



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 04:17 PM UTC

PDB ID : 8OW5 / pdb_00008ow5
Title : Crystal structure of Mycobacterium smegmatis CoaB in complex with CTP and 5-methoxy-1H-indole-2-carboxylic acid
Authors : Mendes, V.; Blundell, T.L.
Deposited on : 2023-04-27
Resolution : 2.84 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 : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

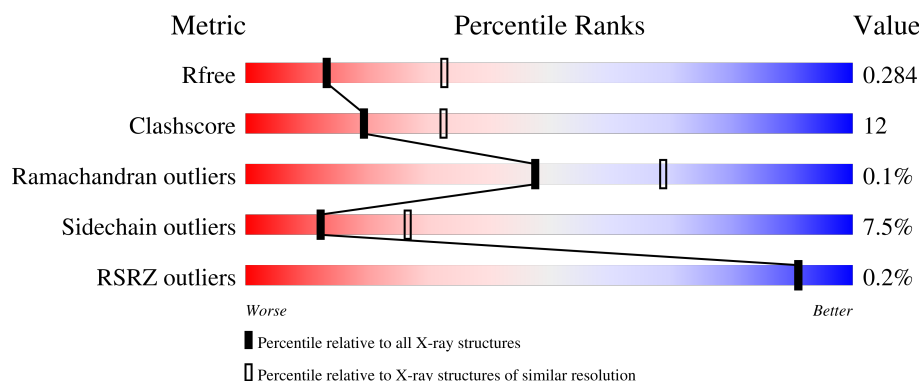
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.84 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}	180053	1520 (2.86-2.82)
Clashscore	190562	1559 (2.86-2.82)
Ramachandran outliers	187476	1517 (2.86-2.82)
Sidechain outliers	187428	1518 (2.86-2.82)
RSRZ outliers	180081	1521 (2.86-2.82)

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	237	
1	B	237	
1	C	237	
1	D	237	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6102 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 Coenzyme A biosynthesis bifunctional protein CoaBC.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	209	Total	C	N	O	S	0	0	0
			1466	913	266	281	6			
1	B	224	Total	C	N	O	S	0	0	0
			1539	954	284	295	6			
1	C	213	Total	C	N	O	S	0	0	0
			1477	916	274	282	5			
1	D	213	Total	C	N	O	S	0	0	0
			1486	925	279	276	6			

There are 32 discrepancies between the modelled and reference sequences:

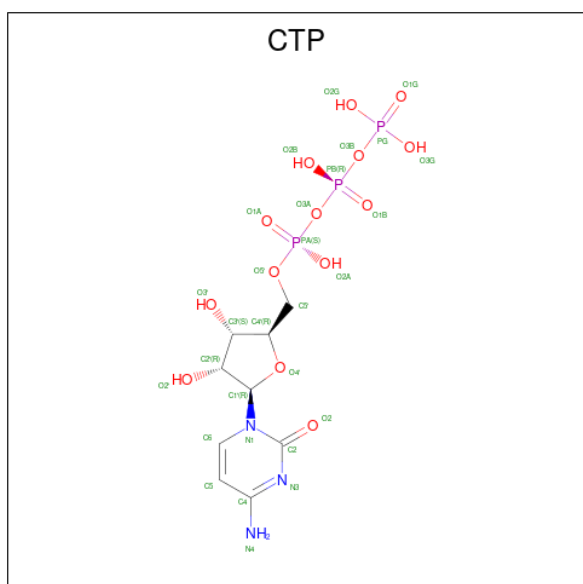
Chain	Residue	Modelled	Actual	Comment	Reference
A	178	MET	-	initiating methionine	UNP A0QWT2
A	179	ALA	-	expression tag	UNP A0QWT2
A	180	HIS	-	expression tag	UNP A0QWT2
A	181	HIS	-	expression tag	UNP A0QWT2
A	182	HIS	-	expression tag	UNP A0QWT2
A	183	HIS	-	expression tag	UNP A0QWT2
A	184	HIS	-	expression tag	UNP A0QWT2
A	185	HIS	-	expression tag	UNP A0QWT2
B	178	MET	-	initiating methionine	UNP A0QWT2
B	179	ALA	-	expression tag	UNP A0QWT2
B	180	HIS	-	expression tag	UNP A0QWT2
B	181	HIS	-	expression tag	UNP A0QWT2
B	182	HIS	-	expression tag	UNP A0QWT2
B	183	HIS	-	expression tag	UNP A0QWT2
B	184	HIS	-	expression tag	UNP A0QWT2
B	185	HIS	-	expression tag	UNP A0QWT2
C	178	MET	-	initiating methionine	UNP A0QWT2
C	179	ALA	-	expression tag	UNP A0QWT2
C	180	HIS	-	expression tag	UNP A0QWT2
C	181	HIS	-	expression tag	UNP A0QWT2
C	182	HIS	-	expression tag	UNP A0QWT2

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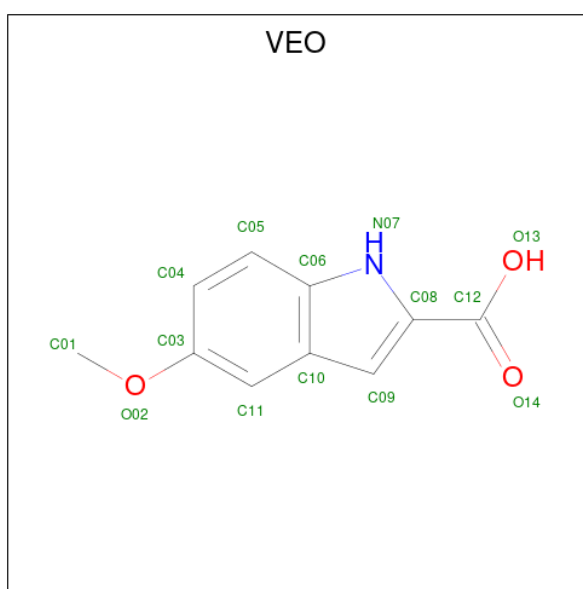
Chain	Residue	Modelled	Actual	Comment	Reference
C	183	HIS	-	expression tag	UNP A0QWT2
C	184	HIS	-	expression tag	UNP A0QWT2
C	185	HIS	-	expression tag	UNP A0QWT2
D	178	MET	-	initiating methionine	UNP A0QWT2
D	179	ALA	-	expression tag	UNP A0QWT2
D	180	HIS	-	expression tag	UNP A0QWT2
D	181	HIS	-	expression tag	UNP A0QWT2
D	182	HIS	-	expression tag	UNP A0QWT2
D	183	HIS	-	expression tag	UNP A0QWT2
D	184	HIS	-	expression tag	UNP A0QWT2
D	185	HIS	-	expression tag	UNP A0QWT2

- Molecule 2 is CYTIDINE-5'-TRIPHOSPHATE (CCD ID: CTP) (formula: $C_9H_{16}N_3O_{14}P_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Ca 1 1	0	0
3	B	1	Total Ca 1 1	0	0
3	C	1	Total Ca 1 1	0	0
3	D	1	Total Ca 1 1	0	0

- Molecule 4 is 5-methoxy-1H-indole-2-carboxylic acid (CCD ID: VEO) (formula: $C_{10}H_9NO_3$) (labeled as "Ligand of Interest" by depositor).

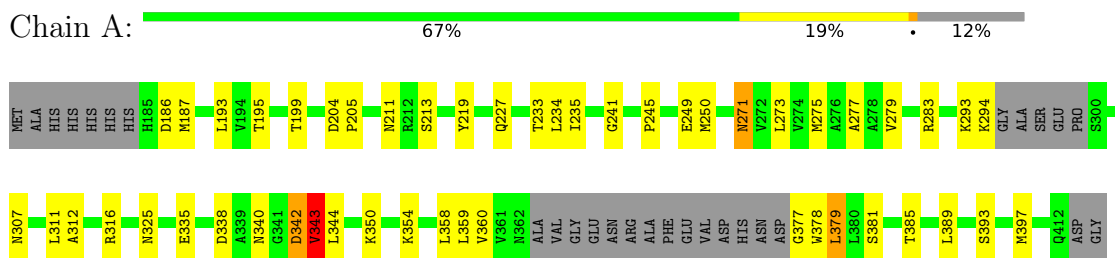


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total C N O 14 10 1 3	0	0

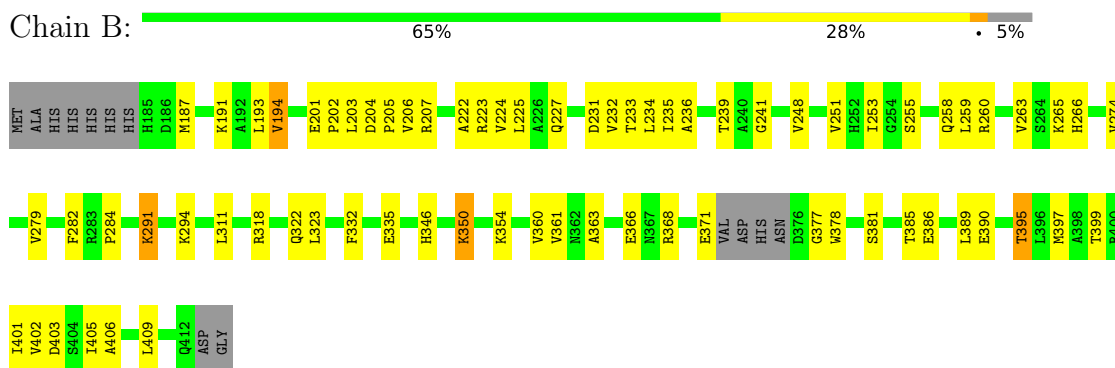
3 Residue-property plots

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.

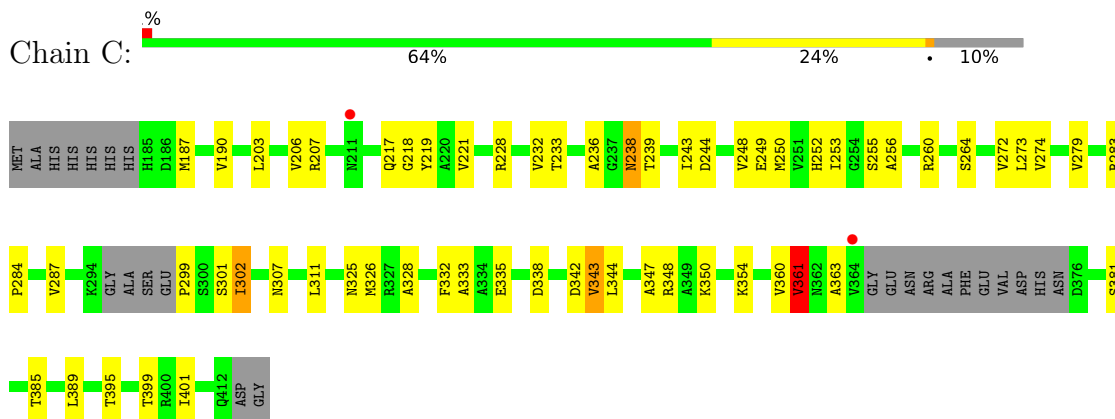
- Molecule 1: Coenzyme A biosynthesis bifunctional protein CoaBC



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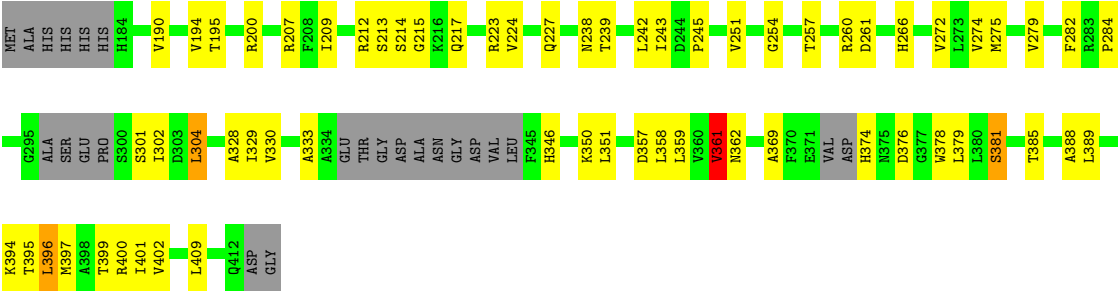
- Molecule 1: Coenzyme A biosynthesis bifunctional protein CoaBC

Chain D:

63%

25%

10%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	76.73Å 77.48Å 144.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	68.31 – 2.84 68.31 – 2.84	Depositor EDS
% Data completeness (in resolution range)	100.0 (68.31-2.84) 97.3 (68.31-2.84)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.12 (at 2.86Å)	Xtriage
Refinement program	PHENIX (1.20.1_4487: ???)	Depositor
R, R_{free}	0.218 , 0.295 0.216 , 0.284	Depositor DCC
R_{free} test set	1036 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	65.3	Xtriage
Anisotropy	0.313	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 144.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.208 for k,h,-l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	6102	wwPDB-VP
Average B, all atoms (Å ²)	72.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.37% of the height of the origin peak. No significant pseudotranslation is detected.*

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: VEO, CTP, CA

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	0.74	2/1483 (0.1%)	0.93	6/2013 (0.3%)
1	B	0.74	3/1557 (0.2%)	0.87	3/2117 (0.1%)
1	C	0.83	2/1495 (0.1%)	0.85	3/2036 (0.1%)
1	D	0.79	4/1503 (0.3%)	0.87	6/2041 (0.3%)
All	All	0.78	11/6038 (0.2%)	0.88	18/8207 (0.2%)

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	338	ASP	C-O	-10.73	1.16	1.24
1	A	340	ASN	CA-C	-6.17	1.44	1.52
1	D	369	ALA	CA-C	-5.79	1.45	1.52
1	D	214	SER	N-CA	-5.79	1.38	1.46
1	D	213	SER	CA-C	-5.65	1.45	1.52
1	A	211	ASN	CA-C	-5.61	1.45	1.52
1	D	376	ASP	C-N	-5.37	1.30	1.33
1	C	238	ASN	CA-C	-5.25	1.46	1.52
1	B	236	ALA	CA-C	-5.19	1.46	1.52
1	B	318	ARG	CA-C	-5.18	1.46	1.52
1	B	239	THR	CA-C	-5.12	1.46	1.52

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	214	SER	N-CA-C	-10.18	100.72	113.55
1	A	340	ASN	N-CA-C	-8.91	102.55	113.50
1	D	215	GLY	N-CA-C	-7.95	104.59	114.92
1	A	344	LEU	N-CA-C	7.58	120.42	111.71
1	D	304	LEU	N-CA-C	7.30	120.51	109.41
1	B	241	GLY	N-CA-C	-7.20	105.37	113.79

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	343	VAL	N-CA-C	-7.02	102.62	110.62
1	B	363	ALA	N-CA-C	6.98	121.72	111.87
1	A	241	GLY	N-CA-C	-6.56	105.25	116.01
1	B	239	THR	CB-CA-C	-6.51	100.17	110.85
1	D	301	SER	CA-CB-OG	6.42	123.94	111.10
1	D	213	SER	N-CA-C	-6.28	98.08	108.07
1	A	271	ASN	N-CA-C	-6.26	105.96	113.97
1	C	361	VAL	N-CA-C	5.57	117.00	108.71
1	C	363	ALA	N-CA-C	5.56	119.40	107.67
1	A	213	SER	N-CA-C	5.38	118.69	110.20
1	D	361	VAL	N-CA-C	5.20	115.39	108.11
1	C	338	ASP	N-CA-C	5.10	115.27	108.23

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	1466	0	1436	29	0
1	B	1539	0	1485	41	0
1	C	1477	0	1424	35	0
1	D	1486	0	1450	37	0
2	A	29	0	12	2	0
2	B	29	0	12	2	0
2	C	29	0	12	1	0
2	D	29	0	12	3	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	B	14	0	0	0	0
All	All	6102	0	5843	138	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 12.

All (138) 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:381:SER:HB3	1:D:385:THR:HB	1.52	0.90
1:D:351:LEU:HB2	1:D:359:LEU:HD12	1.63	0.81
1:A:335:GLU:HB2	1:A:343:VAL:HG13	1.63	0.80
1:B:224:VAL:HG11	1:B:402:VAL:HG11	1.62	0.79
1:A:195:THR:HB	1:A:277:ALA:HB2	1.63	0.78
1:B:191:LYS:HA	1:B:231:ASP:HB3	1.67	0.77
1:B:311:LEU:HD23	1:B:354:LYS:HD3	1.67	0.77
1:C:335:GLU:HG3	1:C:347:ALA:HB2	1.66	0.75
1:B:401:ILE:O	1:B:405:ILE:HG13	1.88	0.74
1:D:359:LEU:HD22	1:D:361:VAL:HG12	1.71	0.73
1:B:381:SER:HB3	1:B:385:THR:HB	1.72	0.72
1:D:260:ARG:NH1	1:D:261:ASP:OD1	2.23	0.71
1:C:232:VAL:HB	1:C:248:VAL:HG22	1.70	0.71
1:B:350:LYS:NZ	2:B:502:CTP:O3A	2.26	0.68
1:D:328:ALA:HA	1:D:357:ASP:OD2	1.94	0.68
1:D:272:VAL:HG22	1:D:328:ALA:HB3	1.77	0.66
1:C:311:LEU:HD23	1:C:354:LYS:HD2	1.79	0.65
1:B:397:MET:O	1:B:401:ILE:HG13	1.97	0.64
1:B:232:VAL:HB	1:B:248:VAL:HG22	1.80	0.63
1:B:193:LEU:HD11	1:B:235:ILE:HD12	1.80	0.63
1:C:332:PHE:O	2:C:501:CTP:H5'1	1.99	0.62
1:D:209:ILE:HD12	1:D:284:PRO:HD3	1.80	0.62
1:B:203:LEU:HD22	1:B:284:PRO:HG3	1.81	0.62
1:D:223:ARG:O	1:D:227:GLN:HG3	1.99	0.62
1:C:360:VAL:HG11	1:C:401:ILE:HD11	1.81	0.62
1:C:381:SER:HB2	1:C:385:THR:HB	1.82	0.61
1:D:350:LYS:NZ	2:D:501:CTP:O2A	2.34	0.60
1:C:244:ASP:HB3	1:C:250:MET:HG3	1.83	0.60
1:A:381:SER:HB2	1:A:385:THR:HB	1.83	0.59
1:D:389:LEU:HD13	1:D:397:MET:HE1	1.84	0.58
1:A:275:MET:HG3	1:A:311:LEU:HD13	1.84	0.58
1:B:222:ALA:HA	1:B:225:LEU:HD12	1.85	0.58
1:A:381:SER:CB	1:A:385:THR:HB	2.34	0.57
1:C:344:LEU:HG	1:C:348:ARG:CZ	2.34	0.57
1:D:379:LEU:HD11	1:D:389:LEU:HD23	1.86	0.57
1:A:358:LEU:HG	1:A:359:LEU:N	2.20	0.56
1:C:283:ARG:NH1	1:C:307:ASN:OD1	2.37	0.56
1:D:397:MET:O	1:D:401:ILE:HG13	2.05	0.55
1:A:193:LEU:HD11	1:A:235:ILE:HD12	1.89	0.55
1:B:204:ASP:HB2	1:B:205:PRO:HD2	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:260:ARG:HH11	1:B:260:ARG:HG2	1.71	0.55
1:B:395:THR:O	1:B:399:THR:OG1	2.20	0.55
1:A:293:LYS:NZ	1:B:350:LYS:HD2	2.21	0.55
1:C:233:THR:HG23	1:C:249:GLU:HB3	1.89	0.55
1:C:187:MET:O	1:C:190:VAL:HG12	2.07	0.54
1:C:395:THR:O	1:C:399:THR:HG23	2.08	0.53
1:B:224:VAL:CG1	1:B:402:VAL:HG11	2.34	0.53
1:D:223:ARG:HG3	1:D:245:PRO:HB3	1.91	0.53
1:B:335:GLU:OE2	1:B:350:LYS:HE2	2.08	0.53
1:D:239:THR:HB	1:D:242:LEU:HB2	1.91	0.53
1:D:224:VAL:HG11	1:D:402:VAL:HG11	1.91	0.53
1:D:330:VAL:HG13	1:D:358:LEU:HB2	1.90	0.53
1:D:217:GLN:HB2	1:D:394:LYS:HE3	1.91	0.52
1:B:368:ARG:O	1:B:371:GLU:HB2	2.10	0.52
1:A:293:LYS:HZ3	1:B:350:LYS:HD2	1.74	0.52
1:B:206:VAL:HG13	1:B:207:ARG:HG3	1.92	0.52
1:B:274:VAL:HG13	1:B:332:PHE:HE1	1.74	0.52
1:B:399:THR:O	1:B:403:ASP:HB2	2.10	0.51
1:A:273:LEU:HD11	1:A:275:MET:HE3	1.93	0.51
1:A:283:ARG:HD3	1:A:307:ASN:OD1	2.10	0.51
1:A:381:SER:OG	1:A:385:THR:HB	2.10	0.51
1:B:322:GLN:O	1:B:323:LEU:HD23	2.11	0.51
1:B:223:ARG:O	1:B:227:GLN:HG3	2.11	0.50
1:A:377:GLY:C	1:A:389:LEU:HB2	2.37	0.50
1:D:279:VAL:HG22	2:D:501:CTP:C6	2.46	0.50
1:A:245:PRO:HD2	1:A:250:MET:SD	2.53	0.49
1:C:333:ALA:O	1:C:361:VAL:HA	2.12	0.49
1:B:406:ALA:HA	1:B:409:LEU:HD12	1.94	0.49
1:B:204:ASP:OD2	1:B:291:LYS:HA	2.12	0.49
1:A:234:LEU:HD23	1:A:250:MET:SD	2.53	0.48
1:B:259:LEU:HD23	1:B:263:VAL:HG23	1.95	0.48
1:C:326:MET:HB2	1:C:326:MET:HE2	1.64	0.48
1:D:224:VAL:CG1	1:D:402:VAL:HG11	2.44	0.48
1:A:271:ASN:HA	1:A:325:ASN:O	2.13	0.48
1:B:194:VAL:HG22	1:B:234:LEU:HD13	1.94	0.48
1:B:294:LYS:HG2	1:B:294:LYS:O	2.12	0.48
1:D:274:VAL:HG12	1:D:330:VAL:HB	1.95	0.48
1:C:344:LEU:O	1:C:348:ARG:HB2	2.14	0.48
1:B:187:MET:HE3	1:B:402:VAL:HG23	1.96	0.48
1:C:342:ASP:O	1:C:343:VAL:C	2.54	0.48
1:A:195:THR:HA	1:A:235:ILE:O	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:350:LYS:NZ	2:D:501:CTP:PA	2.87	0.47
1:C:187:MET:HB3	1:C:190:VAL:HG11	1.96	0.47
1:C:236:ALA:O	1:C:252:HIS:HA	2.15	0.47
1:C:299:PRO:HG2	1:D:282:PHE:CZ	2.50	0.47
1:A:187:MET:HE3	1:A:187:MET:HB3	1.75	0.47
1:B:378:TRP:HZ3	1:B:386:GLU:O	1.96	0.47
1:B:255:SER:OG	1:B:258:GLN:HG3	2.14	0.47
1:C:273:LEU:HD12	1:C:274:VAL:N	2.30	0.47
1:D:357:ASP:O	1:D:358:LEU:HD23	2.14	0.47
1:C:252:HIS:O	1:C:253:ILE:HD12	2.15	0.47
1:C:335:GLU:OE2	1:C:350:LYS:HE3	2.15	0.47
1:A:393:SER:O	1:A:397:MET:HG3	2.14	0.46
1:D:409:LEU:HD23	1:D:409:LEU:HA	1.64	0.46
1:A:204:ASP:HB2	1:A:205:PRO:HD2	1.97	0.46
1:A:350:LYS:NZ	2:A:501:CTP:O3B	2.46	0.46
1:B:366:GLU:CB	1:B:371:GLU:HB3	2.46	0.46
1:A:312:ALA:O	1:A:316:ARG:HG2	2.15	0.46
1:C:255:SER:OG	1:C:256:ALA:N	2.49	0.46
1:B:390:GLU:O	1:B:397:MET:HE1	2.16	0.46
1:D:195:THR:OG1	1:D:275:MET:HA	2.16	0.46
1:B:202:PRO:O	1:B:284:PRO:HG2	2.16	0.45
1:C:260:ARG:HH11	1:C:260:ARG:HG2	1.81	0.45
1:C:217:GLN:HG2	1:C:218:GLY:N	2.32	0.45
1:D:243:ILE:O	1:D:243:ILE:HG13	2.16	0.45
1:B:377:GLY:C	1:B:389:LEU:H	2.24	0.44
1:C:203:LEU:HD13	1:C:284:PRO:CG	2.48	0.44
1:C:217:GLN:HG2	1:C:218:GLY:H	1.82	0.44
1:D:396:LEU:O	1:D:400:ARG:HG2	2.18	0.44
1:D:194:VAL:HG13	1:D:274:VAL:HG23	1.99	0.44
1:C:302:ILE:HB	1:D:302:ILE:HB	2.00	0.44
1:A:342:ASP:O	1:A:343:VAL:C	2.61	0.44
1:C:273:LEU:HD12	1:C:274:VAL:H	1.83	0.43
1:D:378:TRP:HA	1:D:388:ALA:HA	2.00	0.43
1:D:251:VAL:HG21	1:D:266:HIS:NE2	2.33	0.43
1:A:358:LEU:HD11	1:A:379:LEU:HG	2.01	0.43
1:D:200:ARG:HH12	1:D:238:ASN:ND2	2.17	0.42
1:D:395:THR:O	1:D:399:THR:HG23	2.19	0.42
1:A:279:VAL:HB	2:A:501:CTP:C5	2.55	0.42
1:B:193:LEU:HD12	1:B:233:THR:O	2.20	0.42
1:D:361:VAL:HG23	1:D:362:ASN:OD1	2.19	0.42
1:D:333:ALA:O	1:D:361:VAL:HA	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:219:TYR:CE2	1:C:239:THR:HG21	2.54	0.42
1:D:389:LEU:HD12	1:D:389:LEU:O	2.20	0.42
1:B:265:LYS:HG2	1:B:266:HIS:ND1	2.35	0.41
1:A:294:LYS:HG2	2:B:502:CTP:PG	2.61	0.41
1:C:283:ARG:HH12	1:C:307:ASN:HA	1.84	0.41
1:B:201:GLU:HG3	1:B:282:PHE:O	2.21	0.41
1:B:204:ASP:OD1	1:B:204:ASP:C	2.63	0.41
1:C:206:VAL:HG23	1:C:207:ARG:N	2.35	0.41
1:C:260:ARG:O	1:C:264:SER:N	2.42	0.41
1:B:260:ARG:HG2	1:B:260:ARG:NH1	2.35	0.41
1:C:272:VAL:HG22	1:C:328:ALA:HB3	2.02	0.41
1:A:350:LYS:O	1:A:354:LYS:HB2	2.21	0.41
1:C:187:MET:HB3	1:C:190:VAL:CG1	2.51	0.41
1:A:219:TYR:CE2	1:A:245:PRO:HD3	2.56	0.41
1:A:233:THR:HA	1:A:249:GLU:HB3	2.04	0.40
1:D:329:ILE:O	1:D:357:ASP:HB2	2.21	0.40

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	203/237 (86%)	198 (98%)	5 (2%)	0	100	100
1	B	220/237 (93%)	211 (96%)	9 (4%)	0	100	100
1	C	207/237 (87%)	198 (96%)	9 (4%)	0	100	100
1	D	205/237 (86%)	194 (95%)	10 (5%)	1 (0%)	24	43
All	All	835/948 (88%)	801 (96%)	33 (4%)	1 (0%)	48	69

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	254	GLY

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	138/173 (80%)	129 (94%)	9 (6%)	15	32
1	B	139/173 (80%)	129 (93%)	10 (7%)	13	28
1	C	136/173 (79%)	124 (91%)	12 (9%)	9	21
1	D	137/173 (79%)	127 (93%)	10 (7%)	13	28
All	All	550/692 (80%)	509 (92%)	41 (8%)	12	27

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	186	ASP
1	A	199	THR
1	A	227	GLN
1	A	338	ASP
1	A	342	ASP
1	A	343	VAL
1	A	360	VAL
1	A	378	TRP
1	A	379	LEU
1	B	194	VAL
1	B	251	VAL
1	B	253	ILE
1	B	279	VAL
1	B	291	LYS
1	B	346	HIS
1	B	350	LYS
1	B	360	VAL
1	B	361	VAL
1	B	395	THR
1	C	221	VAL
1	C	228	ARG

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Mol	Chain	Res	Type
1	C	238	ASN
1	C	243	ILE
1	C	279	VAL
1	C	287	VAL
1	C	301	SER
1	C	302	ILE
1	C	325	ASN
1	C	343	VAL
1	C	361	VAL
1	C	389	LEU
1	D	190	VAL
1	D	207	ARG
1	D	212	ARG
1	D	257	THR
1	D	304	LEU
1	D	346	HIS
1	D	361	VAL
1	D	374	HIS
1	D	381	SER
1	D	396	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	286	HIS
1	A	362	ASN
1	B	227	GLN
1	B	271	ASN
1	C	238	ASN
1	C	271	ASN
1	D	211	ASN
1	D	238	ASN

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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 4 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
2	CTP	D	501	3	29,30,30	1.19	4 (13%)	43,47,47	1.15	3 (6%)
2	CTP	B	502	3	29,30,30	0.98	1 (3%)	43,47,47	0.91	3 (6%)
4	VEO	B	501	-	15,15,15	1.67	2 (13%)	19,21,21	1.52	4 (21%)
2	CTP	A	501	3	29,30,30	1.17	4 (13%)	43,47,47	1.31	5 (11%)
2	CTP	C	501	3	29,30,30	0.96	0	43,47,47	1.09	5 (11%)

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
2	CTP	D	501	3	-	7/22/38/38	0/2/2/2
2	CTP	B	502	3	-	5/22/38/38	0/2/2/2
4	VEO	B	501	-	-	6/6/14/14	0/2/2/2
2	CTP	A	501	3	-	9/22/38/38	0/2/2/2
2	CTP	C	501	3	-	5/22/38/38	0/2/2/2

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	501	VEO	C08-N07	-4.89	1.31	1.38
2	D	501	CTP	PB-O3B	3.01	1.62	1.59
2	A	501	CTP	PA-O3A	2.74	1.62	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	501	VEO	C06-N07	-2.50	1.34	1.38
2	A	501	CTP	C2-N1	-2.32	1.35	1.40
2	D	501	CTP	PA-O3A	2.25	1.61	1.59
2	D	501	CTP	C4-N3	2.19	1.38	1.34
2	B	502	CTP	PA-O3A	2.14	1.61	1.59
2	A	501	CTP	C5-C4	-2.07	1.38	1.42
2	D	501	CTP	PB-O3A	2.06	1.61	1.59
2	A	501	CTP	C4-N3	2.01	1.38	1.34

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	CTP	C1'-N1-C2	-3.56	110.57	118.44
4	B	501	VEO	C01-O02-C03	-3.03	111.01	117.50
4	B	501	VEO	C10-C09-C08	-3.02	104.71	107.41
4	B	501	VEO	C05-C06-C10	-2.68	119.70	122.13
2	A	501	CTP	O2B-PB-O3A	2.66	114.47	107.27
2	A	501	CTP	O3'-C3'-C2'	-2.65	103.32	111.82
2	C	501	CTP	O3G-PG-O3B	2.63	113.47	104.64
2	D	501	CTP	C2'-C3'-C4'	2.39	107.22	102.61
2	C	501	CTP	O2B-PB-O3A	2.36	113.64	107.27
2	B	502	CTP	C2'-C1'-N1	-2.35	106.70	113.25
2	C	501	CTP	O3A-PB-O1B	-2.25	103.92	110.70
2	D	501	CTP	C2'-C1'-N1	-2.19	107.15	113.25
2	D	501	CTP	O2-C2-N3	-2.18	118.89	122.33
2	A	501	CTP	C2'-C3'-C4'	2.17	106.80	102.61
2	A	501	CTP	O2B-PB-O3B	2.14	113.07	107.27
2	C	501	CTP	C5-C6-N1	-2.14	118.36	121.84
2	C	501	CTP	O2A-PA-O3A	2.10	112.95	107.27
4	B	501	VEO	C09-C08-N07	2.08	111.21	109.15
2	B	502	CTP	O2G-PG-O3B	2.03	111.46	104.64
2	B	502	CTP	C4-N3-C2	-2.00	117.11	120.26

There are no chirality outliers.

All (32) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	CTP	C5'-O5'-PA-O1A
2	A	501	CTP	C5'-O5'-PA-O2A
2	A	501	CTP	C5'-O5'-PA-O3A
2	B	502	CTP	C5'-O5'-PA-O1A
2	B	502	CTP	C5'-O5'-PA-O2A

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Mol	Chain	Res	Type	Atoms
2	B	502	CTP	C5'-O5'-PA-O3A
2	D	501	CTP	C5'-O5'-PA-O1A
2	D	501	CTP	C5'-O5'-PA-O2A
2	D	501	CTP	C5'-O5'-PA-O3A
4	B	501	VEO	N07-C08-C12-O13
4	B	501	VEO	C04-C03-O02-C01
4	B	501	VEO	C11-C03-O02-C01
2	C	501	CTP	O4'-C4'-C5'-O5'
2	A	501	CTP	C3'-C4'-C5'-O5'
2	A	501	CTP	O4'-C4'-C5'-O5'
2	B	502	CTP	C3'-C4'-C5'-O5'
2	B	502	CTP	O4'-C4'-C5'-O5'
2	D	501	CTP	C3'-C4'-C5'-O5'
2	D	501	CTP	O4'-C4'-C5'-O5'
2	C	501	CTP	C3'-C4'-C5'-O5'
4	B	501	VEO	N07-C08-C12-O14
2	C	501	CTP	PB-O3B-PG-O1G
2	C	501	CTP	PA-O3A-PB-O1B
2	A	501	CTP	PG-O3B-PB-O2B
2	D	501	CTP	PG-O3B-PB-O1B
2	A	501	CTP	PB-O3B-PG-O2G
2	A	501	CTP	PB-O3B-PG-O3G
4	B	501	VEO	C09-C08-C12-O13
4	B	501	VEO	C09-C08-C12-O14
2	C	501	CTP	PA-O3A-PB-O2B
2	A	501	CTP	O4'-C1'-N1-C2
2	D	501	CTP	PB-O3A-PA-O2A

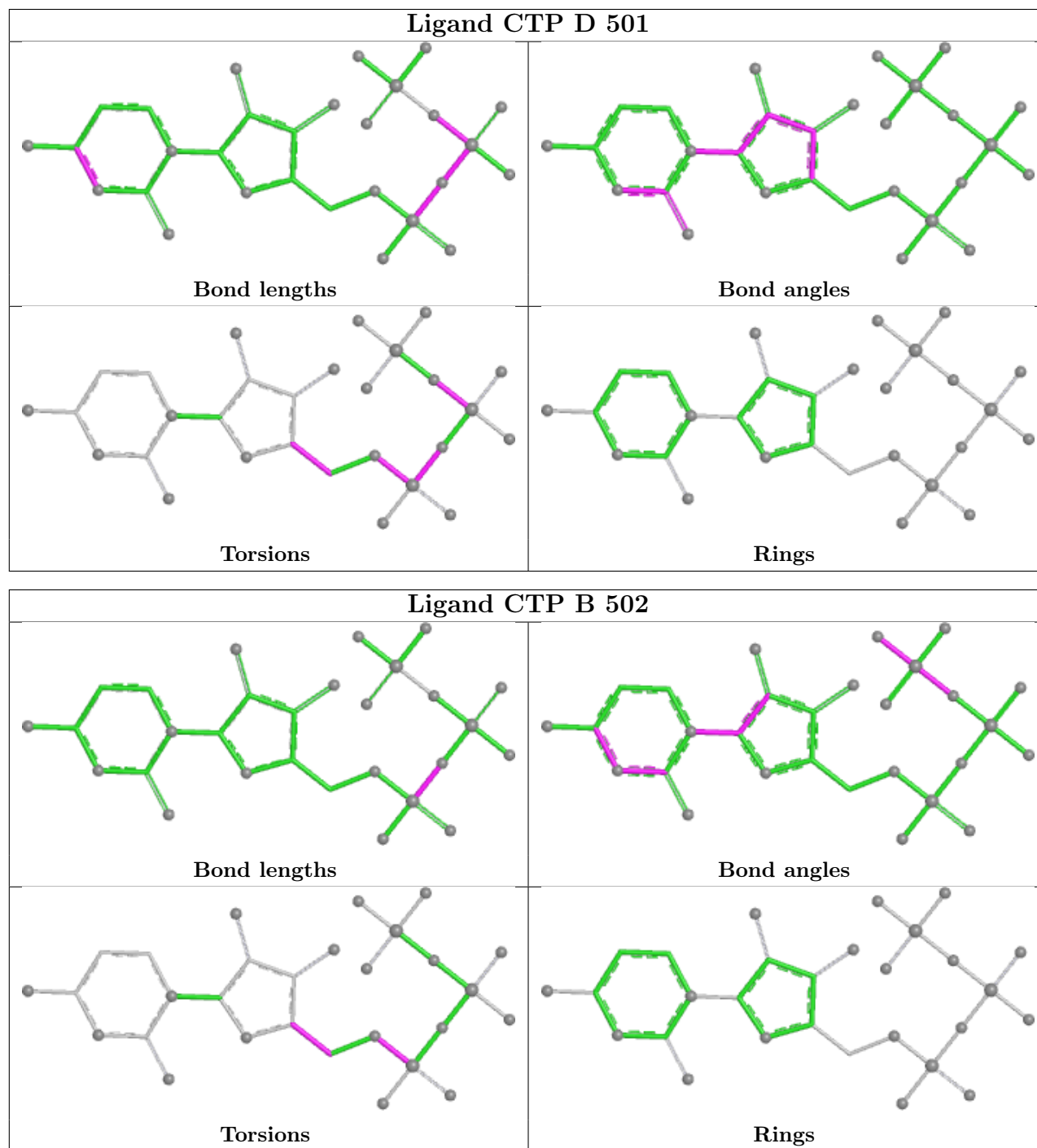
There are no ring outliers.

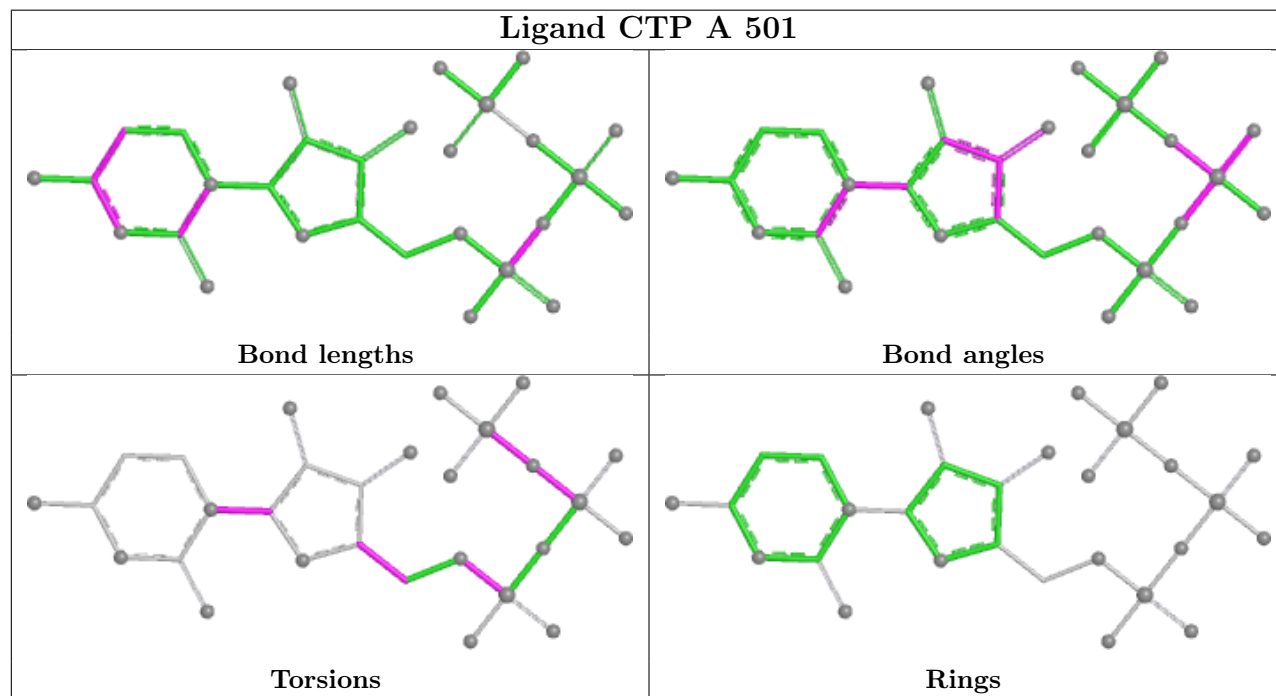
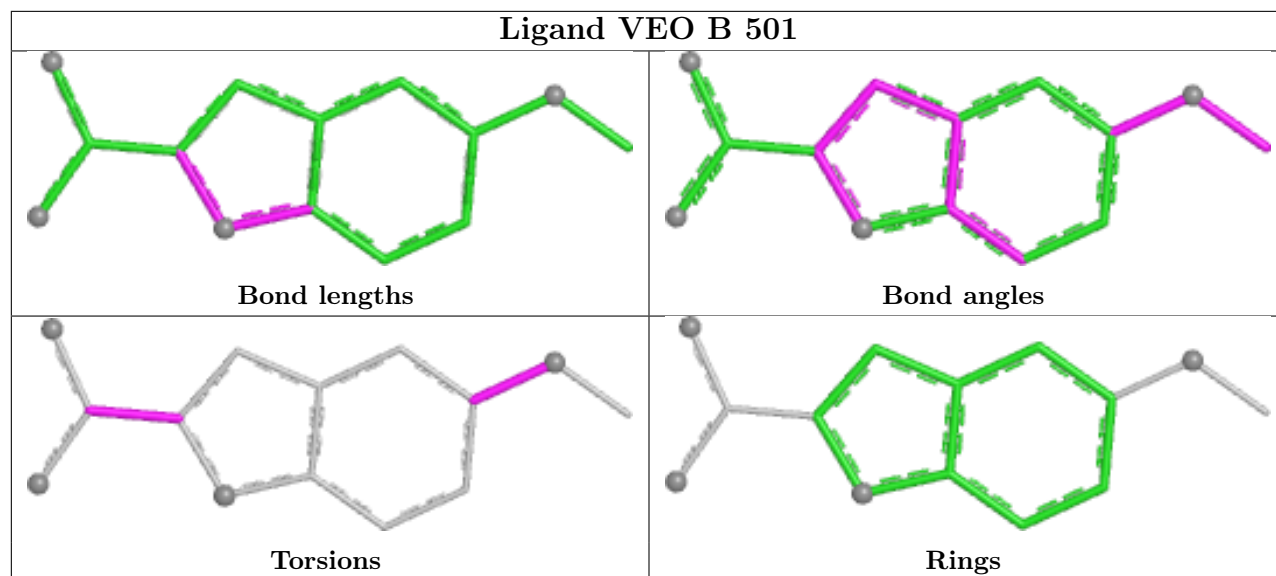
4 monomers are involved in 8 short contacts:

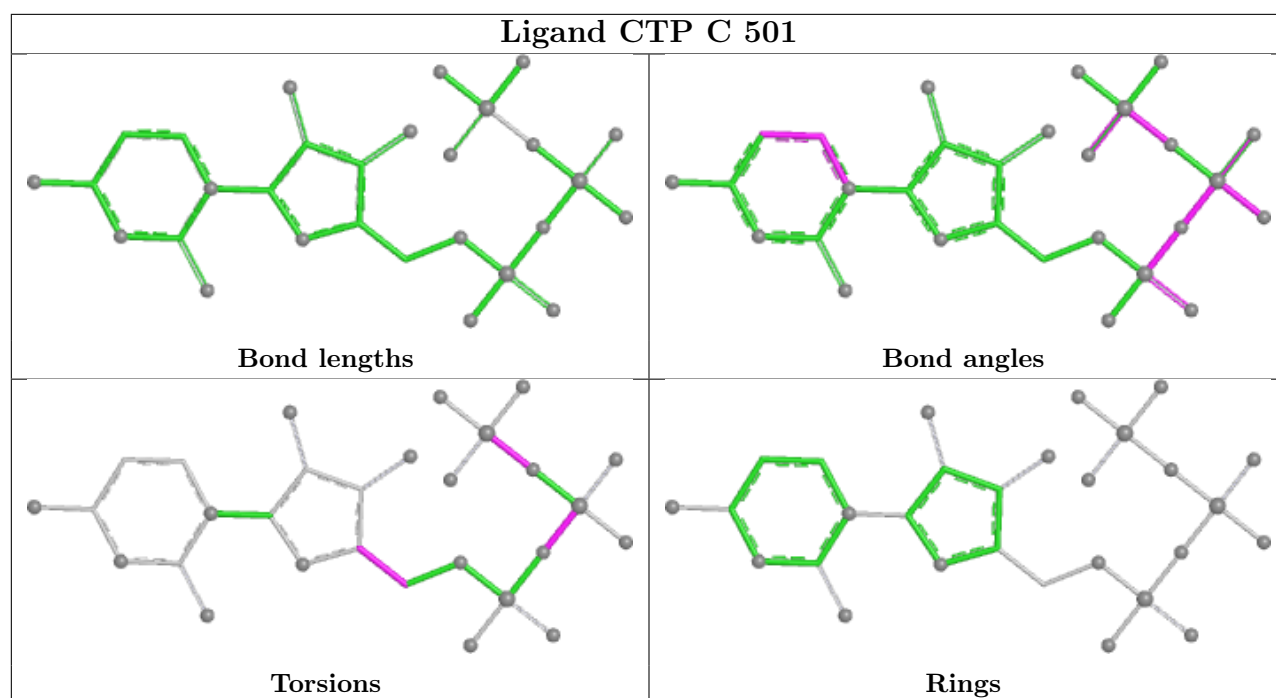
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	501	CTP	3	0
2	B	502	CTP	2	0
2	A	501	CTP	2	0
2	C	501	CTP	1	0

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	209/237 (88%)	-0.21	0 100 100	46, 61, 95, 117	0
1	B	224/237 (94%)	-0.01	0 100 100	48, 70, 115, 153	0
1	C	213/237 (89%)	-0.06	2 (0%) 81 76	50, 67, 110, 155	0
1	D	213/237 (89%)	-0.08	0 100 100	52, 70, 101, 134	0
All	All	859/948 (90%)	-0.09	2 (0%) 91 91	46, 68, 107, 155	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	364	VAL	3.0
1	C	211	ASN	2.2

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](#)

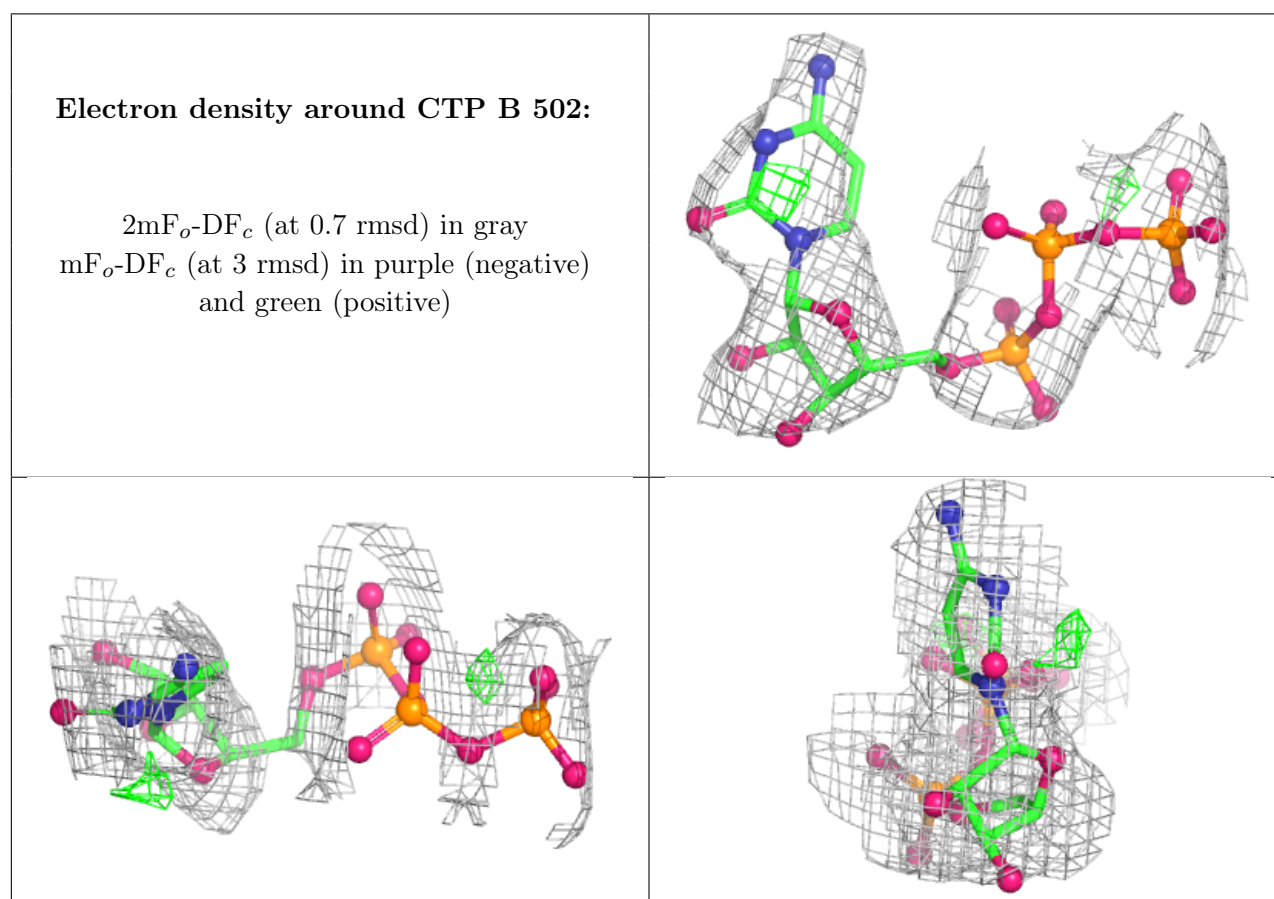
There are no oligosaccharides in this entry.

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.

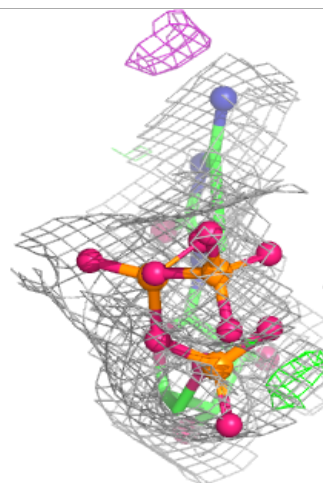
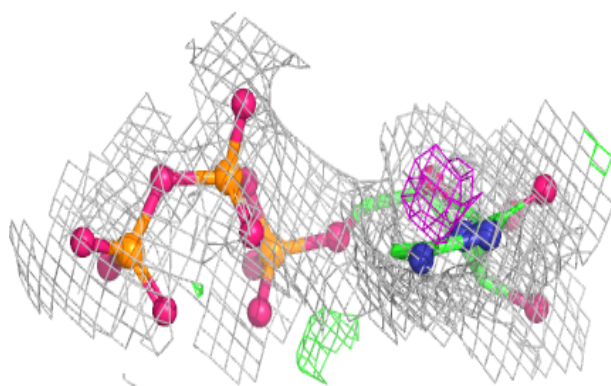
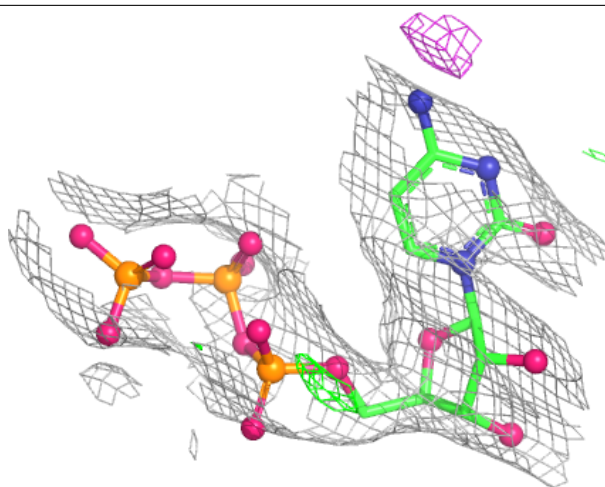
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	CTP	B	502	29/29	0.98	0.07	61,69,79,82	0
2	CTP	C	501	29/29	0.98	0.07	48,53,59,67	0
2	CTP	D	501	29/29	0.98	0.07	56,61,70,73	0
3	CA	A	502	1/1	0.98	0.04	63,63,63,63	0
3	CA	C	502	1/1	0.98	0.04	59,59,59,59	0
4	VEO	B	501	14/14	0.98	0.08	55,63,67,69	0
2	CTP	A	501	29/29	0.99	0.06	48,54,66,68	0
3	CA	D	502	1/1	0.99	0.05	67,67,67,67	0
3	CA	B	503	1/1	0.99	0.04	69,69,69,69	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.



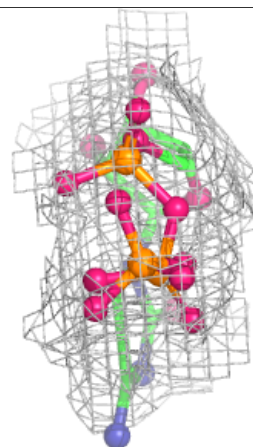
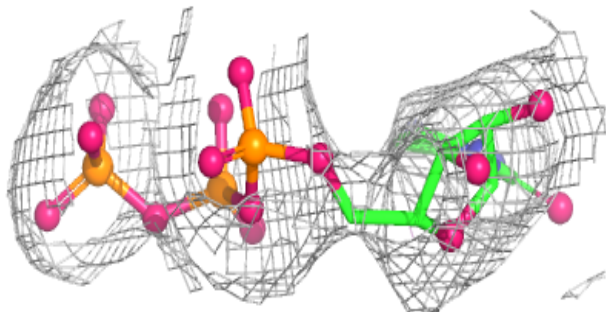
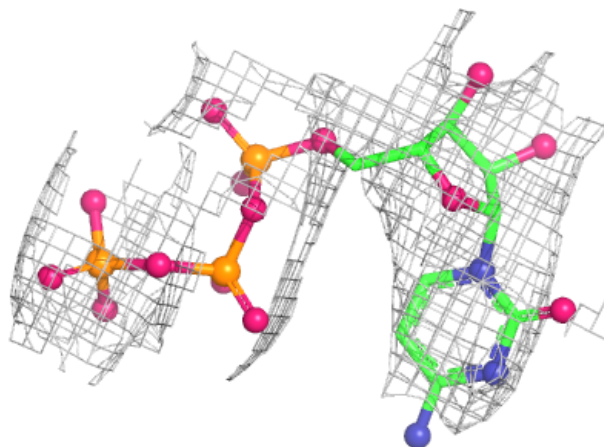
Electron density around CTP C 501:

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



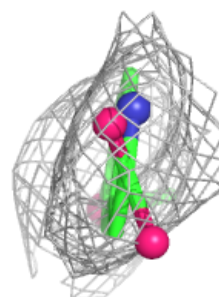
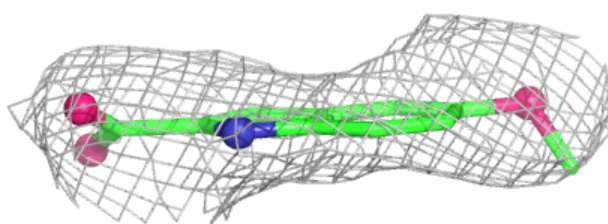
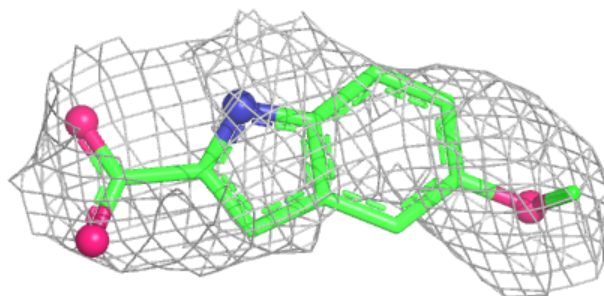
Electron density around CTP D 501:

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

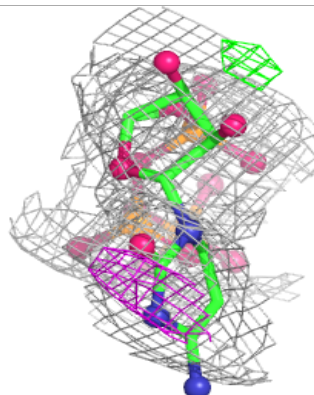
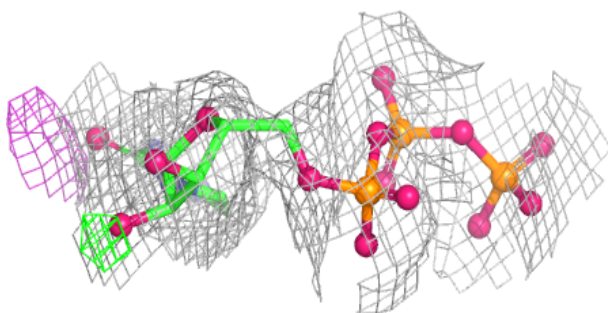
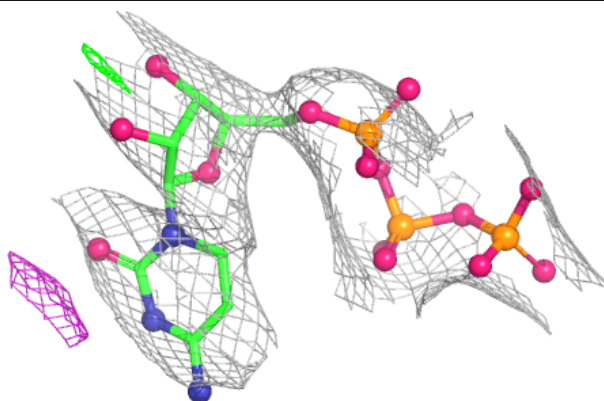


Electron density around VEO B 501:

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

**Electron density around CTP A 501:**

$2mF_o-DF_c$ (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.