



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:41 AM UTC

PDB ID : 6O0J / pdb_00006o0j
Title : M.tb MenD with ThDP and Inhibitor bound
Authors : Johnston, J.M.; Bashiri, G.; Bulloch, E.M.M.; Jirgis, E.M.N.; Nigon, L.V.;
Chuang, H.; Ho, N.A.T.; Baker, E.N.
Deposited on : 2019-02-16
Resolution : 2.35 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

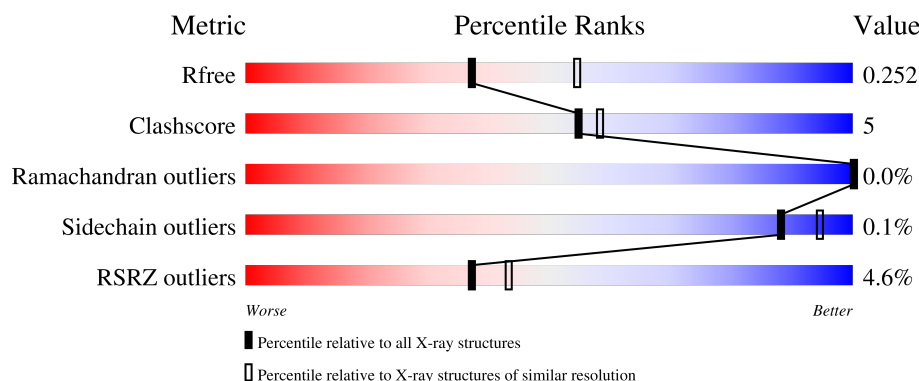
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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	1596 (2.36-2.36)
Clashscore	190562	1663 (2.36-2.36)
Ramachandran outliers	187476	1646 (2.36-2.36)
Sidechain outliers	187428	1646 (2.36-2.36)
RSRZ outliers	180081	1598 (2.36-2.36)

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	574	5% 80% 11% 8%
1	B	574	4% 82% 9% 9%
1	C	574	4% 82% 11% 7%
1	D	574	4% 83% 9% 7%

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 15922 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 2-succinyl-5-enolpyruvyl-6-hydroxy-3-cyclohexene-1-carboxylate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	527	Total	C	N	O	S	0	1	0
			3814	2390	699	715	10			
1	D	532	Total	C	N	O	S	0	2	0
			3855	2415	703	727	10			
1	B	524	Total	C	N	O	S	0	0	0
			3773	2361	688	713	11			
1	C	536	Total	C	N	O	S	0	3	0
			3862	2421	704	727	10			

There are 80 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP P9WK11
A	-18	GLY	-	expression tag	UNP P9WK11
A	-17	SER	-	expression tag	UNP P9WK11
A	-16	SER	-	expression tag	UNP P9WK11
A	-15	HIS	-	expression tag	UNP P9WK11
A	-14	HIS	-	expression tag	UNP P9WK11
A	-13	HIS	-	expression tag	UNP P9WK11
A	-12	HIS	-	expression tag	UNP P9WK11
A	-11	HIS	-	expression tag	UNP P9WK11
A	-10	HIS	-	expression tag	UNP P9WK11
A	-9	SER	-	expression tag	UNP P9WK11
A	-8	SER	-	expression tag	UNP P9WK11
A	-7	GLY	-	expression tag	UNP P9WK11
A	-6	LEU	-	expression tag	UNP P9WK11
A	-5	VAL	-	expression tag	UNP P9WK11
A	-4	PRO	-	expression tag	UNP P9WK11
A	-3	ARG	-	expression tag	UNP P9WK11
A	-2	GLY	-	expression tag	UNP P9WK11
A	-1	SER	-	expression tag	UNP P9WK11
A	0	HIS	-	expression tag	UNP P9WK11

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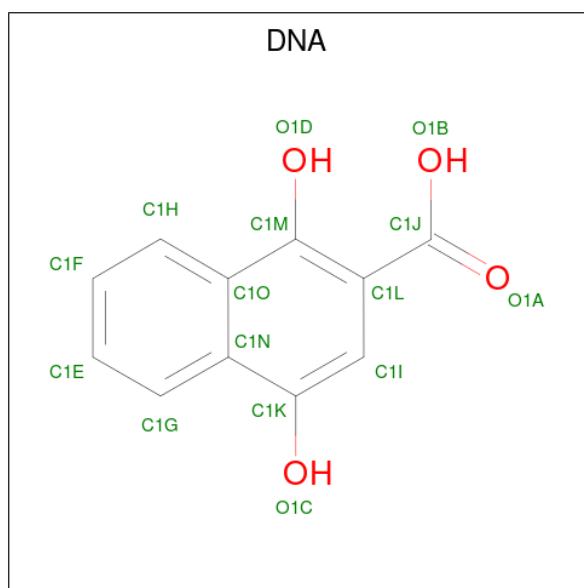
Chain	Residue	Modelled	Actual	Comment	Reference
D	-19	MET	-	initiating methionine	UNP P9WK11
D	-18	GLY	-	expression tag	UNP P9WK11
D	-17	SER	-	expression tag	UNP P9WK11
D	-16	SER	-	expression tag	UNP P9WK11
D	-15	HIS	-	expression tag	UNP P9WK11
D	-14	HIS	-	expression tag	UNP P9WK11
D	-13	HIS	-	expression tag	UNP P9WK11
D	-12	HIS	-	expression tag	UNP P9WK11
D	-11	HIS	-	expression tag	UNP P9WK11
D	-10	HIS	-	expression tag	UNP P9WK11
D	-9	SER	-	expression tag	UNP P9WK11
D	-8	SER	-	expression tag	UNP P9WK11
D	-7	GLY	-	expression tag	UNP P9WK11
D	-6	LEU	-	expression tag	UNP P9WK11
D	-5	VAL	-	expression tag	UNP P9WK11
D	-4	PRO	-	expression tag	UNP P9WK11
D	-3	ARG	-	expression tag	UNP P9WK11
D	-2	GLY	-	expression tag	UNP P9WK11
D	-1	SER	-	expression tag	UNP P9WK11
D	0	HIS	-	expression tag	UNP P9WK11
B	-19	MET	-	initiating methionine	UNP P9WK11
B	-18	GLY	-	expression tag	UNP P9WK11
B	-17	SER	-	expression tag	UNP P9WK11
B	-16	SER	-	expression tag	UNP P9WK11
B	-15	HIS	-	expression tag	UNP P9WK11
B	-14	HIS	-	expression tag	UNP P9WK11
B	-13	HIS	-	expression tag	UNP P9WK11
B	-12	HIS	-	expression tag	UNP P9WK11
B	-11	HIS	-	expression tag	UNP P9WK11
B	-10	HIS	-	expression tag	UNP P9WK11
B	-9	SER	-	expression tag	UNP P9WK11
B	-8	SER	-	expression tag	UNP P9WK11
B	-7	GLY	-	expression tag	UNP P9WK11
B	-6	LEU	-	expression tag	UNP P9WK11
B	-5	VAL	-	expression tag	UNP P9WK11
B	-4	PRO	-	expression tag	UNP P9WK11
B	-3	ARG	-	expression tag	UNP P9WK11
B	-2	GLY	-	expression tag	UNP P9WK11
B	-1	SER	-	expression tag	UNP P9WK11
B	0	HIS	-	expression tag	UNP P9WK11
C	-19	MET	-	initiating methionine	UNP P9WK11
C	-18	GLY	-	expression tag	UNP P9WK11

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-17	SER	-	expression tag	UNP P9WK11
C	-16	SER	-	expression tag	UNP P9WK11
C	-15	HIS	-	expression tag	UNP P9WK11
C	-14	HIS	-	expression tag	UNP P9WK11
C	-13	HIS	-	expression tag	UNP P9WK11
C	-12	HIS	-	expression tag	UNP P9WK11
C	-11	HIS	-	expression tag	UNP P9WK11
C	-10	HIS	-	expression tag	UNP P9WK11
C	-9	SER	-	expression tag	UNP P9WK11
C	-8	SER	-	expression tag	UNP P9WK11
C	-7	GLY	-	expression tag	UNP P9WK11
C	-6	LEU	-	expression tag	UNP P9WK11
C	-5	VAL	-	expression tag	UNP P9WK11
C	-4	PRO	-	expression tag	UNP P9WK11
C	-3	ARG	-	expression tag	UNP P9WK11
C	-2	GLY	-	expression tag	UNP P9WK11
C	-1	SER	-	expression tag	UNP P9WK11
C	0	HIS	-	expression tag	UNP P9WK11

- Molecule 2 is 1,4-dihydroxy-2-naphthoic acid (CCD ID: DNA) (formula: C₁₁H₈O₄) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			15	11	4		
2	D	1	Total	C	O	0	0
			15	11	4		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	C	O	0	0
			15	11	4		
2	C	1	Total	C	O	0	0
			15	11	4		

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



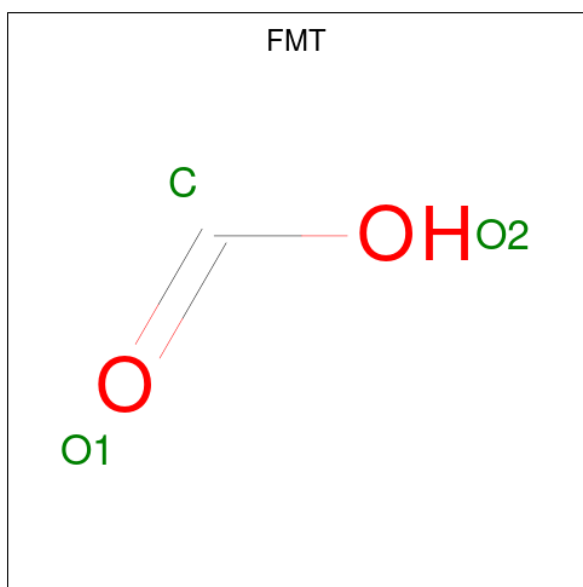
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is ACETATE ION (CCD ID: ACT) (formula: $C_2H_3O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			4	2	2		
4	D	1	Total	C	O	0	0
			4	2	2		
4	C	1	Total	C	O	0	0
			4	2	2		

- Molecule 5 is FORMIC ACID (CCD ID: FMT) (formula: CH_2O_2).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			3	1	2		

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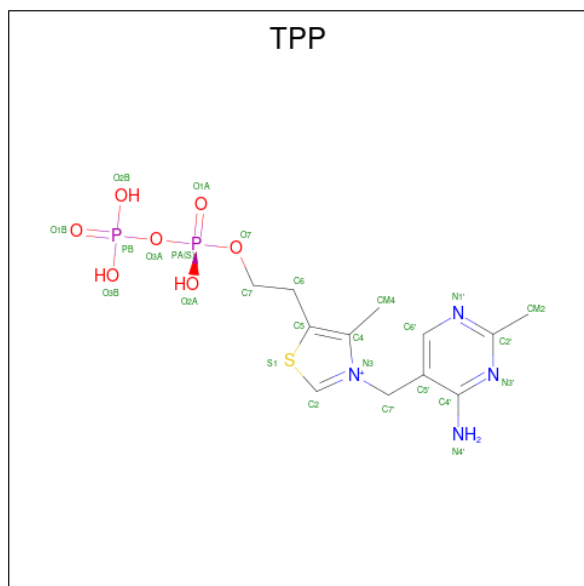
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	D	1	Total	C	O	0	0
			3	1	2		
5	B	1	Total	C	O	0	0
			3	1	2		

- Molecule 6 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	D	1	Total	Mg	0	0
			1	1		
6	C	1	Total	Mg	0	0
			1	1		

- Molecule 7 is THIAMINE DIPHOSPHATE (CCD ID: TPP) (formula: C₁₂H₁₉N₄O₇P₂S) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
7	D	1	Total 26	C 12	N 4	O 7	P 2	S 1	0	0
7	C	1	Total 26	C 12	N 4	O 7	P 2	S 1	0	0

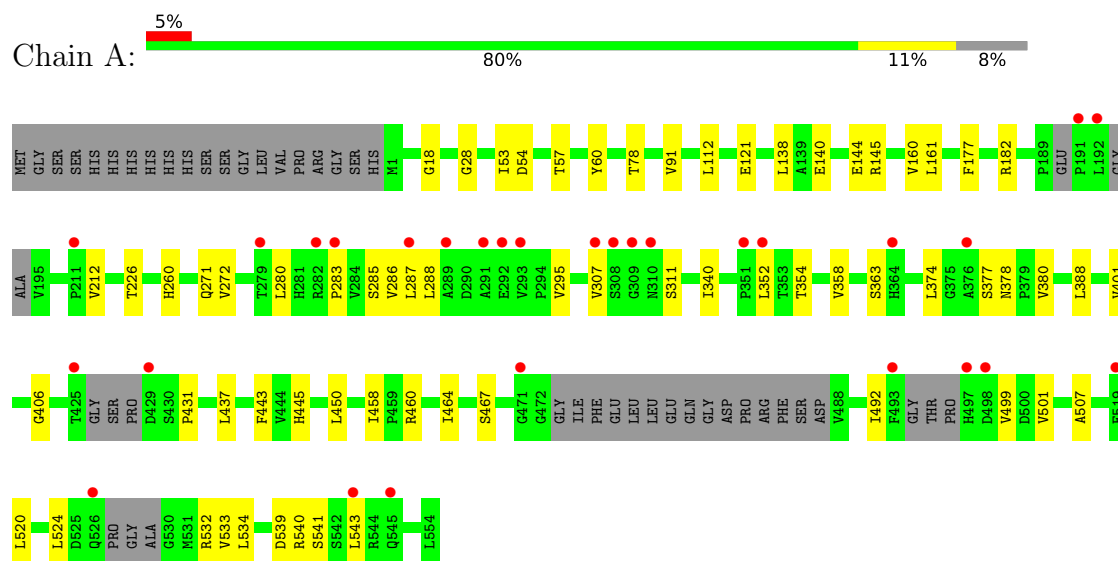
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	141	Total 141	O 141	0	0
8	D	105	Total 105	O 105	0	0
8	B	117	Total 117	O 117	0	0
8	C	108	Total 108	O 108	0	0

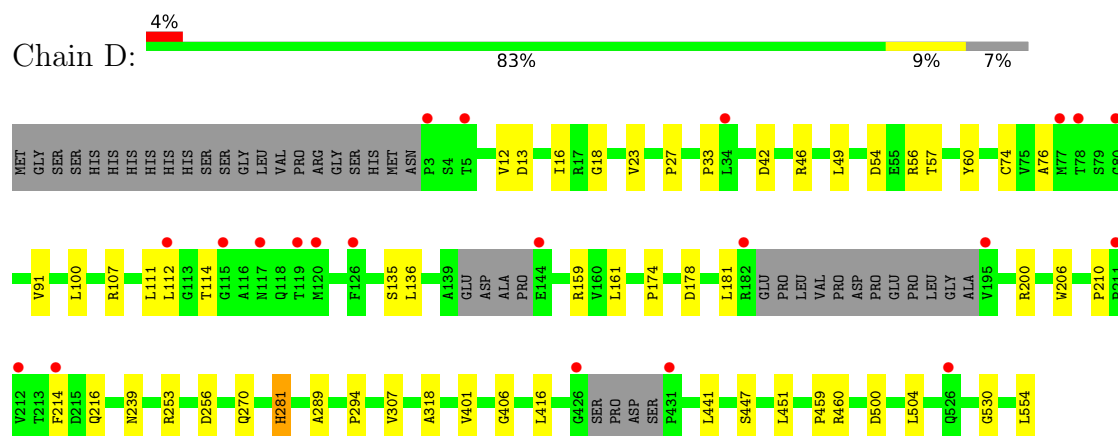
3 Residue-property plots [i](#)

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

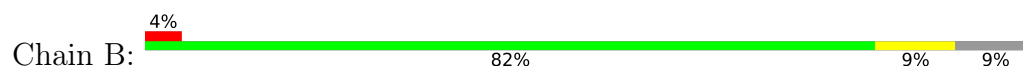
- Molecule 1: 2-succinyl-5-enolpyruvyl-6-hydroxy-3-cyclohexene-1-carboxylate synthase

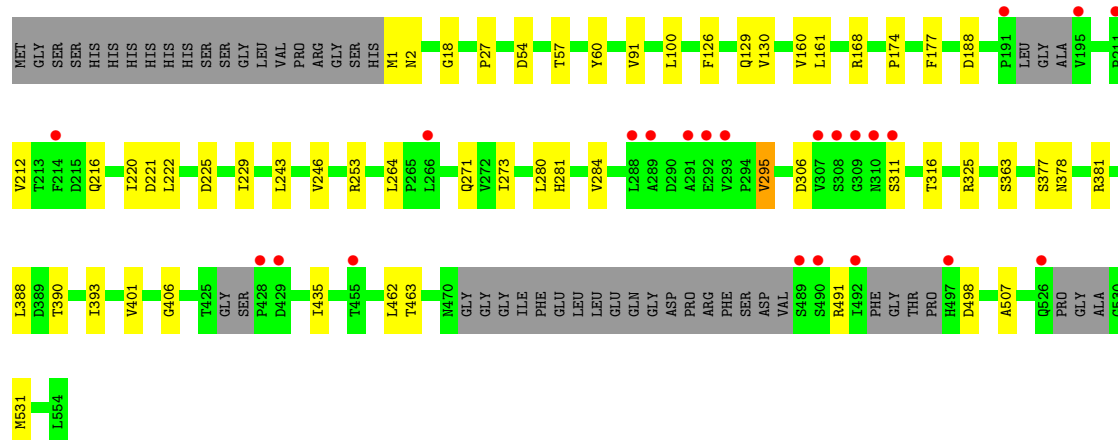


- Molecule 1: 2-succinyl-5-enolpyruvyl-6-hydroxy-3-cyclohexene-1-carboxylate synthase

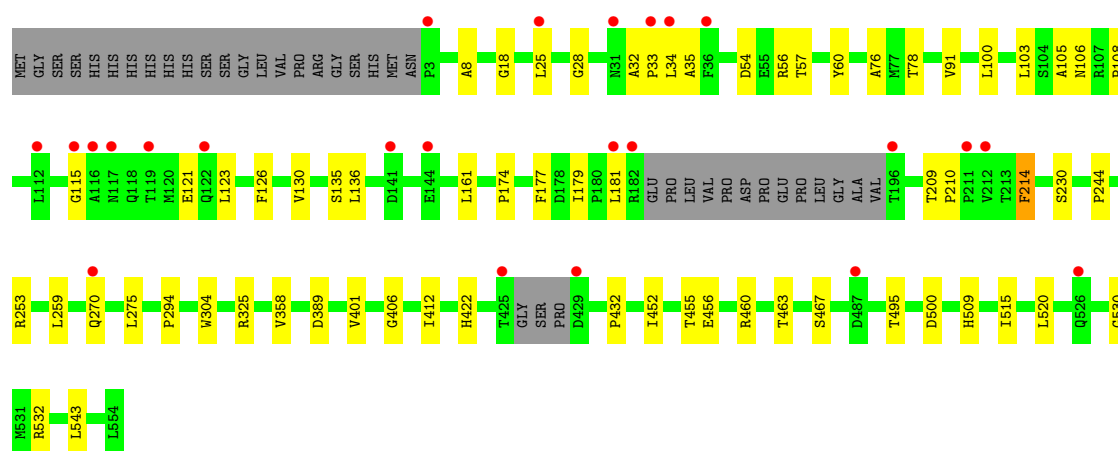
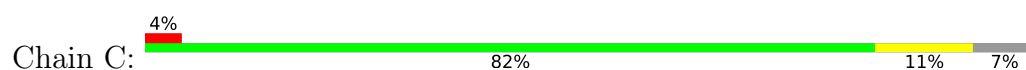


- Molecule 1: 2-succinyl-5-enolpyruvyl-6-hydroxy-3-cyclohexene-1-carboxylate synthase





- Molecule 1: 2-succinyl-5-enolpyruvyl-6-hydroxy-3-cyclohexene-1-carboxylate synthase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	101.55Å 143.67Å 176.11Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.79 – 2.35 48.79 – 2.35	Depositor EDS
% Data completeness (in resolution range)	99.9 (48.79-2.35) 99.9 (48.79-2.35)	Depositor EDS
R_{merge}	0.28	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.32 (at 2.34Å)	Xtrriage
Refinement program	PHENIX 1.14_3260, REFMAC	Depositor
R, R_{free}	0.205 , 0.251 0.208 , 0.252	Depositor DCC
R_{free} test set	5373 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	44.2	Xtrriage
Anisotropy	0.401	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 47.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	15922	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.98% 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: GOL, FMT, MG, DNA, TPP, ACT

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	0.26	0/3892	0.44	0/5337
1	B	0.23	0/3848	0.41	0/5283
1	C	0.24	0/3948	0.42	0/5418
1	D	0.22	0/3940	0.41	0/5400
All	All	0.24	0/15628	0.42	0/21438

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 [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	A	3814	0	3809	43	0
1	B	3773	0	3746	40	0
1	C	3862	0	3831	44	0
1	D	3855	0	3841	38	0
2	A	15	0	5	0	0
2	B	15	0	5	1	0
2	C	15	0	7	1	0
2	D	15	0	5	0	0
3	A	12	0	16	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	4	0	3	0	0
4	C	4	0	3	0	0
4	D	4	0	3	0	0
5	A	3	0	1	0	0
5	B	3	0	1	0	0
5	D	3	0	1	0	0
6	C	1	0	0	0	0
6	D	1	0	0	0	0
7	C	26	0	16	2	0
7	D	26	0	16	0	0
8	A	141	0	0	1	0
8	B	117	0	0	0	0
8	C	108	0	0	1	0
8	D	105	0	0	2	0
All	All	15922	0	15309	146	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 5.

All (146) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1:MET:HE1	1:B:188:ASP:H	1.33	0.94
1:A:307:VAL:HG11	1:A:311:SER:HB2	1.54	0.90
1:C:32:ALA:HB3	1:C:35:ALA:HB2	1.68	0.75
1:B:435:ILE:HD12	1:B:463:THR:HB	1.72	0.72
1:B:229:ILE:HD13	1:B:284:VAL:HG13	1.72	0.71
1:B:253:ARG:NH2	1:B:390:THR:H	1.90	0.70
1:B:253:ARG:HH21	1:B:390:THR:H	1.40	0.69
1:A:285:SER:HA	1:A:288:LEU:HD12	1.75	0.68
1:C:54:ASP:HB3	1:C:57:THR:HG22	1.76	0.67
1:C:33:PRO:HB3	1:C:105:ALA:HB2	1.78	0.65
1:B:1:MET:HE3	1:B:2:ASN:H	1.60	0.65
1:B:498:ASP:O	1:C:509:HIS:NE2	2.32	0.61
1:A:18:GLY:HA3	1:A:161:LEU:HD13	1.81	0.61
1:D:23:VAL:HG22	1:D:74:CYS:HB2	1.81	0.60
1:C:105:ALA:HB1	1:C:181:LEU:HD12	1.85	0.59
1:A:140:GLU:OE2	1:A:145:ARG:NH1	2.37	0.58
1:D:54:ASP:HB3	1:D:57:THR:HG22	1.85	0.58
1:B:1:MET:HE1	1:B:188:ASP:N	2.13	0.57
1:B:295:VAL:HG13	1:B:311:SER:HA	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:60:TYR:CG	1:A:406:GLY:HA3	2.40	0.57
1:C:8:ALA:HA	1:C:34:LEU:CD1	2.34	0.57
1:C:91:VAL:HG12	1:C:401:VAL:HG21	1.87	0.56
1:D:216:GLN:HE22	1:B:168:ARG:HG2	1.69	0.56
1:B:221:ASP:O	1:B:271:GLN:NE2	2.39	0.56
1:C:18:GLY:HA3	1:C:161:LEU:HD13	1.86	0.56
1:D:239:ASN:ND2	1:D:318:ALA:O	2.37	0.56
7:C:604:TPP:HN42	7:C:604:TPP:H2	1.71	0.55
1:B:507:ALA:HA	1:C:500:ASP:HB3	1.89	0.55
1:C:515:ILE:HD11	1:C:520:LEU:HD13	1.88	0.55
1:D:214[B]:PHE:HD2	1:B:212:VAL:HG21	1.72	0.55
1:D:18:GLY:HA3	1:D:161:LEU:HD13	1.89	0.54
1:B:246:VAL:HG21	1:B:264:LEU:HD22	1.89	0.54
1:D:107:ARG:HH12	1:D:181:LEU:HD11	1.72	0.54
1:A:53:ILE:HD12	1:D:441:LEU:HG	1.90	0.53
1:C:103:LEU:HD23	1:C:177:PHE:HB3	1.91	0.53
1:A:431:PRO:HD3	1:A:460:ARG:HH21	1.74	0.53
1:A:354:THR:HB	1:A:543:LEU:HD21	1.90	0.53
1:C:8:ALA:HA	1:C:34:LEU:HD12	1.89	0.53
1:B:280:LEU:O	1:B:381:ARG:NH1	2.42	0.52
1:C:34:LEU:HD21	1:C:179:ILE:HD13	1.92	0.51
1:A:260:HIS:CD2	1:A:340:ILE:HD13	2.46	0.51
1:B:220:ILE:HG21	1:B:273:ILE:HD11	1.92	0.51
1:B:225:ASP:OD2	1:B:325:ARG:NH2	2.44	0.51
1:C:25:LEU:HD23	1:C:76:ALA:HB3	1.91	0.51
1:D:281:HIS:NE2	8:D:702:HOH:O	2.35	0.51
1:B:216:GLN:HB3	1:B:316:THR:HG23	1.93	0.51
1:A:112:LEU:HD21	1:A:121:GLU:HG3	1.92	0.50
1:B:126:PHE:HA	1:C:123:LEU:HD23	1.93	0.50
1:C:304:TRP:C	2:C:601:DNA:H1H	2.36	0.50
1:C:455:THR:OG1	1:C:456:GLU:OE2	2.29	0.50
1:D:33:PRO:HD2	1:D:76:ALA:HB1	1.92	0.50
1:A:54:ASP:HB3	1:A:57:THR:HG22	1.93	0.50
1:A:507:ALA:HA	1:D:500:ASP:HB3	1.94	0.50
1:A:212:VAL:HG21	1:C:214[B]:PHE:CD2	2.47	0.49
7:C:604:TPP:HN42	7:C:604:TPP:C2	2.24	0.49
1:B:100:LEU:O	1:B:174:PRO:HA	2.11	0.49
1:A:144:GLU:H	1:A:144:GLU:CD	2.20	0.49
1:D:289:ALA:HB2	1:D:307:VAL:HG12	1.95	0.48
1:B:363:SER:HB3	1:B:388:LEU:HD13	1.94	0.48
1:B:377:SER:OG	1:B:378:ASN:N	2.31	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:463:THR:HG23	1:C:532:ARG:HG3	1.94	0.48
1:B:60:TYR:CG	1:B:406:GLY:HA3	2.48	0.48
1:A:260:HIS:HD2	1:A:340:ILE:HD13	1.79	0.47
1:D:60:TYR:CD2	1:D:406:GLY:HA3	2.49	0.47
1:A:352:LEU:HD13	1:A:539:ASP:HB3	1.95	0.47
1:D:253:ARG:NH2	8:D:706:HOH:O	2.46	0.47
1:B:129:GLN:NE2	1:C:121:GLU:HG3	2.29	0.47
1:D:91:VAL:HG12	1:D:401:VAL:HG21	1.95	0.47
1:D:100:LEU:O	1:D:174:PRO:HA	2.14	0.47
1:D:159:ARG:HH11	1:B:306:ASP:HB3	1.78	0.47
2:B:601:DNA:H1F	1:C:115:GLY:HA2	1.96	0.47
1:A:272:VAL:HG22	1:A:295:VAL:HG22	1.96	0.47
1:C:126:PHE:O	1:C:130:VAL:HG22	2.15	0.47
1:A:283:PRO:HA	1:A:286:VAL:HG22	1.97	0.46
1:C:100:LEU:O	1:C:174:PRO:HA	2.15	0.46
1:B:18:GLY:HA3	1:B:161:LEU:HD13	1.97	0.46
1:B:54:ASP:OD1	1:C:56:ARG:NH2	2.48	0.46
1:D:214[B]:PHE:CD2	1:B:212:VAL:HG21	2.51	0.46
1:A:499:VAL:HG21	1:D:451:LEU:HD12	1.98	0.46
1:A:91:VAL:HG12	1:A:401:VAL:HG11	1.98	0.45
1:A:520:LEU:O	1:A:524:LEU:HG	2.15	0.45
1:B:160:VAL:HG21	1:B:177:PHE:CD2	2.52	0.45
1:A:54:ASP:OD1	1:D:56:ARG:NH2	2.48	0.45
1:C:412:ILE:HD13	1:C:452:ILE:HD11	1.99	0.45
1:A:363:SER:HB3	1:A:388:LEU:HD13	1.99	0.45
1:A:374:LEU:HD23	1:A:437:LEU:HB3	1.97	0.45
1:D:12:VAL:O	1:D:16:ILE:HG12	2.17	0.45
1:D:135:SER:HB2	1:D:178:ASP:HB3	1.99	0.45
1:C:358:VAL:HG11	1:C:467:SER:HB2	1.98	0.45
1:A:443:PHE:CE2	1:A:501:VAL:HG13	2.52	0.45
1:B:91:VAL:HG12	1:B:401:VAL:HG11	1.98	0.45
1:A:272:VAL:HG11	1:A:287:LEU:HD21	1.98	0.45
1:C:460:ARG:O	1:C:530:GLY:HA2	2.18	0.44
1:C:28:GLY:HA3	1:C:78:THR:HB	1.99	0.44
1:D:210:PRO:HG2	1:B:216:GLN:HG2	1.98	0.44
1:D:111:LEU:HD13	1:D:114:THR:HG21	1.99	0.44
1:A:160:VAL:HG21	1:A:177:PHE:CG	2.53	0.44
1:C:325:ARG:HG2	8:C:712:HOH:O	2.16	0.44
1:D:112:LEU:HB3	1:C:135:SER:HB3	1.99	0.44
1:B:54:ASP:HB3	1:B:57:THR:HG22	2.00	0.44
1:C:60:TYR:CG	1:C:406:GLY:HA3	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:136:LEU:HD23	1:D:136:LEU:HA	1.87	0.43
1:A:492:ILE:HD11	1:D:27:PRO:HG3	1.99	0.43
1:C:60:TYR:CD2	1:C:406:GLY:HA3	2.53	0.43
1:C:136:LEU:HD12	1:C:136:LEU:HA	1.76	0.43
1:A:445:HIS:HE1	1:D:54:ASP:HA	1.84	0.43
1:A:532:ARG:NH2	1:A:534:LEU:HD11	2.34	0.43
1:D:200:ARG:HG3	1:D:206:TRP:HA	1.99	0.43
1:A:464:ILE:HB	1:A:533:VAL:HG22	2.00	0.43
1:D:554:LEU:HD23	1:D:554:LEU:HA	1.91	0.43
1:A:28:GLY:HA3	1:A:78:THR:HB	2.01	0.43
1:D:42:ASP:HB2	1:D:49:LEU:HG	2.01	0.43
1:D:60:TYR:CG	1:D:406:GLY:HA3	2.54	0.43
1:C:106:ASN:O	1:C:108:PRO:HD3	2.19	0.43
1:D:416:LEU:HG	1:D:459:PRO:HG3	2.01	0.43
1:A:283:PRO:O	1:A:286:VAL:HG22	2.19	0.42
1:A:358:VAL:HG11	1:A:467:SER:HB2	2.00	0.42
1:C:230:SER:HB3	1:C:275:LEU:HD12	2.01	0.42
1:A:226:THR:HA	1:A:271:GLN:O	2.19	0.42
1:A:458:ILE:HD11	8:A:705:HOH:O	2.20	0.42
1:D:256:ASP:OD1	1:D:256:ASP:N	2.53	0.42
1:C:543:LEU:HD23	1:C:543:LEU:HA	1.72	0.42
1:A:539:ASP:OD1	1:A:541:SER:OG	2.31	0.42
1:A:443:PHE:CZ	1:A:450:LEU:HD11	2.55	0.42
1:B:130:VAL:HA	1:B:174:PRO:HG2	2.02	0.42
1:C:270[A]:GLN:O	1:C:294:PRO:HD2	2.20	0.42
1:A:377:SER:OG	1:A:378:ASN:N	2.48	0.41
1:D:460:ARG:O	1:D:530:GLY:HA2	2.19	0.41
1:A:539:ASP:CG	1:A:541:SER:H	2.28	0.41
1:B:222:LEU:HD22	1:B:243:LEU:HD11	2.02	0.41
1:B:27:PRO:HB3	1:C:495:THR:HG21	2.02	0.41
1:C:244:PRO:HB2	1:C:259:LEU:HD22	2.02	0.41
1:B:281:HIS:HB2	1:B:284:VAL:HG23	2.03	0.41
1:B:491:ARG:NH1	1:C:456:GLU:OE1	2.43	0.41
1:C:209:THR:HA	1:C:210:PRO:HD3	1.92	0.41
1:C:422:HIS:CG	1:C:432:PRO:HD3	2.56	0.41
1:D:270:GLN:O	1:D:294:PRO:HD2	2.21	0.41
1:D:447:SER:HB2	1:D:504:LEU:HD21	2.03	0.40
1:A:280:LEU:HD11	1:A:380:VAL:HG13	2.04	0.40
1:B:462:LEU:O	1:B:531:MET:HA	2.21	0.40
1:A:540:ARG:HA	1:A:543:LEU:CD1	2.51	0.40
1:D:13:ASP:OD1	1:D:46:ARG:NH1	2.40	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:129:GLN:HE22	1:C:121:GLU:HG3	1.85	0.40
1:B:390:THR:HA	1:B:393:ILE:HD11	2.04	0.40
1:A:138:LEU:HD11	1:A:182:ARG:HB2	2.03	0.40
1:C:253:ARG:HH12	1:C:389:ASP:HA	1.86	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	514/574 (90%)	503 (98%)	11 (2%)	0	100	100
1	B	512/574 (89%)	499 (98%)	13 (2%)	0	100	100
1	C	533/574 (93%)	520 (98%)	13 (2%)	0	100	100
1	D	526/574 (92%)	509 (97%)	16 (3%)	1 (0%)	43	52
All	All	2085/2296 (91%)	2031 (97%)	53 (2%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	281	HIS

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	393/445 (88%)	393 (100%)	0	100	100
1	B	387/445 (87%)	386 (100%)	1 (0%)	86	92
1	C	393/445 (88%)	391 (100%)	2 (0%)	81	89
1	D	397/445 (89%)	397 (100%)	0	100	100
All	All	1570/1780 (88%)	1567 (100%)	3 (0%)	88	94

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	295	VAL
1	C	214[A]	PHE
1	C	214[B]	PHE

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

Mol	Chain	Res	Type
1	A	39	GLN
1	A	117	ASN
1	A	118	GLN
1	A	281	HIS
1	A	346	GLN
1	A	547	HIS
1	D	106	ASN
1	B	39	GLN
1	B	117	ASN
1	B	310	ASN
1	B	547	HIS
1	C	171	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 16 ligands modelled in this entry, 2 are monoatomic - leaving 14 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
7	TPP	D	604	6	26,27,27	1.72	3 (11%)	38,40,40	1.58	8 (21%)
2	DNA	D	601	-	16,16,16	2.26	4 (25%)	23,23,23	1.00	0
4	ACT	C	603	-	3,3,3	0.87	0	3,3,3	0.74	0
3	GOL	A	603	-	5,5,5	0.24	0	5,5,5	0.47	0
5	FMT	B	602	-	2,2,2	0.76	0	1,1,1	0.49	0
3	GOL	A	602	-	5,5,5	0.39	0	5,5,5	0.47	0
5	FMT	A	605	-	2,2,2	0.59	0	1,1,1	0.70	0
5	FMT	D	605	-	2,2,2	0.70	0	1,1,1	0.60	0
4	ACT	A	604	-	3,3,3	0.81	0	3,3,3	0.77	0
2	DNA	C	601	-	16,16,16	2.25	4 (25%)	23,23,23	1.10	1 (4%)
4	ACT	D	603	-	3,3,3	0.81	0	3,3,3	0.71	0
2	DNA	B	601	-	16,16,16	2.17	4 (25%)	23,23,23	1.15	2 (8%)
2	DNA	A	601	-	16,16,16	2.19	5 (31%)	23,23,23	1.36	3 (13%)
7	TPP	C	604	6	26,27,27	1.65	3 (11%)	38,40,40	1.60	10 (26%)

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
7	TPP	D	604	6	-	3/17/17/17	0/2/2/2
2	DNA	D	601	-	-	0/4/4/4	0/2/2/2
3	GOL	A	603	-	-	4/4/4/4	-
3	GOL	A	602	-	-	2/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	DNA	C	601	-	-	0/4/4/4	0/2/2/2
2	DNA	B	601	-	-	2/4/4/4	0/2/2/2
2	DNA	A	601	-	-	0/4/4/4	0/2/2/2
7	TPP	C	604	6	-	2/17/17/17	0/2/2/2

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	601	DNA	C1L-C1M	6.71	1.49	1.39
2	C	601	DNA	C1L-C1M	6.59	1.48	1.39
2	B	601	DNA	C1L-C1M	6.32	1.48	1.39
2	A	601	DNA	C1L-C1M	6.22	1.48	1.39
7	C	604	TPP	C5-C4	6.04	1.47	1.35
7	D	604	TPP	C5-C4	6.04	1.47	1.35
7	D	604	TPP	PA-O3A	3.42	1.63	1.59
7	C	604	TPP	C5'-C4'	3.32	1.48	1.42
7	D	604	TPP	C5'-C4'	3.28	1.48	1.42
2	C	601	DNA	C1K-C1N	3.02	1.48	1.42
2	A	601	DNA	C1K-C1N	2.88	1.48	1.42
2	D	601	DNA	C1K-C1N	2.83	1.48	1.42
7	C	604	TPP	PA-O3A	2.73	1.62	1.59
2	B	601	DNA	C1O-C1N	2.72	1.48	1.43
2	D	601	DNA	C1O-C1N	2.71	1.48	1.43
2	A	601	DNA	C1O-C1N	2.68	1.48	1.43
2	C	601	DNA	C1O-C1N	2.66	1.48	1.43
2	A	601	DNA	C1M-C1O	2.50	1.48	1.43
2	B	601	DNA	C1K-C1N	2.48	1.47	1.42
2	C	601	DNA	C1M-C1O	2.46	1.48	1.43
2	B	601	DNA	C1M-C1O	2.45	1.48	1.43
2	D	601	DNA	C1M-C1O	2.35	1.48	1.43
2	A	601	DNA	C1H-C1O	-2.23	1.37	1.42

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	D	604	TPP	C2-S1-C5	-3.79	88.71	91.22
7	C	604	TPP	C6'-N1'-C2'	3.74	122.21	116.07
7	D	604	TPP	C6'-N1'-C2'	3.62	122.02	116.07
7	D	604	TPP	CM2-C2'-N1'	3.50	120.93	117.20
7	C	604	TPP	C2-S1-C5	-3.47	88.91	91.22
7	C	604	TPP	C5'-C6'-N1'	-3.29	118.47	123.83
7	C	604	TPP	CM4-C4-N3	2.55	126.45	120.57

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	D	604	TPP	S1-C2-N3	2.54	115.52	112.30
7	D	604	TPP	C5'-C6'-N1'	-2.52	119.72	123.83
7	C	604	TPP	CM2-C2'-N1'	2.52	119.88	117.20
7	C	604	TPP	C6'-C5'-C4'	2.51	118.63	115.55
7	D	604	TPP	N1'-C2'-N3'	-2.49	121.38	125.53
2	A	601	DNA	C1F-C1H-C1O	-2.41	117.66	120.91
2	A	601	DNA	C1E-C1G-C1N	-2.41	117.66	120.91
7	D	604	TPP	CM4-C4-N3	2.39	126.09	120.57
2	A	601	DNA	C1I-C1L-C1M	2.39	120.95	119.05
2	C	601	DNA	C1I-C1L-C1M	2.29	120.87	119.05
2	B	601	DNA	C1I-C1L-C1M	2.28	120.86	119.05
7	C	604	TPP	N4'-C4'-N3'	2.28	120.10	117.03
7	D	604	TPP	CM4-C4-C5	-2.18	122.15	127.75
7	C	604	TPP	C6-C5-S1	-2.17	115.83	120.90
7	C	604	TPP	S1-C2-N3	2.15	115.03	112.30
2	B	601	DNA	C1F-C1H-C1O	-2.07	118.12	120.91
7	C	604	TPP	CM4-C4-C5	-2.01	122.61	127.75

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	602	GOL	C1-C2-C3-O3
3	A	603	GOL	O1-C1-C2-C3
3	A	603	GOL	C1-C2-C3-O3
3	A	603	GOL	O2-C2-C3-O3
7	D	604	TPP	PA-O3A-PB-O2B
7	C	604	TPP	PA-O3A-PB-O2B
3	A	602	GOL	O2-C2-C3-O3
3	A	603	GOL	O1-C1-C2-O2
7	D	604	TPP	PA-O3A-PB-O1B
2	B	601	DNA	O1B-C1J-C1L-C1M
7	D	604	TPP	C5-C6-C7-O7
7	C	604	TPP	C5-C6-C7-O7
2	B	601	DNA	O1A-C1J-C1L-C1M

There are no ring outliers.

3 monomers are involved in 4 short contacts:

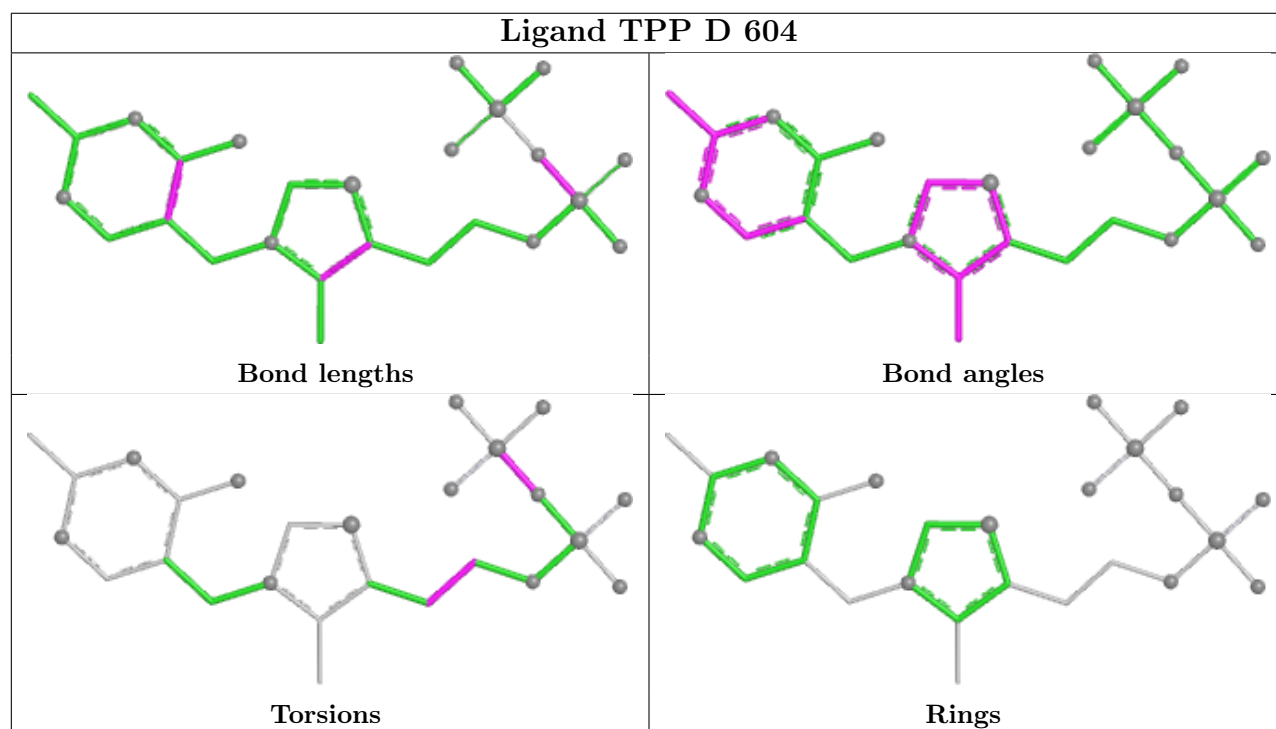
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	601	DNA	1	0

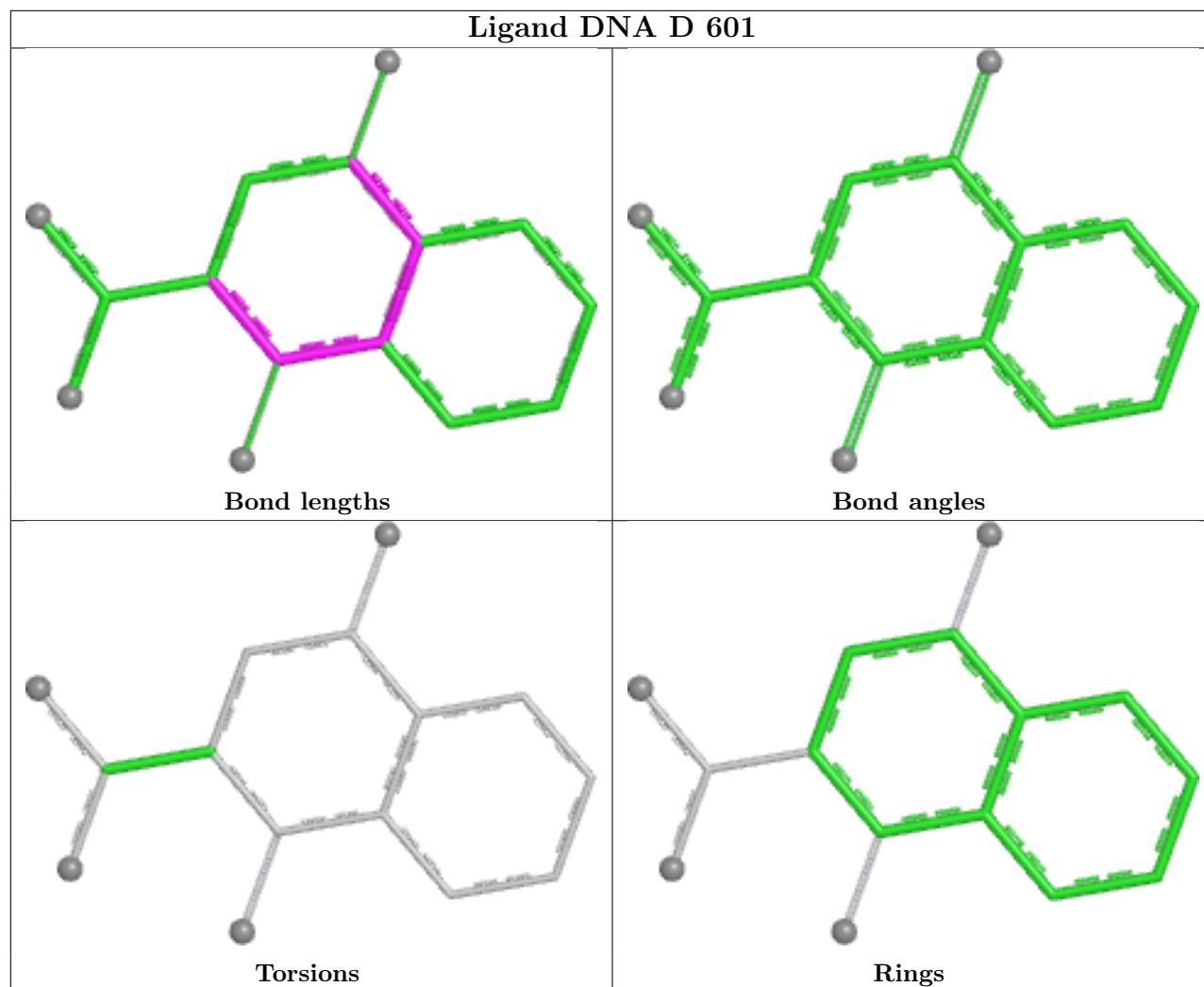
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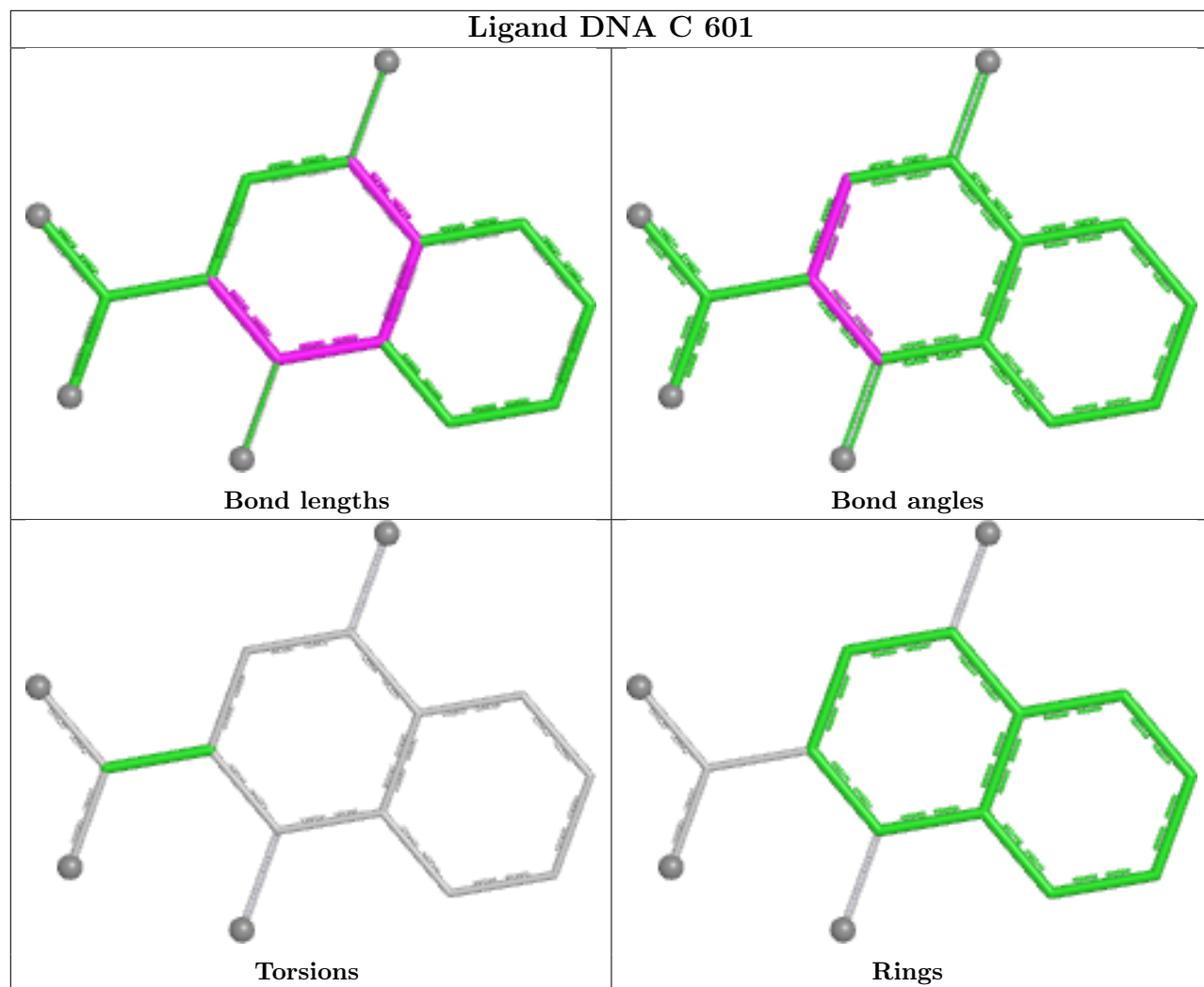
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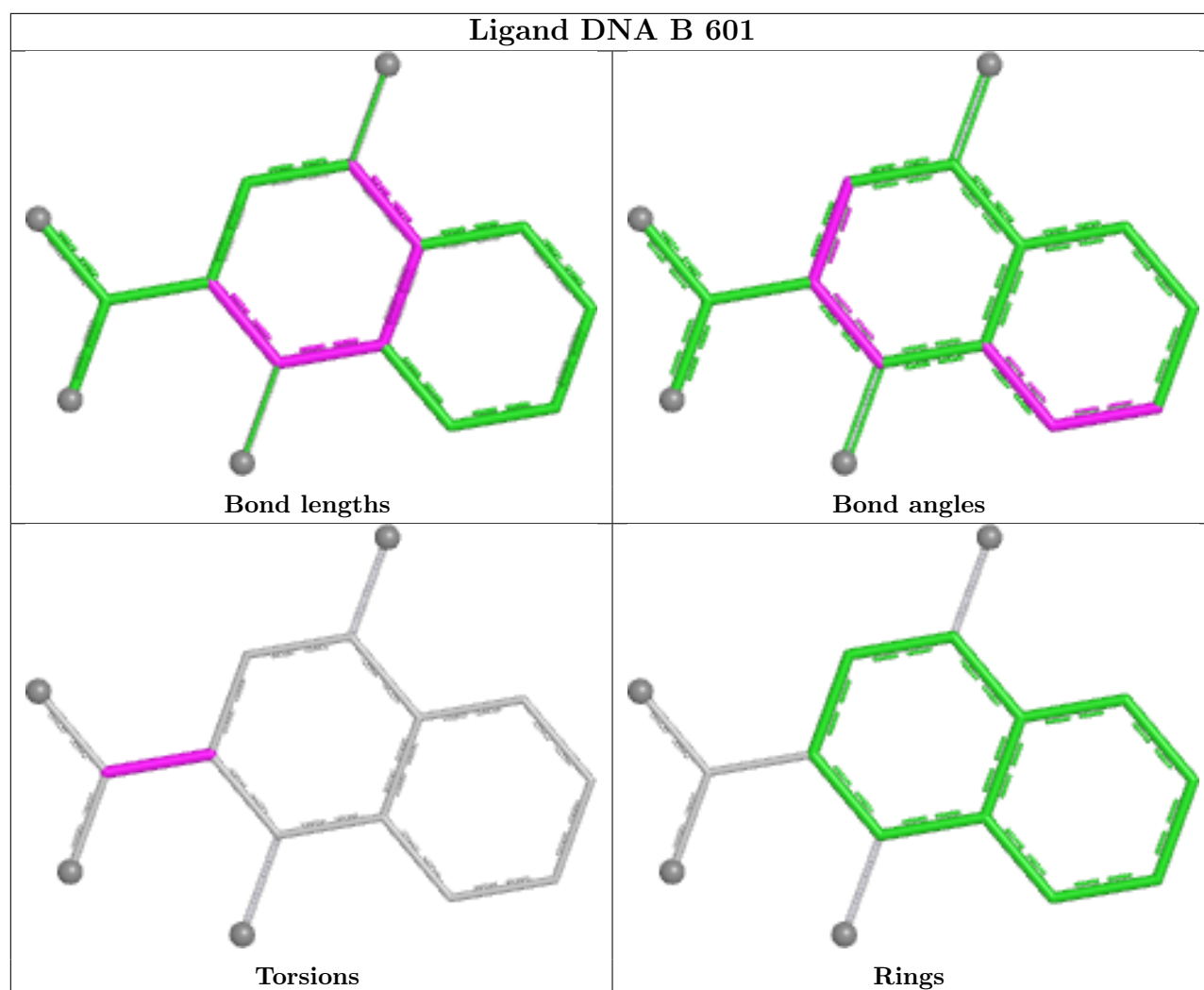
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	601	DNA	1	0
7	C	604	TPP	2	0

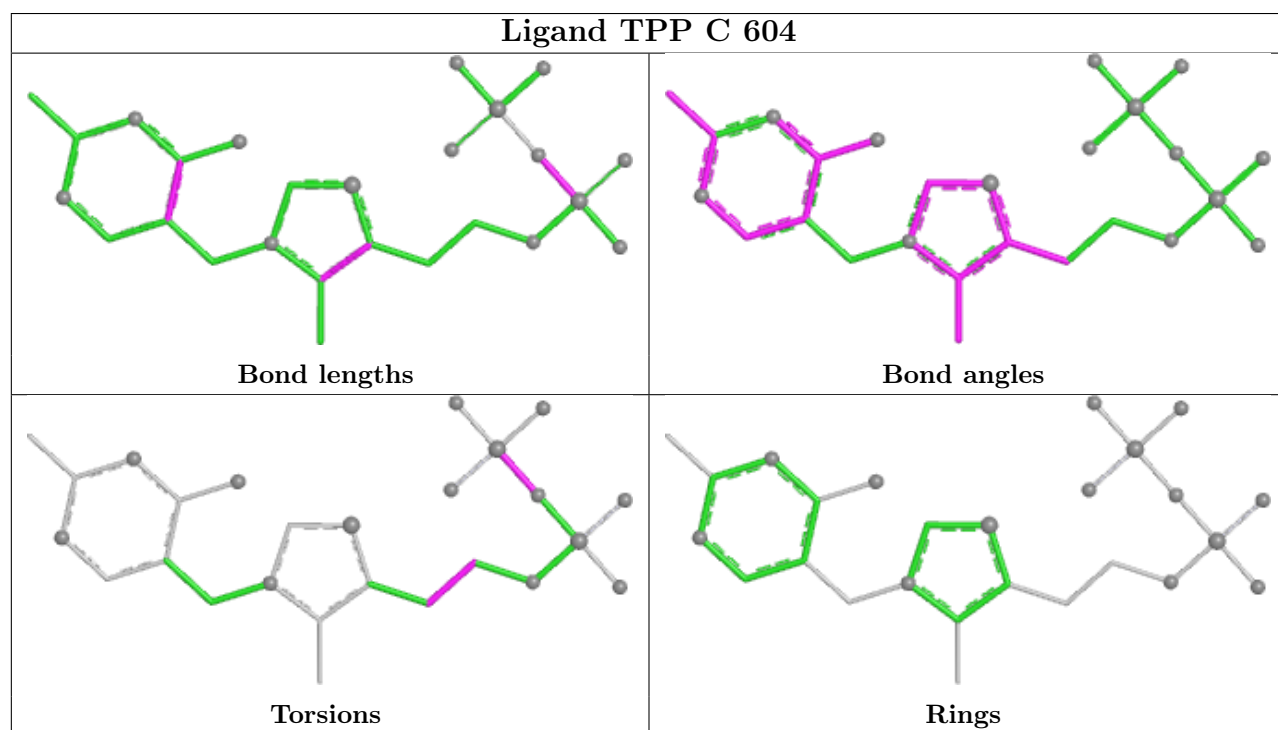
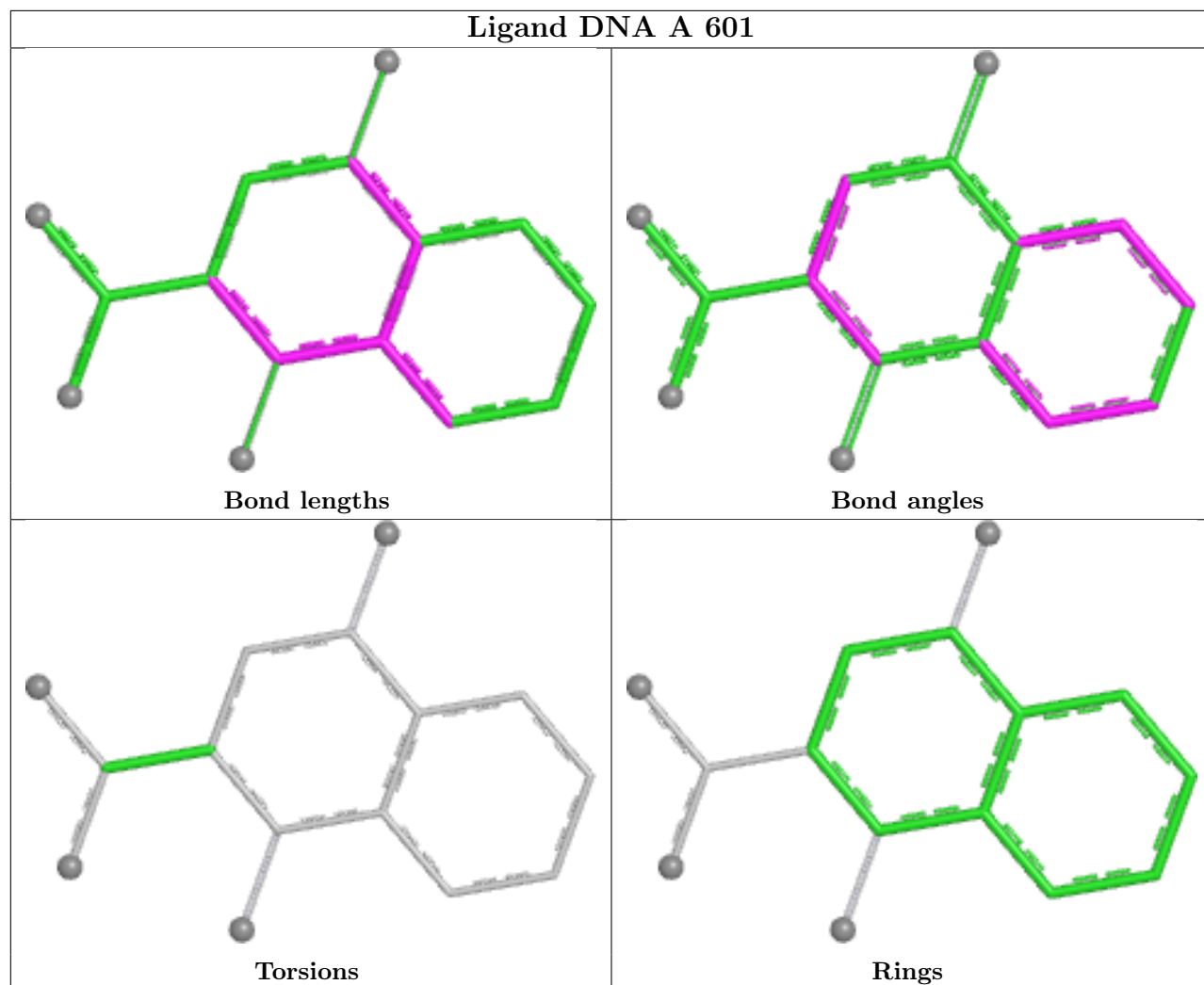
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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

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(Å ²)	Q<0.9
1	A	527/574 (91%)	0.37	29 (5%)	30 36	30, 55, 89, 109	1 (0%)
1	B	524/574 (91%)	0.24	23 (4%)	39 45	32, 53, 84, 105	0
1	C	536/574 (93%)	0.35	24 (4%)	38 44	29, 56, 88, 115	3 (0%)
1	D	532/574 (92%)	0.32	21 (3%)	43 49	31, 57, 83, 106	2 (0%)
All	All	2119/2296 (92%)	0.32	97 (4%)	37 43	29, 55, 86, 115	6 (0%)

All (97) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	426	GLY	5.8
1	B	490	SER	5.1
1	C	211	PRO	4.9
1	D	144	GLU	4.9
1	C	429	ASP	4.7
1	D	211	PRO	4.4
1	C	3	PRO	4.3
1	B	211	PRO	4.2
1	C	31	ASN	4.2
1	A	307	VAL	4.1
1	B	526	GLN	4.1
1	D	119	THR	4.0
1	B	428	PRO	3.9
1	B	429	ASP	3.9
1	A	211	PRO	3.9
1	D	195	VAL	3.9
1	C	115	GLY	3.9
1	B	191	PRO	3.8
1	D	214[A]	PHE	3.7
1	C	182	ARG	3.7
1	B	310	ASN	3.6

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Mol	Chain	Res	Type	RSRZ
1	A	291	ALA	3.6
1	C	33	PRO	3.6
1	D	182	ARG	3.3
1	C	117	ASN	3.3
1	A	376	ALA	3.3
1	C	487	ASP	3.2
1	A	526	GLN	3.2
1	D	77	MET	3.2
1	C	141	ASP	3.2
1	B	492	ILE	3.2
1	A	425	THR	3.2
1	C	181	LEU	3.1
1	B	307	VAL	3.1
1	C	119	THR	3.0
1	A	191	PRO	3.0
1	A	543	LEU	2.9
1	D	431	PRO	2.9
1	B	195	VAL	2.9
1	D	120	MET	2.9
1	A	293	VAL	2.9
1	C	34	LEU	2.8
1	B	308	SER	2.8
1	A	283	PRO	2.7
1	A	310	ASN	2.7
1	D	117	ASN	2.7
1	C	112	LEU	2.7
1	A	292	GLU	2.7
1	B	489	SER	2.7
1	A	309	GLY	2.7
1	C	116	ALA	2.6
1	A	282	ARG	2.6
1	B	292	GLU	2.6
1	D	115	GLY	2.6
1	D	112	LEU	2.6
1	D	34	LEU	2.5
1	A	289	ALA	2.5
1	C	196	THR	2.5
1	A	429	ASP	2.5
1	C	36	PHE	2.5
1	B	309	GLY	2.5
1	A	287	LEU	2.4
1	A	497	HIS	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	455	THR	2.4
1	A	471	GLY	2.4
1	A	351	PRO	2.4
1	D	78	THR	2.4
1	B	293	VAL	2.3
1	A	279	THR	2.3
1	C	212	VAL	2.3
1	B	289	ALA	2.3
1	A	364	HIS	2.3
1	C	122	GLN	2.3
1	C	270[A]	GLN	2.3
1	C	425	THR	2.2
1	B	266	LEU	2.2
1	D	126	PHE	2.2
1	D	212	VAL	2.2
1	A	498	ASP	2.1
1	C	144	GLU	2.1
1	D	526	GLN	2.1
1	A	308	SER	2.1
1	A	192	LEU	2.1
1	A	352	LEU	2.1
1	A	545	GLN	2.1
1	B	497	HIS	2.1
1	D	5	THR	2.1
1	B	288	LEU	2.1
1	C	526	GLN	2.1
1	C	25	LEU	2.1
1	D	3	PRO	2.1
1	A	493	PHE	2.1
1	B	214	PHE	2.1
1	A	519	GLU	2.0
1	B	291	ALA	2.0
1	D	80	GLY	2.0
1	B	311	SER	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 oligosaccharides in this entry.

6.4 Ligands ⓘ

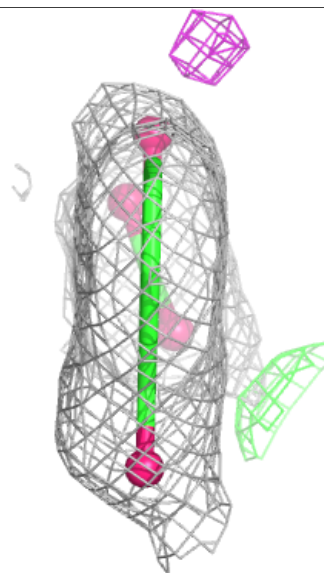
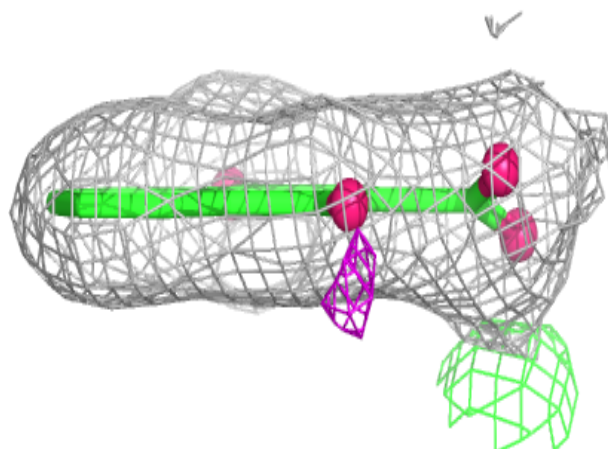
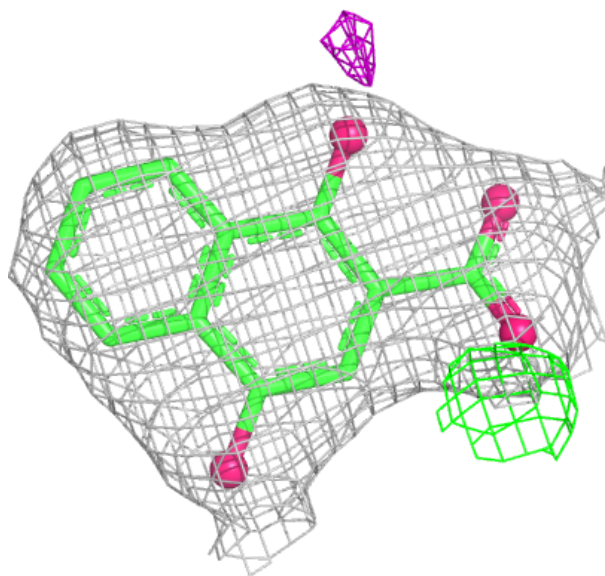
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	ACT	C	603	4/4	0.78	0.17	49,50,52,54	0
5	FMT	D	605	3/3	0.81	0.17	68,68,71,73	0
5	FMT	A	605	3/3	0.84	0.16	60,60,62,64	0
4	ACT	A	604	4/4	0.84	0.18	61,61,61,62	0
4	ACT	D	603	4/4	0.85	0.15	47,50,50,52	0
3	GOL	A	603	6/6	0.88	0.17	53,55,57,60	0
3	GOL	A	602	6/6	0.88	0.18	53,55,61,63	0
5	FMT	B	602	3/3	0.91	0.12	58,58,58,58	0
2	DNA	B	601	15/15	0.95	0.07	38,45,47,49	0
2	DNA	A	601	15/15	0.96	0.07	39,43,45,46	0
2	DNA	C	601	15/15	0.96	0.06	32,38,43,43	0
2	DNA	D	601	15/15	0.97	0.06	34,38,41,42	0
7	TPP	D	604	26/26	0.97	0.07	28,40,47,47	0
7	TPP	C	604	26/26	0.97	0.07	43,47,50,50	0
6	MG	D	602	1/1	0.99	0.02	37,37,37,37	0
6	MG	C	602	1/1	0.99	0.03	41,41,41,41	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.

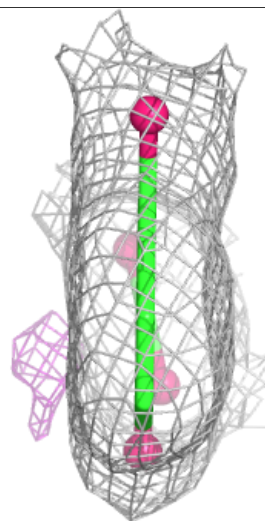
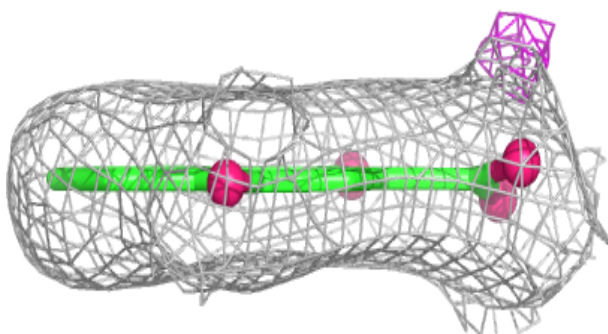
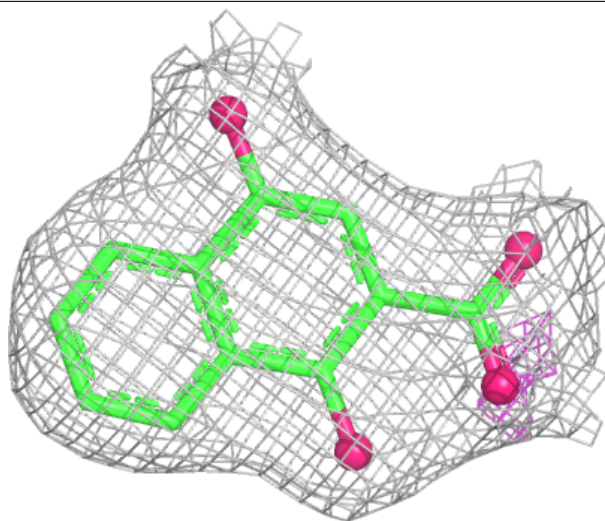
Electron density around DNA B 601:

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



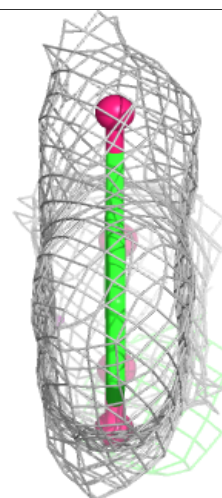
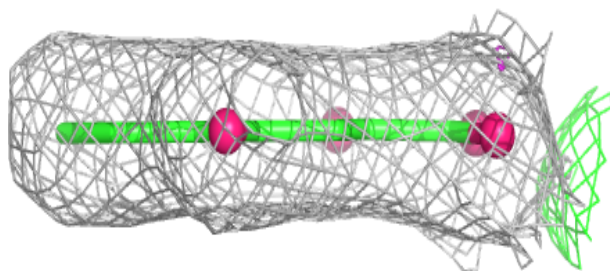
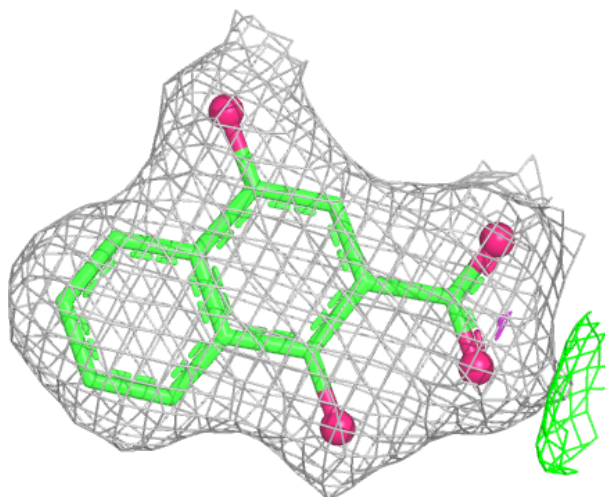
Electron density around DNA A 601:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



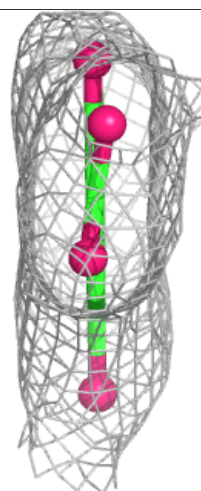
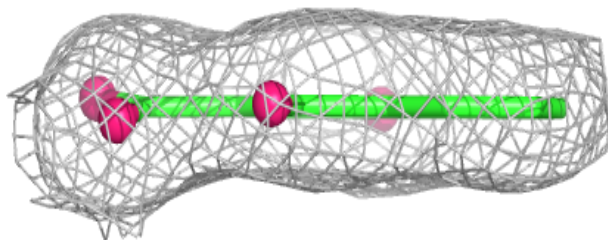
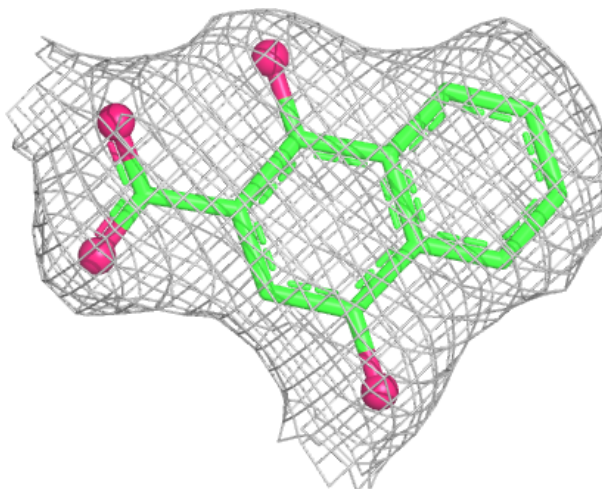
Electron density around DNA C 601:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



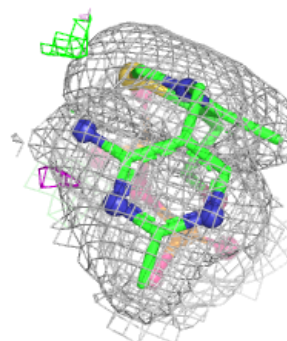
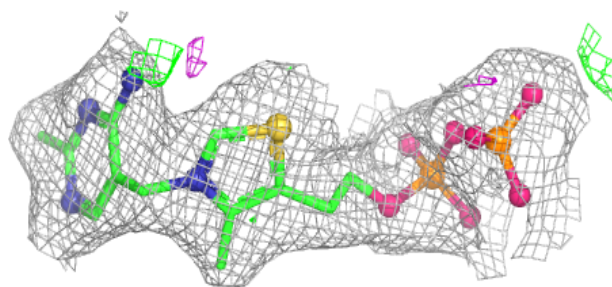
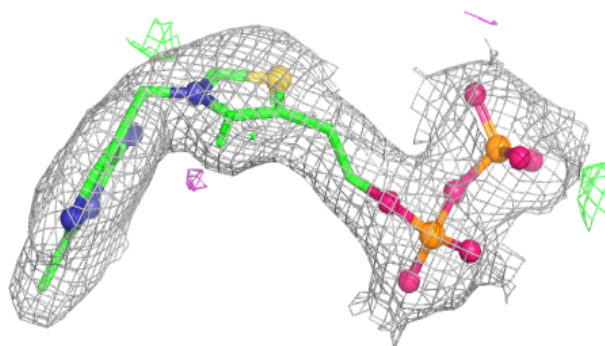
Electron density around DNA D 601:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

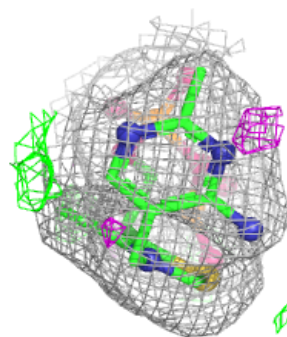
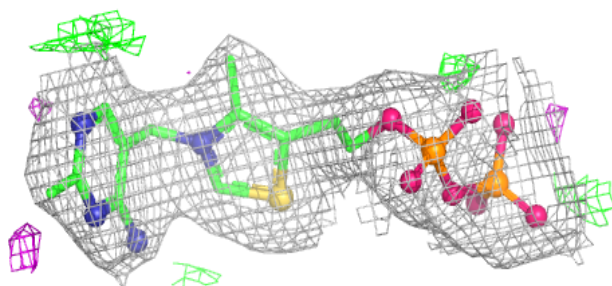
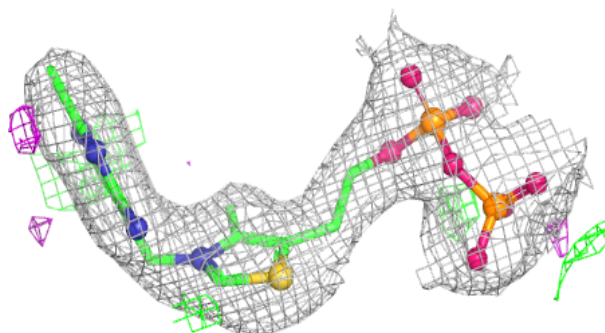


Electron density around TPP D 604:

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

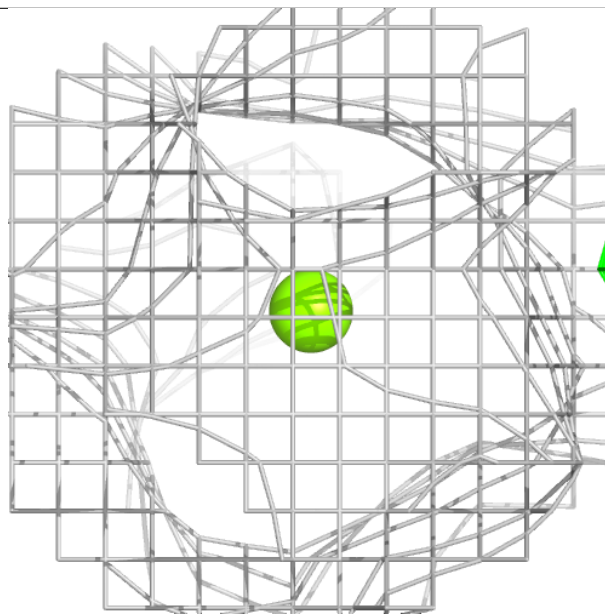
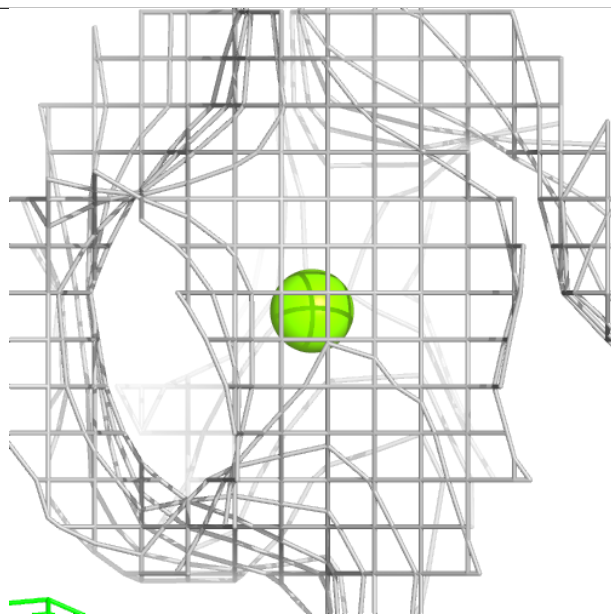
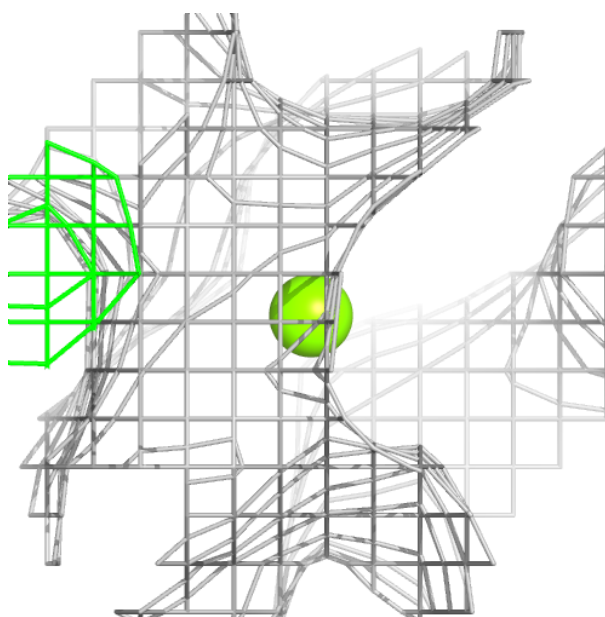
**Electron density around TPP C 604:**

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



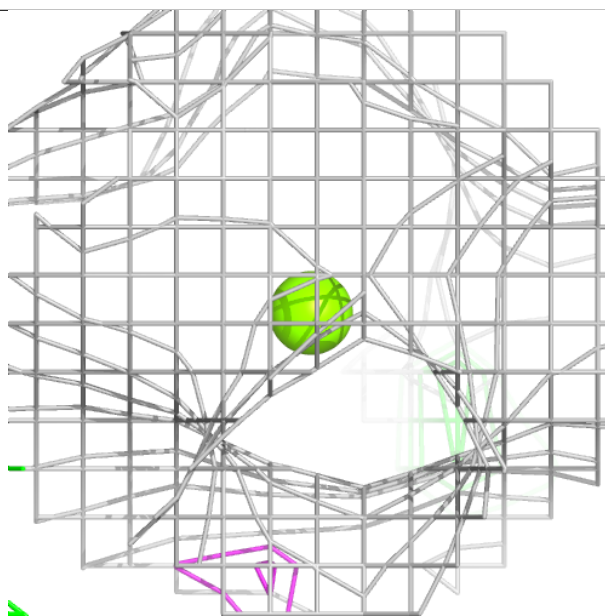
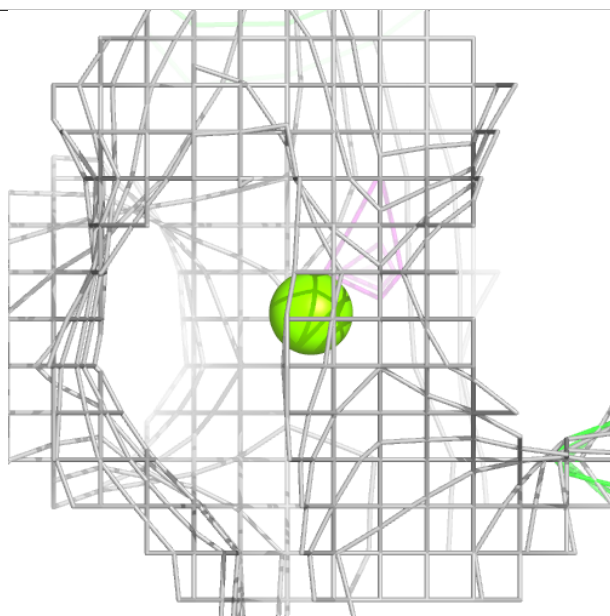
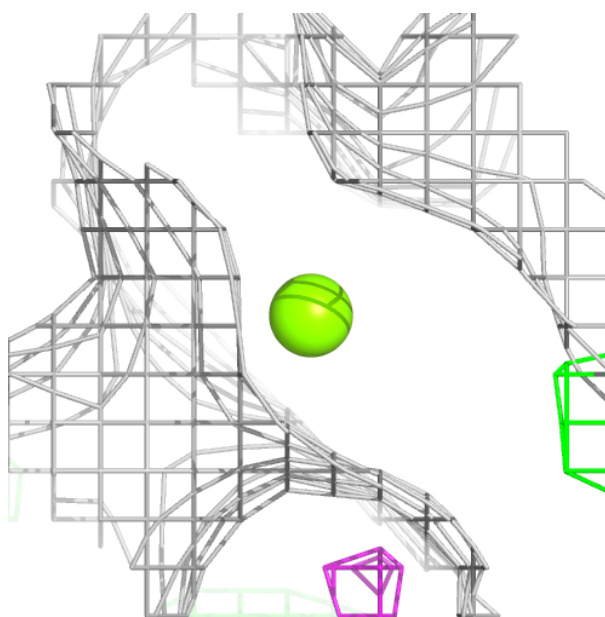
Electron density around MG D 602:

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



Electron density around MG C 602:

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



6.5 Other polymers ⓘ

There are no such residues in this entry.