6/16/16/16	-
12	EDO	J	102	7,9	-	1/1/1/1	-
12	EDO	B	711	-	-	1/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
12	EDO	A	711	-	-	0/1/1/1	-
10	TRS	B	709	-	-	0/9/9/9	-
6	DUP	B	701	7,8	-	3/16/34/34	0/2/2/2
10	TRS	J	101[A]	-	-	6/9/9/9	-
6	DUP	A	701	7,8	-	3/16/34/34	0/2/2/2

The worst 5 of 10 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	701	DUP	PA-O2A	-3.44	1.47	1.56
6	A	701	DUP	C6-N1	2.80	1.39	1.35
6	B	701	DUP	PA-O2A	-2.62	1.49	1.56
6	A	701	DUP	PB-O2B	-2.55	1.49	1.56
6	B	701	DUP	C6-N1	2.31	1.38	1.35

The worst 5 of 15 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
11	B	710[A]	CIT	O6-C6-C3	4.23	120.40	113.05
6	B	701	DUP	O2A-PA-O1A	4.14	118.60	109.92
11	A	710	CIT	O6-C6-C3	4.11	120.19	113.05
6	B	701	DUP	O2B-PB-O1B	3.98	118.27	109.92
6	A	701	DUP	O2B-PB-O1B	3.62	117.50	109.92

There are no chirality outliers.

5 of 29 torsion outliers are listed below:

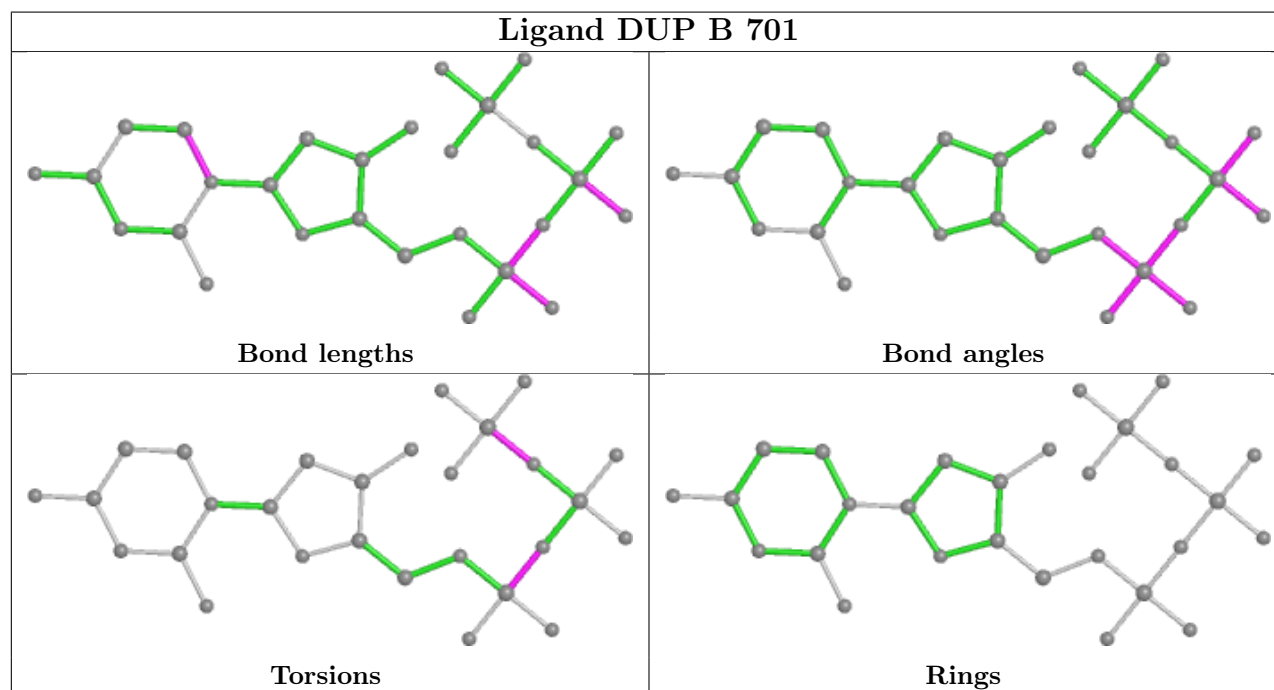
Mol	Chain	Res	Type	Atoms
6	A	701	DUP	PA-N3A-PB-O1B
6	A	701	DUP	PB-O3B-PG-O1G
6	B	701	DUP	PB-N3A-PA-O1A
6	B	701	DUP	PB-O3B-PG-O1G
10	A	709[A]	TRS	C2-C-C1-O1

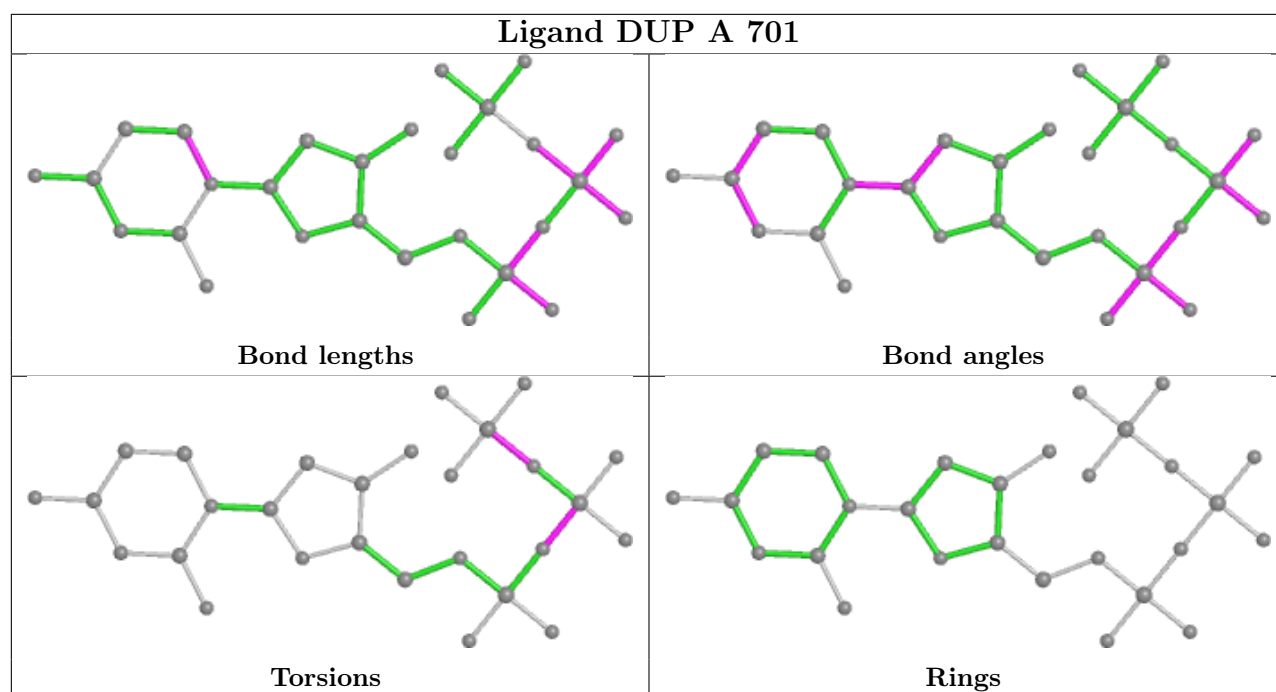
There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
11	B	710[A]	CIT	1	0
12	B	712	EDO	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 [i](#)

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	339/346 (97%)	-0.48	3 (0%) 84 82	10, 19, 35, 60	0
1	B	340/346 (98%)	-0.47	4 (1%) 79 76	10, 18, 34, 61	0
2	G	7/7 (100%)	-0.35	0 100 100	16, 22, 39, 44	0
2	K	7/7 (100%)	-0.48	0 100 100	15, 20, 38, 42	0
3	F	6/6 (100%)	-0.55	0 100 100	14, 16, 36, 40	0
3	J	6/6 (100%)	-0.58	0 100 100	13, 15, 33, 36	0
4	H	4/4 (100%)	-0.76	0 100 100	19, 20, 29, 38	0
4	L	4/4 (100%)	-0.83	0 100 100	20, 22, 29, 40	0
5	E	4/4 (100%)	-0.65	0 100 100	22, 23, 23, 32	0
5	I	4/4 (100%)	-0.72	0 100 100	19, 20, 22, 34	0
All	All	721/734 (98%)	-0.48	7 (0%) 82 80	10, 19, 36, 61	0

The worst 5 of 7 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	540	THR	4.3
1	B	236	LEU	3.7
1	A	540	THR	3.5
1	B	541	HIS	3.2
1	B	542	GLY	3.1

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

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

6.3 Carbohydrates [i](#)

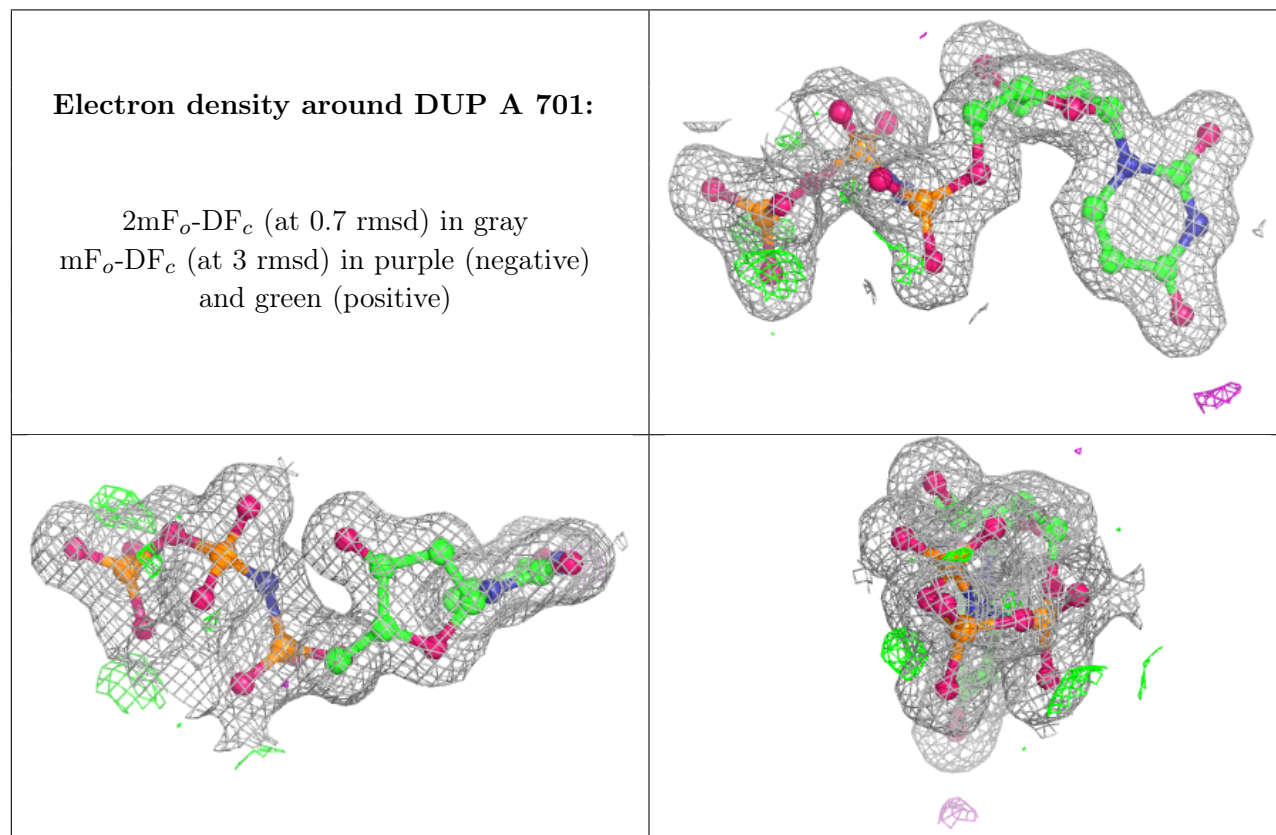
There are no monosaccharides in this entry.

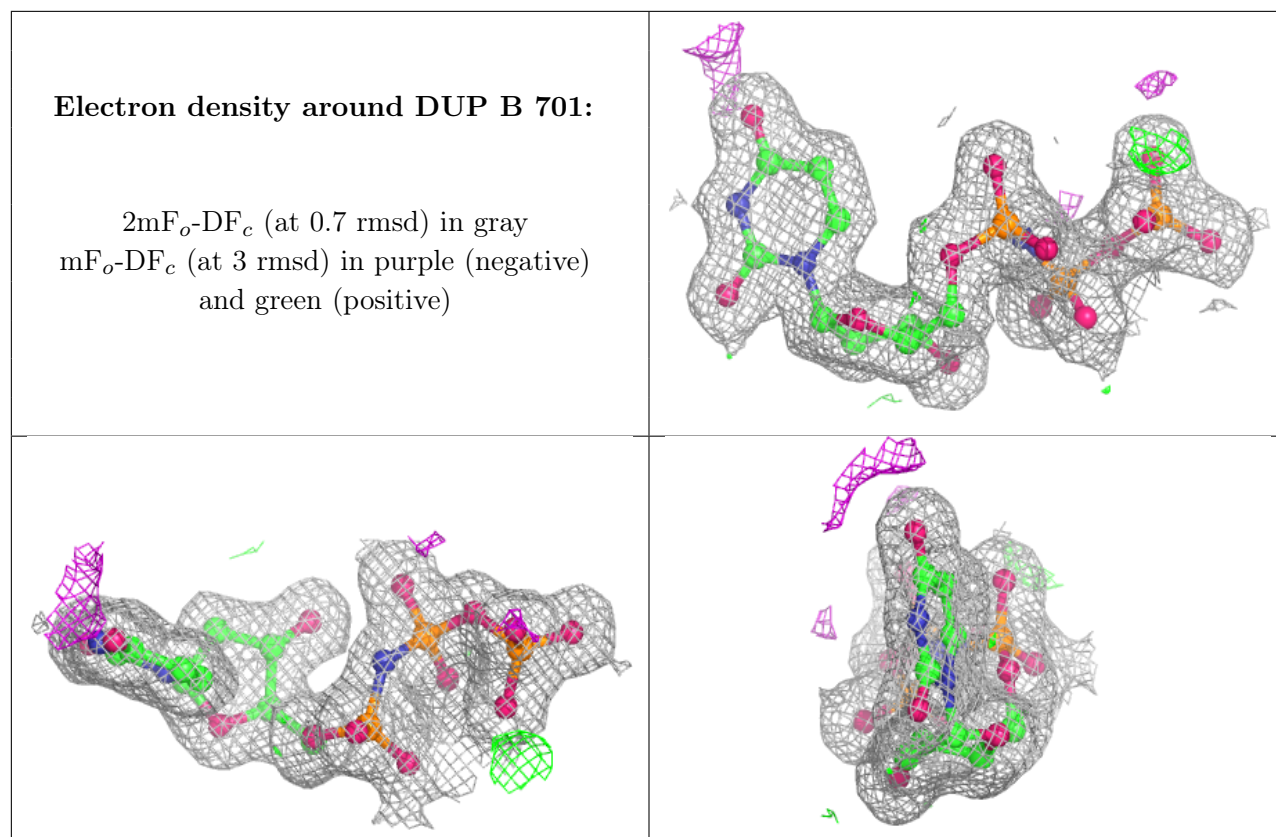
6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
10	TRS	B	709	8/8	0.84	0.15	13,19,22,29	8
14	FMT	B	713	3/3	0.84	0.13	29,29,36,40	0
10	TRS	A	708	8/8	0.87	0.14	15,21,26,29	0
14	FMT	B	714	3/3	0.87	0.21	39,39,44,48	0
12	EDO	B	711	4/4	0.88	0.15	20,29,32,33	0
12	EDO	F	101	4/4	0.90	0.16	19,22,28,31	0
10	TRS	J	101[A]	8/8	0.90	0.18	15,17,21,22	8
11	CIT	A	710	13/13	0.91	0.12	13,23,29,34	13
12	EDO	A	713	4/4	0.91	0.16	22,26,29,33	0
10	TRS	A	709[A]	8/8	0.91	0.19	14,20,25,28	8
11	CIT	B	710[A]	13/13	0.92	0.14	19,22,31,31	13
12	EDO	A	714	4/4	0.92	0.11	27,27,28,39	0
12	EDO	A	712	4/4	0.93	0.13	22,23,30,32	0
12	EDO	J	102	4/4	0.94	0.12	22,22,23,27	0
13	CL	B	708	1/1	0.94	0.09	25,25,25,25	1
12	EDO	A	711	4/4	0.94	0.12	23,29,30,39	0
7	NA	K	101	1/1	0.94	0.09	29,29,29,29	0
12	EDO	B	712	4/4	0.96	0.16	25,27,32,45	0
6	DUP	A	701	28/28	0.99	0.09	10,12,14,19	0
8	MG	A	703	1/1	0.99	0.08	12,12,12,12	0
8	MG	B	703	1/1	0.99	0.09	12,12,12,12	0
9	K	A	705	1/1	0.99	0.06	20,20,20,20	0
9	K	A	706	1/1	0.99	0.04	18,18,18,18	1
9	K	A	707	1/1	0.99	0.04	17,17,17,17	0
9	K	B	705	1/1	0.99	0.04	20,20,20,20	0
9	K	B	706	1/1	0.99	0.04	21,21,21,21	1
9	K	B	707	1/1	0.99	0.04	16,16,16,16	0
6	DUP	B	701	28/28	0.99	0.09	8,11,14,15	0
7	NA	A	702	1/1	0.99	0.07	15,15,15,15	0
7	NA	B	702	1/1	0.99	0.07	14,14,14,14	0
7	NA	B	704	1/1	0.99	0.06	10,10,10,10	0
7	NA	A	704	1/1	1.00	0.09	10,10,10,10	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.