W O R L D W I D E
wwPDB
PROTEIN DATA BANK

Full wwPDB X-ray Structure Validation Report ⓘ

Oct 14, 2023 – 02:26 PM EDT

PDB ID	:	8DQI
Title	:	Crystal structure of pyrrolysyl-tRNA synthetase from Methanomethylophilus alvus engineered for acridone amino acid (RS1) bound to ATP and acridone after 2- weeks of crystal growth
Authors	:	Gottfried-Lee, I.; Karplus, P.A.; Mehl, R.A.; Cooley, R.B.
Deposited on	:	2022-07-19
Resolution	:	1.54 Å(reported)

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

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

The 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:

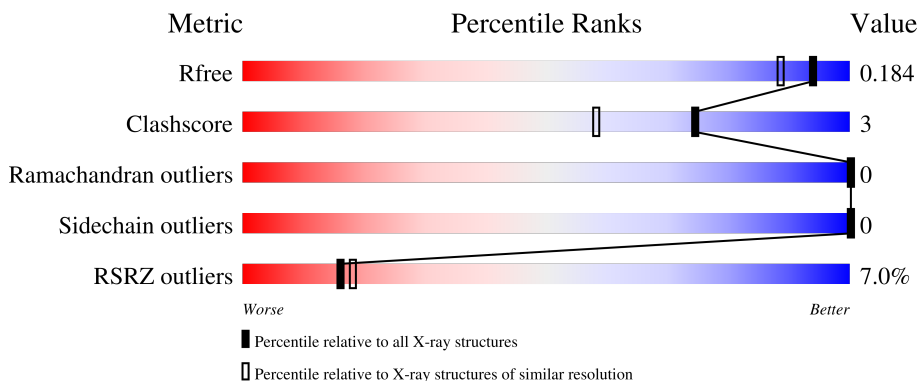|		
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & 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 1.54 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	Percentile Ranks	Value
Rfree		0.184
Clashscore		3
Ramachandran outliers		0
Sidechain outliers		0
RSRZ outliers		7.0%

Worse　　　　　　　　　　　　　　　*Better*

▌ Percentile relative to all X-ray structures

▯ Percentile relative to X-ray structures of similar resolution

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	2556 (1.56-1.52)
Clashscore	141614	2634 (1.56-1.52)
Ramachandran outliers	138981	2580 (1.56-1.52)
Sidechain outliers	138945	2577 (1.56-1.52)
RSRZ outliers	127900	2524 (1.56-1.52)

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	A	276	10% / 87% / 7% / 7%
1	D	276	4% / 95% / 5% ·

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:

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	PG4	D	303	-	-	-	X

2 Entry composition ⓘ

There are 7 unique types of molecules in this entry. The entry contains 8841 atoms, of which 4133 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 AA_TRNA_LIGASE_II domain-containing protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	D	274	Total 4194	C 1334	H 2074	N 358	O 412	S 16	0	6	0
1	A	258	Total 4009	C 1272	H 1998	N 337	O 385	S 17	0	11	0

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	0	SER	-	expression tag	UNP A0A3G3IHP7
D	1	GLY	-	expression tag	UNP A0A3G3IHP7
D	166	ALA	ASN	conflict	UNP A0A3G3IHP7
D	168	GLY	VAL	conflict	UNP A0A3G3IHP7
D	239	CYS	TRP	conflict	UNP A0A3G3IHP7
A	0	SER	-	expression tag	UNP A0A3G3IHP7
A	1	GLY	-	expression tag	UNP A0A3G3IHP7
A	166	ALA	ASN	conflict	UNP A0A3G3IHP7
A	168	GLY	VAL	conflict	UNP A0A3G3IHP7
A	239	CYS	TRP	conflict	UNP A0A3G3IHP7

- Molecule 2 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$) (labeled as "Ligand of Interest" by depositor).

ATP

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	D	1	Total	C	H	N	O	P	0	0
			42	10	11	5	13	3		
2	A	1	Total	C	H	N	O	P	0	0
			43	10	12	5	13	3		

- Molecule 3 is (2 {S})-2-azanyl-3-(9-oxidanylidene-10 {H}-acridin-2-yl)propanoic acid (three-letter code: T7Q) (formula: $C_{16}H_{15}N_2O_3$) (labeled as "Ligand of Interest" by depositor).

T7Q
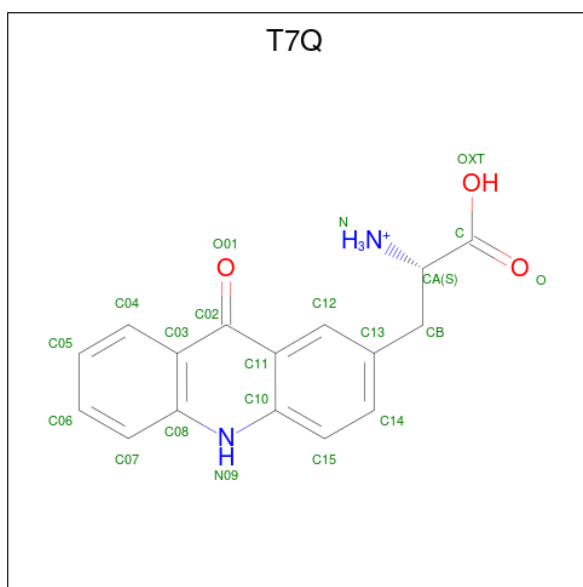
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	D	1	Total	C	H	N	O	0	0
			34	16	13	2	3		

- Molecule 4 is TETRAETHYLENE GLYCOL (three-letter code: PG4) (formula: $C_8H_{18}O_5$).
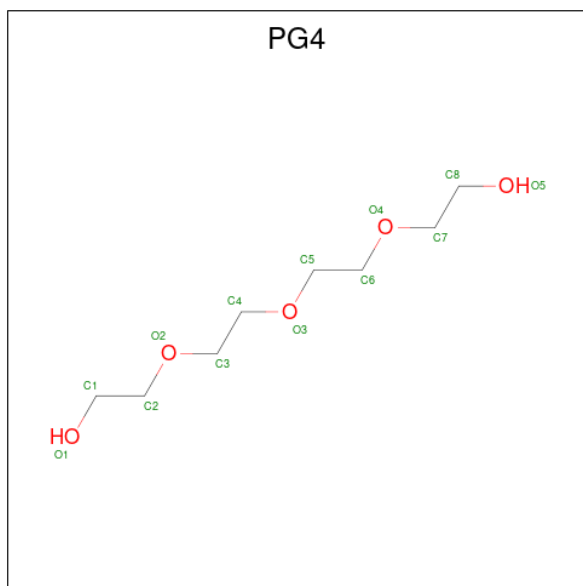
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	D	1	Total 30	C 8	H 17	O 5	0	0

- Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	D	3	Total 3	Mg 3	0	0
5	A	2	Total 2	Mg 2	0	0

- Molecule 6 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).

GOL

C3

OH O3

C1 C2

HO

O1

OH

O2

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total C H O				0	0
			14 3 8 3					
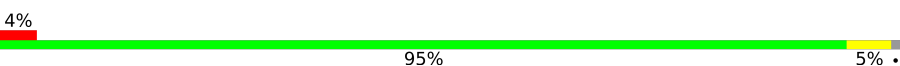
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	D	269	Total O		0	0
			269 269			
7	A	201	Total O		0	0
			201 201			

3 Residue-property plots ⓘ

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.
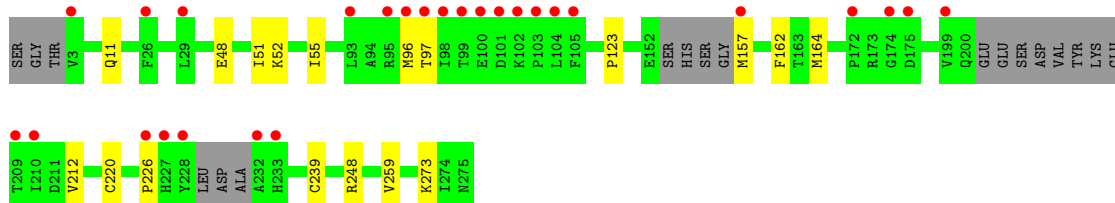
• Molecule 1: AA_TRNA_LIGASE_II domain-containing protein
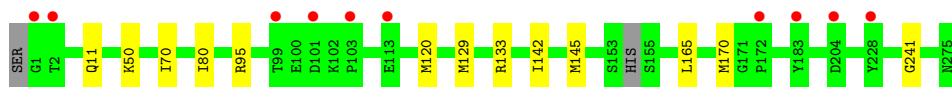
• Molecule 1: AA_TRNA_LIGASE_II domain-containing protein

4 Data and refinement statistics (i)

Property	Value	Source
Space group	I 4	Depositor
Cell constants a, b, c, α, β, γ	110.80Å 110.80Å 113.77Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.43 – 1.54 45.43 – 1.54	Depositor EDS
% Data completeness (in resolution range)	100.0 (45.43-1.54) 100.0 (45.43-1.54)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.30 (at 1.54Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.167 , 0.186 0.165 , 0.184	Depositor DCC
R_{free} test set	5083 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å^2)	24.2	Xtriage
Anisotropy	0.023	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.41 , 49.5	EDS
L-test for twinning[2]	$< |L| > = 0.50, < L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.000 for l,-k,h 0.007 for -l,-k,-h 0.007 for -h,-l,-k 0.000 for -h,l,k 0.024 for -h,k,-l	Xtriage
F_o,F_c correlation	0.97	EDS
Total number of atoms	8841	wwPDB-VP
Average B, all atoms (Å^2)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.12% of the height of the origin peak. No significant pseudotranslation is detected.*

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[1]Intensities estimated from amplitudes.
[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: GOL, ATP, T7Q, PG4, 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	0.45	0/2075	0.65	0/2792
1	D	0.48	0/2176	0.68	0/2935
All	All	0.46	0/4251	0.67	0/5727

There are no bond length outliers.

There are no bond angle outliers.

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	2011	1998	1998	13	0
1	D	2120	2074	2066	15	0
2	A	31	12	12	0	0
2	D	31	11	12	0	0
3	D	21	13	0	3	0
4	D	13	17	18	0	0
5	A	2	0	0	0	0
5	D	3	0	0	0	0
6	A	6	8	8	0	0
7	A	201	0	0	2	0
7	D	269	0	0	4	0
All	All	4708	4133	4114	27	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 3.

All (27) 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:D:133:ARG:HB2	1:D:170:MET:HE1	1.59	0.85
1:D:95:ARG:NH2	7:D:402:HOH:O	2.25	0.69
1:D:11:GLN:NE2	7:D:401:HOH:O	2.20	0.68
1:D:129:MET:HE3	1:D:170:MET:HB2	1.78	0.65
1:A:96:MET:O	1:A:97:THR:OG1	2.14	0.64
1:A:11:GLN:NE2	7:A:401:HOH:O	2.31	0.60
1:D:50:LYS:HD2	7:D:484:HOH:O	2.02	0.58
1:D:133:ARG:HB2	1:D:170:MET:CE	2.31	0.55
1:A:51:ILE:HD11	1:A:273:LYS:O	2.06	0.55
1:D:129:MET:CE	1:D:170:MET:HB2	2.40	0.52
1:A:162:PHE:HE2	1:A:164[B]:MET:SD	2.33	0.52
1:A:226:PRO:HD2	1:A:239[B]:CYS:SG	2.51	0.51
1:A:48:GLU:OE2	1:A:52[B]:LYS:HE3	2.11	0.50
1:D:145[A]:MET:CE	7:D:450:HOH:O	2.60	0.49
1:A:157:MET:HG2	1:A:248:ARG:HG2	1.94	0.49
1:A:157:MET:HE3	7:A:581:HOH:O	2.14	0.48
1:D:241:GLY:HA3	3:D:302:T7Q:C08	2.44	0.47
1:D:120:MET:HE2	3:D:302:T7Q:OXT	2.15	0.47
1:D:70:ILE:HG21	1:D:165:LEU:HD22	1.98	0.46
1:A:96:MET:HG3	1:A:123:PRO:HG3	1.98	0.46
1:D:129:MET:HG3	1:D:142:ILE:CD1	2.46	0.45
1:D:80:ILE:HD11	1:A:55:ILE:HA	1.99	0.44
1:A:162:PHE:CE2	1:A:164[B]:MET:SD	3.12	0.43
1:A:212:VAL:HB	1:A:220:CYS:HB3	1.99	0.43
1:D:241:GLY:HA3	3:D:302:T7Q:N09	2.35	0.42
1:A:259:VAL:O	1:A:259:VAL:HG22	2.20	0.41
1:D:129:MET:HG3	1:D:142:ILE:HD11	2.02	0.41

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	261/276 (95%)	257 (98%)	4 (2%)	0	100	100
1	D	276/276 (100%)	273 (99%)	3 (1%)	0	100	100
All	All	537/552 (97%)	530 (99%)	7 (1%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2   Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	215/231 (93%)	215 (100%)	0	100	100
1	D	225/231 (97%)	225 (100%)	0	100	100
All	All	440/462 (95%)	440 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

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

5.3.3   RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5   Carbohydrates ⓘ

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 10 ligands modelled in this entry, 5 are monoatomic - leaving 5 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	#$|Z| > 2$
6	GOL	A	301	-	5,5,5	0.71	0	5,5,5	1.34	0
3	T7Q	D	302	-	22,23,23	1.91	3 (13%)	32,33,33	1.29	4 (12%)
2	ATP	A	302	5	26,33,33	0.58	0	31,52,52	1.12	3 (9%)
4	PG4	D	303	-	12,12,12	0.11	0	11,11,11	0.20	0
2	ATP	D	301	5	26,33,33	0.69	0	31,52,52	0.91	1 (3%)

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
6	GOL	A	301	-	-	0/4/4/4	-
3	T7Q	D	302	-	-	2/8/8/8	0/3/3/3
2	ATP	A	302	5	-	3/18/38/38	0/3/3/3
4	PG4	D	303	-	-	6/10/10/10	-
2	ATP	D	301	5	-	4/18/38/38	0/3/3/3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	302	T7Q	C08-N09	5.78	1.51	1.38
3	D	302	T7Q	C10-N09	5.23	1.49	1.38
3	D	302	T7Q	C11-C10	-2.15	1.37	1.41

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	302	T7Q	CB-C13-C12	-3.70	114.08	120.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	302	ATP	C5-C6-N6	2.51	124.16	120.35
3	D	302	T7Q	OXT-C-CA	2.42	121.63	113.38
2	D	301	ATP	C5-C6-N6	2.37	123.96	120.35
2	A	302	ATP	O3'-C3'-C2'	-2.24	104.57	111.82
2	A	302	ATP	O5'-PA-O1A	2.19	117.62	109.07
3	D	302	T7Q	O01-C02-C11	-2.11	117.81	120.91
3	D	302	T7Q	CB-C13-C14	2.07	125.01	120.91

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	301	ATP	C5'-O5'-PA-O2A
2	A	302	ATP	O4'-C4'-C5'-O5'
4	D	303	PG4	O2-C3-C4-O3
4	D	303	PG4	O1-C1-C2-O2
4	D	303	PG4	O3-C5-C6-O4
2	A	302	ATP	C3'-C4'-C5'-O5'
3	D	302	T7Q	OXT-C-CA-N
3	D	302	T7Q	O-C-CA-N
4	D	303	PG4	C3-C4-O3-C5
2	D	301	ATP	C5'-O5'-PA-O3A
2	D	301	ATP	C5'-O5'-PA-O1A
4	D	303	PG4	C1-C2-O2-C3
2	D	301	ATP	C4'-C5'-O5'-PA
4	D	303	PG4	C5-C6-O4-C7
2	A	302	ATP	C5'-O5'-PA-O3A

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	302	T7Q	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 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.

Ligand T7Q D 302

Bond lengths

Bond angles

Torsions

Rings

Ligand ATP A 302

Bond lengths

Bond angles

Torsions

Rings

Ligand ATP D 301

Bond lengths

Bond angles

Torsions

Rings
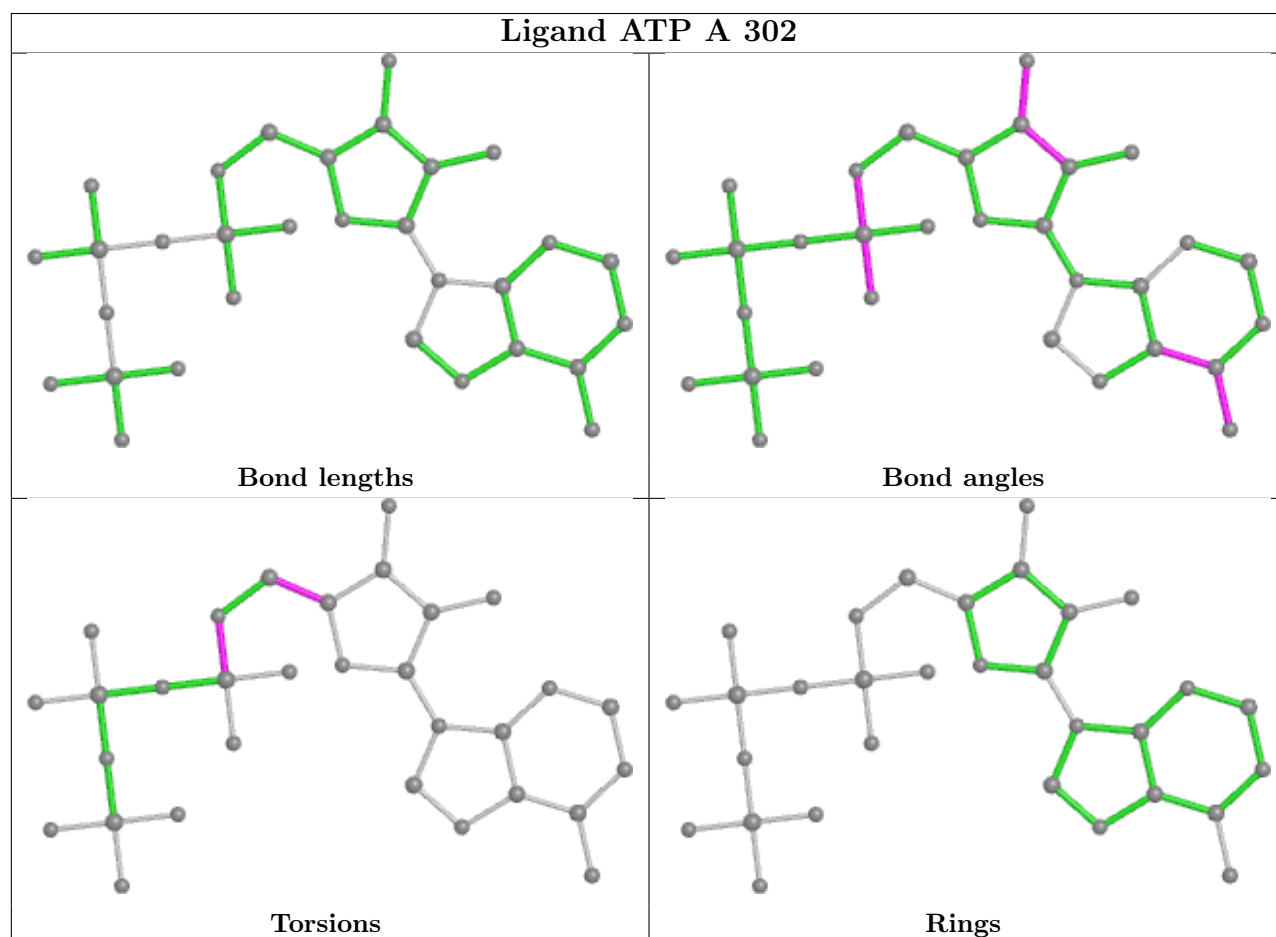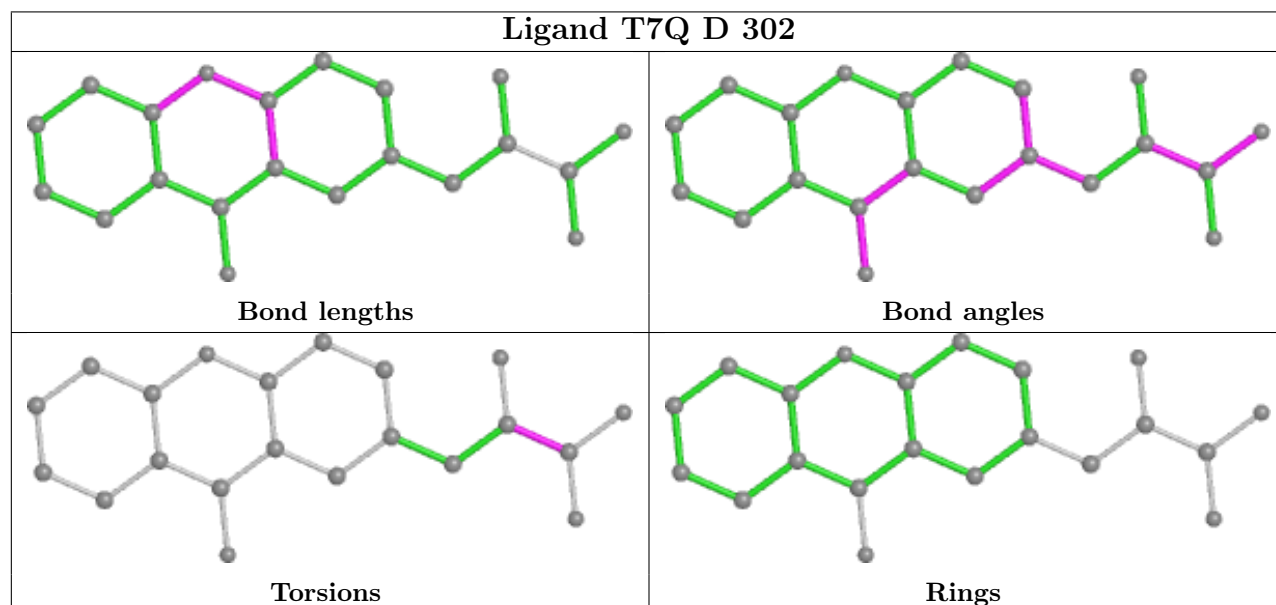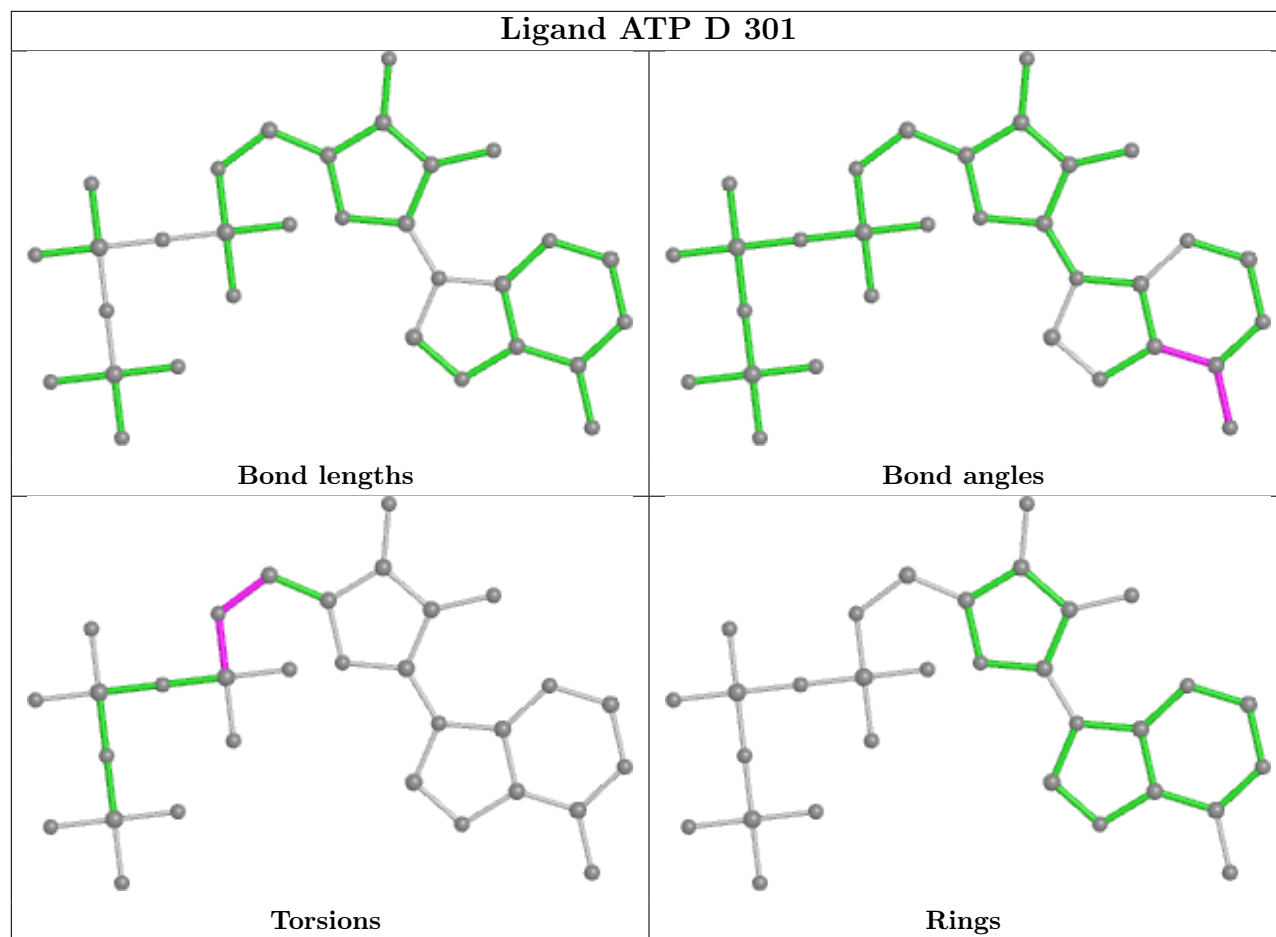
5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å²)	Q<0.9
1	A	258/276 (93%)	0.43	27 (10%)	6	6	20, 32, 67, 93	0
1	D	274/276 (99%)	0.04	10 (3%)	42	49	18, 26, 52, 75	0
All	All	532/552 (96%)	0.23	37 (6%)	16	18	18, 30, 60, 93	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	98	ILE	5.9
1	A	104	LEU	5.1
1	A	228	TYR	4.8
1	A	172	PRO	4.5
1	A	101	ASP	4.4
1	A	103	PRO	4.1
1	A	232	ALA	4.1
1	A	105	PHE	4.0
1	A	99	THR	3.8
1	A	226	PRO	3.8
1	D	101	ASP	3.7
1	A	93	LEU	3.7
1	A	174	GLY	3.6
1	A	175	ASP	3.6
1	A	102	LYS	3.5
1	D	204	ASP	3.5
1	D	99	THR	3.1
1	A	97	THR	3.1
1	D	228	TYR	3.1
1	A	199	VAL	3.1
1	D	103	PRO	2.8
1	A	227	HIS	2.8
1	D	2	THR	2.7
1	A	209	THR	2.7

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Mol	Chain	Res	Type	RSRZ
1	D	113	GLU	2.6
1	D	1	GLY	2.5
1	A	157	MET	2.5
1	D	172	PRO	2.5
1	A	233	HIS	2.5
1	A	3	VAL	2.3
1	A	210	ILE	2.3
1	A	96	MET	2.3
1	A	100	GLU	2.2
1	A	26	PHE	2.1
1	A	29	LEU	2.1
1	D	183	TYR	2.0
1	A	95	ARG	2.0

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

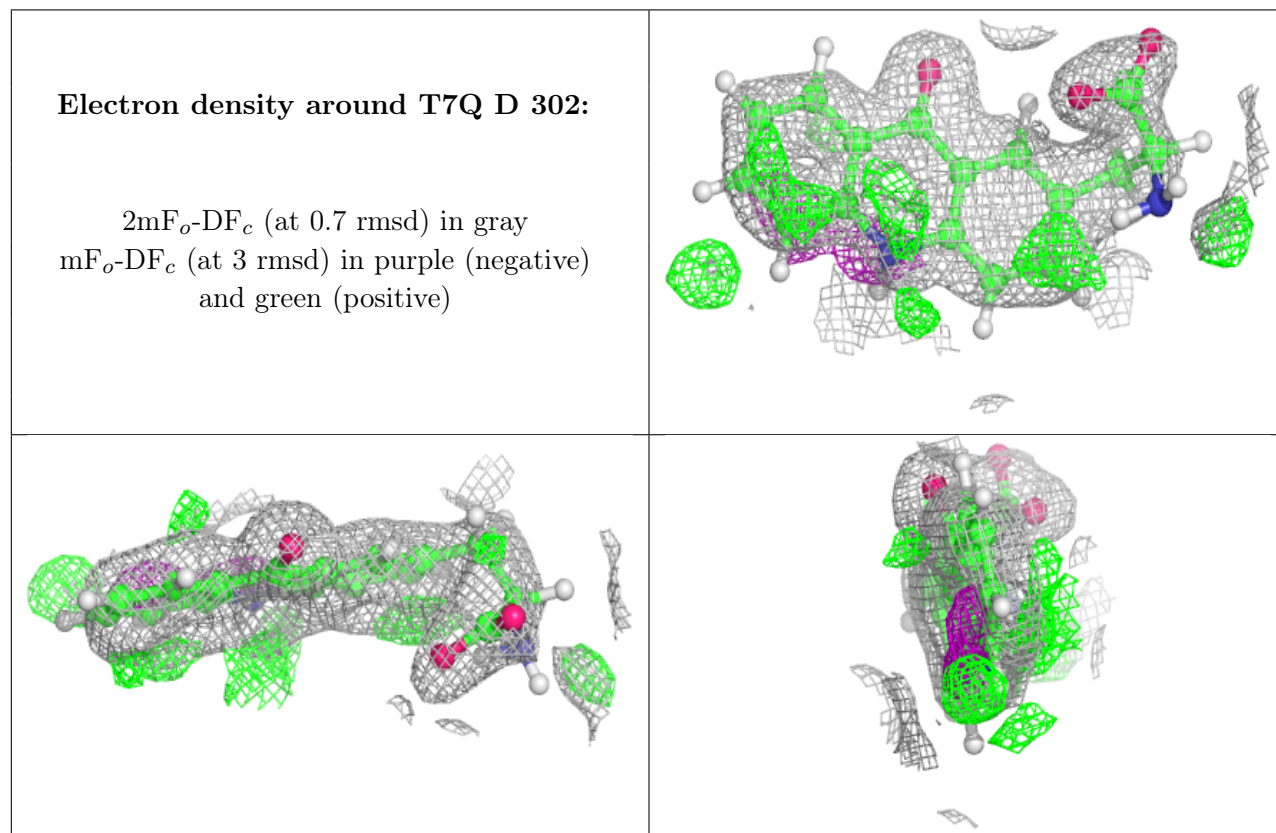There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no 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, 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.
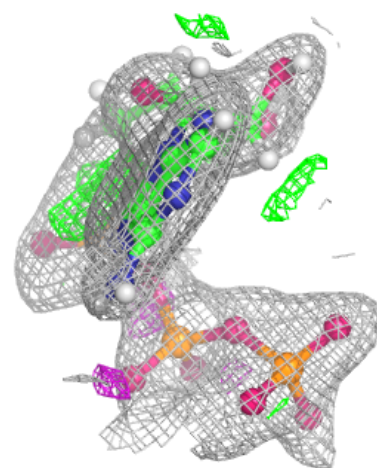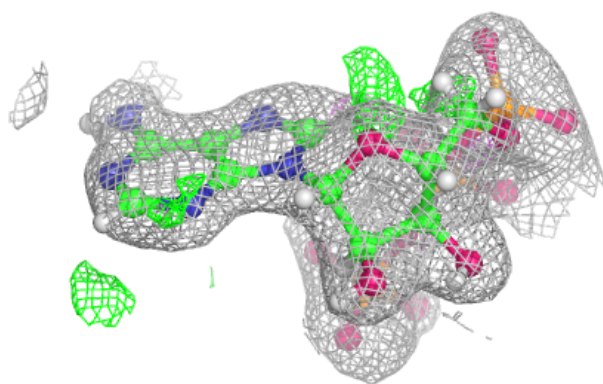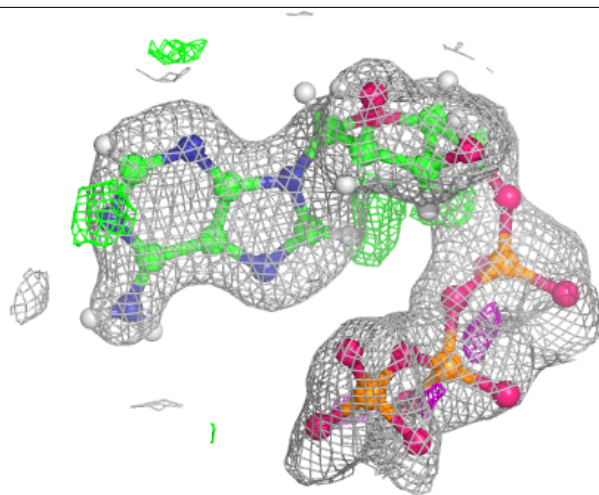
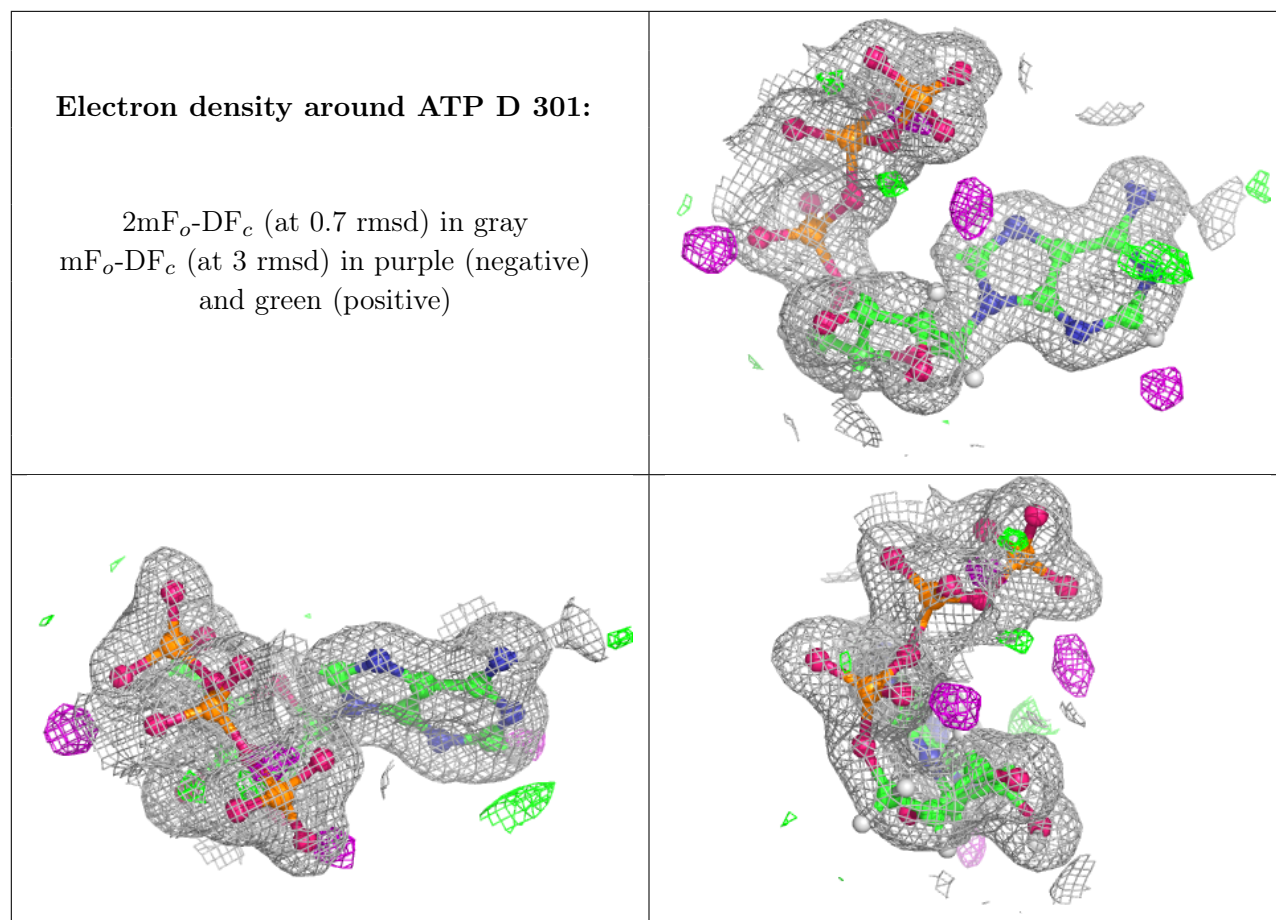| Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
4	PG4	D	303	13/13	0.73	0.50	20,20,20,20	0
3	T7Q	D	302	21/21	0.80	0.17	27,36,59,70	34
2	ATP	A	302	31/31	0.86	0.15	30,41,62,72	43
5	MG	A	304	1/1	0.87	0.20	57,57,57,57	0
5	MG	D	306	1/1	0.93	0.07	32,32,32,32	1
5	MG	D	305	1/1	0.93	0.06	27,27,27,27	1
5	MG	A	303	1/1	0.94	0.12	56,56,56,56	1
6	GOL	A	301	6/6	0.94	0.12	21,34,52,56	0
2	ATP	D	301	31/31	0.95	0.08	20,26,31,32	42
5	MG	D	304	1/1	0.96	0.07	27,27,27,27	1

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 T7Q D 302:

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

Electron density around ATP A 302:

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 ATP D 301:

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

6.5 Other polymers ⓘ

There are no such residues in this entry.