



Full wwPDB X-ray Structure Validation Report ⓘ

May 25, 2020 – 11:40 pm BST

PDB ID : 3FSC
Title : Crystal structure of QdtC, the dTDP-3-amino-3,6-dideoxy-D-glucose N-acetyl transferase from *Thermoanaerobacterium thermosaccharolyticum* in complex with CoA and dTDP-3-amino-fucose
Authors : Holden, H.M.; Thoden, J.B.
Deposited on : 2009-01-09
Resolution : 1.80 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

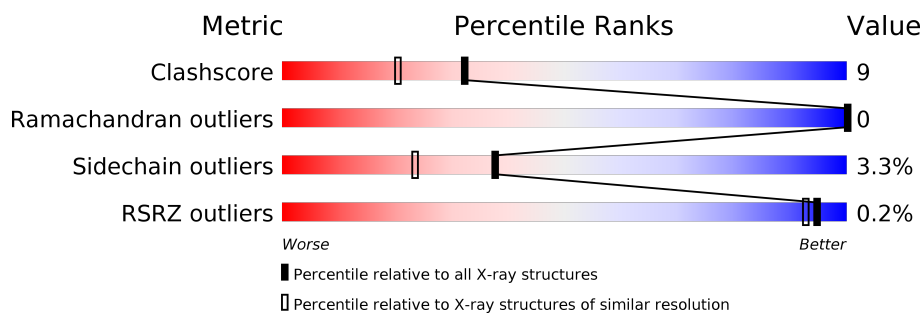
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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))
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 on 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	273	
1	B	273	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4689 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.

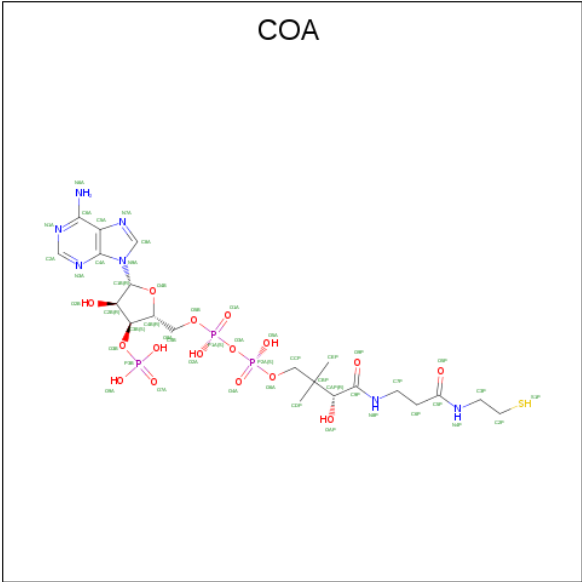
- Molecule 1 is a protein called QdtC.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	259	Total	C	N	O	S	0	2	0
			2060	1321	353	383	3			
1	B	257	Total	C	N	O	S	0	3	0
			2044	1312	350	379	3			

There are 16 discrepancies between the modelled and reference sequences:

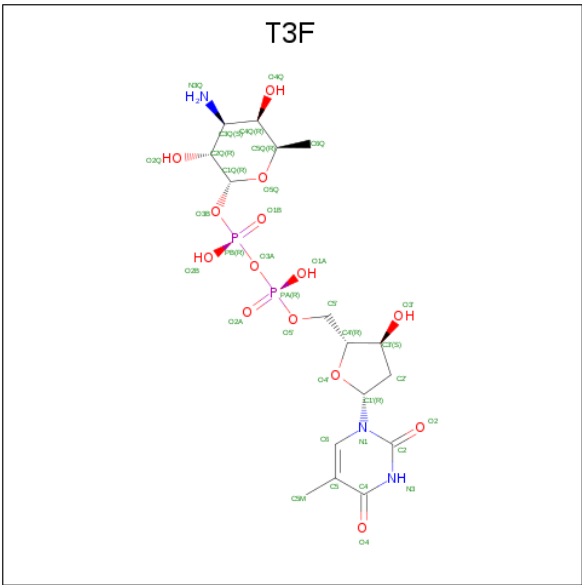
Chain	Residue	Modelled	Actual	Comment	Reference
A	266	LEU	-	EXPRESSION TAG	UNP Q6TFC6
A	267	GLU	-	EXPRESSION TAG	UNP Q6TFC6
A	268	HIS	-	EXPRESSION TAG	UNP Q6TFC6
A	269	HIS	-	EXPRESSION TAG	UNP Q6TFC6
A	270	HIS	-	EXPRESSION TAG	UNP Q6TFC6
A	271	HIS	-	EXPRESSION TAG	UNP Q6TFC6
A	272	HIS	-	EXPRESSION TAG	UNP Q6TFC6
A	273	HIS	-	EXPRESSION TAG	UNP Q6TFC6
B	266	LEU	-	EXPRESSION TAG	UNP Q6TFC6
B	267	GLU	-	EXPRESSION TAG	UNP Q6TFC6
B	268	HIS	-	EXPRESSION TAG	UNP Q6TFC6
B	269	HIS	-	EXPRESSION TAG	UNP Q6TFC6
B	270	HIS	-	EXPRESSION TAG	UNP Q6TFC6
B	271	HIS	-	EXPRESSION TAG	UNP Q6TFC6
B	272	HIS	-	EXPRESSION TAG	UNP Q6TFC6
B	273	HIS	-	EXPRESSION TAG	UNP Q6TFC6

- Molecule 2 is COENZYME A (three-letter code: COA) (formula: $C_{21}H_{36}N_7O_{16}P_3S$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	S	0	0
			48	21	7	16	3	1		
2	B	1	Total	C	N	O	P	S	0	0
			48	21	7	16	3	1		

- Molecule 3 is (3R,4S,5R,6R)-4-amino-3,5-dihydroxy-6-methyloxan-2-yl[[hydroxy-[[[(2R,3S,5R)-3-hydroxy-5-(5-methyl-2,4-dioxypyrimidin-1-yl)oxolan-2-yl]methoxy]phosphoryl]hydrogen phosphate (three-letter code: T3F) (formula: C₁₆H₂₇N₃O₁₄P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			35	16	3	14	2		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	B	1	Total	C	N	O	P	0	0
			35	16	3	14	2		

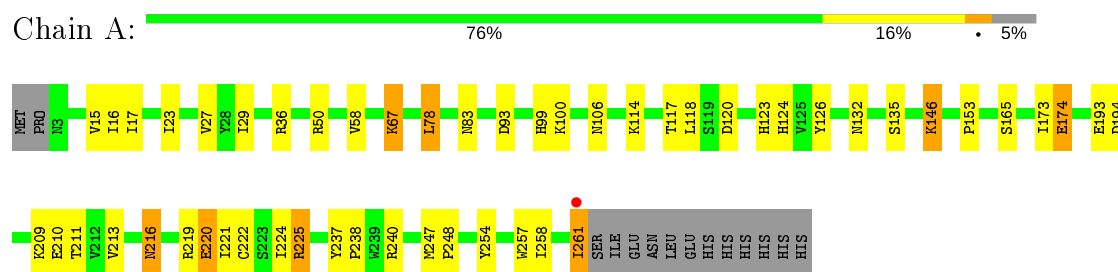
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	180	Total	O	0	0
			180	180		
4	B	239	Total	O	0	0
			239	239		

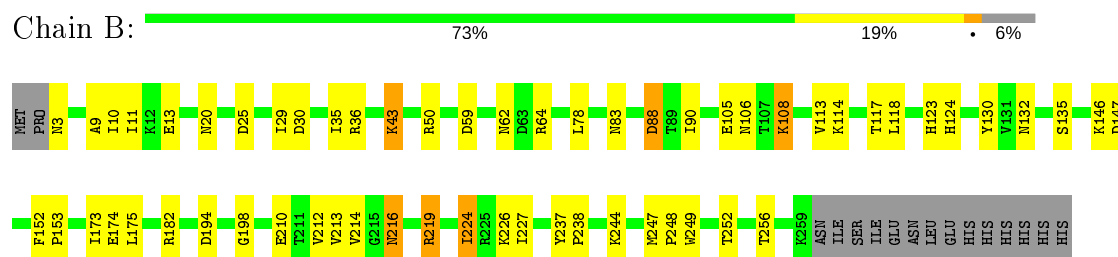
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: QdtC



• Molecule 1: QdtC



4 Data and refinement statistics

Property	Value	Source
Space group	P 3	Depositor
Cell constants a, b, c, α , β , γ	67.10Å 67.10Å 112.00Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 1.80 40.32 – 1.80	Depositor EDS
% Data completeness (in resolution range)	95.9 (30.00-1.80) 96.1 (40.32-1.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.65 (at 1.79Å)	Xtriage
Refinement program	TNT	Depositor
R, R_{free}	0.162 , 0.231 0.161 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	9.8	Xtriage
Anisotropy	0.174	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 83.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.020 for -h,-k,l 0.053 for h,-h-k,-l 0.027 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4689	wwPDB-VP
Average B, all atoms (Å ²)	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 15.16% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² 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: COA, T3F

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	0.75	0/2113	1.34	15/2873 (0.5%)
1	B	0.78	1/2102 (0.0%)	1.32	14/2858 (0.5%)
All	All	0.76	1/4215 (0.0%)	1.33	29/5731 (0.5%)

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
1	B	1	0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	113	VAL	CB-CG1	5.20	1.63	1.52

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	50	ARG	NE-CZ-NH2	-9.79	115.41	120.30
1	A	225	ARG	NE-CZ-NH1	-8.14	116.23	120.30
1	B	64	ARG	NE-CZ-NH2	-8.05	116.28	120.30
1	B	36	ARG	NE-CZ-NH2	7.30	123.95	120.30
1	A	36	ARG	NE-CZ-NH2	7.24	123.92	120.30
1	B	182	ARG	NE-CZ-NH2	-7.17	116.71	120.30
1	A	225	ARG	NE-CZ-NH2	6.89	123.74	120.30
1	A	78	LEU	CB-CG-CD1	-6.78	99.47	111.00
1	A	36	ARG	NE-CZ-NH1	-6.63	116.99	120.30
1	B	224	ILE	CG1-CB-CG2	6.62	125.96	111.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	50	ARG	NE-CZ-NH2	-6.52	117.04	120.30
1	B	219	ARG	NE-CZ-NH1	-6.45	117.08	120.30
1	B	30	ASP	CB-CG-OD1	6.24	123.91	118.30
1	A	174	GLU	OE1-CD-OE2	-5.74	116.42	123.30
1	B	50	ARG	NE-CZ-NH1	5.59	123.10	120.30
1	A	50	ARG	NE-CZ-NH1	5.54	123.07	120.30
1	B	25	ASP	CB-CG-OD1	5.53	123.28	118.30
1	B	64	ARG	NE-CZ-NH1	5.51	123.05	120.30
1	A	93	ASP	CB-CG-OD2	5.37	123.13	118.30
1	A	120	ASP	CB-CG-OD1	5.33	123.10	118.30
1	B	88	ASP	CB-CG-OD2	5.31	123.08	118.30
1	B	59	ASP	CB-CG-OD1	5.28	123.05	118.30
1	A	120	ASP	CB-CG-OD2	-5.22	113.60	118.30
1	A	165	SER	N-CA-CB	5.20	118.31	110.50
1	B	256	THR	CB-CA-C	-5.09	97.85	111.60
1	A	220[A]	GLU	CB-CA-C	-5.09	100.23	110.40
1	A	220[B]	GLU	CB-CA-C	-5.09	100.23	110.40
1	A	17	ILE	CG1-CB-CG2	5.04	122.50	111.40
1	B	252	THR	CA-CB-CG2	-5.00	105.39	112.40

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	B	224	ILE	CB

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	2060	0	2081	35	0
1	B	2044	0	2070	38	0
2	A	48	0	32	3	0
2	B	48	0	32	0	0
3	A	35	0	25	0	0
3	B	35	0	25	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	180	0	0	4	0
4	B	239	0	0	7	0
All	All	4689	0	4265	75	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 9.

All (75) 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:123:HIS:HD2	1:A:124:HIS:HD2	1.16	0.88
1:B:123:HIS:ND1	1:B:124:HIS:HD2	1.80	0.79
2:A:300:COA:O5A	2:A:300:COA:O1A	2.07	0.72
1:A:123:HIS:HD2	1:A:124:HIS:CD2	2.05	0.71
1:B:29:ILE:HD13	1:B:35[A]:ILE:HD11	1.72	0.70
1:A:216:ASN:HD22	1:A:216:ASN:C	1.91	0.69
1:A:123:HIS:CD2	1:A:124:HIS:HD2	2.04	0.69
1:A:67:LYS:N	1:A:67:LYS:HD3	2.09	0.68
1:A:224:ILE:HD12	1:A:237:TYR:HB3	1.75	0.68
1:A:210:GLU:OE2	4:A:407:HOH:O	2.11	0.68
1:A:237:TYR:HA	1:A:238:PRO:C	2.16	0.66
1:B:43:LYS:NZ	4:B:275:HOH:O	2.30	0.65
1:B:29:ILE:HD13	1:B:35[A]:ILE:CD1	2.28	0.63
1:B:3:ASN:OD1	1:B:20:ASN:N	2.33	0.62
1:A:211:THR:CG2	1:A:220[B]:GLU:HG2	2.30	0.61
1:A:221:ILE:O	1:A:222:CYS:HB3	2.01	0.60
1:A:194:ASP:O	1:A:224:ILE:HD13	2.03	0.59
1:A:211:THR:HG21	1:A:220[B]:GLU:HG2	1.82	0.59
1:A:146:LYS:HG3	1:A:174:GLU:HA	1.85	0.59
1:B:62:ASN:ND2	4:B:443:HOH:O	2.36	0.58
1:B:244:LYS:NZ	4:B:359:HOH:O	2.37	0.58
1:A:114:LYS:HD3	1:A:132:ASN:HD21	1.68	0.58
1:B:212:VAL:HG21	1:B:227:ILE:CD1	2.34	0.57
1:A:114:LYS:HD3	1:A:132:ASN:ND2	2.19	0.57
1:B:198:GLY:HA3	1:B:214:VAL:HG12	1.86	0.57
1:A:146:LYS:HG2	1:A:173:ILE:O	2.05	0.56
1:A:123:HIS:HE1	4:A:324:HOH:O	1.89	0.56
1:B:135[A]:SER:OG	4:B:492:HOH:O	2.18	0.56
1:B:152:PHE:HB3	1:B:153:PRO:CD	2.37	0.55
1:B:88:ASP:O	1:B:106:ASN:HA	2.06	0.54
1:B:216:ASN:C	1:B:216:ASN:HD22	2.11	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:117:THR:HG22	1:A:118:LEU:HG	1.89	0.54
1:A:240:ARG:HD2	1:A:254:TYR:CE2	2.42	0.54
1:A:247:MET:HB3	1:A:248:PRO:HD2	1.90	0.54
1:B:219:ARG:HG2	4:B:457:HOH:O	2.08	0.53
1:B:114:LYS:HD3	1:B:132:ASN:HD21	1.74	0.53
1:B:210:GLU:HB3	1:B:224:ILE:HG12	1.89	0.53
1:B:152:PHE:HB3	1:B:153:PRO:HD2	1.91	0.53
1:B:117:THR:HG22	1:B:118:LEU:HG	1.91	0.52
1:A:257:TRP:O	1:A:261:ILE:HG12	2.10	0.52
1:B:123:HIS:ND1	1:B:124:HIS:CD2	2.72	0.49
1:B:146:LYS:HG2	1:B:173:ILE:O	2.13	0.49
1:A:219:ARG:NH2	4:A:366:HOH:O	2.38	0.49
1:B:237:TYR:HA	1:B:238:PRO:C	2.33	0.49
1:A:225:ARG:HD2	1:B:10:ILE:CD1	2.44	0.48
1:B:212:VAL:HG21	1:B:227:ILE:HD13	1.95	0.48
1:B:194:ASP:O	1:B:224:ILE:HD13	2.14	0.47
1:B:212:VAL:HG21	1:B:227:ILE:HD11	1.95	0.47
1:B:90:ILE:HB	1:B:108:LYS:HD2	1.95	0.47
1:A:15:VAL:O	1:A:16:ILE:HD13	2.14	0.47
1:B:147:ASP:HB2	1:B:175:LEU:CD2	2.45	0.47
1:B:114:LYS:HD3	1:B:132:ASN:ND2	2.30	0.47
1:A:15:VAL:C	1:A:16:ILE:HD13	2.37	0.46
1:A:135:SER:O	1:A:153:PRO:HA	2.16	0.46
1:B:213:VAL:HA	1:B:219:ARG:O	2.16	0.46
1:B:118:LEU:HD12	4:B:492:HOH:O	2.15	0.45
1:B:146:LYS:HG3	1:B:174:GLU:HA	1.97	0.45
1:A:216:ASN:ND2	1:A:216:ASN:C	2.63	0.45
1:A:193:GLU:HG3	4:A:388:HOH:O	2.16	0.44
1:A:23:ILE:HG23	1:A:27:VAL:HG11	1.99	0.44
1:B:130:TYR:CE2	1:B:248:PRO:HD3	2.53	0.43
2:A:300:COA:C2B	2:A:300:COA:O5B	2.67	0.43
1:B:247:MET:HG3	1:B:249:TRP:CZ2	2.54	0.43
1:B:10:ILE:HD13	1:B:10:ILE:N	2.34	0.42
1:B:226:LYS:NZ	4:B:410:HOH:O	2.35	0.42
1:B:105:GLU:O	1:B:123:HIS:HA	2.20	0.41
1:B:9:ALA:C	1:B:10:ILE:HD13	2.41	0.41
1:A:106:ASN:O	1:A:124:HIS:HA	2.20	0.41
1:A:126:TYR:C	1:A:126:TYR:CD2	2.94	0.41
1:A:194:ASP:O	1:A:210:GLU:HA	2.21	0.41
2:A:300:COA:H2B	2:A:300:COA:O5B	2.21	0.41
1:A:99:HIS:O	1:A:100:LYS:HB2	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	A	259/273 (95%)	250 (96%)	9 (4%)	0	100	100
1	B	258/273 (94%)	247 (96%)	11 (4%)	0	100	100
All	All	517/546 (95%)	497 (96%)	20 (4%)	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	230/242 (95%)	221 (96%)	9 (4%)	32	17
1	B	229/242 (95%)	223 (97%)	6 (3%)	46	32
All	All	459/484 (95%)	444 (97%)	15 (3%)	38	23

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	29	ILE
1	A	58	VAL
1	A	67	LYS
1	A	83	ASN
1	A	146	LYS

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Mol	Chain	Res	Type
1	A	209	LYS
1	A	216	ASN
1	A	258	ILE
1	A	261	ILE
1	B	11	ILE
1	B	13	GLU
1	B	43	LYS
1	B	83	ASN
1	B	108	LYS
1	B	216	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	94	ASN
1	A	112	ASN
1	A	123	HIS
1	A	124	HIS
1	A	132	ASN
1	A	216	ASN
1	B	62	ASN
1	B	94	ASN
1	B	112	ASN
1	B	124	HIS
1	B	132	ASN
1	B	216	ASN

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 carbohydrates in this entry.

5.6 Ligand geometry

4 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	T3F	A	301	-	31,37,37	1.15	3 (9%)	42,57,57	1.95	6 (14%)
3	T3F	B	301	-	31,37,37	1.23	2 (6%)	42,57,57	1.53	5 (11%)
2	COA	B	300	-	41,50,50	1.17	5 (12%)	52,75,75	1.70	9 (17%)
2	COA	A	300	-	41,50,50	1.14	3 (7%)	52,75,75	1.98	9 (17%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	T3F	A	301	-	-	2/18/53/53	0/3/3/3
3	T3F	B	301	-	-	5/18/53/53	0/3/3/3
2	COA	B	300	-	-	10/44/64/64	0/3/3/3
2	COA	A	300	-	-	5/44/64/64	0/3/3/3

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	301	T3F	PA-O2A	3.20	1.62	1.50
3	B	301	T3F	C4-N3	3.02	1.38	1.33
2	B	300	COA	O4B-C1B	2.87	1.45	1.41
3	A	301	T3F	PA-O2A	2.85	1.61	1.50
2	A	300	COA	C2B-C1B	2.83	1.58	1.53
2	A	300	COA	O4B-C1B	2.71	1.44	1.41
3	A	301	T3F	C4-N3	2.61	1.37	1.33
2	B	300	COA	C8A-N7A	-2.49	1.30	1.34
2	B	300	COA	P3B-O9A	2.27	1.63	1.54
2	A	300	COA	P3B-O3B	2.27	1.63	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	300	COA	C6A-N6A	2.12	1.41	1.34
2	B	300	COA	P3B-O8A	2.07	1.62	1.54
3	A	301	T3F	C2'-C1'	2.03	1.58	1.52

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	301	T3F	C4-N3-C2	8.47	122.29	115.14
2	A	300	COA	P2A-O3A-P1A	-7.49	107.12	132.83
2	B	300	COA	C7P-C6P-C5P	-5.60	103.02	112.36
3	B	301	T3F	C4-N3-C2	5.50	119.78	115.14
2	A	300	COA	C7P-C6P-C5P	-5.12	103.83	112.36
2	A	300	COA	C7P-N8P-C9P	-4.37	114.80	122.59
2	B	300	COA	P2A-O3A-P1A	-4.30	118.07	132.83
2	A	300	COA	O6A-CCP-CBP	-4.15	103.88	110.55
3	A	301	T3F	O5Q-C1Q-O3B	-4.13	105.97	111.36
3	B	301	T3F	O3A-PB-O3B	-3.91	94.60	102.48
3	B	301	T3F	PA-O3A-PB	-3.87	119.55	132.83
2	B	300	COA	N3A-C2A-N1A	-3.72	122.87	128.68
3	A	301	T3F	PA-O3A-PB	-3.55	120.64	132.83
3	A	301	T3F	C2'-C1'-N1	-3.39	106.44	114.27
2	A	300	COA	N3A-C2A-N1A	-3.39	123.39	128.68
2	B	300	COA	O4B-C4B-C5B	-3.11	99.15	109.37
3	B	301	T3F	C2'-C1'-N1	-2.84	107.72	114.27
2	A	300	COA	P1A-O5B-C5B	-2.78	105.37	121.68
2	A	300	COA	O9P-C9P-N8P	-2.70	117.19	122.99
3	A	301	T3F	O3B-C1Q-C2Q	2.65	113.23	108.38
3	A	301	T3F	C6Q-C5Q-C4Q	-2.58	108.31	113.07
2	B	300	COA	C6P-C5P-N4P	2.35	120.38	116.42
2	A	300	COA	CAP-C9P-N8P	2.23	121.01	116.58
2	A	300	COA	N6A-C6A-N1A	2.22	123.19	118.57
2	B	300	COA	P1A-O5B-C5B	-2.18	108.91	121.68
2	B	300	COA	O4B-C1B-C2B	-2.17	103.75	106.93
2	B	300	COA	N6A-C6A-N1A	2.12	122.97	118.57
2	B	300	COA	C2A-N1A-C6A	2.08	122.32	118.75
3	B	301	T3F	C5Q-C4Q-C3Q	-2.02	107.44	110.51

There are no chirality outliers.

All (22) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	300	COA	O4B-C4B-C5B-O5B

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Mol	Chain	Res	Type	Atoms
2	B	300	COA	C3B-C4B-C5B-O5B
2	B	300	COA	C3B-O3B-P3B-O9A
2	B	300	COA	CCP-O6A-P2A-O3A
2	A	300	COA	C3B-C4B-C5B-O5B
3	B	301	T3F	C1Q-O3B-PB-O1B
3	B	301	T3F	PB-O3A-PA-O1A
2	B	300	COA	CCP-O6A-P2A-O5A
2	B	300	COA	CAP-CBP-CCP-O6A
2	A	300	COA	CAP-CBP-CCP-O6A
2	A	300	COA	CEP-CBP-CCP-O6A
3	A	301	T3F	PB-O3A-PA-O1A
2	B	300	COA	O5P-C5P-C6P-C7P
3	B	301	T3F	C1Q-O3B-PB-O2B
2	A	300	COA	CDP-CBP-CCP-O6A
3	B	301	T3F	C1Q-O3B-PB-O3A
2	B	300	COA	N4P-C5P-C6P-C7P
2	B	300	COA	CDP-CBP-CCP-O6A
2	B	300	COA	CEP-CBP-CCP-O6A
2	B	300	COA	O4B-C4B-C5B-O5B
3	A	301	T3F	PB-O3A-PA-O2A
3	B	301	T3F	PB-O3A-PA-O2A

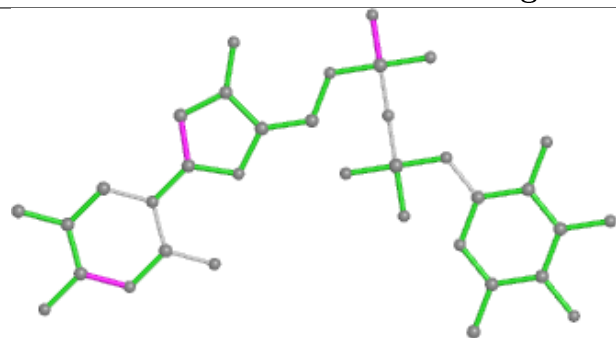
There are no ring outliers.

1 monomer is involved in 3 short contacts:

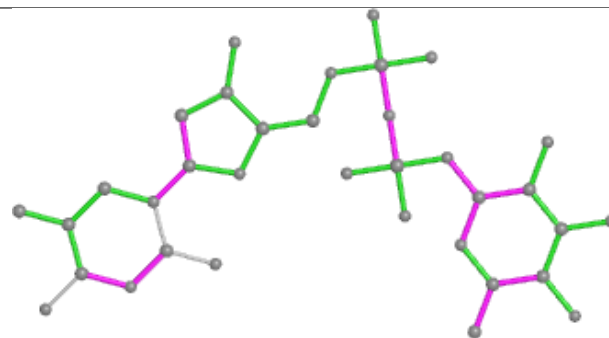
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	300	COA	3	0

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.

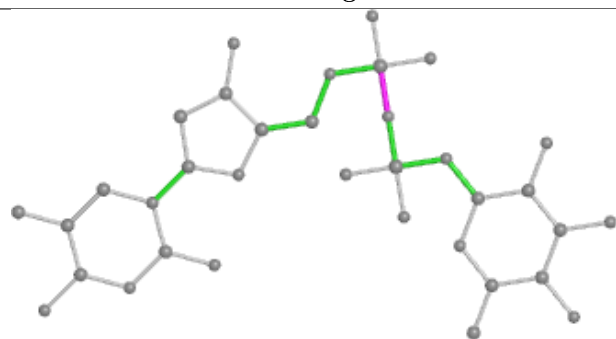
Ligand T3F A 301



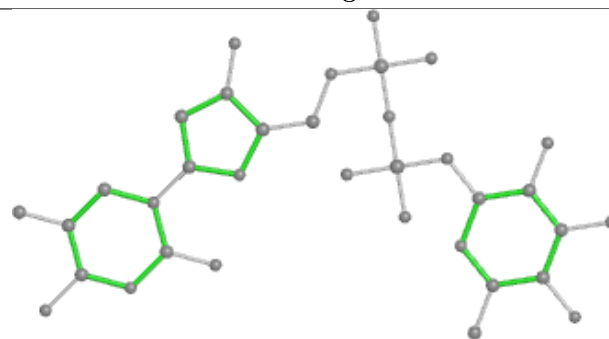
Bond lengths



Bond angles

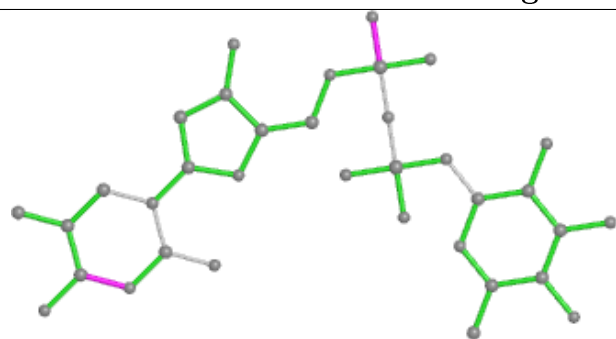


Torsions

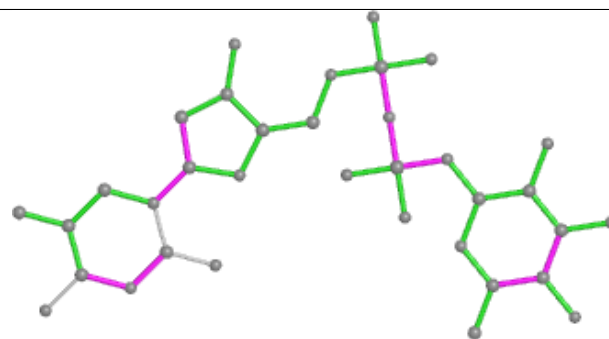


Rings

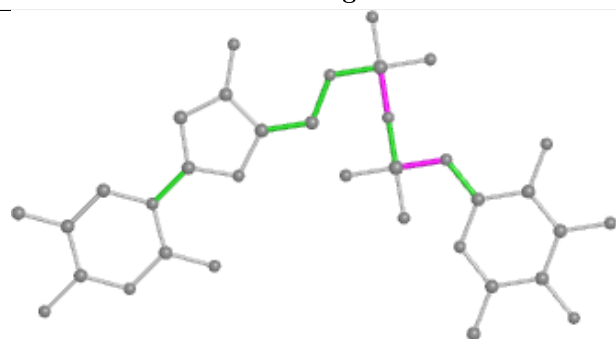
Ligand T3F B 301



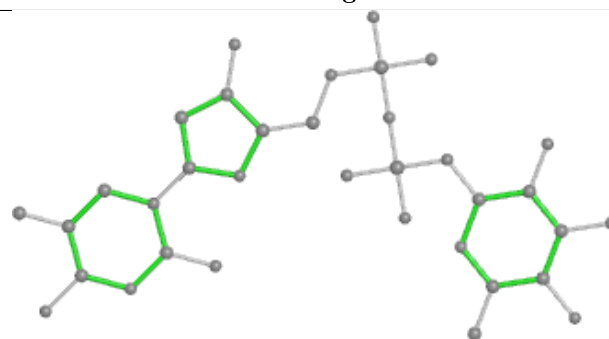
Bond lengths



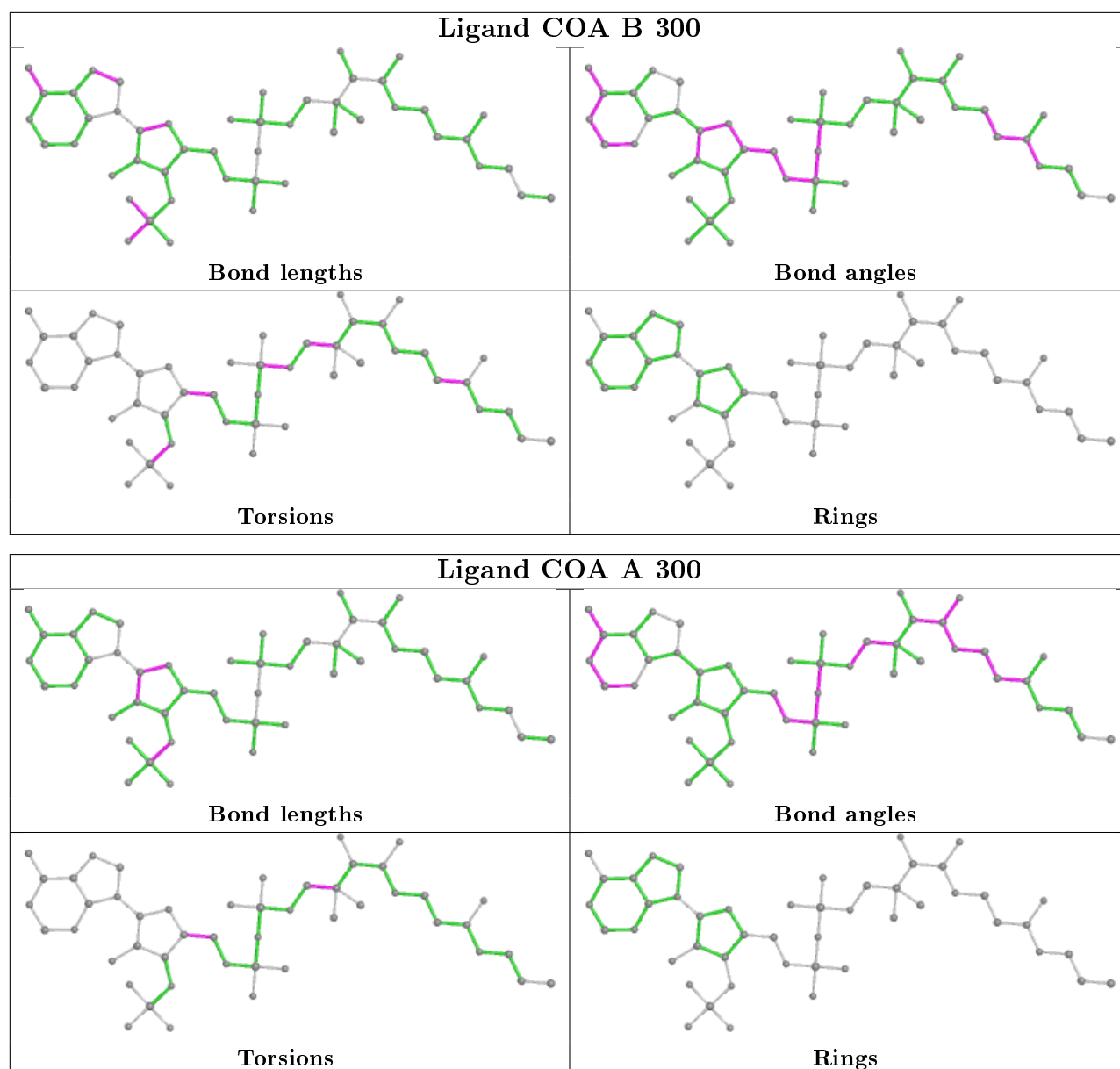
Bond angles



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 [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	259/273 (94%)	-0.40	1 (0%) 92 90	5, 15, 43, 74	0
1	B	257/273 (94%)	-0.55	0 100 100	5, 12, 40, 60	0
All	All	516/546 (94%)	-0.47	1 (0%) 95 93	5, 13, 43, 74	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	261	ILE	3.2

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

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

6.3 Carbohydrates [i](#)

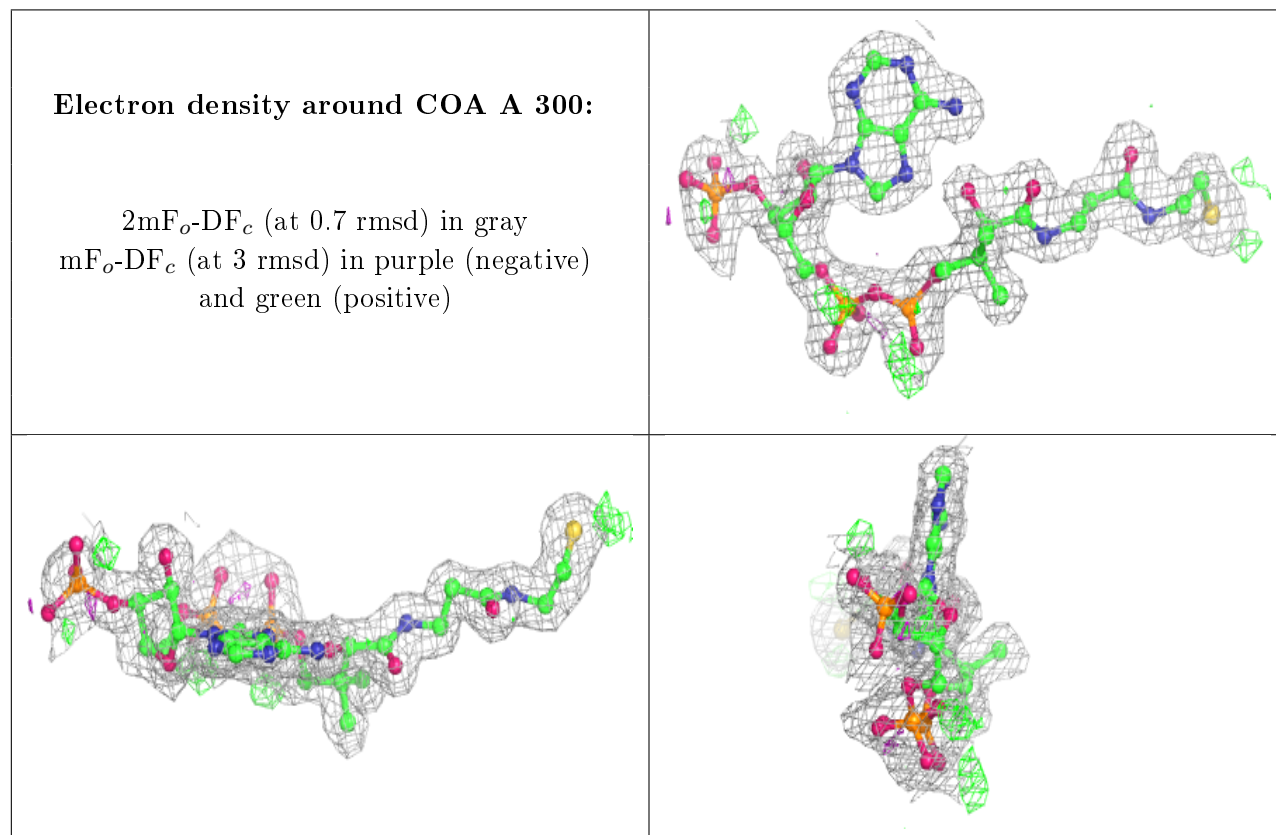
There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 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.

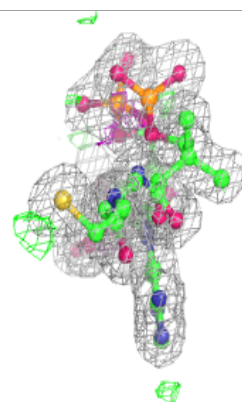
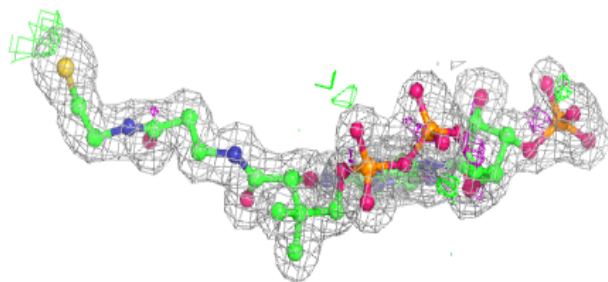
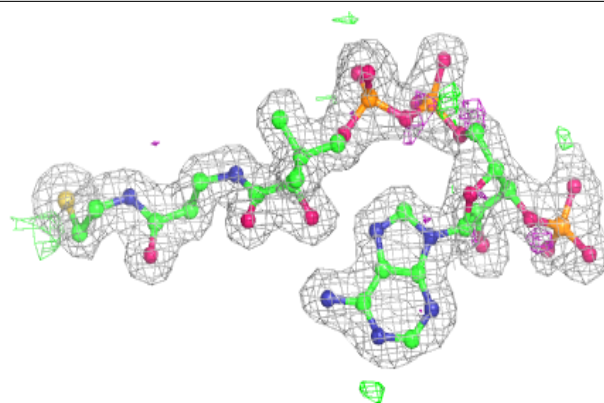
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	COA	A	300	48/48	0.95	0.09	6,21,63,100	0
2	COA	B	300	48/48	0.96	0.08	6,13,38,51	0
3	T3F	A	301	35/35	0.97	0.07	9,15,23,27	0
3	T3F	B	301	35/35	0.97	0.07	6,13,26,31	0

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

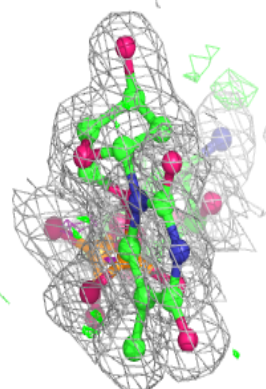
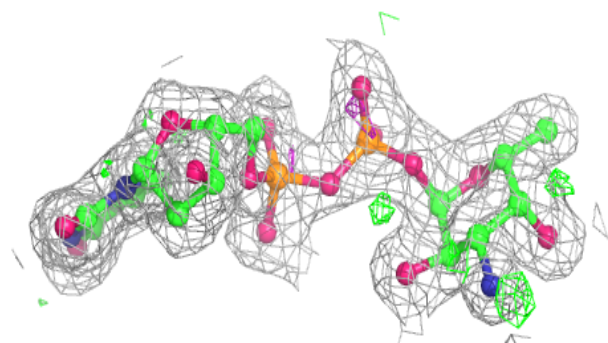
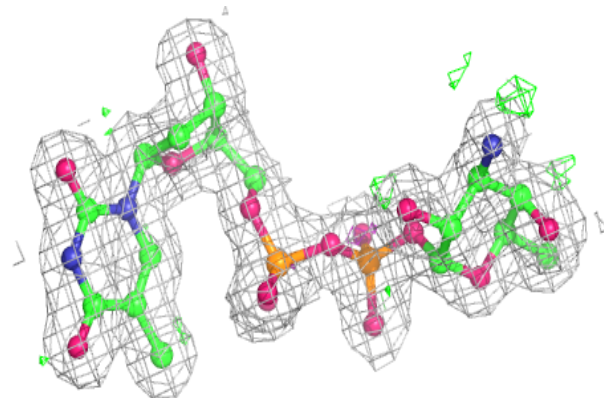


Electron density around COA B 300:

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

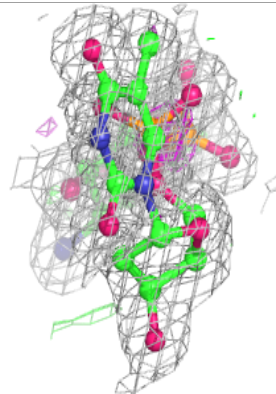
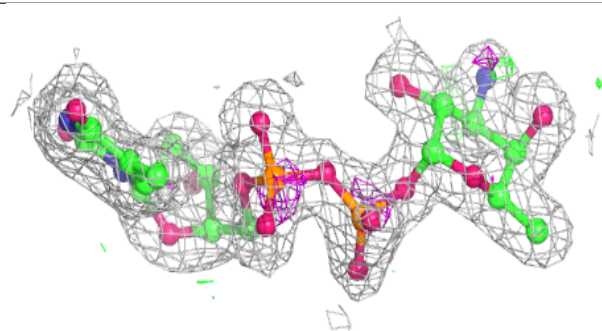
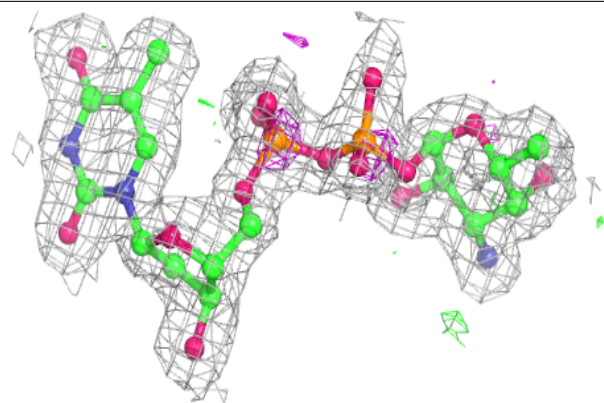
**Electron density around T3F A 301:**

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



Electron density around T3F B 301:

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



6.5 Other polymers [i](#)

There are no such residues in this entry.