



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 08:20 AM UTC

PDB ID : 8VZJ / pdb_00008vzj
Title : Crystal Structure of 2-Hydroxyacyl-CoA Lyase/Synthase DhcHACS from Dehalococcoidia bacterium in the Complex with THDP and ADP
Authors : Kim, Y.; Gade, P.; Endres, M.; Lee, S.; Yoshikuni, Y.; Gonzalez, R.; Michalska, K.; Joachimiak, A.
Deposited on : 2024-02-11
Resolution : 2.70 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

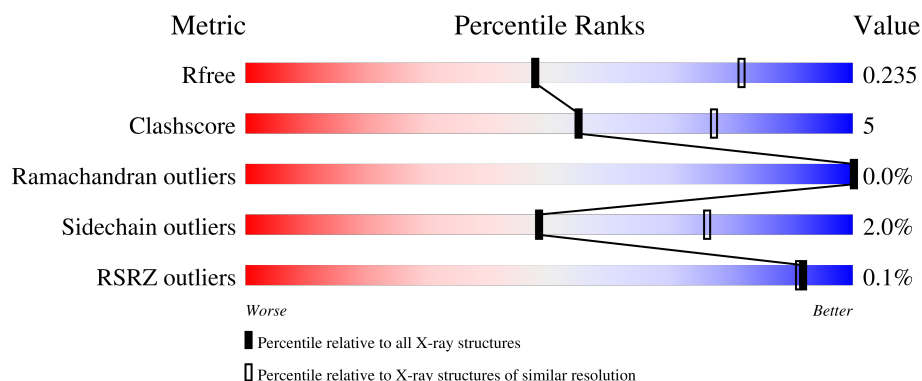
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	554	
1	B	554	
1	C	554	
1	D	554	

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Mol	Chain	Length	Quality of chain
1	E	554	 85% 12% ..
1	F	554	 84% 13% ..
1	G	554	 86% 10% ..
1	H	554	 85% 11% ..

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 33287 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 Oxalyl-CoA decarboxylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	542	Total	C	N	O	S	0	0	0
			4111	2604	712	765	30			
1	B	539	Total	C	N	O	S	0	0	0
			4084	2587	705	762	30			
1	C	542	Total	C	N	O	S	0	0	0
			4102	2597	708	766	31			
1	D	546	Total	C	N	O	S	0	1	0
			4149	2626	719	772	32			
1	E	541	Total	C	N	O	S	0	0	0
			4104	2599	711	764	30			
1	F	542	Total	C	N	O	S	0	0	0
			4102	2597	708	766	31			
1	G	540	Total	C	N	O	S	0	0	0
			4090	2591	706	762	31			
1	H	538	Total	C	N	O	S	0	0	0
			4080	2585	704	761	30			

There are 24 discrepancies between the modelled and reference sequences:

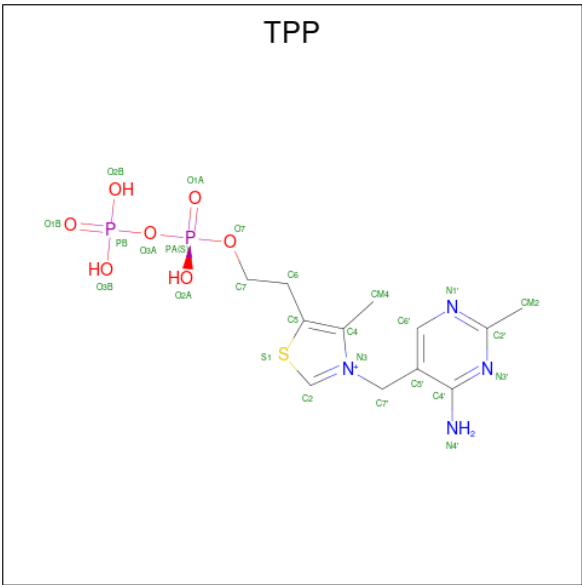
Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	UNP A0A315XEK8
A	-1	ASN	-	expression tag	UNP A0A315XEK8
A	0	ALA	-	expression tag	UNP A0A315XEK8
B	-2	SER	-	expression tag	UNP A0A315XEK8
B	-1	ASN	-	expression tag	UNP A0A315XEK8
B	0	ALA	-	expression tag	UNP A0A315XEK8
C	-2	SER	-	expression tag	UNP A0A315XEK8
C	-1	ASN	-	expression tag	UNP A0A315XEK8
C	0	ALA	-	expression tag	UNP A0A315XEK8
D	-2	SER	-	expression tag	UNP A0A315XEK8
D	-1	ASN	-	expression tag	UNP A0A315XEK8
D	0	ALA	-	expression tag	UNP A0A315XEK8
E	-2	SER	-	expression tag	UNP A0A315XEK8

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Chain	Residue	Modelled	Actual	Comment	Reference
E	-1	ASN	-	expression tag	UNP A0A315XEK8
E	0	ALA	-	expression tag	UNP A0A315XEK8
F	-2	SER	-	expression tag	UNP A0A315XEK8
F	-1	ASN	-	expression tag	UNP A0A315XEK8
F	0	ALA	-	expression tag	UNP A0A315XEK8
G	-2	SER	-	expression tag	UNP A0A315XEK8
G	-1	ASN	-	expression tag	UNP A0A315XEK8
G	0	ALA	-	expression tag	UNP A0A315XEK8
H	-2	SER	-	expression tag	UNP A0A315XEK8
H	-1	ASN	-	expression tag	UNP A0A315XEK8
H	0	ALA	-	expression tag	UNP A0A315XEK8

- Molecule 2 is THIAMINE DIPHOSPHATE (CCD ID: TPP) (formula: C₁₂H₁₉N₄O₇P₂S) (labeled as "Ligand of Interest" by depositor).



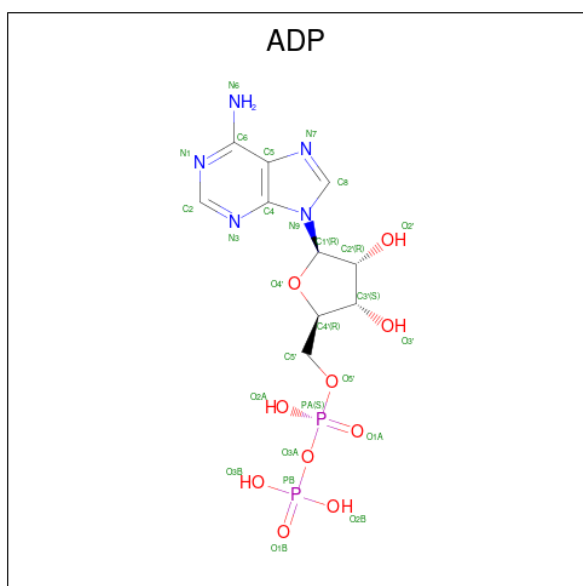
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total 26	C 12	N 4	O 7	P 2	S 1	0	0
2	B	1	Total 26	C 12	N 4	O 7	P 2	S 1	0	0
2	C	1	Total 26	C 12	N 4	O 7	P 2	S 1	0	0
2	D	1	Total 26	C 12	N 4	O 7	P 2	S 1	0	0
2	E	1	Total 26	C 12	N 4	O 7	P 2	S 1	0	0

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

- Molecule 3 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	B	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	C	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	D	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	E	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	F	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	G	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	H	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 5 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Mg	0	0
			1	1		
5	B	1	Total	Mg	0	0
			1	1		
5	C	1	Total	Mg	0	0
			1	1		
5	D	1	Total	Mg	0	0
			1	1		
5	E	1	Total	Mg	0	0
			1	1		
5	F	1	Total	Mg	0	0
			1	1		
5	G	1	Total	Mg	0	0
			1	1		
5	H	1	Total	Mg	0	0
			1	1		

- Molecule 6 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	C	1	Total	C	O	0	0
			4	2	2		
6	F	1	Total	C	O	0	0
			4	2	2		
6	G	1	Total	C	O	0	0
			4	2	2		
6	H	1	Total	C	O	0	0
			4	2	2		

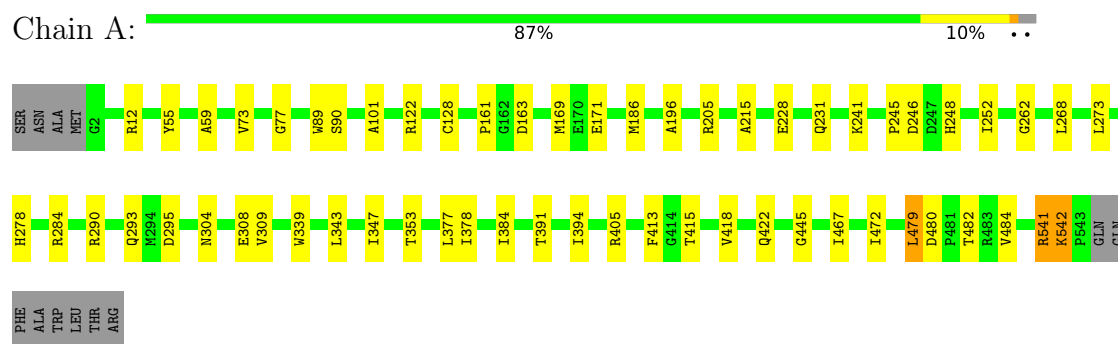
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	C	1	Total	O	0	0
			1	1		
7	E	1	Total	O	0	0
			1	1		
7	F	2	Total	O	0	0
			2	2		
7	G	1	Total	O	0	0
			1	1		

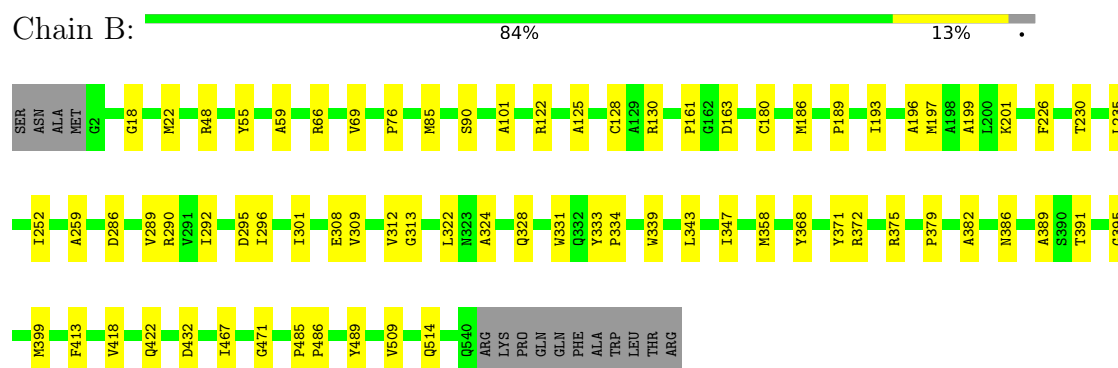
3 Residue-property plots

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.

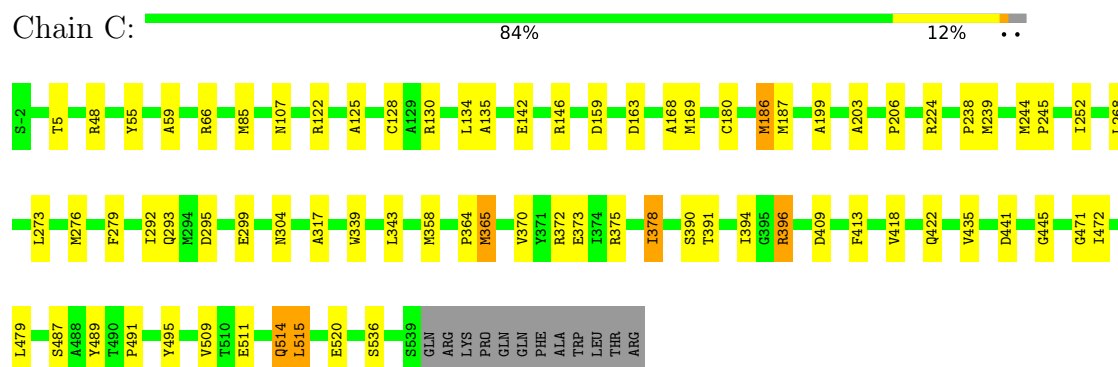
- Molecule 1: Oxalyl-CoA decarboxylase



- Molecule 1: Oxalyl-CoA decarboxylase

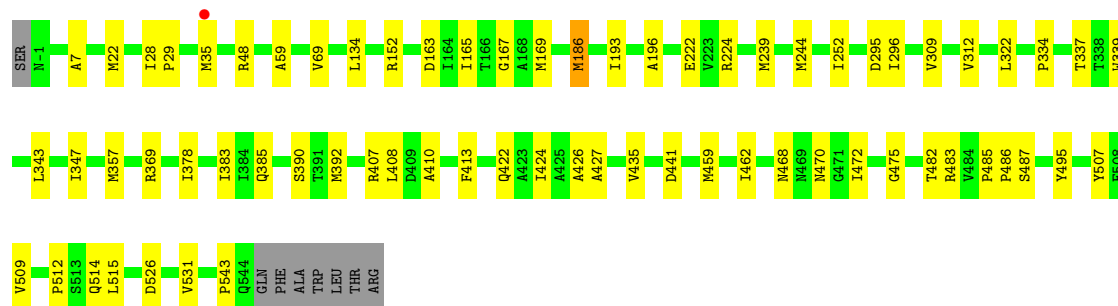


- Molecule 1: Oxalyl-CoA decarboxylase



- Molecule 1: Oxalyl-CoA decarboxylase

Chain D:  86% 12% .



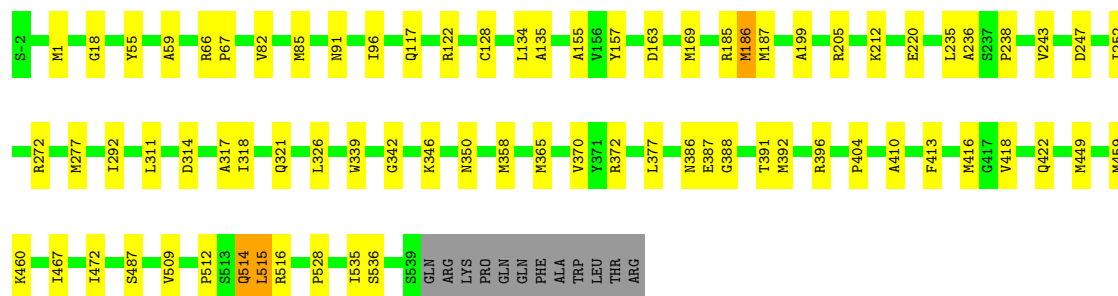
- Molecule 1: Oxalyl-CoA decarboxylase

Chain E:  85% 12% ..



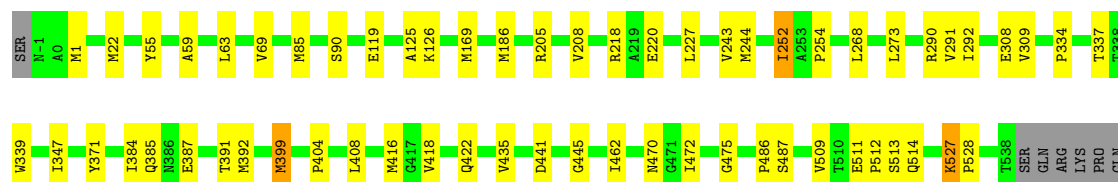
- Molecule 1: Oxalyl-CoA decarboxylase

Chain F:  84% 13% ..



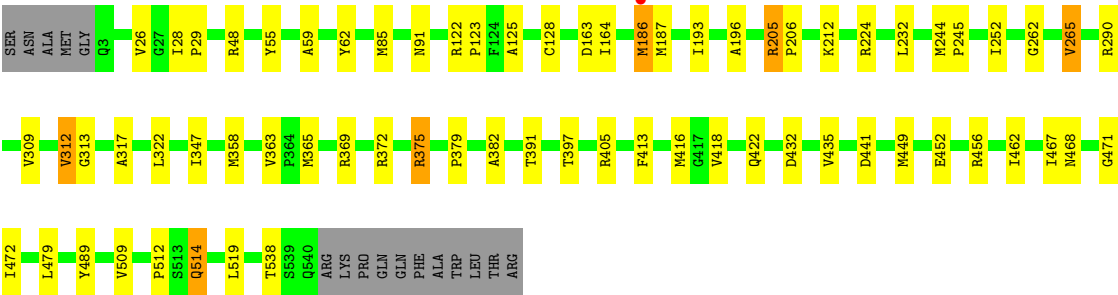
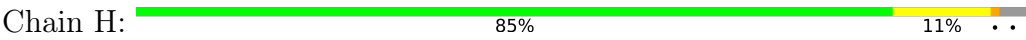
- Molecule 1: Oxalyl-CoA decarboxylase

Chain G:  86% 10% ..



GLN
PHE
ALA
TRP
LEU
THR
ARG

● Molecule 1: Oxalyl-CoA decarboxylase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	324.22Å 77.84Å 229.55Å 90.00° 134.53° 90.00°	Depositor
Resolution (Å)	40.53 – 2.70 40.53 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.6 (40.53-2.70) 99.6 (40.53-2.70)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.42 (at 2.69Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.197 , 0.235 0.197 , 0.235	Depositor DCC
R_{free} test set	5493 reflections (4.87%)	wwPDB-VP
Wilson B-factor (Å ²)	66.0	Xtriage
Anisotropy	0.253	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 44.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.016 for h+2*l,k,-h-l 0.003 for h,-k,-h-l 0.010 for -h-2*l,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	33287	wwPDB-VP
Average B, all atoms (Å ²)	80.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.71% 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: EDO, GOL, TPP, ADP, MG

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	0.08	0/4211	0.24	0/5722
1	B	0.08	0/4183	0.24	0/5685
1	C	0.08	0/4201	0.24	0/5709
1	D	0.08	0/4249	0.25	0/5772
1	E	0.08	0/4203	0.23	0/5710
1	F	0.09	0/4201	0.24	0/5709
1	G	0.08	0/4189	0.23	0/5693
1	H	0.08	0/4179	0.24	0/5680
All	All	0.08	0/33616	0.24	0/45680

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	4111	0	4060	34	0
1	B	4084	0	4027	46	0
1	C	4102	0	4047	43	0
1	D	4149	0	4099	42	0
1	E	4104	0	4053	38	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	F	4102	0	4047	45	0
1	G	4090	0	4037	37	0
1	H	4080	0	4024	40	0
2	A	26	0	16	1	0
2	B	26	0	16	1	0
2	C	26	0	16	1	0
2	D	26	0	16	1	0
2	E	26	0	16	3	0
2	F	26	0	16	1	0
2	G	26	0	16	3	0
2	H	26	0	16	2	0
3	A	27	0	12	1	0
3	B	27	0	12	0	0
3	C	27	0	12	1	0
3	D	27	0	12	1	0
3	E	27	0	12	0	0
3	F	27	0	12	1	0
3	G	27	0	12	0	0
3	H	27	0	12	0	0
4	A	6	0	8	2	0
4	B	6	0	8	1	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
5	C	1	0	0	0	0
5	D	1	0	0	0	0
5	E	1	0	0	0	0
5	F	1	0	0	0	0
5	G	1	0	0	0	0
5	H	1	0	0	0	0
6	C	4	0	6	0	0
6	F	4	0	6	0	0
6	G	4	0	6	1	0
6	H	4	0	6	0	0
7	C	1	0	0	0	0
7	E	1	0	0	0	0
7	F	2	0	0	0	0
7	G	1	0	0	0	0
All	All	33287	0	32658	314	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 (314) 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:85:MET:HE1	1:B:125:ALA:HB2	1.77	0.67
1:C:85:MET:HE1	1:C:125:ALA:HB2	1.78	0.65
1:H:252:ILE:HA	1:H:347:ILE:HD11	1.82	0.62
1:C:479:LEU:HD12	1:C:491:PRO:HD3	1.84	0.59
1:F:220:GLU:HG2	1:F:404:PRO:HG3	1.84	0.59
1:F:387:GLU:HA	1:F:392:MET:HE2	1.84	0.59
1:F:311:LEU:HB3	1:F:318:ILE:HD13	1.84	0.59
1:C:48:ARG:NH2	1:D:495:TYR:OH	2.30	0.59
1:G:509:VAL:HG13	1:G:514:GLN:HB2	1.84	0.58
1:B:193:ILE:HG23	1:B:322:LEU:HD23	1.85	0.58
1:F:235:LEU:HB2	1:F:252:ILE:HD11	1.85	0.58
1:C:59:ALA:HB2	1:C:422:GLN:HG2	1.86	0.58
1:G:59:ALA:HB2	1:G:422:GLN:HG2	1.86	0.58
1:H:391:THR:HG23	1:H:467:ILE:HD12	1.85	0.58
1:G:511:GLU:HG2	1:G:512:PRO:HD2	1.84	0.58
1:A:391:THR:HG23	1:A:467:ILE:HD12	1.86	0.57
1:F:59:ALA:HB2	1:F:422:GLN:HG2	1.86	0.57
1:G:252:ILE:HG22	1:G:347:ILE:HD11	1.87	0.57
1:D:59:ALA:HB2	1:D:422:GLN:HG2	1.86	0.57
1:B:59:ALA:HB2	1:B:422:GLN:HG2	1.87	0.57
1:B:130:ARG:HE	1:B:161:PRO:HG3	1.69	0.57
1:C:365:MET:HE2	1:C:370:VAL:HG22	1.87	0.57
1:E:369:ARG:HH12	1:E:373:GLU:HG2	1.70	0.56
1:E:379:PRO:HG2	1:E:382:ALA:HB2	1.88	0.56
1:F:377:LEU:HD11	1:F:516:ARG:HG3	1.86	0.56
1:F:391:THR:HG23	1:F:467:ILE:HD12	1.88	0.56
1:H:91:ASN:OD1	1:H:212:LYS:NZ	2.38	0.56
1:A:378:ILE:HG21	1:A:384:ILE:HD11	1.87	0.56
1:B:358:MET:HE1	1:B:375:ARG:HH21	1.71	0.56
1:C:358:MET:HB3	1:C:372:ARG:HD2	1.88	0.55
1:C:472:ILE:HG23	2:C:601:TPP:H71	1.89	0.55
1:E:472:ILE:HG23	2:E:601:TPP:H62	1.87	0.55
1:F:472:ILE:HB	1:F:487:SER:HB2	1.88	0.55
1:E:134:LEU:HB2	1:E:169:MET:HE3	1.89	0.55
1:C:66:ARG:NH2	1:C:180:CYS:SG	2.80	0.55
1:D:369:ARG:HE	1:D:512:PRO:HB3	1.72	0.55
1:D:435:VAL:HB	1:D:462:ILE:HD13	1.88	0.55
1:C:495:TYR:OH	1:D:48:ARG:NH2	2.35	0.54
1:F:117:GLN:NE2	1:F:157:TYR:OH	2.37	0.54
1:H:471:GLY:HA2	1:H:489:TYR:HD2	1.73	0.54
1:C:273:LEU:O	1:C:304:ASN:ND2	2.39	0.54
1:G:472:ILE:HG23	2:G:601:TPP:H71	1.88	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:346:LYS:O	1:F:350:ASN:ND2	2.33	0.53
1:A:241:LYS:NZ	1:A:246:ASP:OD2	2.41	0.53
1:A:273:LEU:O	1:A:304:ASN:ND2	2.41	0.53
1:B:391:THR:HG23	1:B:467:ILE:HD12	1.89	0.53
1:E:509:VAL:HG13	1:E:514:GLN:HB2	1.90	0.53
1:G:371:TYR:OH	1:G:391:THR:O	2.24	0.53
1:H:196:ALA:HB2	1:H:309:VAL:HG11	1.91	0.53
1:A:472:ILE:HG23	2:A:601:TPP:H71	1.91	0.53
1:B:252:ILE:HA	1:B:347:ILE:HD11	1.91	0.53
1:B:163:ASP:OD1	1:B:163:ASP:N	2.38	0.52
1:B:371:TYR:HE1	1:B:399:MET:HE2	1.74	0.52
1:E:358:MET:HE2	1:E:372:ARG:HG3	1.91	0.52
1:E:59:ALA:HB2	1:E:422:GLN:HG2	1.91	0.52
1:F:365:MET:HE3	1:F:512:PRO:HG3	1.92	0.52
1:F:509:VAL:HG13	1:F:514:GLN:HB2	1.91	0.52
1:H:265:VAL:HG23	1:H:290:ARG:HB2	1.92	0.52
1:G:387:GLU:HA	1:G:392:MET:HE2	1.92	0.52
1:E:472:ILE:HB	1:E:487:SER:HB2	1.91	0.52
1:A:228:GLU:O	1:A:231:GLN:NE2	2.39	0.51
1:F:243:VAL:HA	1:F:404:PRO:HB3	1.92	0.51
1:C:199:ALA:HB3	1:C:292:ILE:HD11	1.93	0.51
1:H:472:ILE:HG23	2:H:601:TPP:H71	1.93	0.51
1:E:441:ASP:OD2	1:E:471:GLY:N	2.42	0.51
1:G:441:ASP:O	1:H:48:ARG:NH2	2.44	0.51
1:H:452:GLU:OE1	1:H:456:ARG:NH2	2.43	0.51
1:A:196:ALA:HB2	1:A:309:VAL:HG11	1.93	0.51
1:C:142:GLU:OE1	1:C:146:ARG:NH1	2.44	0.51
1:D:196:ALA:HB2	1:D:309:VAL:HG11	1.92	0.51
1:H:206:PRO:HA	1:H:265:VAL:HG12	1.93	0.51
1:A:415:THR:HG23	1:B:76:PRO:HB3	1.93	0.50
1:H:187:MET:HG2	1:H:317:ALA:HB1	1.92	0.50
1:A:163:ASP:OD1	1:A:163:ASP:N	2.41	0.50
1:C:509:VAL:HG13	1:C:514:GLN:HB2	1.93	0.50
1:G:384:ILE:HD13	1:G:399:MET:HG2	1.92	0.50
1:G:445:GLY:HA2	1:H:449:MET:HE2	1.92	0.50
1:C:441:ASP:OD1	1:C:441:ASP:N	2.45	0.50
1:F:18:GLY:HA3	1:F:66:ARG:HD3	1.93	0.50
1:D:193:ILE:HG23	1:D:322:LEU:HD23	1.94	0.50
1:F:187:MET:HG2	1:F:317:ALA:HB1	1.94	0.50
1:F:358:MET:HB3	1:F:372:ARG:HD3	1.94	0.50
1:C:134:LEU:HB2	1:C:169:MET:HE3	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:383:ILE:HB	1:D:435:VAL:HG22	1.93	0.50
1:A:122:ARG:NH2	1:A:128:CYS:HB2	2.26	0.50
1:C:187:MET:HG2	1:C:317:ALA:HB1	1.94	0.50
1:E:442:SER:HB3	2:E:601:TPP:O7	2.11	0.50
1:D:472:ILE:HG23	2:D:601:TPP:H71	1.92	0.49
1:G:55:TYR:CG	1:G:418:VAL:HG12	2.47	0.49
1:H:55:TYR:CG	1:H:418:VAL:HG12	2.47	0.49
1:F:163:ASP:OD1	1:F:163:ASP:N	2.43	0.49
1:F:199:ALA:HB3	1:F:292:ILE:HD11	1.94	0.49
1:H:379:PRO:HG2	1:H:382:ALA:HB2	1.93	0.49
1:F:122:ARG:NH2	1:F:128:CYS:HB2	2.27	0.49
1:D:482:THR:HG23	1:D:483:ARG:HG3	1.94	0.49
1:D:392:MET:HE1	1:D:410:ALA:HB2	1.94	0.49
1:F:186:MET:SD	1:F:186:MET:N	2.83	0.49
1:F:85:MET:HG3	1:F:155:ALA:HB1	1.95	0.49
1:B:371:TYR:CE1	1:B:399:MET:HE2	2.48	0.48
1:A:479:LEU:HD12	1:A:484:VAL:HG22	1.95	0.48
1:F:509:VAL:HG11	1:F:515:LEU:HD22	1.94	0.48
1:A:101:ALA:HB3	1:A:161:PRO:HA	1.94	0.48
1:A:205:ARG:HD2	1:A:339:TRP:CE3	2.48	0.48
1:E:445:GLY:HA2	1:F:449:MET:HE2	1.94	0.48
1:A:59:ALA:HB2	1:A:422:GLN:HG2	1.96	0.48
1:C:268:LEU:HB2	1:C:293:GLN:HA	1.95	0.48
1:F:55:TYR:CG	1:F:418:VAL:HG12	2.49	0.48
1:A:55:TYR:CG	1:A:418:VAL:HG12	2.49	0.48
1:D:468:ASN:HD22	1:D:495:TYR:HE2	1.62	0.48
1:E:55:TYR:CG	1:E:418:VAL:HG12	2.49	0.48
1:F:67:PRO:HB2	1:F:96:ILE:HG13	1.96	0.48
1:B:509:VAL:HG13	1:B:514:GLN:HB2	1.94	0.48
1:G:85:MET:HE1	1:G:125:ALA:HB2	1.96	0.47
1:H:59:ALA:HB2	1:H:422:GLN:HG2	1.96	0.47
1:F:236:ALA:O	1:F:396:ARG:NH2	2.45	0.47
1:B:18:GLY:HA3	1:B:66:ARG:HD3	1.96	0.47
1:C:295:ASP:OD1	3:C:602:ADP:O2'	2.21	0.47
1:H:435:VAL:HB	1:H:462:ILE:HG12	1.96	0.47
1:D:427:ALA:HB3	1:D:459:MET:HE2	1.95	0.47
1:E:252:ILE:HG21	1:E:343:LEU:HD13	1.96	0.47
1:F:392:MET:HE1	1:F:410:ALA:HB2	1.97	0.47
1:H:358:MET:HE1	1:H:375:ARG:HH12	1.80	0.47
1:B:196:ALA:HB2	1:B:309:VAL:HG11	1.96	0.47
1:D:252:ILE:HA	1:D:347:ILE:HD11	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:241:LYS:HD2	1:E:396:ARG:HD3	1.96	0.47
1:E:440:GLY:HA3	2:E:601:TPP:O1A	2.14	0.47
1:C:55:TYR:CG	1:C:418:VAL:HG12	2.50	0.47
1:F:91:ASN:OD1	1:F:212:LYS:NZ	2.48	0.47
1:H:193:ILE:HG23	1:H:322:LEU:HD23	1.97	0.47
1:H:224:ARG:HA	1:H:244:MET:HE2	1.96	0.47
1:H:509:VAL:HG13	1:H:514:GLN:HB2	1.97	0.47
1:A:90:SER:HA	4:A:603:GOL:H31	1.96	0.47
1:C:373:GLU:HG2	1:C:515:LEU:HB3	1.97	0.47
1:A:295:ASP:OD1	3:A:602:ADP:O2'	2.26	0.47
1:H:186:MET:HG2	1:H:313:GLY:HA2	1.96	0.47
1:F:247:ASP:OD1	1:F:247:ASP:N	2.45	0.46
1:G:254:PRO:HG2	1:G:347:ILE:HG12	1.96	0.46
1:G:408:LEU:HD13	1:G:422:GLN:HB3	1.96	0.46
1:C:445:GLY:HA3	1:D:48:ARG:HD2	1.97	0.46
1:E:367:TYR:HE1	1:E:391:THR:HG22	1.79	0.46
1:B:226:PHE:O	1:B:230:THR:HG22	2.16	0.46
1:B:259:ALA:HA	1:B:339:TRP:HH2	1.79	0.46
1:B:286:ASP:HB3	1:B:289:VAL:HB	1.98	0.46
1:F:134:LEU:HB2	1:F:169:MET:HE3	1.97	0.46
1:G:334:PRO:HB2	1:G:337:THR:HG23	1.97	0.46
1:B:85:MET:HE2	1:B:85:MET:HB3	1.67	0.46
1:D:334:PRO:HB2	1:D:337:THR:HG23	1.96	0.46
1:E:149:ILE:HG13	1:E:150:TYR:HD1	1.80	0.46
1:G:220:GLU:HB3	1:G:404:PRO:HG3	1.96	0.46
1:C:163:ASP:OD1	1:C:163:ASP:N	2.49	0.46
1:C:375:ARG:HA	1:C:378:ILE:HG12	1.98	0.46
1:B:90:SER:HA	4:B:603:GOL:H32	1.97	0.46
1:C:299:GLU:OE1	1:C:299:GLU:N	2.48	0.46
1:G:385:GLN:HG2	1:G:408:LEU:HB2	1.97	0.46
1:A:480:ASP:OD2	1:A:482:THR:OG1	2.30	0.46
1:E:85:MET:HA	1:E:95:MET:HE1	1.97	0.46
1:C:203:ALA:HB3	1:C:206:PRO:HG3	1.98	0.46
1:C:130:ARG:HB2	1:C:159:ASP:HB3	1.97	0.46
1:E:199:ALA:HB3	1:E:292:ILE:HD11	1.97	0.46
1:H:122:ARG:NH2	1:H:128:CYS:HB2	2.31	0.46
1:B:197:MET:HE2	1:B:322:LEU:HD22	1.99	0.45
1:G:227:LEU:HD12	1:G:244:MET:HG3	1.98	0.45
1:B:379:PRO:HG2	1:B:382:ALA:HB2	1.97	0.45
1:D:509:VAL:HG13	1:D:514:GLN:HB2	1.97	0.45
1:E:287:PRO:HG2	1:E:288:LYS:HD2	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:339:TRP:CZ2	1:B:343:LEU:HD11	2.51	0.45
1:H:471:GLY:HA2	1:H:489:TYR:CD2	2.50	0.45
1:B:55:TYR:CG	1:B:418:VAL:HG12	2.52	0.45
1:C:122:ARG:NH2	1:C:128:CYS:HB2	2.31	0.45
1:D:7:ALA:HB2	1:D:165:ILE:HG23	1.98	0.45
1:B:186:MET:HB2	1:B:313:GLY:HA2	1.98	0.45
1:E:387:GLU:HA	1:E:392:MET:HE2	1.99	0.45
1:C:107:ASN:HD21	1:C:130:ARG:HH12	1.65	0.45
1:G:470:ASN:O	1:G:475:GLY:HA3	2.17	0.45
1:E:107:ASN:HD21	1:E:130:ARG:HH12	1.64	0.45
1:F:187:MET:HB3	1:F:321:GLN:NE2	2.32	0.45
1:B:471:GLY:HA2	1:B:489:TYR:CD2	2.52	0.45
1:F:82:VAL:HA	1:F:85:MET:HE2	1.98	0.45
1:H:244:MET:SD	1:H:245:PRO:HD2	2.56	0.45
1:B:22:MET:HE2	1:B:69:VAL:HG11	1.99	0.45
2:G:601:TPP:HN42	2:G:601:TPP:C2	2.30	0.45
1:H:205:ARG:HD2	1:H:262:GLY:O	2.17	0.45
1:B:324:ALA:O	1:B:328:GLN:HG2	2.18	0.44
1:H:369:ARG:CZ	1:H:512:PRO:HB3	2.47	0.44
1:B:358:MET:HB3	1:B:372:ARG:HD3	1.99	0.44
1:D:134:LEU:HD23	1:D:169:MET:HG2	2.00	0.44
1:E:254:PRO:O	1:E:350:ASN:ND2	2.51	0.44
1:A:215:ALA:O	1:A:405:ARG:NH1	2.47	0.44
1:H:363:VAL:HG22	1:H:512:PRO:HD3	2.00	0.44
1:B:358:MET:HE2	1:B:372:ARG:HB2	1.99	0.44
1:B:122:ARG:NH2	1:B:128:CYS:HB2	2.33	0.44
1:A:268:LEU:HB2	1:A:293:GLN:HA	2.00	0.44
1:D:222:GLU:N	1:D:222:GLU:OE1	2.50	0.44
1:B:66:ARG:NH2	1:B:180:CYS:O	2.47	0.44
1:B:101:ALA:HB3	1:B:161:PRO:HA	2.00	0.44
1:B:189:PRO:HA	1:E:187:MET:SD	2.58	0.44
1:B:333:TYR:CD1	1:B:334:PRO:HD2	2.52	0.44
1:B:358:MET:HG2	1:B:368:TYR:HB3	2.00	0.44
1:B:371:TYR:OH	1:B:391:THR:O	2.36	0.43
1:E:273:LEU:O	1:E:304:ASN:ND2	2.48	0.43
1:F:386:ASN:CG	1:F:392:MET:HA	2.43	0.43
1:A:73:VAL:HG22	1:A:77:GLY:HA3	2.00	0.43
1:E:201:LYS:HG2	1:E:331:TRP:CE2	2.53	0.43
1:F:135:ALA:HB2	1:F:169:MET:HE1	2.00	0.43
1:F:365:MET:O	1:F:536:SER:N	2.43	0.43
1:B:386:ASN:HD21	1:B:395:GLY:HA3	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:134:LEU:HD21	1:D:167:GLY:C	2.43	0.43
1:G:435:VAL:HB	1:G:462:ILE:HG12	2.00	0.43
1:G:472:ILE:HB	1:G:487:SER:HB2	2.00	0.43
1:D:295:ASP:OD1	1:D:296:ILE:N	2.52	0.43
1:E:28:ILE:HA	1:E:29:PRO:HA	1.87	0.43
1:A:55:TYR:CD1	1:A:418:VAL:HG12	2.54	0.43
1:C:339:TRP:CZ2	1:C:343:LEU:HD11	2.52	0.43
1:D:186:MET:H	1:D:186:MET:HG3	1.67	0.43
1:D:509:VAL:HG11	1:D:515:LEU:HD23	2.00	0.43
1:B:485:PRO:HA	1:B:486:PRO:HD3	1.92	0.43
1:H:28:ILE:HD12	1:H:28:ILE:H	1.84	0.43
1:E:442:SER:HB2	1:E:489:TYR:HE2	1.84	0.43
1:F:238:PRO:HD3	1:F:396:ARG:HH21	1.84	0.43
1:F:272:ARG:HG3	3:F:602:ADP:O5'	2.19	0.43
1:F:416:MET:HE3	2:F:601:TPP:S1	2.58	0.43
1:A:252:ILE:HG21	1:A:343:LEU:HD13	2.01	0.43
1:A:394:ILE:HG12	1:A:541:ARG:HE	1.84	0.43
1:E:265:VAL:HG22	1:E:290:ARG:HB2	2.01	0.43
1:E:471:GLY:HA2	1:E:489:TYR:CD2	2.54	0.43
1:G:208:VAL:HG11	1:G:227:LEU:HD21	1.99	0.43
1:H:416:MET:HE3	2:H:601:TPP:S1	2.59	0.43
1:B:389:ALA:HB2	2:B:601:TPP:S1	2.59	0.42
1:D:152:ARG:NH2	3:D:602:ADP:O3'	2.50	0.42
1:A:278:HIS:CG	1:A:284:ARG:HD3	2.54	0.42
1:E:193:ILE:HG23	1:E:322:LEU:HD23	2.01	0.42
1:E:363:VAL:HA	1:E:364:PRO:C	2.44	0.42
1:A:290:ARG:HD2	1:A:308:GLU:OE1	2.18	0.42
1:D:470:ASN:O	1:D:475:GLY:HA3	2.19	0.42
1:G:416:MET:HE3	2:G:601:TPP:S1	2.59	0.42
1:H:62:TYR:CZ	1:H:405:ARG:HG2	2.54	0.42
1:A:205:ARG:NE	1:A:262:GLY:O	2.30	0.42
1:D:163:ASP:OD1	1:D:163:ASP:N	2.51	0.42
1:G:268:LEU:HD23	1:G:291:VAL:HG13	2.01	0.42
1:G:511:GLU:OE1	1:G:513:SER:OG	2.32	0.42
1:C:244:MET:SD	1:C:245:PRO:HD2	2.59	0.42
1:D:339:TRP:CZ2	1:D:343:LEU:HD11	2.55	0.42
1:E:392:MET:HE1	1:E:410:ALA:HB2	2.01	0.42
1:E:515:LEU:HD23	1:E:515:LEU:HA	1.83	0.42
1:G:90:SER:HA	6:G:603:EDO:H22	2.02	0.42
1:H:252:ILE:H	1:H:252:ILE:HG13	1.70	0.42
1:C:364:PRO:HB2	1:C:536:SER:HB3	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:239:MET:HE3	1:D:407:ARG:O	2.19	0.42
1:E:343:LEU:O	1:E:347:ILE:HD12	2.20	0.42
1:D:485:PRO:HA	1:D:486:PRO:HD3	1.93	0.42
1:F:205:ARG:HB3	1:F:339:TRP:CD2	2.54	0.42
1:A:186:MET:SD	1:A:186:MET:N	2.92	0.42
1:D:28:ILE:HA	1:D:29:PRO:HA	1.88	0.42
1:D:239:MET:HE2	1:D:239:MET:HB3	1.96	0.42
1:D:441:ASP:OD1	1:D:441:ASP:N	2.52	0.42
1:G:63:LEU:O	1:G:218:ARG:NH2	2.52	0.42
1:G:186:MET:SD	1:G:186:MET:N	2.93	0.42
1:G:290:ARG:HD2	1:G:308:GLU:OE1	2.19	0.42
1:C:239:MET:HG3	1:C:409:ASP:OD2	2.20	0.42
1:C:472:ILE:HB	1:C:487:SER:HB2	2.02	0.42
1:F:185:ARG:NH1	1:F:314:ASP:OD2	2.45	0.42
1:F:459:MET:O	1:F:528:PRO:HG3	2.20	0.42
1:B:199:ALA:HB3	1:B:292:ILE:HD11	2.01	0.41
1:C:252:ILE:H	1:C:252:ILE:HG13	1.72	0.41
1:D:383:ILE:HG21	1:D:426:ALA:HB1	2.02	0.41
1:C:224:ARG:HG2	1:C:244:MET:SD	2.61	0.41
1:D:424:ILE:O	1:D:459:MET:HE1	2.20	0.41
1:F:370:VAL:HG21	1:F:535:ILE:HG22	2.02	0.41
1:E:441:ASP:OD1	1:E:441:ASP:N	2.50	0.41
1:B:290:ARG:HD2	1:B:308:GLU:OE1	2.20	0.41
1:B:301:ILE:HG13	1:B:312:VAL:HG22	2.02	0.41
1:G:292:ILE:HG23	1:G:309:VAL:HB	2.03	0.41
1:A:245:PRO:O	1:A:248:HIS:HB2	2.21	0.41
1:B:201:LYS:HA	1:B:331:TRP:CZ2	2.55	0.41
1:C:471:GLY:HA2	1:C:489:TYR:CD2	2.56	0.41
1:H:441:ASP:HB3	1:H:468:ASN:HA	2.03	0.41
1:A:12:ARG:NH1	1:A:171:GLU:OE2	2.52	0.41
1:C:186:MET:HG3	1:H:312:VAL:HG13	2.03	0.41
1:C:238:PRO:HG3	1:C:396:ARG:HD2	2.03	0.41
1:D:526:ASP:OD1	1:D:526:ASP:N	2.53	0.41
1:G:119:GLU:HB3	1:H:123:PRO:HG3	2.02	0.41
1:G:205:ARG:HB3	1:G:339:TRP:CD2	2.55	0.41
1:H:28:ILE:HA	1:H:29:PRO:HA	1.90	0.41
1:C:5:THR:HG22	1:C:168:ALA:HB2	2.03	0.41
1:D:22:MET:HE2	1:D:69:VAL:HG11	2.01	0.41
1:F:388:GLY:H	1:F:392:MET:HE3	1.85	0.41
1:F:388:GLY:O	1:F:392:MET:HE3	2.21	0.41
1:A:445:GLY:HA3	1:B:48:ARG:HH21	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:135:ALA:HB2	1:C:169:MET:HE1	2.03	0.41
1:D:357:MET:HE1	1:D:543:PRO:HD2	2.03	0.41
1:D:507:TYR:HD2	1:D:531:VAL:HG12	1.86	0.41
1:E:85:MET:HE1	1:E:125:ALA:HB2	2.02	0.41
1:G:126:LYS:HE2	1:G:126:LYS:HB3	1.87	0.41
1:G:527:LYS:HG3	1:G:528:PRO:HD2	2.03	0.41
1:H:432:ASP:OD1	1:H:432:ASP:N	2.46	0.41
1:A:252:ILE:HA	1:A:347:ILE:HD11	2.02	0.41
1:A:353:THR:HB	1:A:542:LYS:HE3	2.02	0.41
1:C:186:MET:SD	1:C:186:MET:N	2.82	0.41
1:D:224:ARG:HG2	1:D:244:MET:SD	2.61	0.41
1:F:342:GLY:O	1:F:346:LYS:HG2	2.21	0.41
1:D:385:GLN:HG2	1:D:408:LEU:HB2	2.03	0.40
1:E:171:GLU:O	1:E:174:VAL:HG22	2.21	0.40
1:H:479:LEU:HD23	1:H:479:LEU:HA	1.85	0.40
1:B:235:LEU:HB2	1:B:252:ILE:HD11	2.03	0.40
1:B:295:ASP:OD1	1:B:296:ILE:N	2.54	0.40
1:C:273:LEU:O	1:C:279:PHE:HA	2.21	0.40
1:H:85:MET:HE1	1:H:125:ALA:HB2	2.03	0.40
1:C:365:MET:HE1	1:C:515:LEU:HD11	2.03	0.40
1:D:472:ILE:HB	1:D:487:SER:HB2	2.02	0.40
1:G:22:MET:HE2	1:G:69:VAL:HG11	2.03	0.40
1:A:89:TRP:CZ2	4:A:603:GOL:H32	2.57	0.40
1:G:486:PRO:HB2	1:H:26:VAL:HG11	2.04	0.40
1:H:163:ASP:OD1	1:H:163:ASP:N	2.40	0.40
1:G:220:GLU:HA	1:G:243:VAL:HG13	2.04	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	540/554 (98%)	524 (97%)	16 (3%)	0	100	100
1	B	537/554 (97%)	522 (97%)	15 (3%)	0	100	100
1	C	540/554 (98%)	521 (96%)	18 (3%)	1 (0%)	43	68
1	D	545/554 (98%)	530 (97%)	15 (3%)	0	100	100
1	E	539/554 (97%)	520 (96%)	19 (4%)	0	100	100
1	F	540/554 (98%)	525 (97%)	15 (3%)	0	100	100
1	G	538/554 (97%)	521 (97%)	17 (3%)	0	100	100
1	H	536/554 (97%)	518 (97%)	18 (3%)	0	100	100
All	All	4315/4432 (97%)	4181 (97%)	133 (3%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	515	LEU

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	422/432 (98%)	416 (99%)	6 (1%)	59	82
1	B	419/432 (97%)	417 (100%)	2 (0%)	81	92
1	C	421/432 (98%)	408 (97%)	13 (3%)	35	65
1	D	426/432 (99%)	419 (98%)	7 (2%)	55	80
1	E	421/432 (98%)	409 (97%)	12 (3%)	37	67
1	F	421/432 (98%)	413 (98%)	8 (2%)	50	77
1	G	419/432 (97%)	413 (99%)	6 (1%)	59	82
1	H	419/432 (97%)	405 (97%)	14 (3%)	33	63
All	All	3368/3456 (98%)	3300 (98%)	68 (2%)	48	76

All (68) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	169	MET
1	A	377	LEU
1	A	413	PHE
1	A	479	LEU
1	A	541	ARG
1	A	542	LYS
1	B	413	PHE
1	B	432	ASP
1	C	186	MET
1	C	276	MET
1	C	365	MET
1	C	378	ILE
1	C	390	SER
1	C	391	THR
1	C	394	ILE
1	C	396	ARG
1	C	413	PHE
1	C	435	VAL
1	C	511	GLU
1	C	514	GLN
1	C	520	GLU
1	D	35[A]	MET
1	D	35[B]	MET
1	D	186	MET
1	D	312	VAL
1	D	378	ILE
1	D	390	SER
1	D	413	PHE
1	E	8	GLN
1	E	186	MET
1	E	252	ILE
1	E	338	THR
1	E	347	ILE
1	E	357	MET
1	E	390	SER
1	E	394	ILE
1	E	396	ARG
1	E	413	PHE
1	E	540	GLN
1	E	541	ARG
1	F	1	MET
1	F	186	MET
1	F	277	MET

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Mol	Chain	Res	Type
1	F	326	LEU
1	F	413	PHE
1	F	460	LYS
1	F	514	GLN
1	F	515	LEU
1	G	1	MET
1	G	169	MET
1	G	252	ILE
1	G	273	LEU
1	G	399	MET
1	G	527	LYS
1	H	164	ILE
1	H	186	MET
1	H	205	ARG
1	H	232	LEU
1	H	265	VAL
1	H	312	VAL
1	H	365	MET
1	H	372	ARG
1	H	375	ARG
1	H	397	THR
1	H	413	PHE
1	H	514	GLN
1	H	519	LEU
1	H	538	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (29) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	143	GLN
1	B	102	ASN
1	B	386	ASN
1	B	540	GLN
1	C	143	GLN
1	C	328	GLN
1	C	350	ASN
1	C	385	GLN
1	D	3	GLN
1	D	231	GLN
1	D	386	ASN
1	E	87	ASN
1	E	350	ASN

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Mol	Chain	Res	Type
1	E	385	GLN
1	E	422	GLN
1	E	430	HIS
1	F	321	GLN
1	F	329	GLN
1	F	385	GLN
1	F	386	ASN
1	F	422	GLN
1	G	87	ASN
1	G	143	GLN
1	G	323	ASN
1	G	328	GLN
1	G	386	ASN
1	H	143	GLN
1	H	385	GLN
1	H	422	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 30 ligands modelled in this entry, 8 are monoatomic - leaving 22 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	EDO	G	603	-	3,3,3	0.43	0	2,2,2	0.38	0
3	ADP	D	602	-	28,29,29	1.41	5 (17%)	43,45,45	1.81	10 (23%)
3	ADP	H	602	-	28,29,29	1.41	4 (14%)	43,45,45	1.84	10 (23%)
6	EDO	F	603	-	3,3,3	0.43	0	2,2,2	0.37	0
6	EDO	H	603	-	3,3,3	0.42	0	2,2,2	0.38	0
3	ADP	A	602	-	28,29,29	1.40	4 (14%)	43,45,45	1.85	10 (23%)
2	TPP	B	601	5	26,27,27	0.26	0	38,40,40	0.29	0
2	TPP	E	601	5	26,27,27	0.27	0	38,40,40	0.30	0
3	ADP	B	602	-	28,29,29	1.42	4 (14%)	43,45,45	1.86	10 (23%)
2	TPP	D	601	5	26,27,27	0.25	0	38,40,40	0.29	0
4	GOL	A	603	-	5,5,5	0.90	0	5,5,5	1.03	0
2	TPP	F	601	5	26,27,27	0.25	0	38,40,40	0.28	0
2	TPP	C	601	5	26,27,27	0.25	0	38,40,40	0.29	0
2	TPP	G	601	5	26,27,27	0.26	0	38,40,40	0.29	0
3	ADP	G	602	-	28,29,29	1.41	4 (14%)	43,45,45	1.83	10 (23%)
3	ADP	C	602	-	28,29,29	1.41	4 (14%)	43,45,45	1.84	9 (20%)
3	ADP	F	602	-	28,29,29	1.40	4 (14%)	43,45,45	1.84	10 (23%)
3	ADP	E	602	-	28,29,29	1.41	4 (14%)	43,45,45	1.82	10 (23%)
2	TPP	H	601	5	26,27,27	0.26	0	38,40,40	0.29	0
4	GOL	B	603	-	5,5,5	0.93	0	5,5,5	1.07	0
6	EDO	C	603	-	3,3,3	0.42	0	2,2,2	0.39	0
2	TPP	A	601	5	26,27,27	0.25	0	38,40,40	0.28	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	EDO	G	603	-	-	0/1/1/1	-
3	ADP	D	602	-	-	7/16/32/32	0/3/3/3
3	ADP	H	602	-	-	9/16/32/32	0/3/3/3
6	EDO	F	603	-	-	0/1/1/1	-
6	EDO	H	603	-	-	0/1/1/1	-
3	ADP	A	602	-	-	5/16/32/32	0/3/3/3
2	TPP	B	601	5	-	3/17/17/17	0/2/2/2
2	TPP	E	601	5	-	3/17/17/17	0/2/2/2
3	ADP	B	602	-	-	7/16/32/32	0/3/3/3
2	TPP	D	601	5	-	0/17/17/17	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	GOL	A	603	-	-	2/4/4/4	-
2	TPP	F	601	5	-	0/17/17/17	0/2/2/2
2	TPP	C	601	5	-	1/17/17/17	0/2/2/2
2	TPP	G	601	5	-	3/17/17/17	0/2/2/2
3	ADP	G	602	-	-	7/16/32/32	0/3/3/3
3	ADP	C	602	-	-	5/16/32/32	0/3/3/3
3	ADP	F	602	-	-	8/16/32/32	0/3/3/3
3	ADP	E	602	-	-	7/16/32/32	0/3/3/3
2	TPP	H	601	5	-	1/17/17/17	0/2/2/2
4	GOL	B	603	-	-	0/4/4/4	-
6	EDO	C	603	-	-	0/1/1/1	-
2	TPP	A	601	5	-	0/17/17/17	0/2/2/2

All (33) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	602	ADP	C5-C4	4.71	1.47	1.39
3	E	602	ADP	C5-C4	4.67	1.47	1.39
3	A	602	ADP	C5-C4	4.66	1.47	1.39
3	H	602	ADP	C5-C4	4.66	1.47	1.39
3	G	602	ADP	C5-C4	4.66	1.47	1.39
3	F	602	ADP	C5-C4	4.64	1.47	1.39
3	D	602	ADP	C5-C4	4.62	1.47	1.39
3	B	602	ADP	C5-C4	4.62	1.47	1.39
3	B	602	ADP	C5-C6	2.78	1.48	1.41
3	C	602	ADP	C5-C6	2.77	1.48	1.41
3	G	602	ADP	C5-C6	2.74	1.48	1.41
3	F	602	ADP	C5-C6	2.73	1.48	1.41
3	H	602	ADP	C5-C6	2.73	1.48	1.41
3	A	602	ADP	C5-C6	2.73	1.48	1.41
3	D	602	ADP	C5-C6	2.72	1.48	1.41
3	E	602	ADP	C5-C6	2.72	1.48	1.41
3	B	602	ADP	C8-N7	2.47	1.36	1.31
3	D	602	ADP	C8-N7	2.43	1.36	1.31
3	C	602	ADP	C8-N7	2.41	1.36	1.31
3	F	602	ADP	C8-N7	2.40	1.36	1.31
3	H	602	ADP	C8-N7	2.40	1.36	1.31
3	G	602	ADP	C8-N7	2.39	1.36	1.31
3	A	602	ADP	C8-N7	2.39	1.36	1.31
3	E	602	ADP	C8-N7	2.37	1.36	1.31
3	C	602	ADP	C5-N7	-2.23	1.35	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	G	602	ADP	C5-N7	-2.22	1.35	1.39
3	F	602	ADP	C5-N7	-2.20	1.35	1.39
3	E	602	ADP	C5-N7	-2.20	1.35	1.39
3	H	602	ADP	C5-N7	-2.19	1.35	1.39
3	A	602	ADP	C5-N7	-2.19	1.35	1.39
3	D	602	ADP	C5-N7	-2.18	1.35	1.39
3	B	602	ADP	C5-N7	-2.04	1.35	1.39
3	D	602	ADP	C4-N9	-2.00	1.33	1.37

All (79) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	602	ADP	C5-C4-N3	-5.82	118.70	126.72
3	G	602	ADP	C5-C4-N3	-5.69	118.88	126.72
3	F	602	ADP	C5-C4-N3	-5.68	118.90	126.72
3	A	602	ADP	C5-C4-N3	-5.62	118.98	126.72
3	H	602	ADP	C5-C4-N3	-5.61	118.99	126.72
3	E	602	ADP	C5-C4-N3	-5.57	119.05	126.72
3	B	602	ADP	C5-C4-N3	-5.51	119.12	126.72
3	D	602	ADP	C5-C4-N3	-5.45	119.21	126.72
3	C	602	ADP	N3-C4-N9	4.51	134.84	127.17
3	A	602	ADP	N3-C4-N9	4.46	134.75	127.17
3	F	602	ADP	N3-C4-N9	4.46	134.75	127.17
3	G	602	ADP	N3-C4-N9	4.44	134.72	127.17
3	E	602	ADP	N3-C4-N9	4.41	134.67	127.17
3	B	602	ADP	N3-C4-N9	4.40	134.66	127.17
3	H	602	ADP	N3-C4-N9	4.40	134.65	127.17
3	D	602	ADP	N3-C4-N9	4.22	134.34	127.17
3	C	602	ADP	C2-N3-C4	3.72	120.91	111.83
3	F	602	ADP	C2-N3-C4	3.70	120.87	111.83
3	G	602	ADP	C2-N3-C4	3.68	120.82	111.83
3	B	602	ADP	C2-N3-C4	3.67	120.80	111.83
3	H	602	ADP	C2-N3-C4	3.66	120.77	111.83
3	A	602	ADP	C2-N3-C4	3.66	120.76	111.83
3	E	602	ADP	C2-N3-C4	3.61	120.65	111.83
3	D	602	ADP	C2-N3-C4	3.60	120.61	111.83
3	C	602	ADP	C4-C5-N7	-3.59	106.48	110.58
3	D	602	ADP	C4-C5-N7	-3.57	106.50	110.58
3	G	602	ADP	C4-C5-N7	-3.56	106.51	110.58
3	B	602	ADP	C4-C5-N7	-3.55	106.52	110.58
3	F	602	ADP	C4-C5-N7	-3.52	106.56	110.58
3	A	602	ADP	C4-C5-N7	-3.51	106.57	110.58

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	H	602	ADP	C4-C5-N7	-3.51	106.57	110.58
3	E	602	ADP	C4-C5-N7	-3.50	106.58	110.58
3	F	602	ADP	N3-C2-N1	-3.32	123.55	128.58
3	H	602	ADP	N3-C2-N1	-3.31	123.58	128.58
3	G	602	ADP	N3-C2-N1	-3.30	123.59	128.58
3	A	602	ADP	N3-C2-N1	-3.28	123.62	128.58
3	D	602	ADP	N3-C2-N1	-3.27	123.62	128.58
3	E	602	ADP	N3-C2-N1	-3.27	123.63	128.58
3	C	602	ADP	N3-C2-N1	-3.24	123.67	128.58
3	B	602	ADP	N3-C2-N1	-3.21	123.72	128.58
3	B	602	ADP	C4-N9-C8	3.16	109.05	105.74
3	A	602	ADP	C4-N9-C8	2.85	108.73	105.74
3	E	602	ADP	C4-N9-C8	2.84	108.72	105.74
3	D	602	ADP	C4-N9-C8	2.80	108.68	105.74
3	F	602	ADP	C4-N9-C8	2.75	108.63	105.74
3	H	602	ADP	C4-N9-C8	2.74	108.62	105.74
3	G	602	ADP	C4-N9-C8	2.69	108.56	105.74
3	B	602	ADP	C5-N7-C8	2.61	107.56	103.45
3	A	602	ADP	C5-N7-C8	2.60	107.53	103.45
3	C	602	ADP	C5-N7-C8	2.59	107.52	103.45
3	F	602	ADP	C5-N7-C8	2.58	107.51	103.45
3	G	602	ADP	C5-N7-C8	2.58	107.51	103.45
3	E	602	ADP	C5-N7-C8	2.58	107.50	103.45
3	D	602	ADP	C5-N7-C8	2.57	107.49	103.45
3	H	602	ADP	C5-N7-C8	2.56	107.47	103.45
3	C	602	ADP	C4-N9-C8	2.55	108.42	105.74
3	B	602	ADP	C6-C5-N7	2.45	136.81	132.09
3	A	602	ADP	O4'-C1'-N9	2.39	112.68	108.09
3	B	602	ADP	N9-C8-N7	-2.31	110.65	113.94
3	H	602	ADP	O4'-C1'-N9	2.31	112.53	108.09
3	D	602	ADP	C6-C5-N7	2.31	136.54	132.09
3	E	602	ADP	O4'-C1'-N9	2.27	112.45	108.09
3	D	602	ADP	O4'-C1'-N9	2.26	112.42	108.09
3	H	602	ADP	C6-C5-N7	2.22	136.37	132.09
3	A	602	ADP	C6-C5-N7	2.21	136.35	132.09
3	F	602	ADP	C6-C5-N7	2.20	136.34	132.09
3	E	602	ADP	C6-C5-N7	2.20	136.33	132.09
3	G	602	ADP	C6-C5-N7	2.20	136.32	132.09
3	F	602	ADP	O4'-C1'-N9	2.19	112.29	108.09
3	C	602	ADP	C6-C5-N7	2.17	136.27	132.09
3	B	602	ADP	O2A-PA-O3A	2.16	113.10	107.27
3	D	602	ADP	N9-C8-N7	-2.15	110.89	113.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	G	602	ADP	O4'-C1'-N9	2.14	112.20	108.09
3	A	602	ADP	N9-C8-N7	-2.14	110.90	113.94
3	E	602	ADP	N9-C8-N7	-2.12	110.93	113.94
3	C	602	ADP	O4'-C1'-N9	2.11	112.14	108.09
3	F	602	ADP	N9-C8-N7	-2.10	110.95	113.94
3	H	602	ADP	N9-C8-N7	-2.08	110.98	113.94
3	G	602	ADP	N9-C8-N7	-2.06	111.02	113.94

There are no chirality outliers.

All (68) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	601	TPP	PA-O3A-PB-O2B
2	E	601	TPP	C7-O7-PA-O1A
3	A	602	ADP	C5'-O5'-PA-O1A
3	A	602	ADP	C5'-O5'-PA-O3A
3	A	602	ADP	O4'-C4'-C5'-O5'
3	B	602	ADP	PA-O3A-PB-O2B
3	C	602	ADP	C5'-O5'-PA-O1A
3	C	602	ADP	C5'-O5'-PA-O3A
3	C	602	ADP	C3'-C4'-C5'-O5'
3	D	602	ADP	PA-O3A-PB-O2B
3	D	602	ADP	C5'-O5'-PA-O1A
3	D	602	ADP	C5'-O5'-PA-O3A
3	D	602	ADP	O4'-C4'-C5'-O5'
3	E	602	ADP	PA-O3A-PB-O2B
3	E	602	ADP	C5'-O5'-PA-O2A
3	E	602	ADP	C5'-O5'-PA-O3A
3	E	602	ADP	C3'-C4'-C5'-O5'
3	F	602	ADP	C5'-O5'-PA-O1A
3	G	602	ADP	PA-O3A-PB-O2B
3	G	602	ADP	C5'-O5'-PA-O1A
3	G	602	ADP	C5'-O5'-PA-O3A
3	G	602	ADP	O4'-C4'-C5'-O5'
3	H	602	ADP	PA-O3A-PB-O2B
3	H	602	ADP	C5'-O5'-PA-O1A
3	H	602	ADP	C5'-O5'-PA-O2A
3	H	602	ADP	C5'-O5'-PA-O3A
3	H	602	ADP	O4'-C4'-C5'-O5'
4	A	603	GOL	C1-C2-C3-O3
3	A	602	ADP	C3'-C4'-C5'-O5'
3	C	602	ADP	O4'-C4'-C5'-O5'

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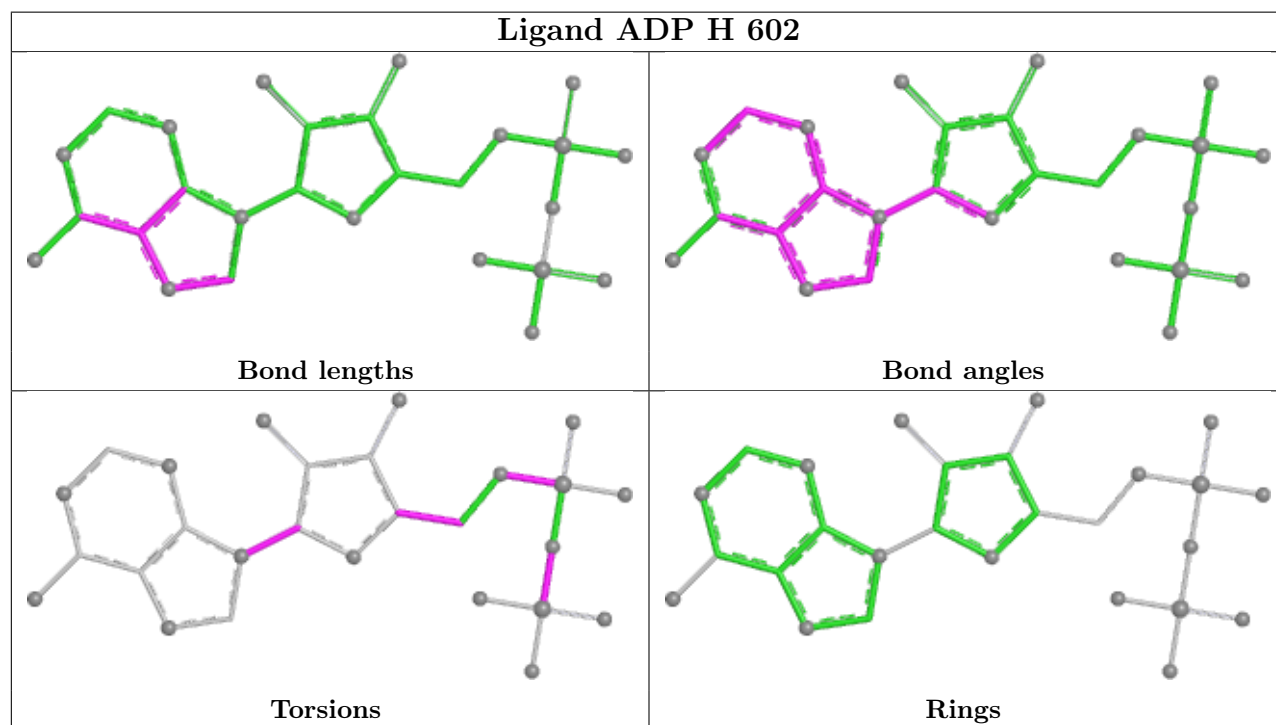
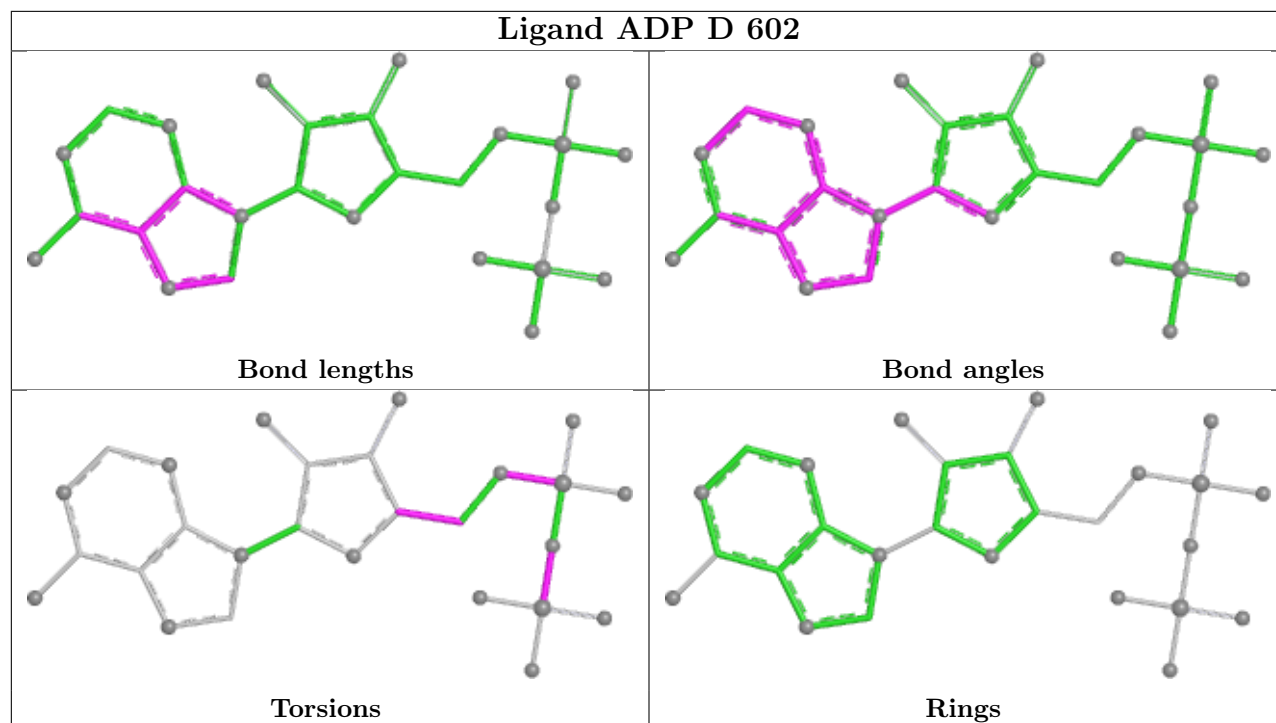
Mol	Chain	Res	Type	Atoms
3	D	602	ADP	C3'-C4'-C5'-O5'
3	G	602	ADP	C3'-C4'-C5'-O5'
3	H	602	ADP	C3'-C4'-C5'-O5'
3	B	602	ADP	O4'-C4'-C5'-O5'
3	E	602	ADP	O4'-C4'-C5'-O5'
3	F	602	ADP	O4'-C4'-C5'-O5'
3	B	602	ADP	C3'-C4'-C5'-O5'
3	F	602	ADP	C3'-C4'-C5'-O5'
4	A	603	GOL	O2-C2-C3-O3
3	B	602	ADP	PB-O3A-PA-O1A
2	G	601	TPP	C4'-C5'-C7'-N3
2	B	601	TPP	PB-O3A-PA-O7
3	E	602	ADP	PA-O3A-PB-O3B
3	F	602	ADP	PA-O3A-PB-O2B
3	F	602	ADP	PA-O3A-PB-O3B
2	E	601	TPP	C5-C6-C7-O7
2	G	601	TPP	C5-C6-C7-O7
3	A	602	ADP	C5'-O5'-PA-O2A
3	C	602	ADP	C5'-O5'-PA-O2A
3	D	602	ADP	C5'-O5'-PA-O2A
3	F	602	ADP	C5'-O5'-PA-O2A
3	F	602	ADP	C5'-O5'-PA-O3A
3	G	602	ADP	C5'-O5'-PA-O2A
3	B	602	ADP	PB-O3A-PA-O2A
2	G	601	TPP	C6'-C5'-C7'-N3
2	E	601	TPP	S1-C5-C6-C7
3	D	602	ADP	PA-O3A-PB-O1B
3	E	602	ADP	PA-O3A-PB-O1B
2	C	601	TPP	PA-O3A-PB-O3B
3	G	602	ADP	PA-O3A-PB-O3B
3	H	602	ADP	PA-O3A-PB-O3B
2	B	601	TPP	C5-C6-C7-O7
2	H	601	TPP	C5-C6-C7-O7
3	B	602	ADP	C2'-C1'-N9-C8
3	H	602	ADP	C2'-C1'-N9-C8
3	B	602	ADP	PA-O3A-PB-O1B
3	H	602	ADP	PA-O3A-PB-O1B
3	F	602	ADP	C2'-C1'-N9-C8

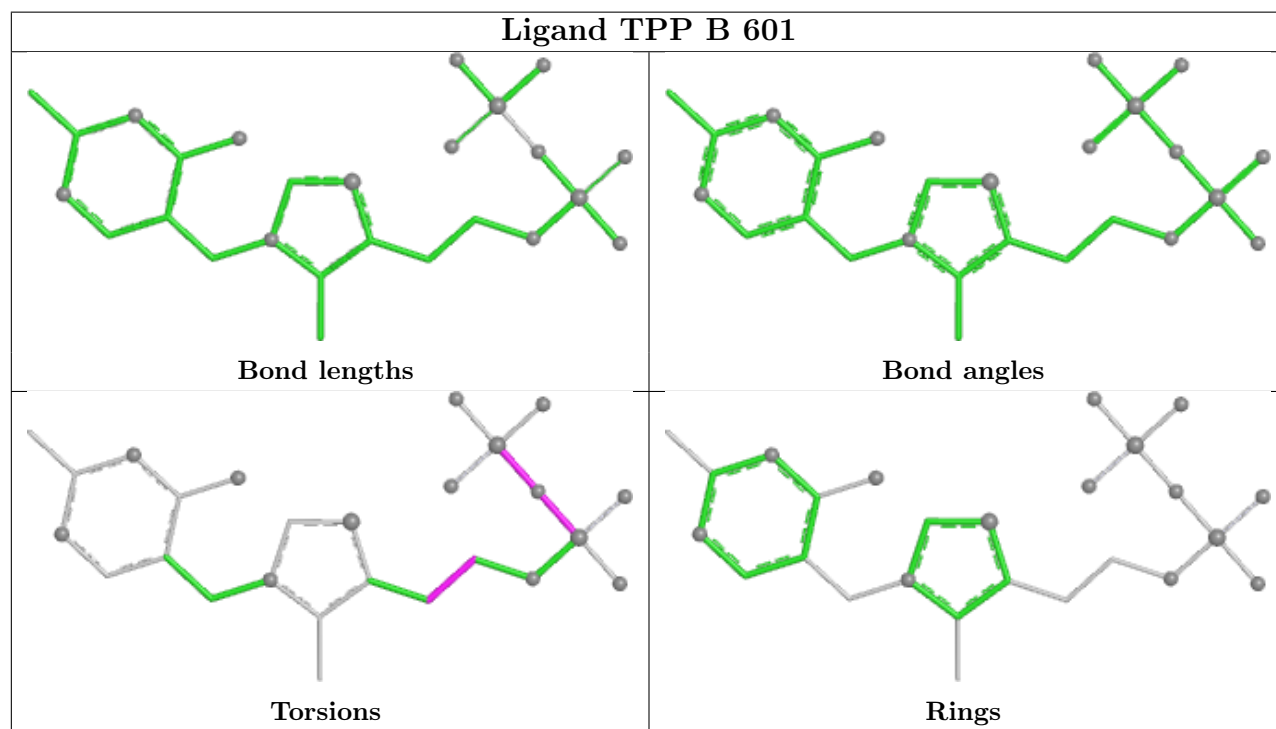
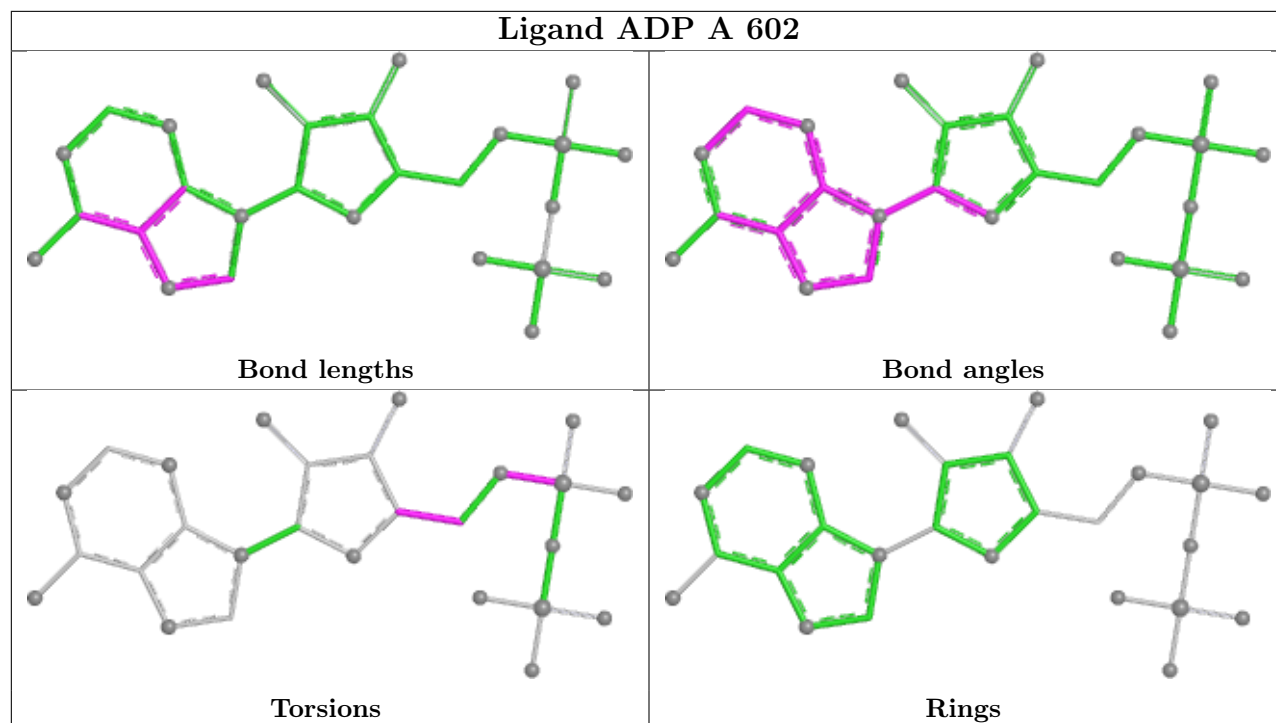
There are no ring outliers.

15 monomers are involved in 21 short contacts:

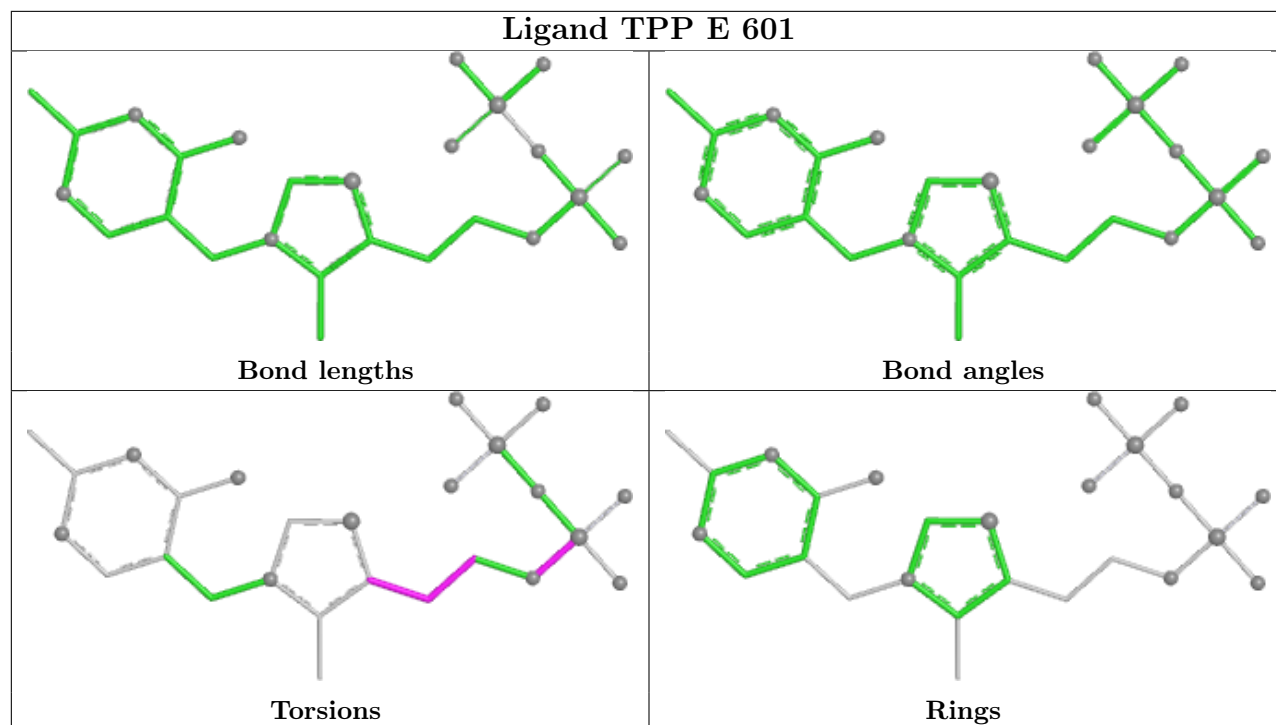
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	G	603	EDO	1	0
3	D	602	ADP	1	0
3	A	602	ADP	1	0
2	B	601	TPP	1	0
2	E	601	TPP	3	0
2	D	601	TPP	1	0
4	A	603	GOL	2	0
2	F	601	TPP	1	0
2	C	601	TPP	1	0
2	G	601	TPP	3	0
3	C	602	ADP	1	0
3	F	602	ADP	1	0
2	H	601	TPP	2	0
4	B	603	GOL	1	0
2	A	601	TPP	1	0

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.

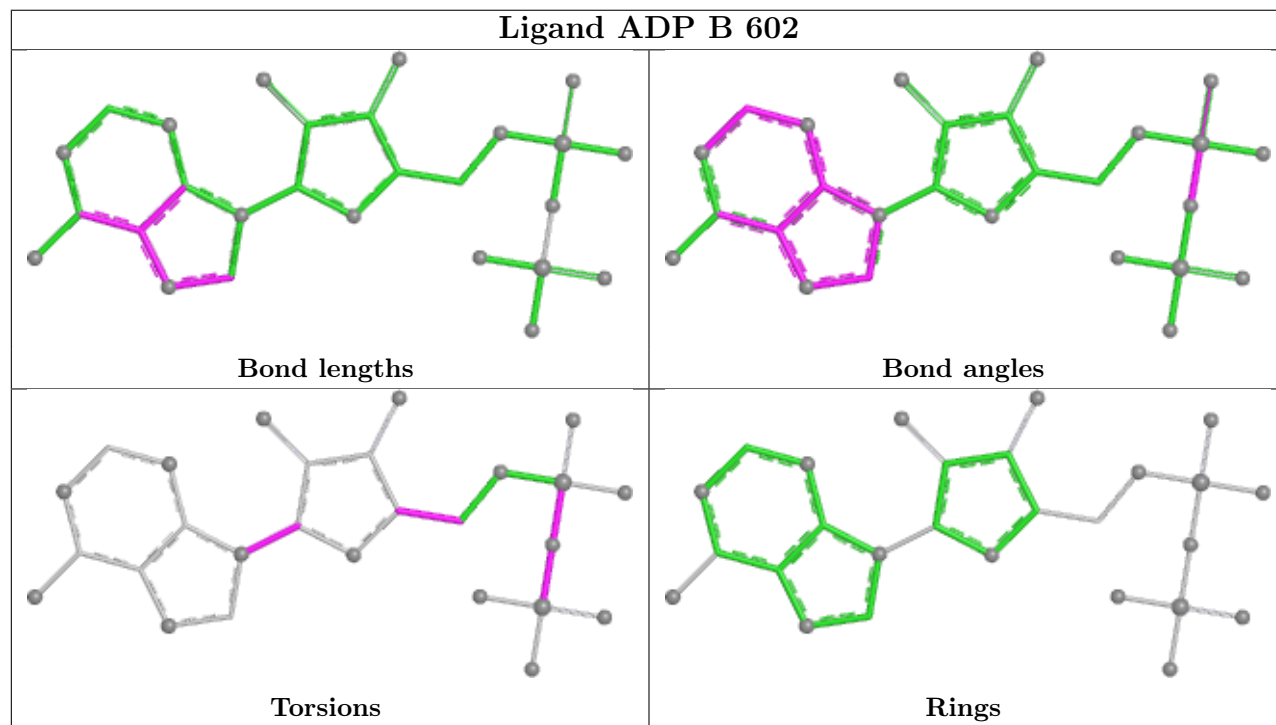




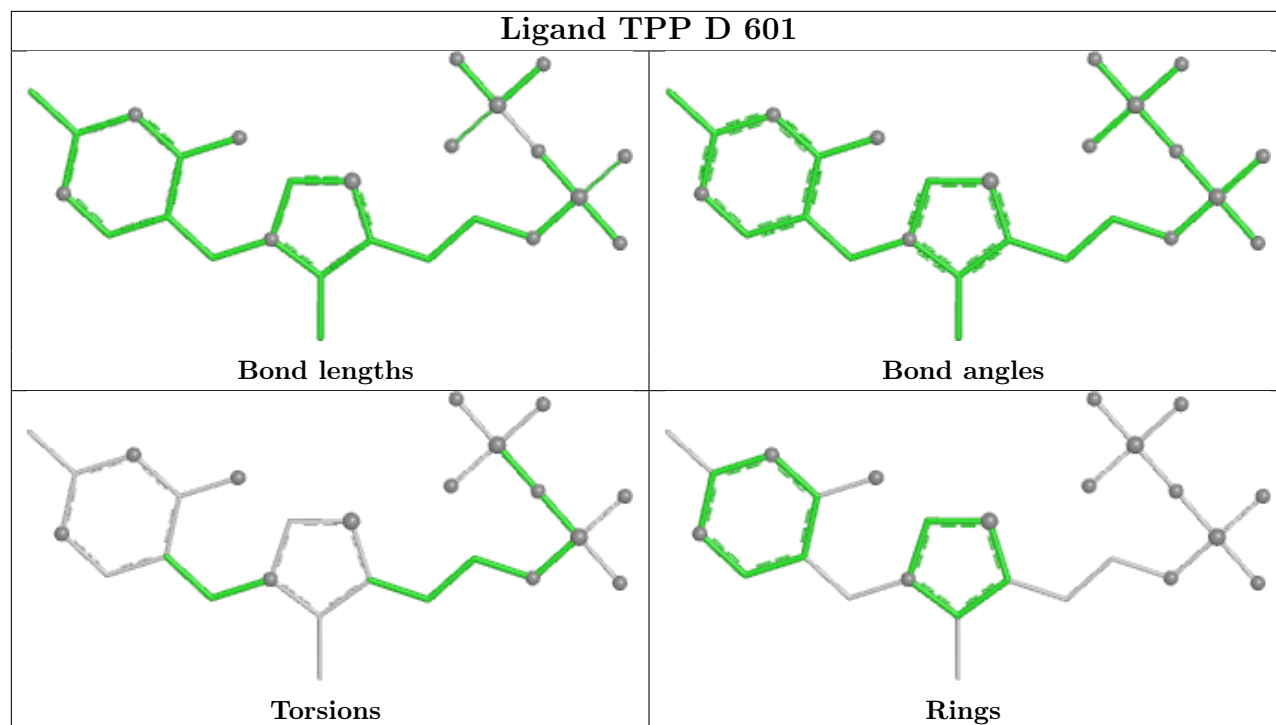
Ligand TPP E 601



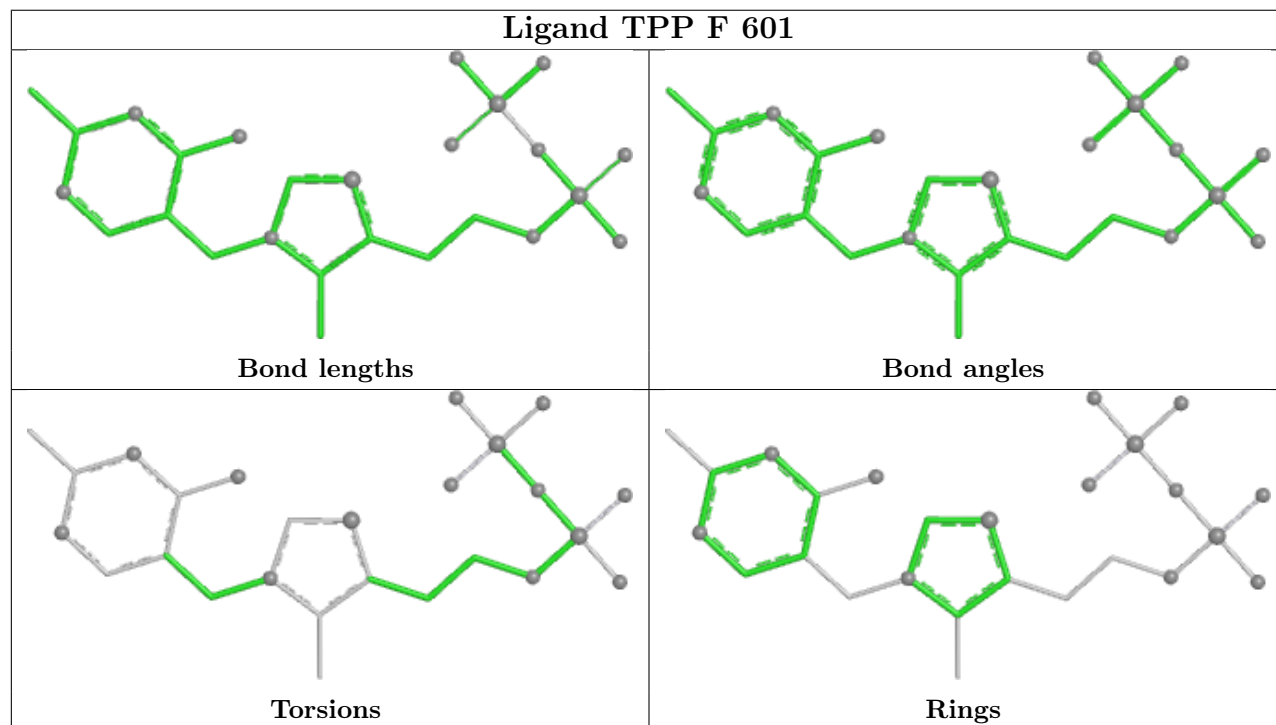
Ligand ADP B 602



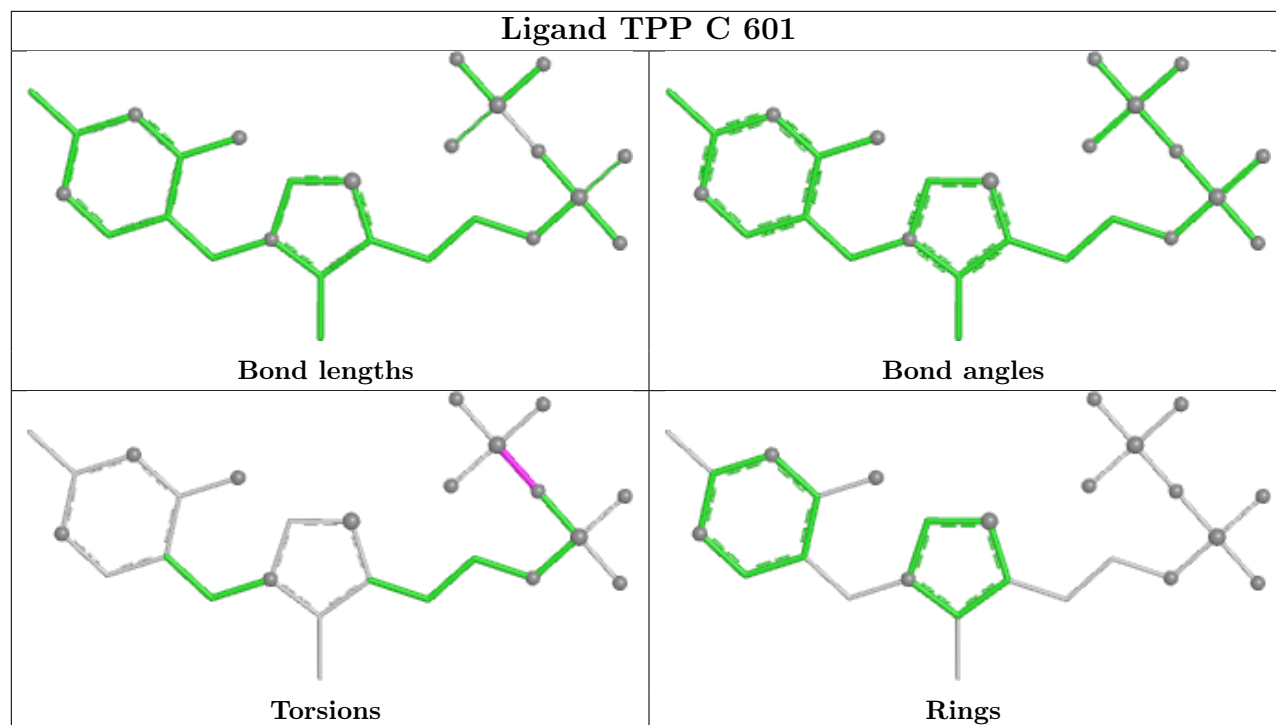
Ligand TPP D 601



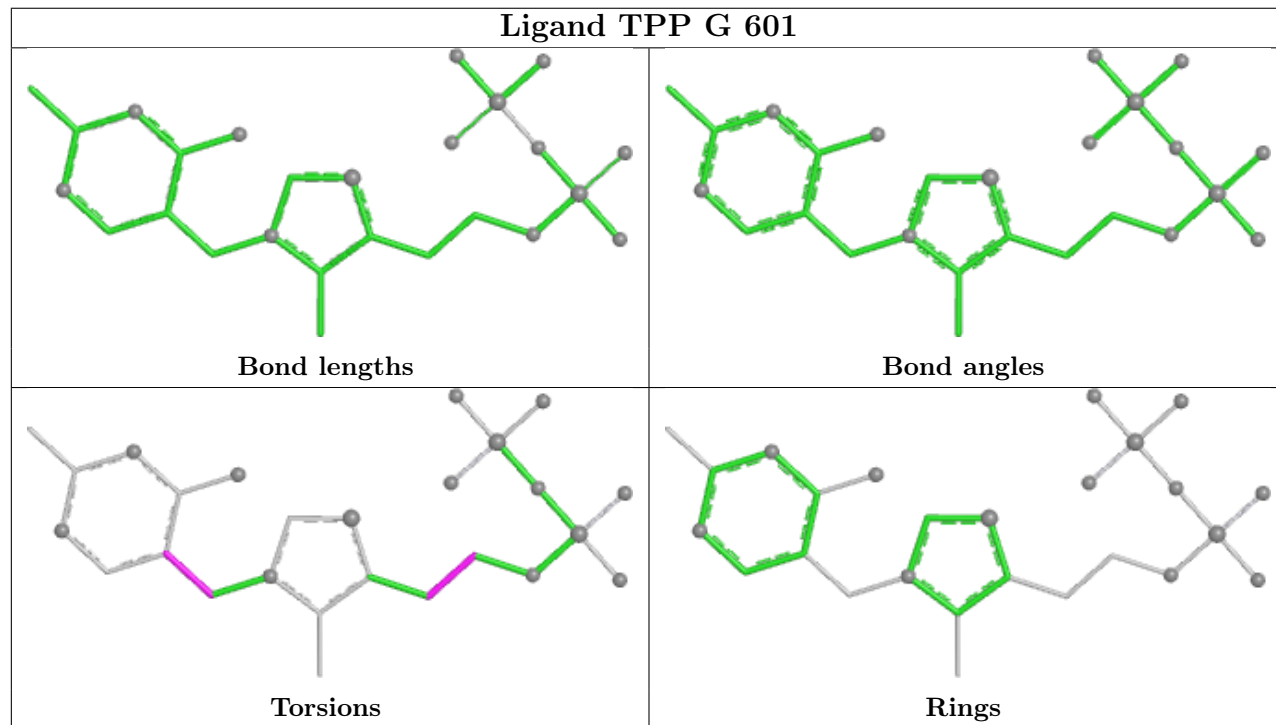
Ligand TPP F 601

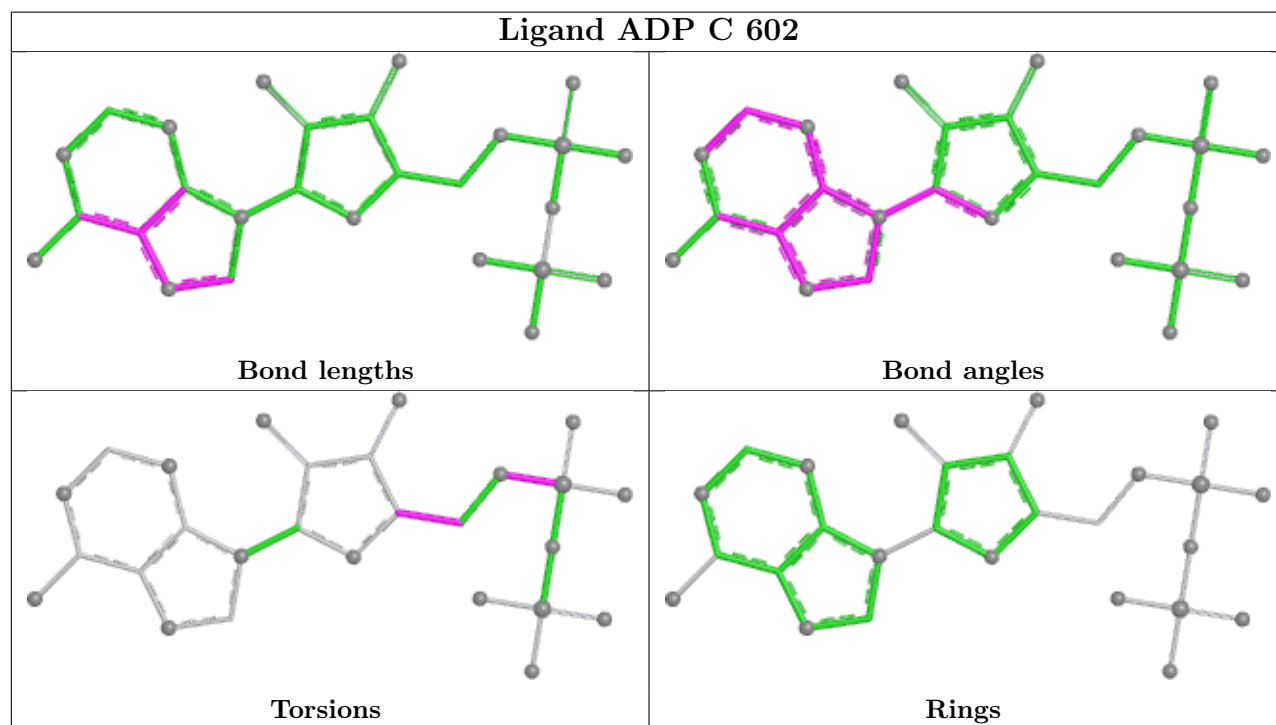
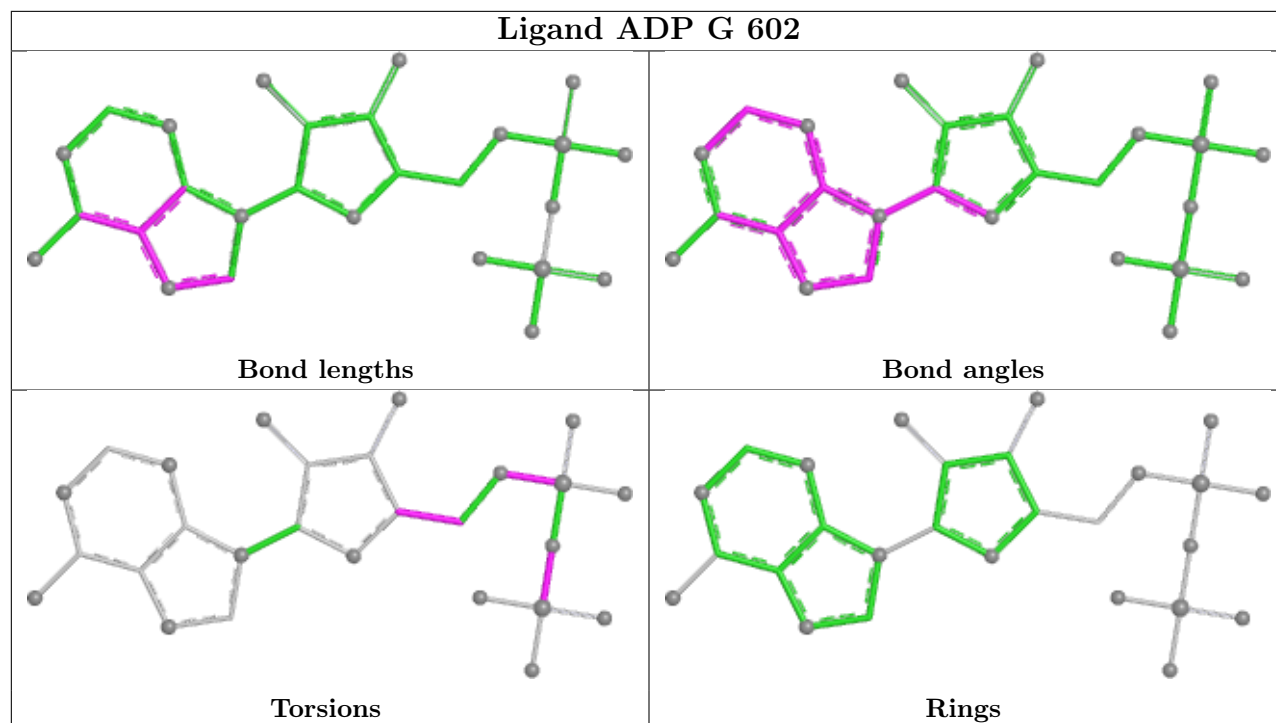


Ligand TPP C 601

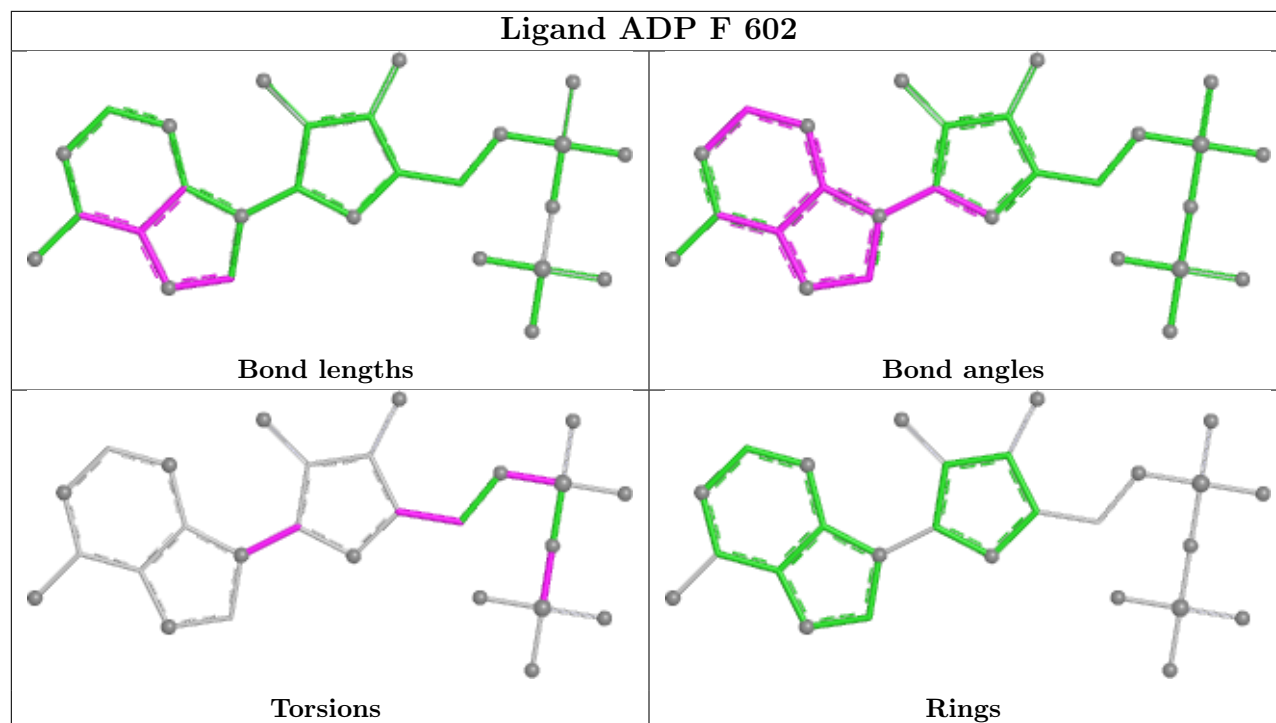


Ligand TPP G 601

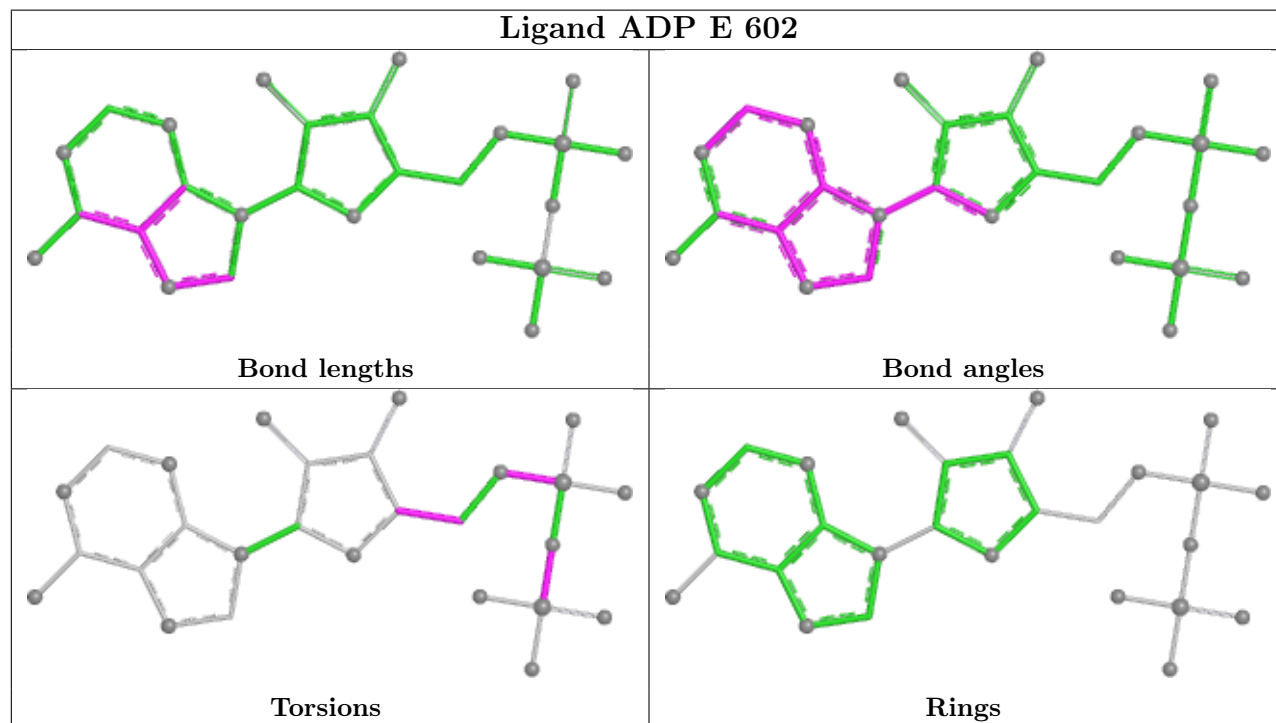


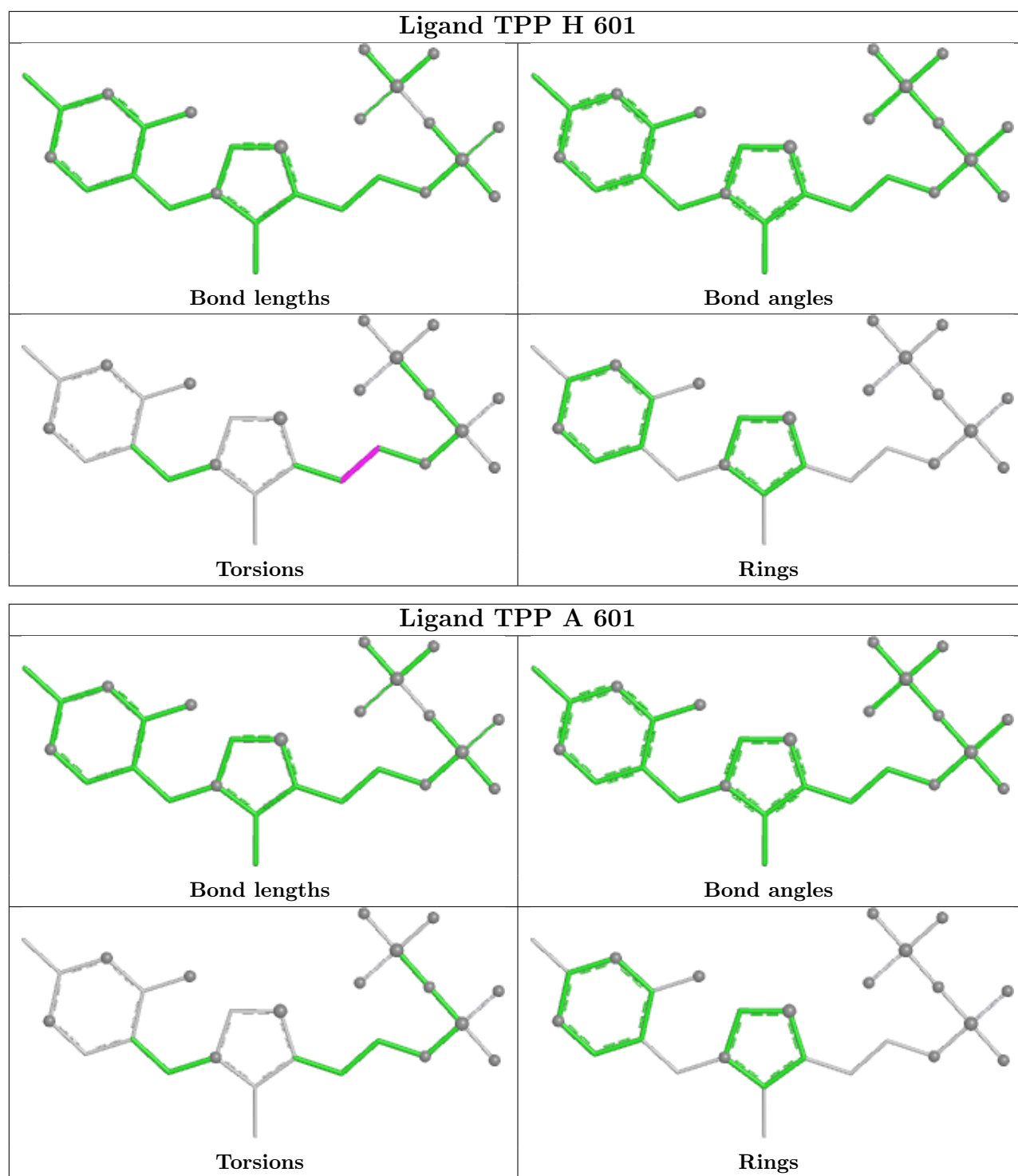


Ligand ADP F 602



Ligand ADP E 602





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	542/554 (97%)	-0.44	0	100	100	47, 62, 89, 124	1 (0%)
1	B	539/554 (97%)	-0.38	0	100	100	47, 69, 106, 137	0
1	C	542/554 (97%)	-0.22	0	100	100	54, 85, 119, 153	1 (0%)
1	D	546/554 (98%)	-0.25	1 (0%)	91	90	47, 80, 109, 146	1 (0%)
1	E	541/554 (97%)	-0.31	1 (0%)	91	90	51, 76, 108, 140	0
1	F	542/554 (97%)	-0.23	0	100	100	49, 79, 118, 144	0
1	G	540/554 (97%)	-0.23	0	100	100	55, 84, 117, 158	0
1	H	538/554 (97%)	-0.20	1 (0%)	91	90	56, 85, 121, 149	0
All	All	4330/4432 (97%)	-0.28	3 (0%)	92	91	47, 77, 114, 158	3 (0%)

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	35[A]	MET	2.7
1	H	186	MET	2.6
1	E	542	LYS	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 ⓘ

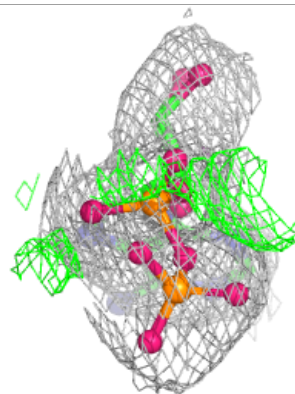
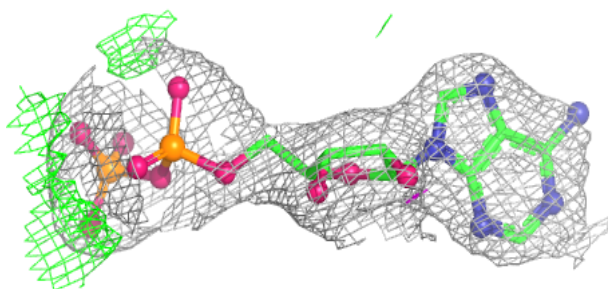
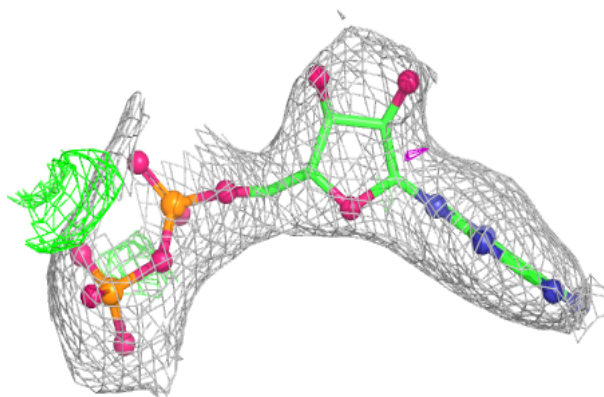
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	EDO	H	603	4/4	0.62	0.16	62,63,64,66	0
6	EDO	G	603	4/4	0.75	0.13	65,65,65,68	0
6	EDO	F	603	4/4	0.81	0.10	61,61,62,62	0
4	GOL	A	603	6/6	0.84	0.12	51,52,55,55	0
6	EDO	C	603	4/4	0.87	0.10	65,66,66,69	0
5	MG	H	604	1/1	0.88	0.13	86,86,86,86	0
4	GOL	B	603	6/6	0.90	0.08	59,61,62,66	0
3	ADP	G	602	27/27	0.92	0.07	72,74,77,79	0
3	ADP	E	602	27/27	0.92	0.08	66,68,72,74	0
3	ADP	F	602	27/27	0.93	0.08	65,68,73,75	0
3	ADP	C	602	27/27	0.93	0.07	71,75,81,83	0
3	ADP	H	602	27/27	0.93	0.08	70,75,81,83	0
3	ADP	D	602	27/27	0.93	0.08	68,72,79,80	0
3	ADP	B	602	27/27	0.93	0.08	62,65,67,67	0
2	TPP	F	601	26/26	0.94	0.08	66,75,86,86	0
5	MG	D	603	1/1	0.95	0.09	72,72,72,72	0
3	ADP	A	602	27/27	0.95	0.07	55,58,62,64	0
2	TPP	E	601	26/26	0.95	0.09	59,65,74,76	0
2	TPP	C	601	26/26	0.95	0.08	63,69,81,84	0
2	TPP	G	601	26/26	0.95	0.08	67,74,82,85	0
2	TPP	H	601	26/26	0.95	0.08	63,69,89,96	0
2	TPP	A	601	26/26	0.96	0.06	48,53,57,64	0
2	TPP	D	601	26/26	0.96	0.07	69,73,77,79	0
5	MG	B	604	1/1	0.96	0.09	58,58,58,58	0
5	MG	F	604	1/1	0.97	0.08	82,82,82,82	0
5	MG	C	604	1/1	0.97	0.06	75,75,75,75	0
2	TPP	B	601	26/26	0.97	0.07	53,58,64,66	0
5	MG	G	604	1/1	0.98	0.07	80,80,80,80	0
5	MG	A	604	1/1	0.98	0.08	54,54,54,54	0
5	MG	E	603	1/1	0.99	0.05	70,70,70,70	0

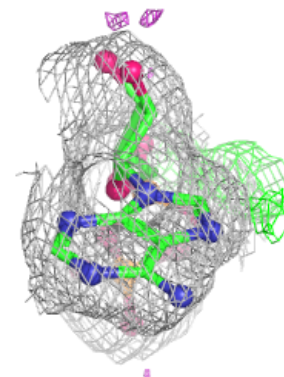
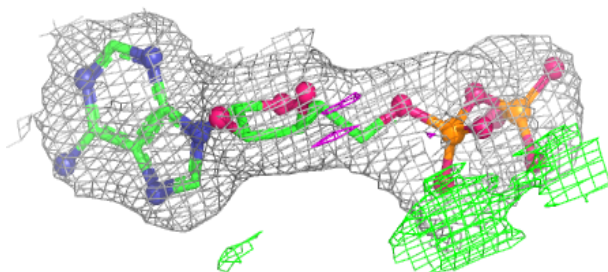
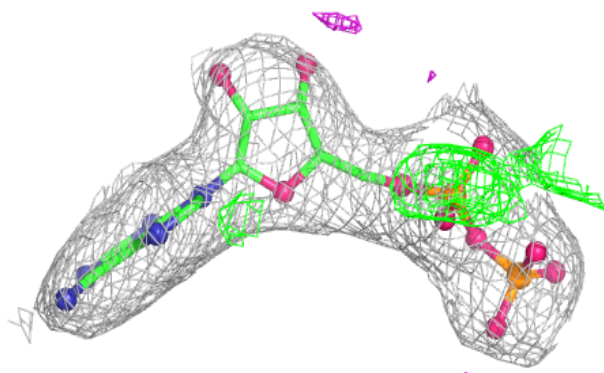
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around ADP G 602:

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

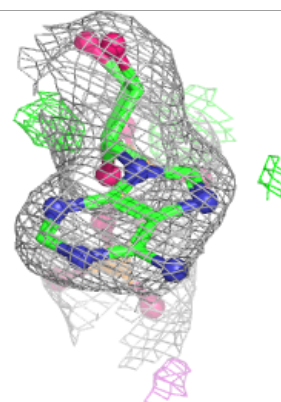
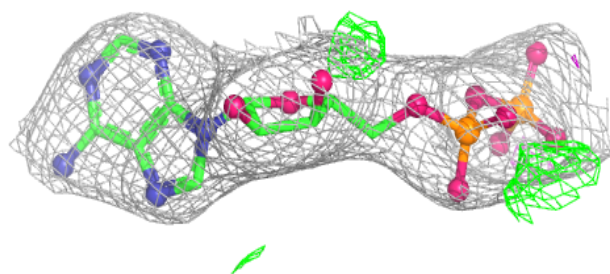
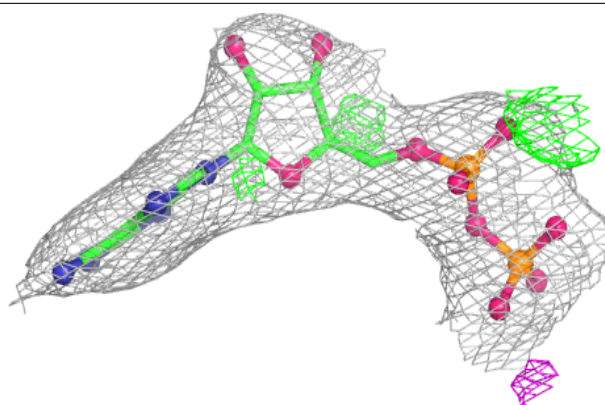
**Electron density around ADP E 602:**

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and green (positive)

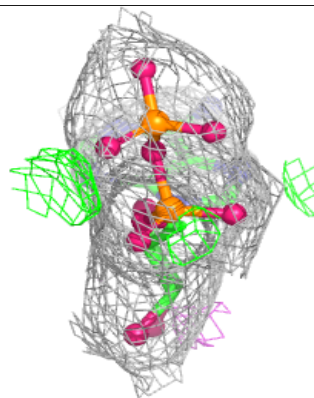
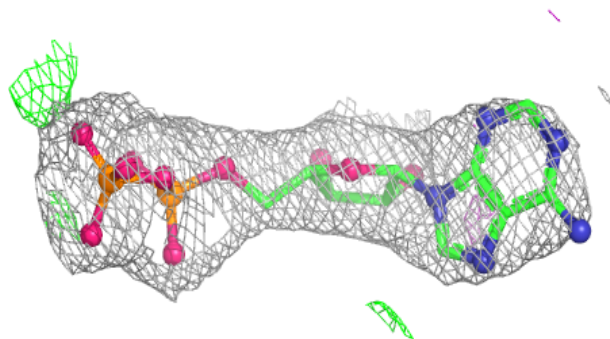
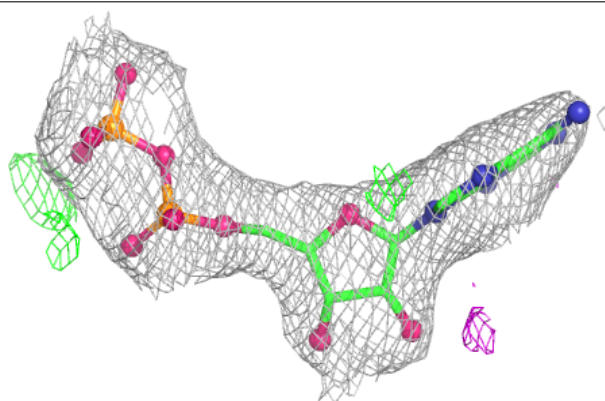


Electron density around ADP F 602:

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

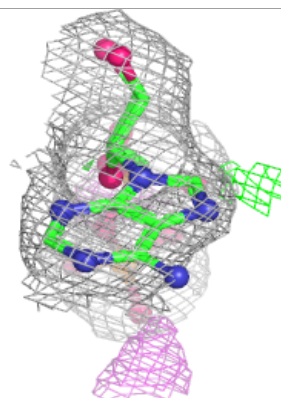
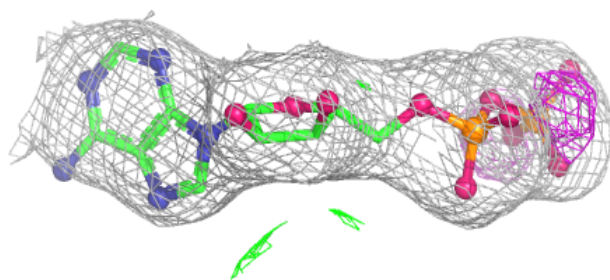
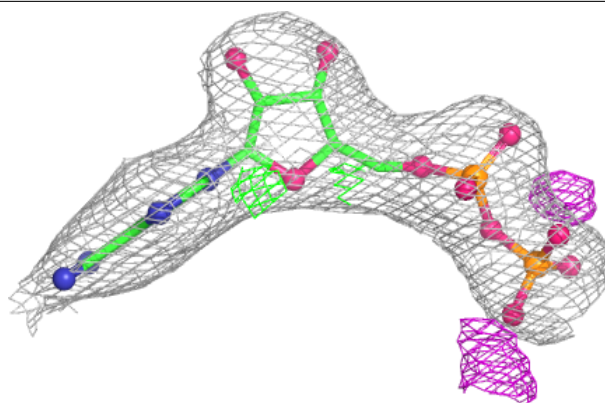
**Electron density around ADP C 602:**

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and green (positive)

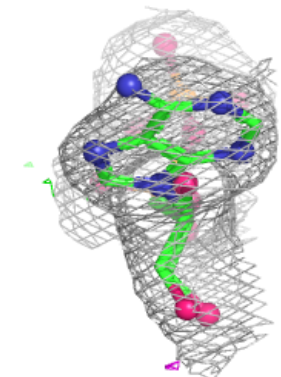
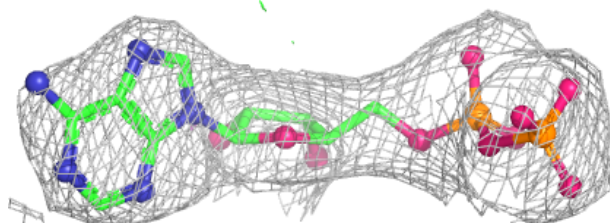
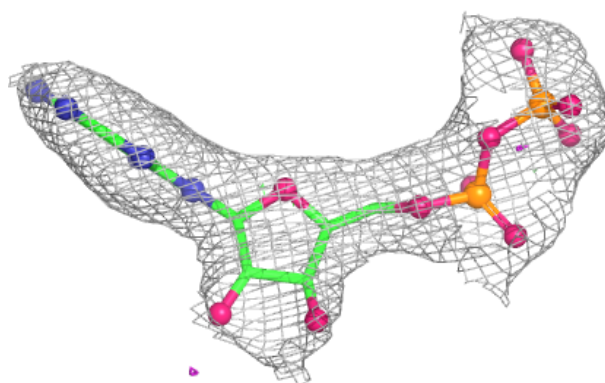


Electron density around ADP H 602:

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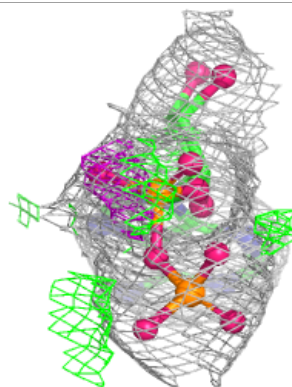
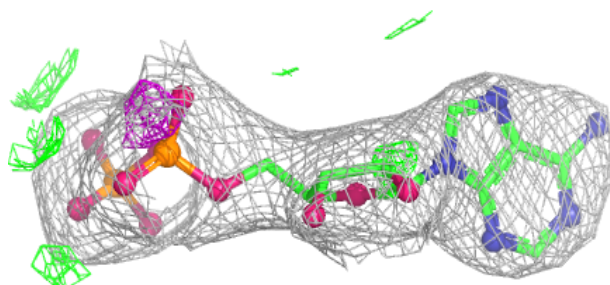
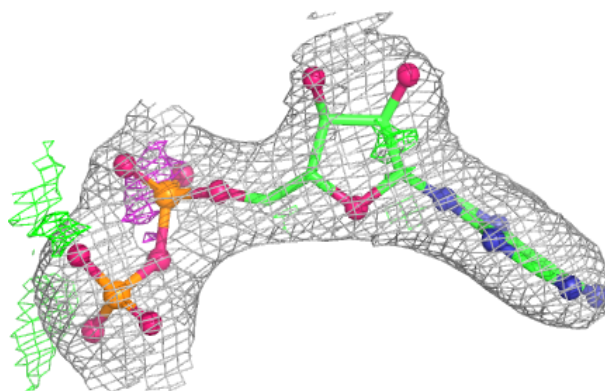
**Electron density around ADP D 602:**

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

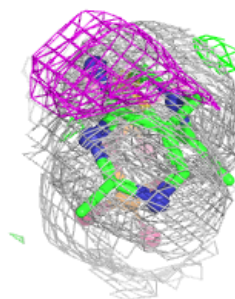
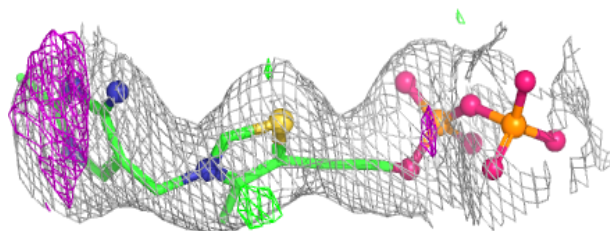
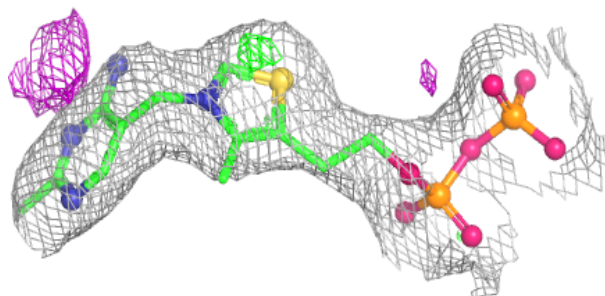


Electron density around ADP B 602:

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and green (positive)

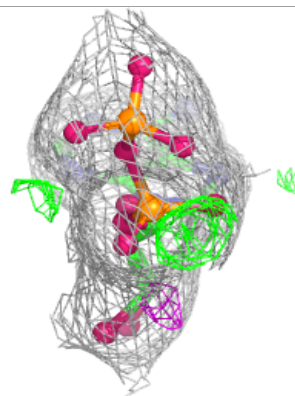
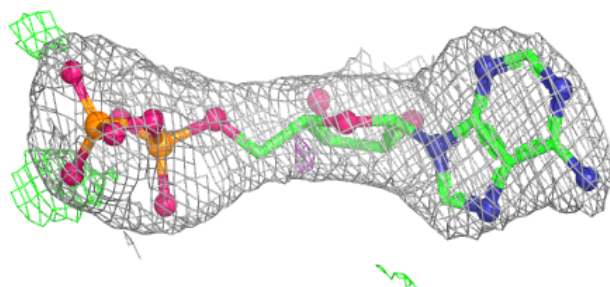
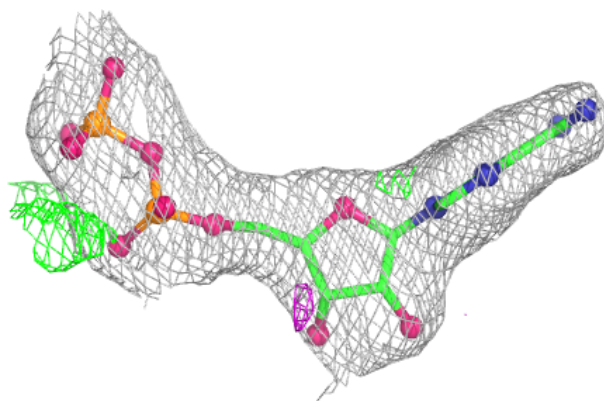
**Electron density around TPP F 601:**

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

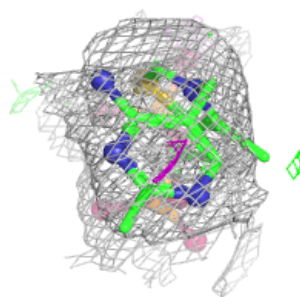
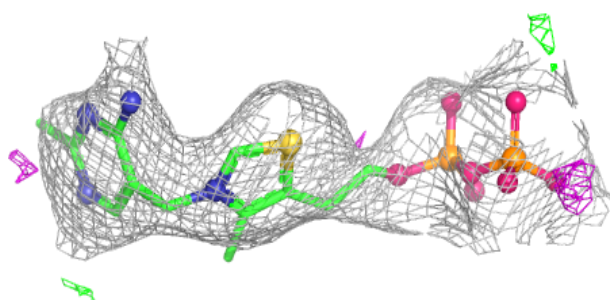
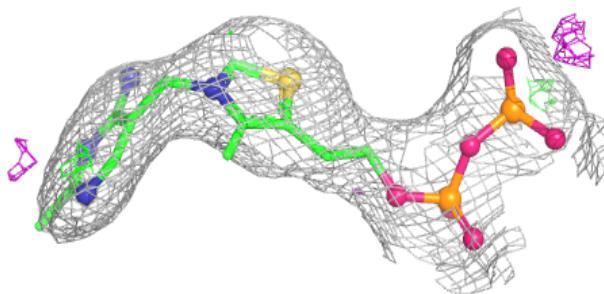


Electron density around ADP A 602:

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

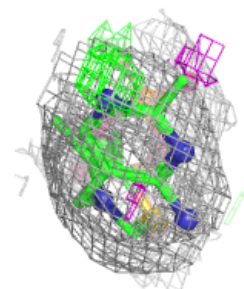
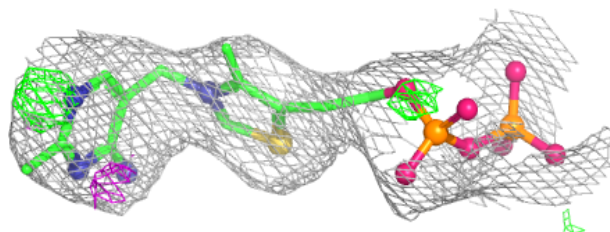
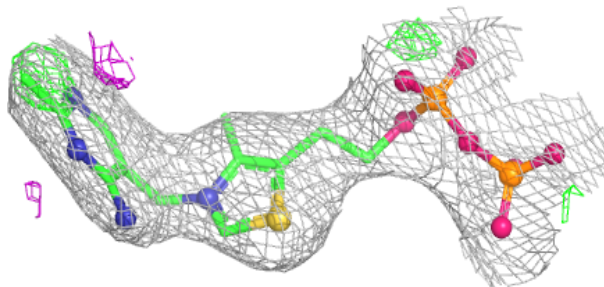
**Electron density around TPP E 601:**

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

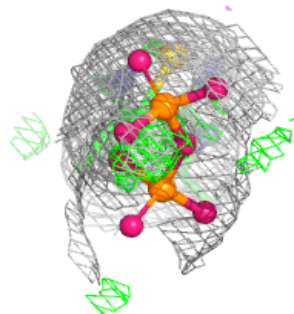
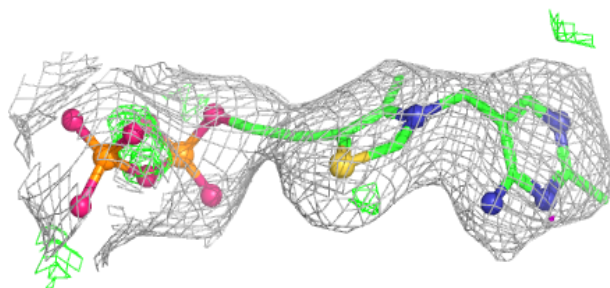
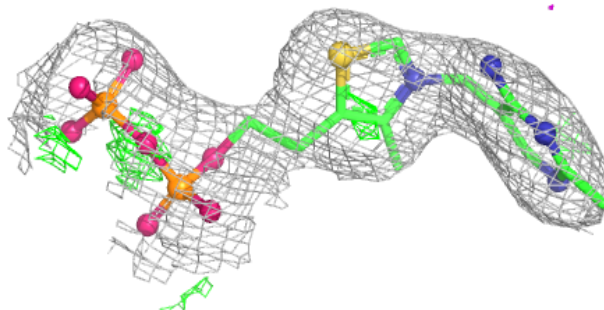


Electron density around TPP C 601:

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

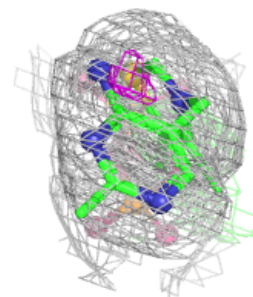
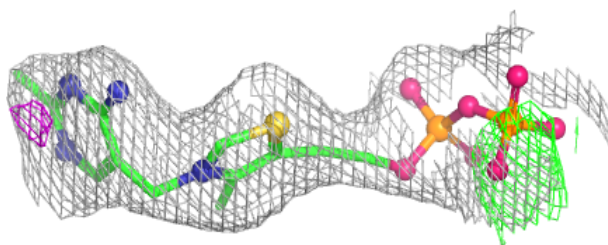
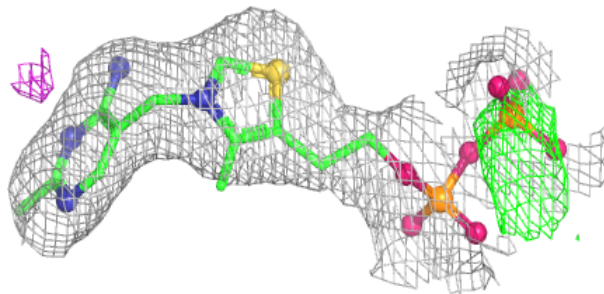
**Electron density around TPP G 601:**

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

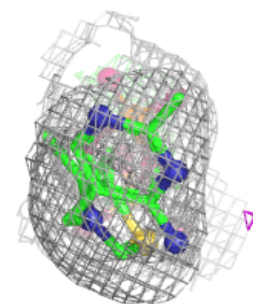
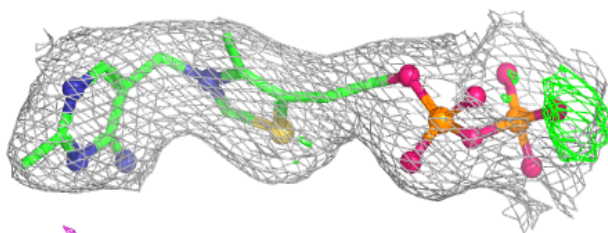
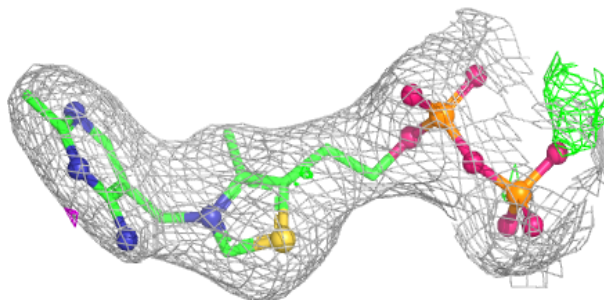


Electron density around TPP H 601:

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

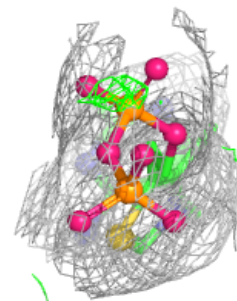
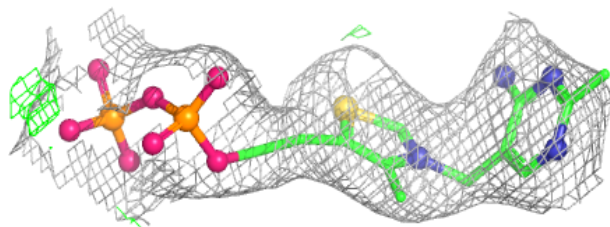
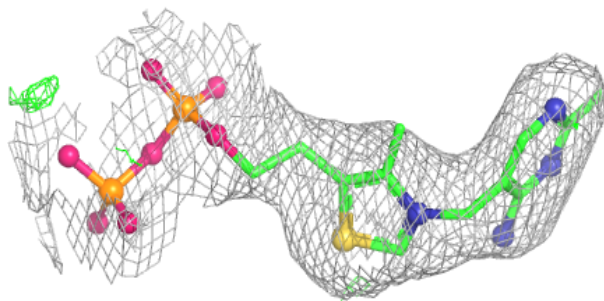
**Electron density around TPP A 601:**

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

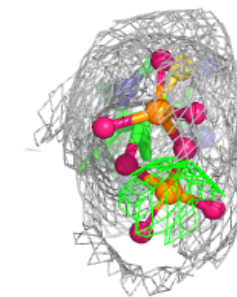
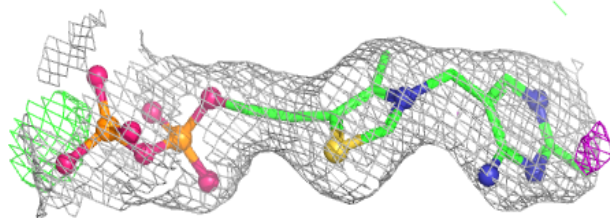
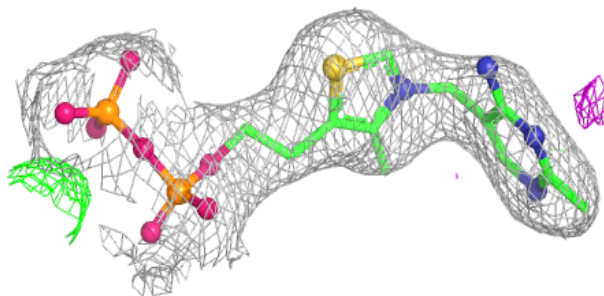


Electron density around TPP D 601:

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

**Electron density around TPP 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)



6.5 Other polymers [i](#)

There are no such residues in this entry.