

Full wwPDB X-ray Structure Validation Report (i)

Oct 10, 2023 – 12:36 AM EDT

PDB ID	:	7M4I
Title	:	DNA Polymerase Lambda, dCTP:At Mn2+ Product State Ternary Complex,
		420 min
Authors	:	Jamsen, J.A.; Wilson, S.H.
Deposited on	:	2021-03-21
Resolution	:	2.00 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org A user guide is available at https://www.wwpdb.org/validation/2017/XrayValidationReportHelp with specific help available everywhere you see the (i) 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 (1)) were used in the production of this report:

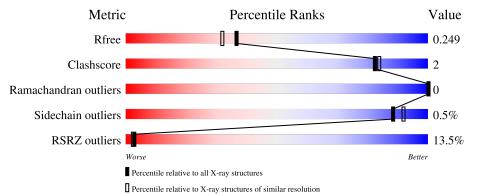
Xtriage (Phenix) EDS buster-report Percentile statistics Refmac CCP4 Ideal geometry (proteins) Ideal geometry (DNA, RNA)	: : : : :	20191225.v01 (using entries in the PDB archive December 25th 2019) 5.8.0158 7.0.044 (Gargrove) Engh & Huber (2001) Parkinson et al. (1996)
Ideal geometry (DNA, RNA) Validation Pipeline (wwPDB-VP)		Parkinson et al. (1996) 2.35.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\text{-}RAY \, DIFFRACTION$

The reported resolution of this entry is 2.00 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	$\begin{array}{c} \textbf{Whole archive} \\ (\#\textbf{Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
R_{free}	130704	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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 <=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	А	329	91%	6% ·
2	Т	11	73%	27%
3	Р	7	100%	
4	D	4	75%	25%



2 Entry composition (i)

There are 9 unique types of molecules in this entry. The entry contains 2938 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	
1	А	318	Total 2303	C 1457	N 412	0 423	S 11	0	4	0

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
А	463	LYS	SER	conflict	UNP Q9UGP5
А	464	GLY	GLN	conflict	UNP Q9UGP5
А	?	-	GLU	deletion	UNP Q9UGP5
А	?	-	ASN	deletion	UNP Q9UGP5
А	?	-	GLY	deletion	UNP Q9UGP5
А	?	-	GLN	deletion	UNP Q9UGP5
А	?	-	GLN	deletion	UNP Q9UGP5
А	471	THR	GLN	conflict	UNP Q9UGP5
А	543	ALA	CYS	engineered mutation	UNP Q9UGP5

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	Т	11	Total 306	C 146	N 58	O 88	Р 14	0	4	0

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

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
3	Р	7	Total 138	C 67	N 26	O 39	P 6	0	0	0

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



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	Л	4	Total	С	Ν	Ο	Р	0	0	Ο
4	D	1	83	38	16	25	4	0	0	0

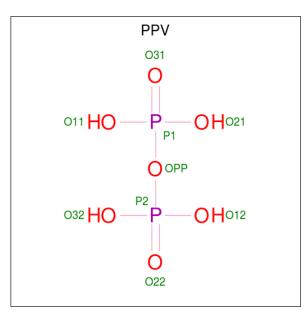
• Molecule 5 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	А	3	Total Mn 3 3	0	0

• Molecule 6 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	А	3	Total Na 3 3	0	0

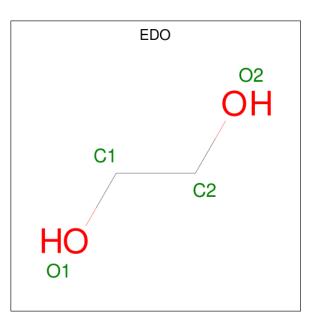
• Molecule 7 is PYROPHOSPHATE (three-letter code: PPV) (formula: $H_4O_7P_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	А	1	Total 9	O 7	Р 2	0	0

• Molecule 8 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).





Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	А	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
8	А	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 3 & 2 & 1 \end{array}$	0	0
8	А	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0

• Molecule 9 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	А	57	$\begin{array}{cc} \text{Total} & \text{O} \\ 57 & 57 \end{array}$	0	1
9	Т	10	Total O 10 10	0	2
9	Р	15	Total O 15 15	0	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.

Chain A:	91%	6% ·
ALA ALA PRO PRO SER SER GLN GLN CVS ALA ASA ASA ASA ASA ASA ASA ASA ASA AS	125 / 125 / 125 / 1226 1226 1228 1228 1228 1228 1228 1228	P303 P304 P304 P310 P317
R419 V430 L431 S439 S439 R452 L457 L457 C464	1474 1490 1491 1491 1492 1503 1504 8510 8511 8531 8531 8531 8531 8535 8531 8535 8531 8536 8531 8536 8531 8536 8537 8536 8536 8536 8537 8536 8536 8536 8536 8536 8536 8536 8536	H541 6542 A543 A543 A543 A544 C546 C548 C548 C548 C548 C548 C548 C548 C565 C565 C565 C565 C565 C565 C565 C56
• Molecule 2: DNA (5'-D(*CP*GP*GP*CP*AP*GP	*TP*AP*CP*TP*G)-3')
Chain T:	73%	27%
C1 17 17 110 111 111 111		
• Molecule 3: DNA (5'-D(*CP*AP*GP*TP*AP*CP	*C)-3')
Chain P:	100%	
There are no outlier	residues recorded for this chain.	
• Molecule 4: DNA ((5'-D(P*GP*CP*CP*G)-3')	
Chain D:	75%	25%
<mark>5</mark> 8		

• Molecule 1: DNA polymerase lambda



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	55.75Å 62.32Å 142.18Å	Deperitor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	37.72 - 2.00	Depositor
Resolution (A)	37.72 - 2.00	EDS
% Data completeness	98.5 (37.72-2.00)	Depositor
(in resolution range)	98.5(37.72-2.00)	EDS
R _{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	$0.70 (at 2.00 \text{\AA})$	Xtriage
Refinement program	PHENIX v1.15.2-3472	Depositor
D D.	0.222 , 0.247	Depositor
R, R_{free}	0.222 , 0.249	DCC
R_{free} test set	1635 reflections (4.87%)	wwPDB-VP
Wilson B-factor $(Å^2)$	46.2	Xtriage
Anisotropy	0.472	Xtriage
Bulk solvent $k_{sol}(e/Å^3), B_{sol}(Å^2)$	0.31,63.3	EDS
L-test for twinning ²	$ < L >=0.48, < L^2>=0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	2938	wwPDB-VP
Average B, all atoms $(Å^2)$	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 5.50% of the height of the origin peak. No significant pseudotranslation is detected.

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.



¹Intensities estimated from amplitudes.

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, PPV, MN, NA

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	Boi	Bond lengths		angles
	Unam	RMSZ	# Z > 5	RMSZ	# Z > 5
1	А	0.25	0/2355	0.44	0/3200
2	Т	0.50	0/343	0.97	0/528
3	Р	0.54	0/154	0.84	0/235
4	D	1.20	1/92~(1.1%)	0.76	0/138
All	All	0.37	1/2944~(0.0%)	0.58	0/4101

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	1	DG	OP3-P	-10.57	1.48	1.61

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	А	2303	0	2087	11	0
2	Т	306	0	170	2	0
3	Р	138	0	80	0	0
4	D	83	0	45	0	0
5	А	3	0	0	0	0
6	А	3	0	0	0	0

Continued on next page...



Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	А	9	0	0	0	0
8	А	11	0	15	0	0
9	А	57	0	0	0	0
9	Р	15	0	0	0	0
9	Т	10	0	0	0	0
All	All	2938	0	2397	12	0

Continued from previous page...

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 2.

All (12) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:519:LEU:CD2	1:A:565:LEU:HD11	2.28	0.64
1:A:269:VAL:O	1:A:346[B]:THR:HG23	2.06	0.56
1:A:531:ALA:HB1	1:A:550:VAL:HG13	1.87	0.55
1:A:434:HIS:ND1	1:A:439:SER:CB	2.70	0.55
1:A:474:LEU:HD22	1:A:490:ASP:CG	2.29	0.53
2:T:6[B]:DG:H2"	2:T:7[B]:DT:O5'	2.10	0.51
1:A:519:LEU:HD22	1:A:565:LEU:HD11	2.00	0.44
1:A:415:CYS:HA	1:A:419:ARG:HB2	2.00	0.43
1:A:407:ASN:OD1	1:A:409:GLY:N	2.45	0.43
1:A:452:ARG:HG3	1:A:457:LEU:HB2	2.01	0.42
1:A:271:GLY:CA	1:A:346[B]:THR:HG21	2.50	0.42
1:A:464:GLY:HA3	2:T:9[B]:DC:O5'	2.21	0.41

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	А	317/329~(96%)	307~(97%)	10 (3%)	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	А	200/274~(73%)	199 (100%)	1 (0%)	88 92

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	А	295	SER

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 10 ligands modelled in this entry, 6 are monoatomic - leaving 4 for Mogul analysis.



7M4I

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).

Mal True		Chain	Chain	Chain	Chain	Chain	Chain	Chain	Chain	Chain	Chain	Chain	Chain	Res	Dec	Dec	Dee	D	Timle	B	Bond lengths			Bond angles		
Mol Type	Type	Unam	nes	Link	Counts	RMSZ	# Z > 2	Counts	RMSZ $ $ $# Z >$	# Z > 2																
8	EDO	А	610	-	3,3,3	0.47	0	2,2,2	0.37	0																
8	EDO	А	608	-	3,3,3	0.45	0	2,2,2	0.35	0																
8	EDO	А	609	-	2,2,3	0.49	0	1,1,2	0.37	0																
7	PPV	А	607	5	6,8,8	0.73	0	13,13,13	1.43	1 (7%)																

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
8	EDO	А	610	-	-	1/1/1/1	-
8	EDO	А	608	-	-	1/1/1/1	-
7	PPV	А	607	5	-	0/6/6/6	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
7	А	607	PPV	P2-OPP-P1	-3.55	120.64	132.83

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
8	А	610	EDO	O1-C1-C2-O2
8	А	608	EDO	O1-C1-C2-O2

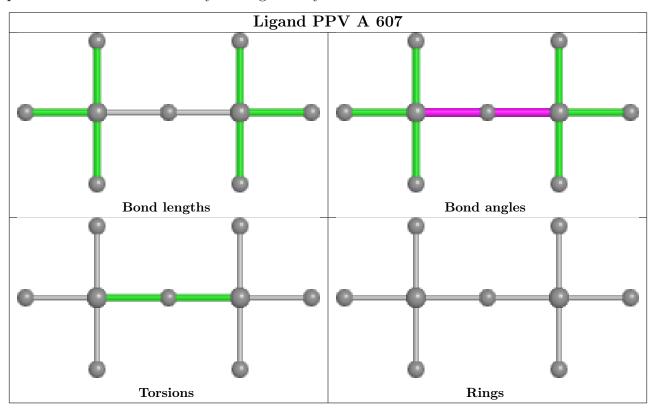
There are no ring outliers.

No monomer is involved in short contacts.

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 then 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^{th} 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 $>$	# RSR	Z>2	$\mathbf{OWAB}(\mathrm{\AA}^2)$	Q<0.9
1	А	318/329~(96%)	0.86	46 (14%)	2 2	37, 56, 99, 129	0
2	Т	11/11~(100%)	0.07	0 100	100	52,56,65,67	0
3	Р	7/7~(100%)	0.40	0 100	100	39, 44, 46, 53	0
4	D	4/4 (100%)	-0.23	0 100	100	63,64,66,71	0
All	All	340/351~(96%)	0.81	46 (13%)	3 2	37, 56, 98, 129	0

All (46) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	А	545	VAL	8.2
1	А	319	SER	6.5
1	А	317	LEU	6.1
1	А	535	ALA	5.7
1	А	536	VAL	4.6
1	А	544	LYS	4.4
1	А	293	VAL	4.2
1	А	537	VAL	4.2
1	А	290	HIS	4.2
1	А	543	ALA	4.0
1	А	542	GLY	3.9
1	А	538	ARG	3.8
1	А	320	GLY	3.8
1	А	255	HIS	3.8
1	А	503	LEU	3.6
1	А	257	THR	3.6
1	А	299	ALA	3.5
1	А	284	ASN	3.3
1	А	321	HIS	3.2
1	А	254	LEU	3.1
1	А	541	HIS	3.1

Continued on next page...



Mol	Chain	Res	Type	RSRZ
1	А	511[A]	HIS	3.0
1	А	547	PRO	3.0
1	А	325	LEU	2.7
1	А	310	ALA	2.7
1	А	548	GLY	2.7
1	А	430	VAL	2.7
1	А	286	LEU	2.6
1	А	289	PHE	2.6
1	А	439	SER	2.5
1	А	309	MET	2.5
1	А	297	GLN	2.5
1	А	292	PRO	2.5
1	А	296	TYR	2.5
1	А	302	ILE	2.5
1	А	304	GLY	2.5
1	А	546	GLY	2.3
1	А	566	PRO	2.3
1	А	415	CYS	2.3
1	А	492	ILE	2.2
1	А	502	ALA	2.2
1	А	431	LEU	2.2
1	А	301	SER	2.2
1	А	504	LEU	2.1
1	А	510	ALA	2.1
1	А	253	ASN	2.0

Continued from previous page...

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)

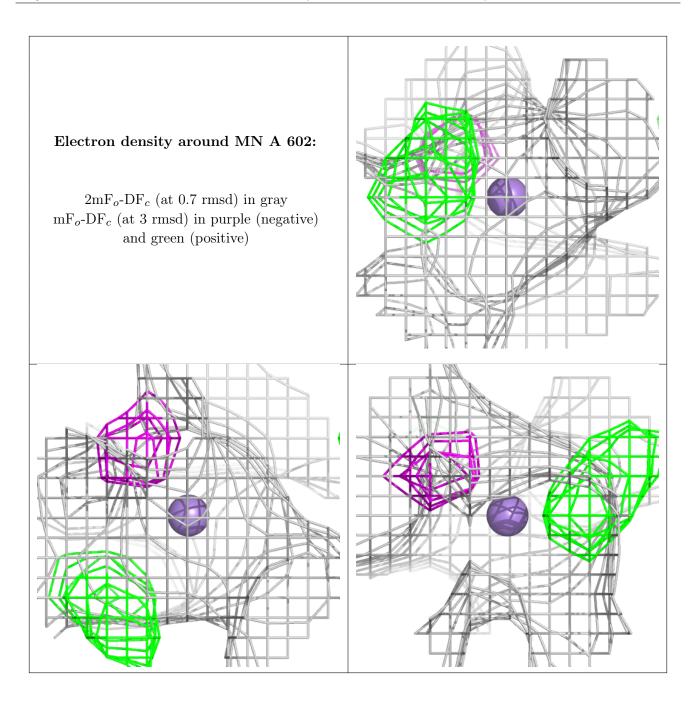
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9'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



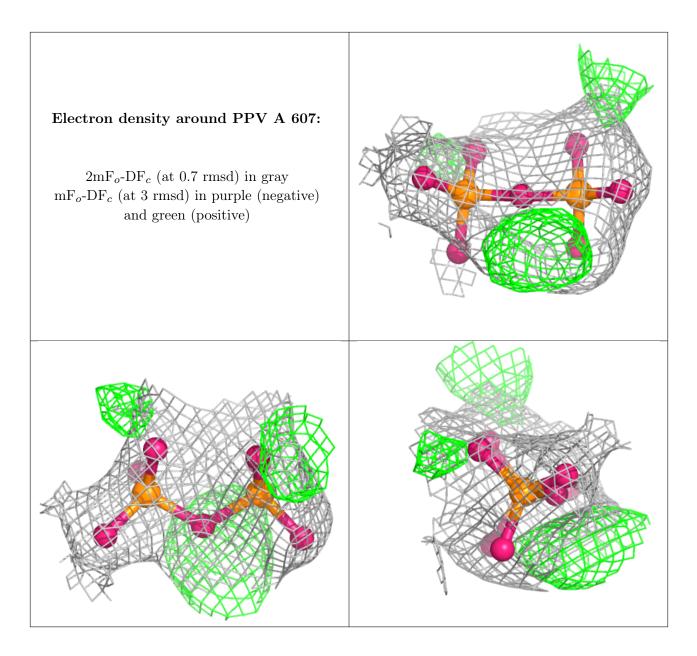
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q < 0.9
6	NA	А	606	1/1	0.19	0.17	$54,\!54,\!54,\!54$	1
5	MN	А	602	1/1	0.72	0.14	56, 56, 56, 56	1
8	EDO	А	609	3/4	0.76	0.23	36, 36, 51, 55	3
7	PPV	А	607	9/9	0.84	0.34	51,58,64,66	6
6	NA	А	605	1/1	0.85	0.12	45,45,45,45	1
8	EDO	А	610	4/4	0.90	0.16	$51,\!51,\!53,\!55$	4
5	MN	А	603	1/1	0.91	0.19	48,48,48,48	1
8	EDO	А	608	4/4	0.94	0.42	39,40,41,64	4
6	NA	А	604	1/1	0.95	0.13	34,34,34,34	0
5	MN	А	601	1/1	0.99	0.19	$50,\!50,\!50,\!50$	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.

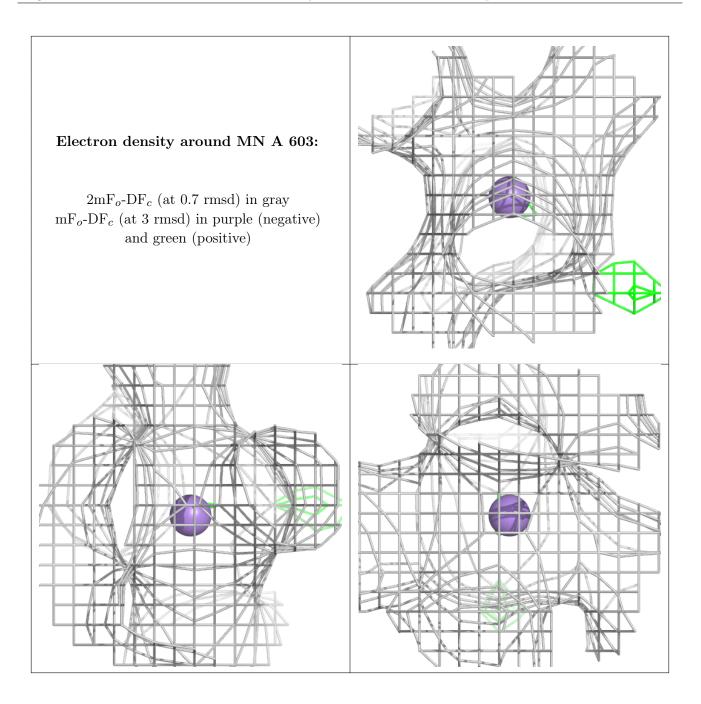




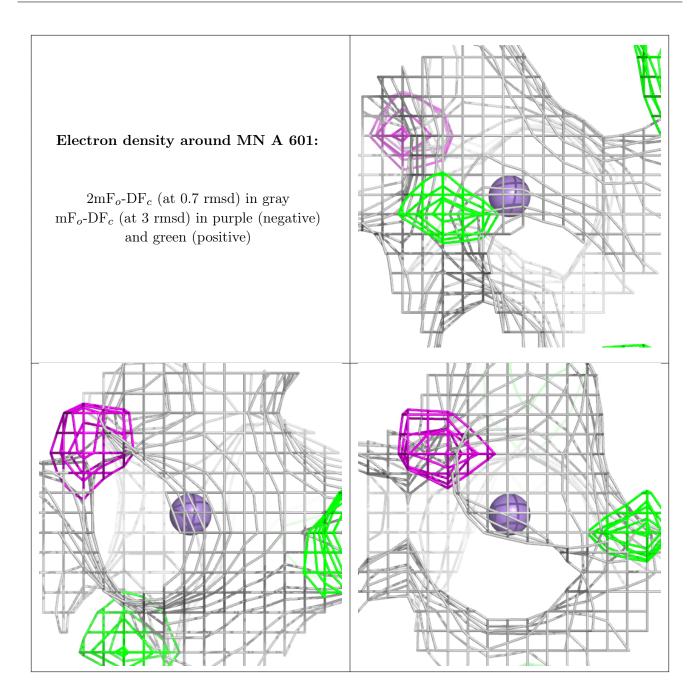












6.5 Other polymers (i)

There are no such residues in this entry.

