



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:20 AM UTC

PDB ID : 9EV6 / pdb_00009ev6
Title : Corynebacterium glutamicum pyruvate:quinone oxidoreductase (PQO), C-terminal truncated construct
Authors : Da Silva Lameira, C.; Muenssinger, S.; Yang, L.; Eikmanns, B.J.; Bellinzoni, M.
Deposited on : 2024-03-28
Resolution : 1.89 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

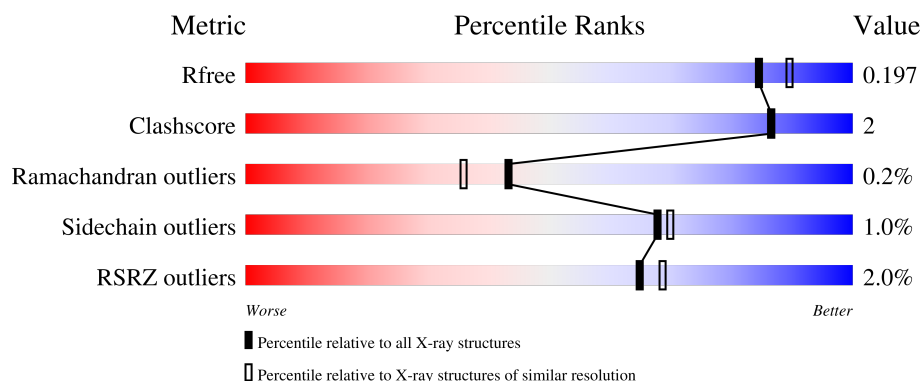
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.89 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	564	2% 90% 5% 5%
1	B	564	% 91% • 5%
1	C	564	% 89% 5% 5%
1	D	564	4% 90% 5% 5%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 17270 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.

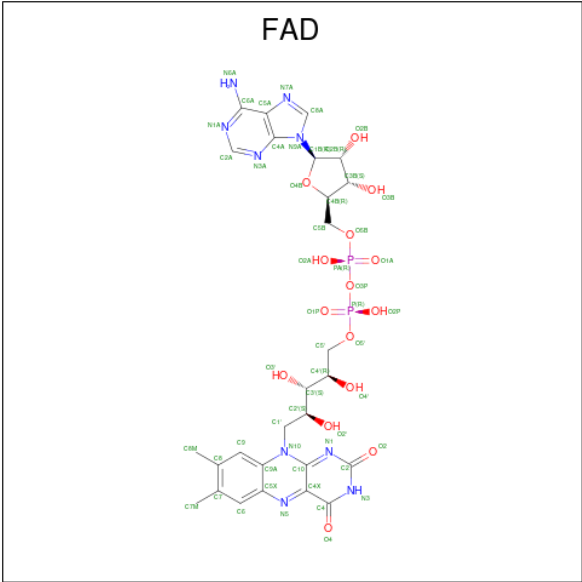
- Molecule 1 is a protein called Thiamine pyrophosphate-requiring enzymes [acetolactate synthase, pyruvate dehydrogenase (Cytochrome), glyoxylate carboligase, phosphonopyruvate decarboxylase].

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	535	Total	C	N	O	S	0	0	0
			3974	2496	690	770	18			
1	B	536	Total	C	N	O	S	0	0	0
			3997	2510	693	776	18			
1	C	535	Total	C	N	O	S	0	0	0
			4000	2512	693	777	18			
1	D	537	Total	C	N	O	S	0	0	0
			3974	2498	690	768	18			

There are 8 discrepancies between the modelled and reference sequences:

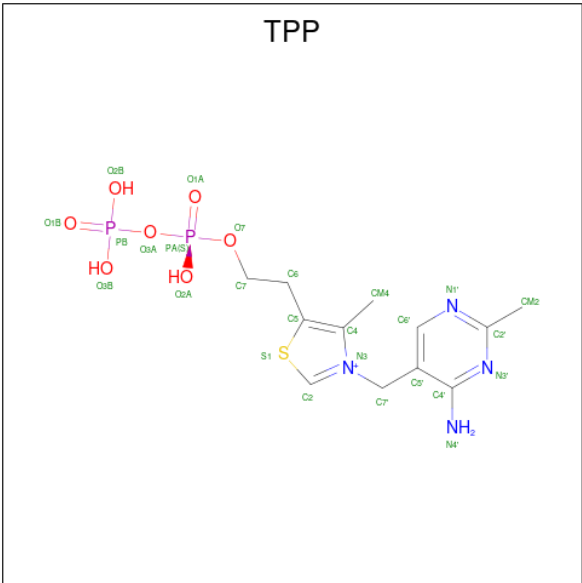
Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP Q8NMG5
A	0	HIS	-	expression tag	UNP Q8NMG5
B	-1	GLY	-	expression tag	UNP Q8NMG5
B	0	HIS	-	expression tag	UNP Q8NMG5
C	-1	GLY	-	expression tag	UNP Q8NMG5
C	0	HIS	-	expression tag	UNP Q8NMG5
D	-1	GLY	-	expression tag	UNP Q8NMG5
D	0	HIS	-	expression tag	UNP Q8NMG5

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	C	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	D	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 3 is THIAMINE DIPHOSPHATE (CCD ID: TPP) (formula: C₁₂H₁₉N₄O₇P₂S) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Mg	0	0
			1	1		
4	B	1	Total	Mg	0	0
			1	1		
4	C	1	Total	Mg	0	0
			1	1		
4	D	1	Total	Mg	0	0
			1	1		

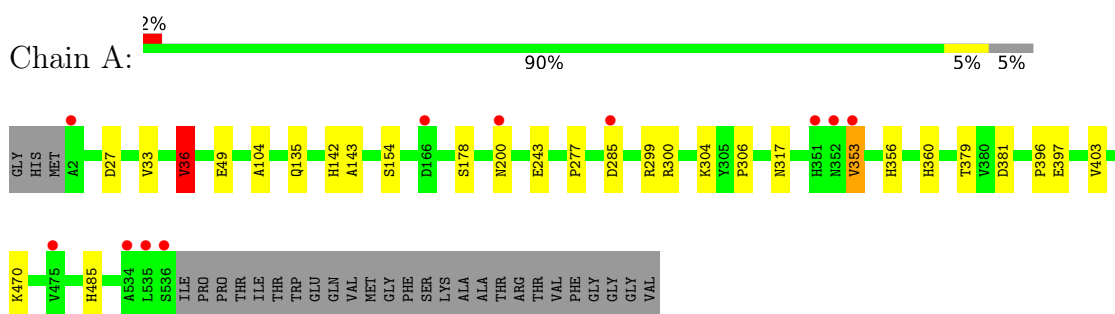
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	262	Total	O	0	0
			262	262		
5	B	276	Total	O	0	0
			276	276		
5	C	238	Total	O	0	0
			238	238		
5	D	229	Total	O	0	0
			229	229		

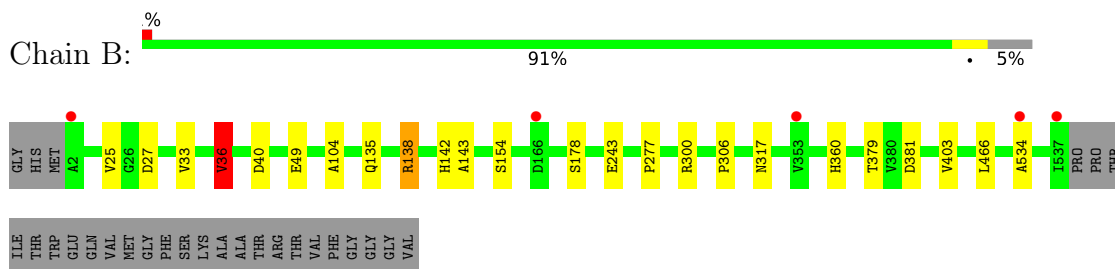
3 Residue-property plots

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.

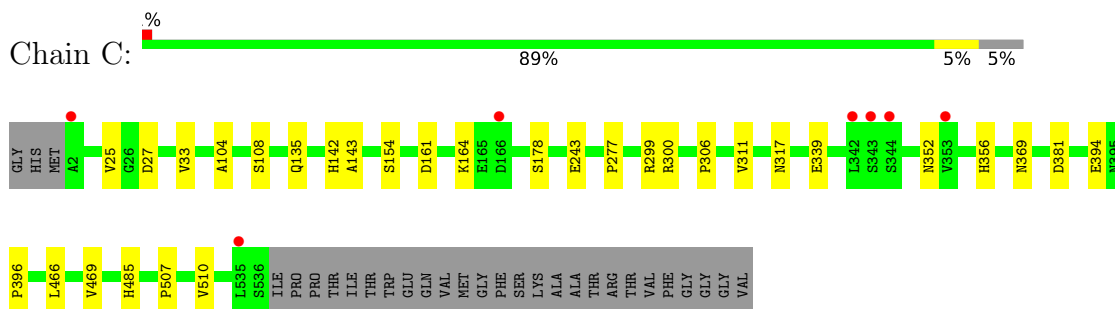
- Molecule 1: Thiamine pyrophosphate-requiring enzymes [acetolactate synthase, pyruvate dehydrogenase (Cytochrome), glyoxylate carboligase, phosphonopyruvate decarboxylase]



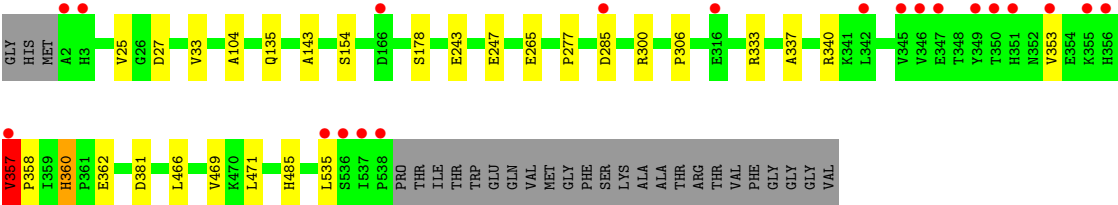
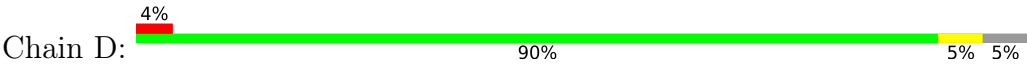
- Molecule 1: Thiamine pyrophosphate-requiring enzymes [acetolactate synthase, pyruvate dehydrogenase (Cytochrome), glyoxylate carboligase, phosphonopyruvate decarboxylase]



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4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	100.55Å 98.48Å 110.52Å 90.00° 90.84° 90.00°	Depositor
Resolution (Å)	37.46 – 1.89 37.46 – 1.89	Depositor EDS
% Data completeness (in resolution range)	77.8 (37.46-1.89) 77.7 (37.46-1.89)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.59 (at 1.89Å)	Xtriage
Refinement program	BUSTER 2.10.4	Depositor
R, R_{free}	0.168 , 0.197 (Not available) , 0.197	Depositor DCC
R_{free} test set	6391 reflections (4.78%)	wwPDB-VP
Wilson B-factor (Å ²)	30.0	Xtriage
Anisotropy	0.100	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 46.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.000 for k,h,-l 0.000 for -k,-h,-l 0.019 for h,-k,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	17270	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.79% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: TPP, MG, FAD

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.75	0/4050	1.00	4/5512 (0.1%)
1	B	0.76	0/4073	1.00	4/5541 (0.1%)
1	C	0.72	0/4076	0.98	4/5543 (0.1%)
1	D	0.73	0/4050	1.00	3/5514 (0.1%)
All	All	0.74	0/16249	1.00	15/22110 (0.1%)

There are no bond length outliers.

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	200	ASN	CA-CB-CG	-7.04	105.56	112.60
1	B	36	VAL	N-CA-CB	6.58	119.49	110.54
1	B	40	ASP	CA-CB-CG	5.93	118.53	112.60
1	B	317	ASN	CA-CB-CG	5.92	118.52	112.60
1	A	36	VAL	N-CA-CB	5.71	119.15	110.58
1	C	317	ASN	CA-CB-CG	5.60	118.20	112.60
1	D	466	LEU	N-CA-C	-5.38	101.54	109.62
1	A	317	ASN	CA-CB-CG	5.36	117.96	112.60
1	C	466	LEU	N-CA-C	-5.26	101.73	109.62
1	A	285	ASP	N-CA-C	5.25	116.69	111.07
1	D	357	VAL	CA-C-O	-5.11	113.10	119.95
1	B	466	LEU	N-CA-C	-5.03	102.08	109.62
1	C	466	LEU	CA-C-N	5.00	125.64	119.99
1	C	466	LEU	C-N-CA	5.00	125.64	119.99
1	D	285	ASP	N-CA-C	5.00	116.42	111.07

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3974	0	3883	16	0
1	B	3997	0	3917	15	0
1	C	4000	0	3930	22	0
1	D	3974	0	3881	18	0
2	A	53	0	31	0	0
2	B	53	0	31	1	0
2	C	53	0	31	1	0
2	D	53	0	31	0	0
3	A	26	0	16	3	0
3	B	26	0	16	3	0
3	C	26	0	16	3	0
3	D	26	0	16	2	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
5	A	262	0	0	1	1
5	B	276	0	0	0	0
5	C	238	0	0	2	0
5	D	229	0	0	4	0
All	All	17270	0	15799	66	1

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 2.

All (66) 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:C:394:GLU:HG3	1:C:396:PRO:HD3	1.68	0.73
1:D:277:PRO:HA	1:D:300:ARG:HD2	1.72	0.71
1:A:277:PRO:HA	1:A:300:ARG:HD2	1.73	0.71
1:B:277:PRO:HA	1:B:300:ARG:HD2	1.73	0.71
1:B:138:ARG:NH1	1:C:108:SER:OG	2.24	0.70
1:D:300:ARG:HA	5:D:801:HOH:O	1.89	0.70
1:D:265:GLU:OE2	1:D:333:ARG:NH2	2.25	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:277:PRO:HA	1:C:300:ARG:HD2	1.76	0.68
1:A:178:SER:HB2	1:C:306:PRO:HB2	1.77	0.67
1:A:306:PRO:HB2	1:C:178:SER:HB2	1.78	0.66
1:D:360:HIS:ND1	1:D:362:GLU:OE2	2.29	0.65
1:B:138:ARG:NH1	1:B:142:HIS:NE2	2.45	0.65
1:A:485:HIS:HD2	5:A:1028:HOH:O	1.82	0.61
1:B:138:ARG:NH2	5:D:801:HOH:O	2.34	0.61
1:D:471:LEU:HD13	1:D:535:LEU:HB3	1.85	0.59
1:C:485:HIS:HD2	5:C:981:HOH:O	1.86	0.58
3:C:702:TPP:HN42	3:C:702:TPP:H2	1.70	0.56
1:B:306:PRO:HB2	1:D:178:SER:HB2	1.88	0.56
1:B:178:SER:HB2	1:D:306:PRO:HB2	1.88	0.56
1:B:104:ALA:HA	1:C:135:GLN:HB2	1.88	0.56
1:D:485:HIS:HD2	5:D:984:HOH:O	1.87	0.55
1:D:353:VAL:HA	1:D:360:HIS:CD2	2.40	0.55
1:D:337:ALA:HA	1:D:340:ARG:HE	1.72	0.54
1:B:33:VAL:HA	1:B:36:VAL:HG13	1.90	0.53
1:D:25:VAL:HG21	1:D:33:VAL:HG21	1.89	0.53
3:A:702:TPP:HN42	3:A:702:TPP:H2	1.75	0.52
1:A:356:HIS:CD2	1:A:360:HIS:CE1	2.98	0.52
1:B:25:VAL:HG21	1:B:33:VAL:HG21	1.91	0.52
1:C:369:ASN:OD1	1:C:394:GLU:HG2	2.11	0.51
1:D:485:HIS:HE1	5:D:1016:HOH:O	1.92	0.51
1:C:25:VAL:HG21	1:C:33:VAL:HG21	1.92	0.51
1:A:33:VAL:HA	1:A:36:VAL:HG13	1.92	0.51
3:B:702:TPP:HN42	3:B:702:TPP:H2	1.75	0.51
1:B:360:HIS:CE1	1:B:534:ALA:HB1	2.46	0.50
1:A:104:ALA:HA	1:D:135:GLN:HB2	1.94	0.50
1:C:507:PRO:HA	1:C:510:VAL:HG23	1.94	0.50
3:D:702:TPP:H2	3:D:702:TPP:HN42	1.77	0.50
3:C:702:TPP:HN42	3:C:702:TPP:C2	2.25	0.48
1:A:396:PRO:O	1:A:397:GLU:HB2	2.13	0.48
1:C:352:ASN:O	1:C:356:HIS:HB2	2.13	0.47
1:C:143:ALA:HB1	1:C:154:SER:HB3	1.97	0.47
1:C:485:HIS:HE1	5:C:1024:HOH:O	1.97	0.46
1:B:135:GLN:HB2	1:C:104:ALA:HA	1.98	0.46
1:C:161:ASP:HA	1:C:164:LYS:NZ	2.30	0.46
1:D:357:VAL:CB	1:D:358:PRO:HD3	2.45	0.46
1:C:469:VAL:HG11	1:D:25:VAL:HG13	1.99	0.45
3:A:702:TPP:HN42	3:A:702:TPP:C2	2.30	0.44
1:A:143:ALA:HB1	1:A:154:SER:HB3	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:138:ARG:NH1	1:C:108:SER:HG	2.13	0.44
1:D:143:ALA:HB1	1:D:154:SER:HB3	2.00	0.44
1:A:356:HIS:CD2	1:A:360:HIS:NE2	2.86	0.44
1:C:25:VAL:HG13	1:D:469:VAL:HG11	1.99	0.43
3:D:702:TPP:HN42	3:D:702:TPP:C2	2.31	0.43
1:A:49:GLU:OE2	3:B:702:TPP:N1'	2.52	0.43
1:B:143:ALA:HB1	1:B:154:SER:HB3	2.01	0.43
1:B:379:THR:HA	1:B:403:VAL:O	2.20	0.42
1:A:135:GLN:HB2	1:D:104:ALA:HA	2.02	0.41
1:A:353:VAL:HG12	1:A:360:HIS:CD2	2.55	0.41
1:A:379:THR:HA	1:A:403:VAL:O	2.21	0.41
1:A:299:ARG:HA	1:C:142:HIS:CE1	2.56	0.41
3:B:702:TPP:HN42	3:B:702:TPP:C2	2.32	0.41
1:C:469:VAL:HG21	3:C:702:TPP:HM42	2.02	0.41
1:A:142:HIS:CE1	1:C:299:ARG:HA	2.56	0.40
3:A:702:TPP:N1'	1:B:49:GLU:OE2	2.54	0.40
2:B:701:FAD:H9	2:B:701:FAD:H1'1	1.91	0.40
1:C:311:VAL:HG23	2:C:701:FAD:C2A	2.51	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:847:HOH:O	5:A:911:HOH:O[2_655]	2.08	0.12

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	533/564 (94%)	526 (99%)	6 (1%)	1 (0%)	43 36
1	B	534/564 (95%)	528 (99%)	5 (1%)	1 (0%)	43 36

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	533/564 (94%)	526 (99%)	6 (1%)	1 (0%)	43	36
1	D	535/564 (95%)	526 (98%)	7 (1%)	2 (0%)	30	22
All	All	2135/2256 (95%)	2106 (99%)	24 (1%)	5 (0%)	43	36

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	357	VAL
1	A	27	ASP
1	B	27	ASP
1	C	27	ASP
1	D	27	ASP

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	416/451 (92%)	410 (99%)	6 (1%)	59	59
1	B	421/451 (93%)	417 (99%)	4 (1%)	68	70
1	C	423/451 (94%)	420 (99%)	3 (1%)	76	78
1	D	414/451 (92%)	410 (99%)	4 (1%)	68	70
All	All	1674/1804 (93%)	1657 (99%)	17 (1%)	68	70

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	36	VAL
1	A	243	GLU
1	A	304	LYS
1	A	353	VAL
1	A	381	ASP
1	A	470	LYS
1	B	36	VAL

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Mol	Chain	Res	Type
1	B	138	ARG
1	B	243	GLU
1	B	381	ASP
1	C	243	GLU
1	C	339	GLU
1	C	381	ASP
1	D	243	GLU
1	D	247	GLU
1	D	360	HIS
1	D	381	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	16	GLN
1	A	264	ASN
1	A	317	ASN
1	A	413	ASN
1	A	453	GLN
1	A	485	HIS
1	A	513	GLN
1	B	16	GLN
1	B	413	ASN
1	B	453	GLN
1	B	513	GLN
1	B	533	ASN
1	C	16	GLN
1	C	453	GLN
1	C	485	HIS
1	C	513	GLN
1	D	16	GLN
1	D	321	HIS
1	D	413	ASN
1	D	453	GLN
1	D	485	HIS
1	D	533	ASN

5.3.3 RNA ⓘ

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 12 ligands modelled in this entry, 4 are monoatomic - leaving 8 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	FAD	D	701	-	58,58,58	0.39	0	85,89,89	0.48	0
2	FAD	A	701	-	58,58,58	0.41	0	85,89,89	0.53	0
2	FAD	B	701	-	58,58,58	0.39	0	85,89,89	0.52	0
3	TPP	A	702	4	26,27,27	0.46	0	38,40,40	0.58	0
3	TPP	D	702	4	26,27,27	0.44	0	38,40,40	0.53	0
3	TPP	B	702	4	26,27,27	0.47	0	38,40,40	0.49	0
2	FAD	C	701	-	58,58,58	0.42	0	85,89,89	0.46	0
3	TPP	C	702	4	26,27,27	0.47	0	38,40,40	0.53	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
2	FAD	D	701	-	-	3/34/50/50	0/6/6/6
2	FAD	A	701	-	-	4/34/50/50	0/6/6/6
2	FAD	B	701	-	-	2/34/50/50	0/6/6/6
3	TPP	A	702	4	-	4/17/17/17	0/2/2/2
3	TPP	D	702	4	-	2/17/17/17	0/2/2/2
3	TPP	B	702	4	-	3/17/17/17	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	FAD	C	701	-	-	3/34/50/50	0/6/6/6
3	TPP	C	702	4	-	4/17/17/17	0/2/2/2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (25) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	702	TPP	PA-O3A-PB-O3B
3	C	702	TPP	PA-O3A-PB-O3B
3	D	702	TPP	PA-O3A-PB-O3B
3	B	702	TPP	PA-O3A-PB-O3B
2	A	701	FAD	C2'-C1'-N10-C10
2	B	701	FAD	C2'-C1'-N10-C10
2	C	701	FAD	C2'-C1'-N10-C10
2	D	701	FAD	C2'-C1'-N10-C10
2	D	701	FAD	PA-O3P-P-O1P
3	C	702	TPP	PA-O3A-PB-O1B
3	D	702	TPP	PA-O3A-PB-O1B
2	A	701	FAD	PA-O3P-P-O1P
2	A	701	FAD	PA-O3P-P-O2P
2	C	701	FAD	PA-O3P-P-O1P
2	C	701	FAD	PA-O3P-P-O2P
2	D	701	FAD	PA-O3P-P-O2P
3	A	702	TPP	PA-O3A-PB-O1B
3	A	702	TPP	C5-C6-C7-O7
3	B	702	TPP	C5-C6-C7-O7
3	C	702	TPP	C5-C6-C7-O7
2	B	701	FAD	PA-O3P-P-O2P
2	A	701	FAD	C2'-C3'-C4'-O4'
3	A	702	TPP	C5'-C7'-N3-C2
3	B	702	TPP	C5'-C7'-N3-C2
3	C	702	TPP	C5'-C7'-N3-C2

There are no ring outliers.

6 monomers are involved in 13 short contacts:

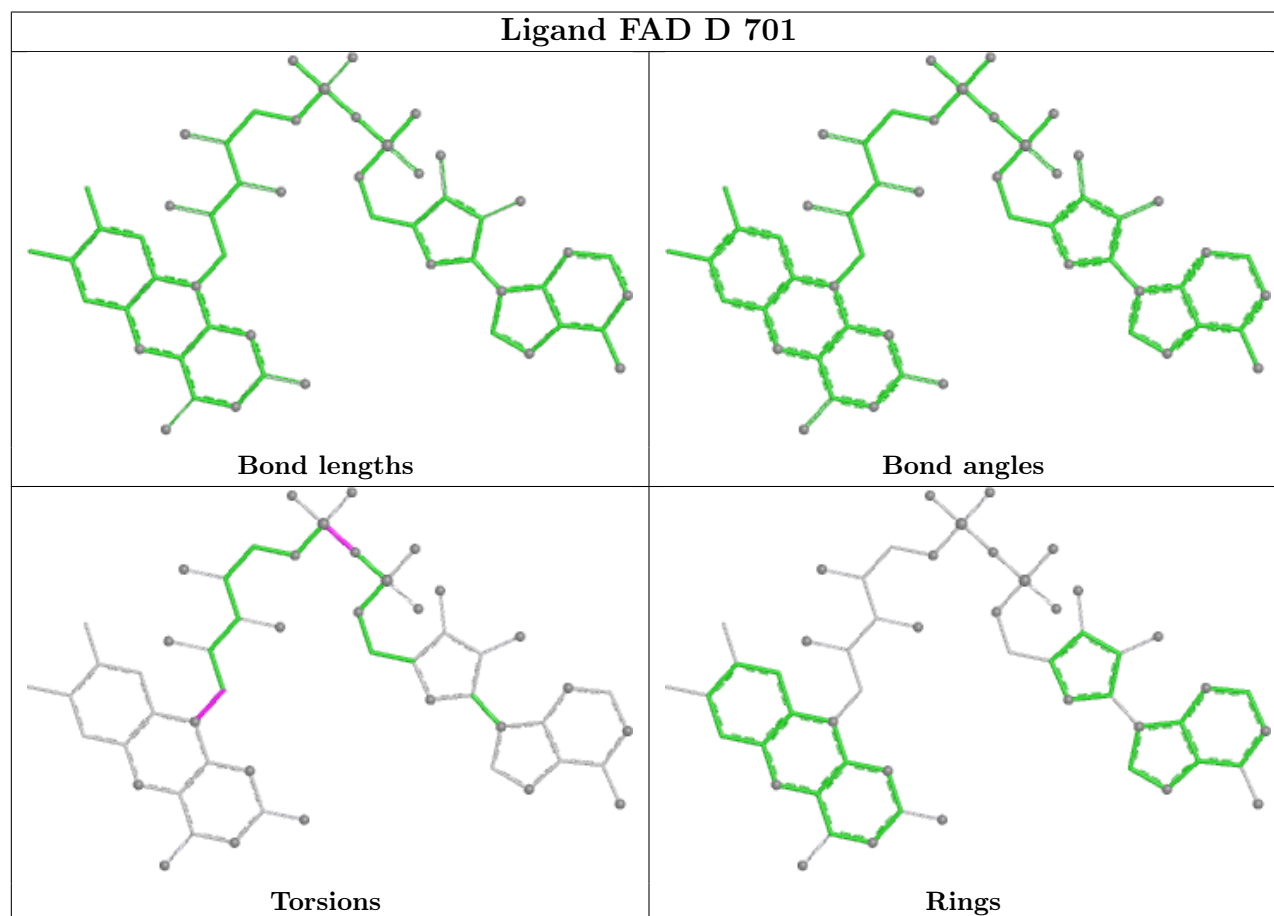
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	701	FAD	1	0

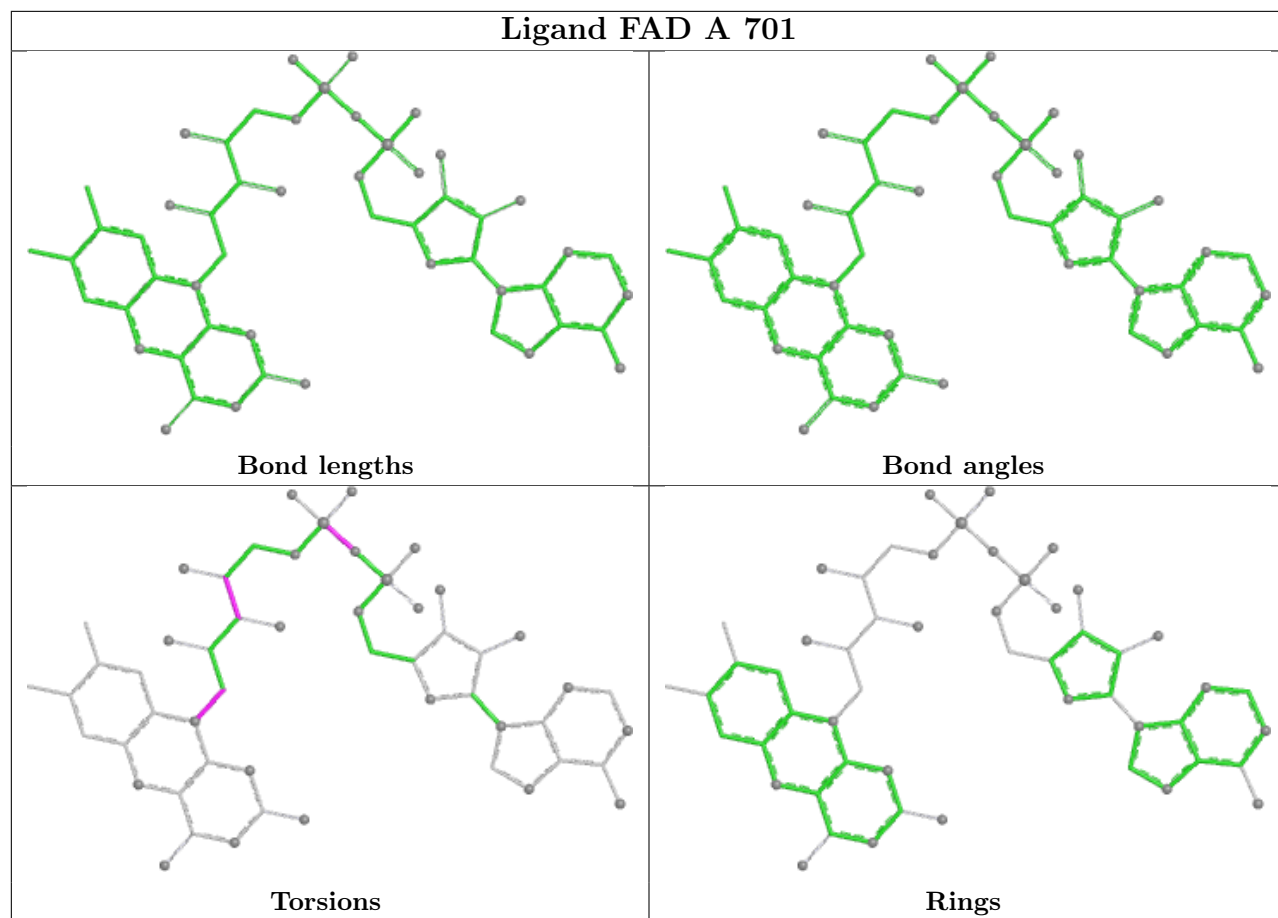
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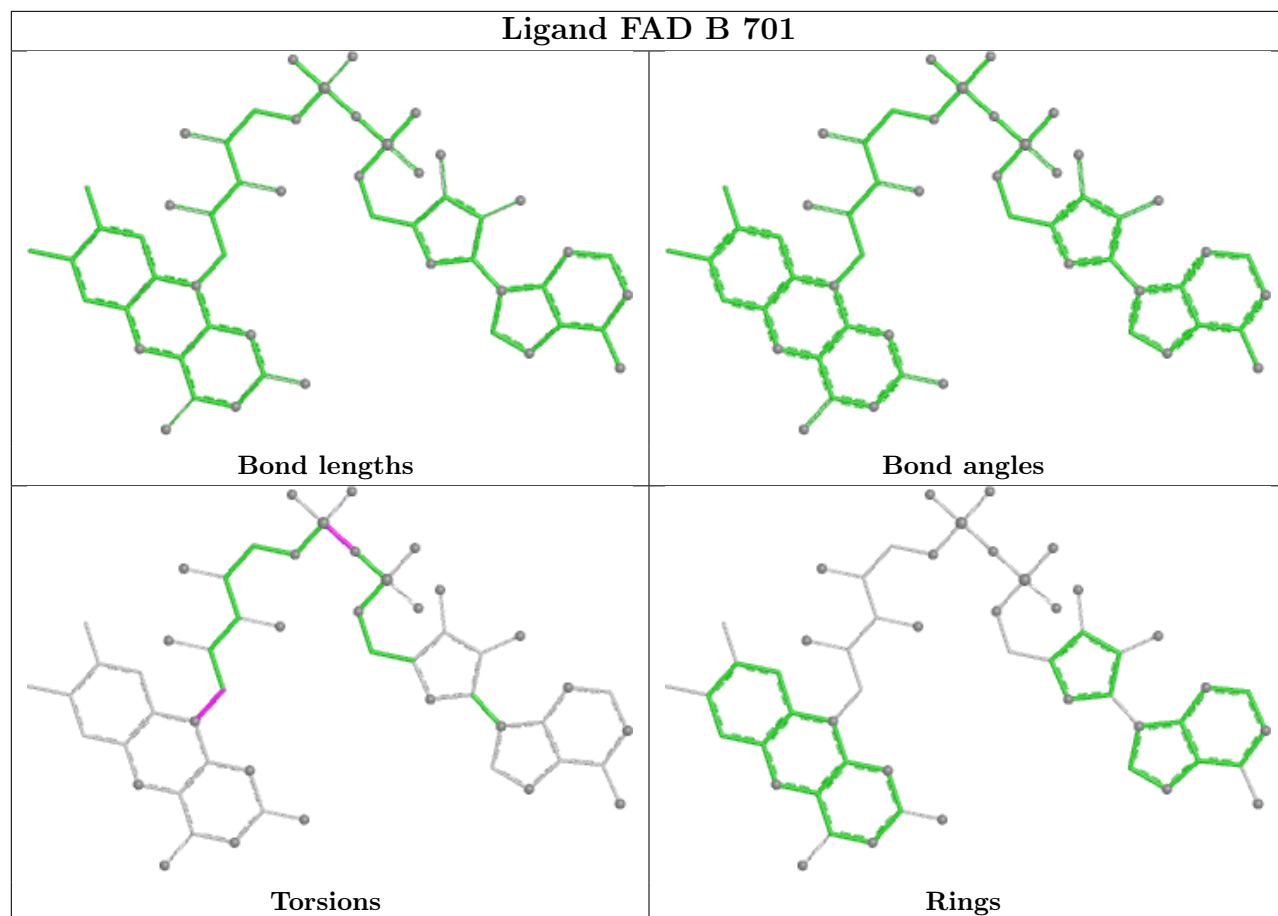
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	702	TPP	3	0
3	D	702	TPP	2	0
3	B	702	TPP	3	0
2	C	701	FAD	1	0
3	C	702	TPP	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

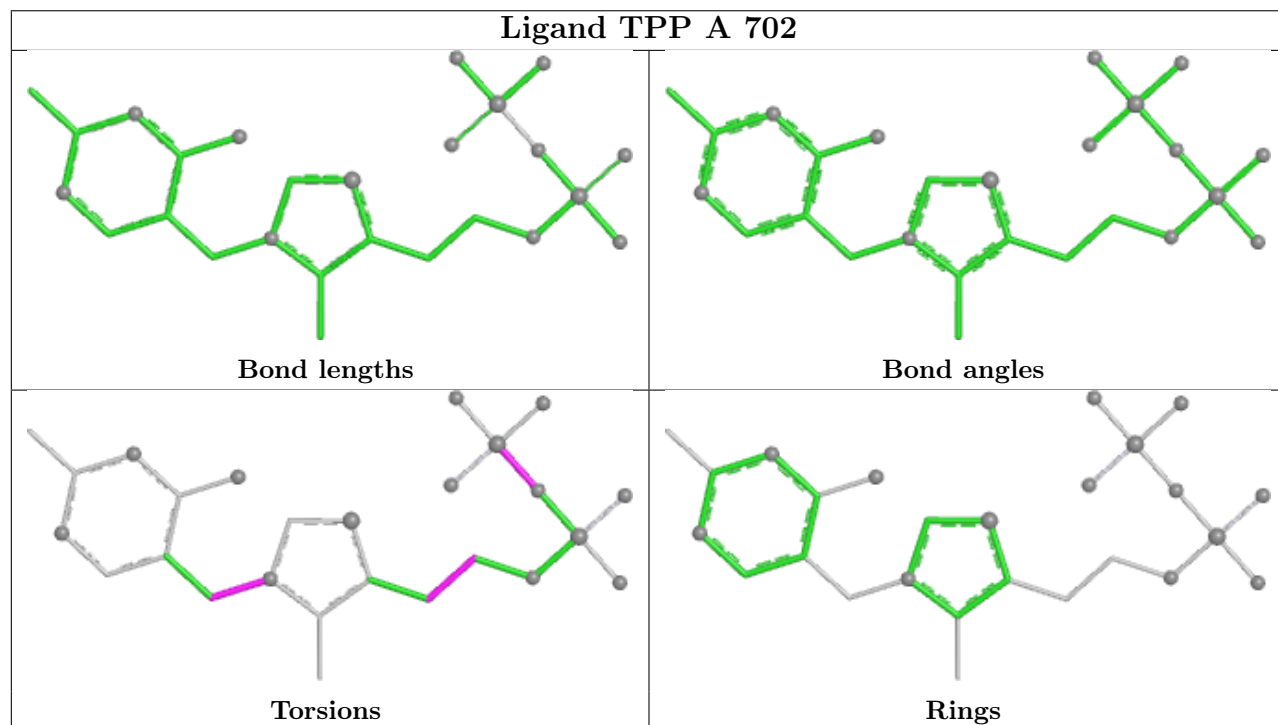




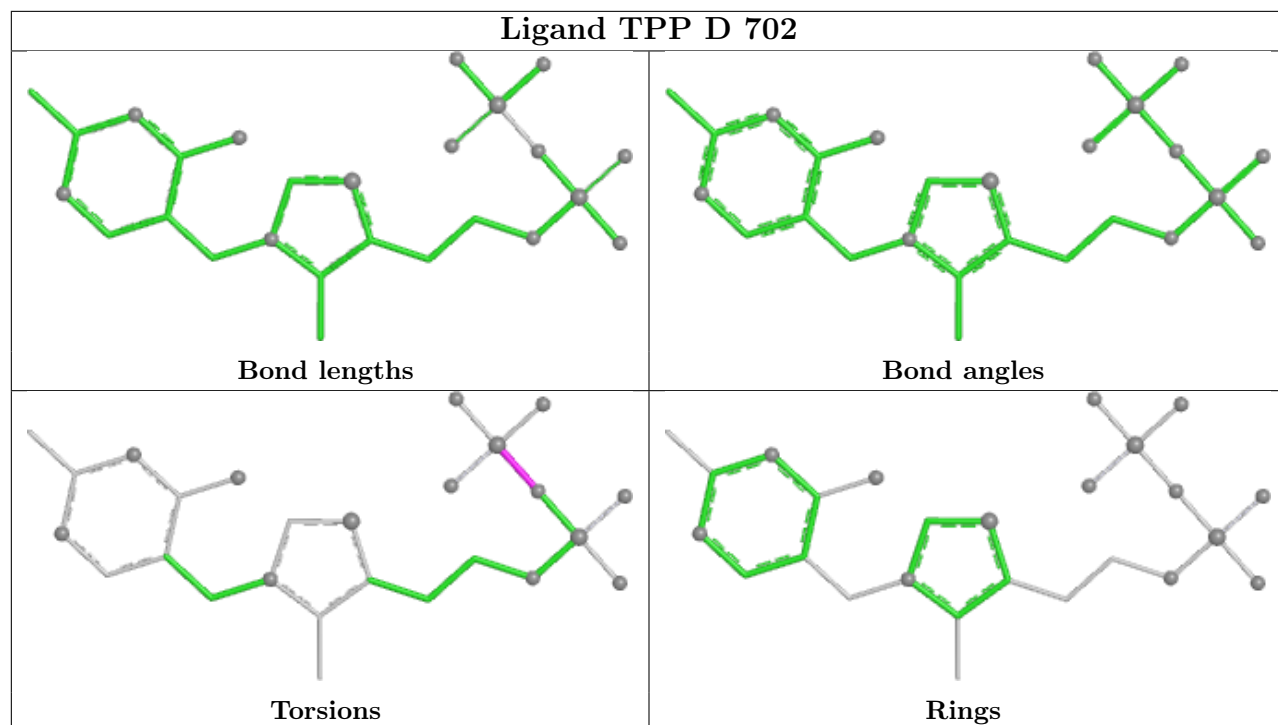
Ligand FAD B 701



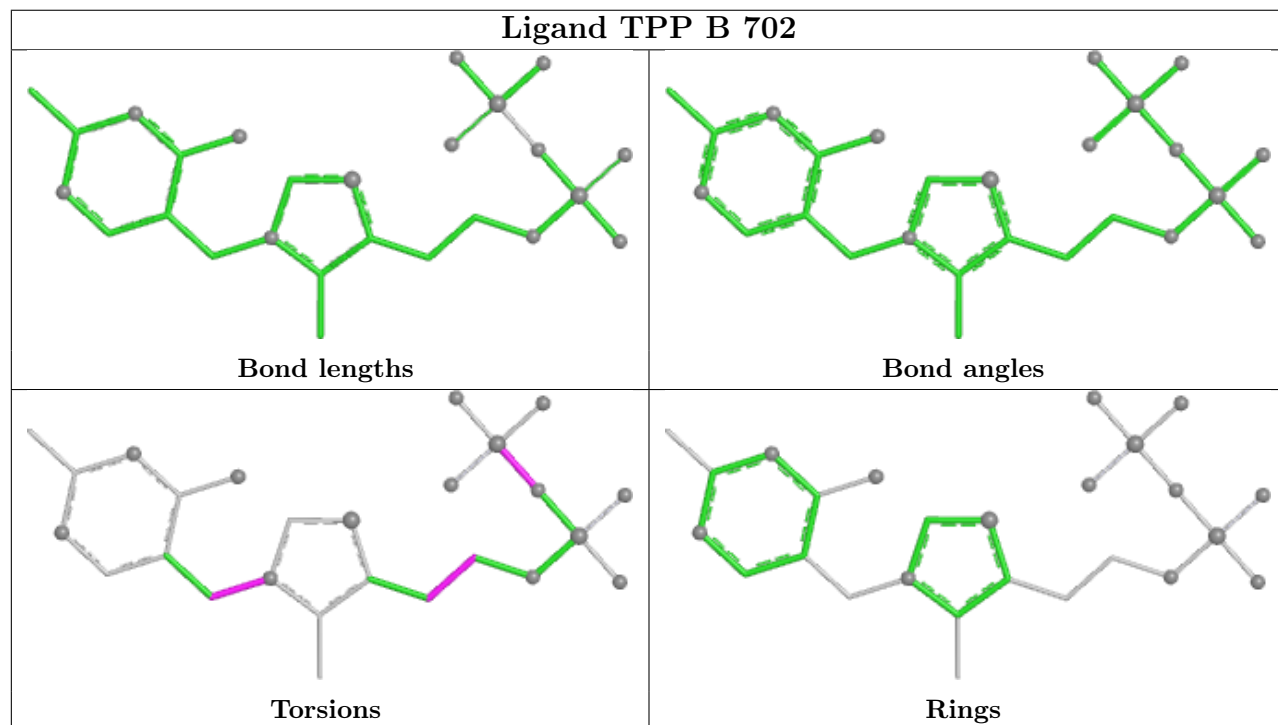
Ligand TPP A 702



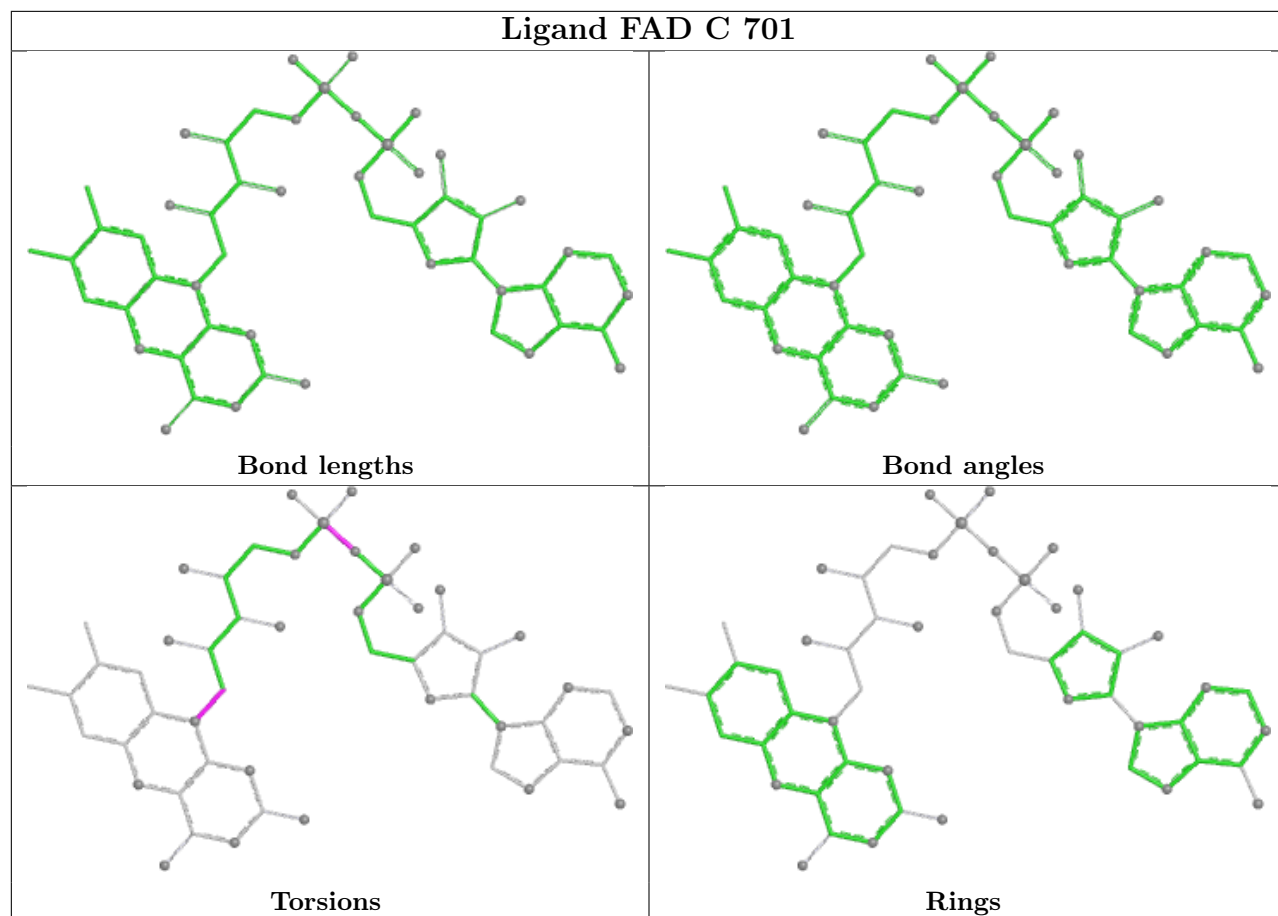
Ligand TPP D 702



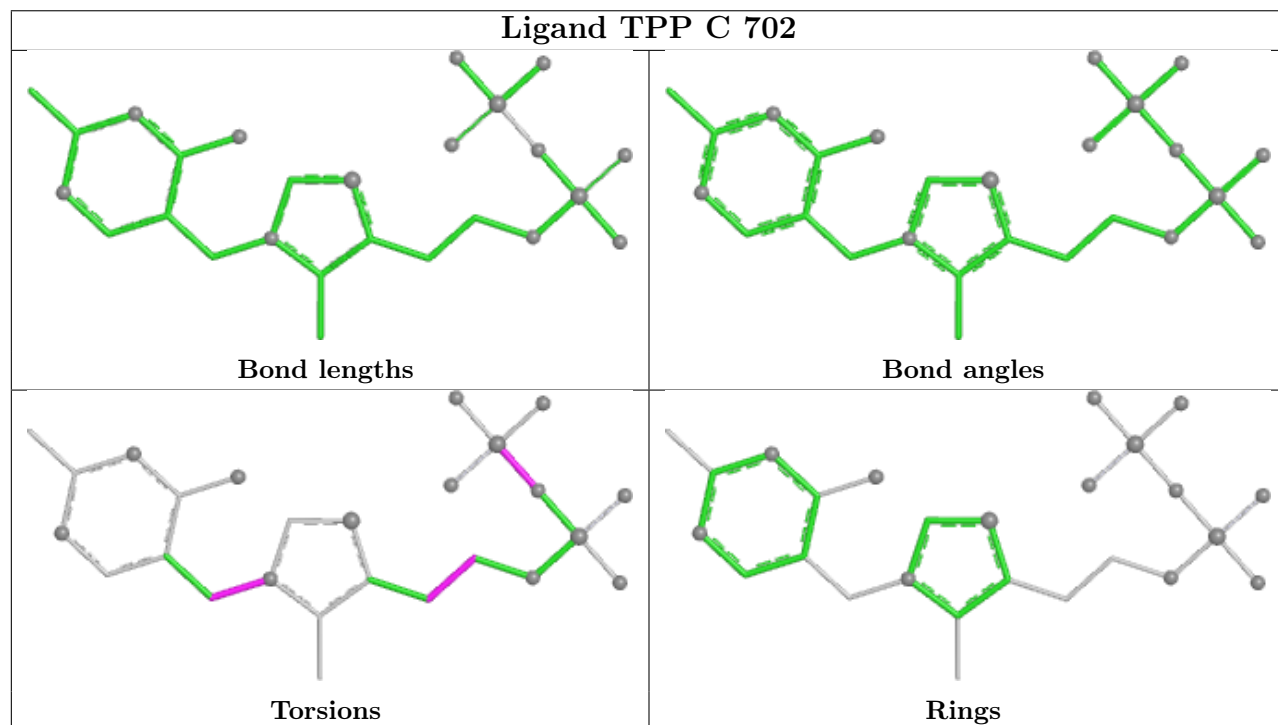
Ligand TPP B 702



Ligand FAD C 701



Ligand TPP C 702



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	535/564 (94%)	-0.10	11 (2%) 63 67	19, 33, 54, 86	0
1	B	536/564 (95%)	-0.19	5 (0%) 81 83	22, 33, 53, 65	0
1	C	535/564 (94%)	-0.07	7 (1%) 75 78	23, 37, 57, 70	0
1	D	537/564 (95%)	0.13	20 (3%) 45 48	19, 41, 77, 110	0
All	All	2143/2256 (94%)	-0.06	43 (2%) 65 69	19, 35, 63, 110	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	353	VAL	6.1
1	D	351	HIS	5.1
1	D	538	PRO	4.9
1	D	357	VAL	4.8
1	D	356	HIS	4.5
1	D	535	LEU	4.1
1	D	2	ALA	4.0
1	D	342	LEU	3.9
1	B	166	ASP	3.7
1	D	355	LYS	3.7
1	D	350	THR	3.6
1	B	2	ALA	3.5
1	A	534	ALA	3.5
1	B	537	ILE	3.5
1	D	166	ASP	3.4
1	C	166	ASP	3.3
1	D	349	TYR	3.0
1	A	166	ASP	3.0
1	D	347	GLU	3.0
1	A	2	ALA	3.0
1	A	353	VAL	3.0

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Mol	Chain	Res	Type	RSRZ
1	D	346	VAL	2.9
1	D	3	HIS	2.9
1	A	200	ASN	2.8
1	D	345	VAL	2.8
1	C	342	LEU	2.8
1	A	351	HIS	2.8
1	D	537	ILE	2.7
1	C	2	ALA	2.7
1	D	536	SER	2.5
1	C	353	VAL	2.5
1	A	535	LEU	2.4
1	A	536	SER	2.4
1	A	285	ASP	2.3
1	C	535	LEU	2.3
1	B	534	ALA	2.2
1	C	344	SER	2.2
1	A	352	ASN	2.1
1	B	353	VAL	2.1
1	D	285	ASP	2.1
1	C	343	SER	2.1
1	A	475	VAL	2.0
1	D	316	GLU	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	TPP	A	702	26/26	0.96	0.06	25,32,36,38	0

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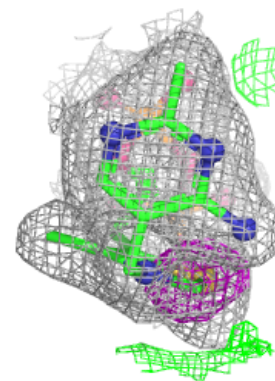
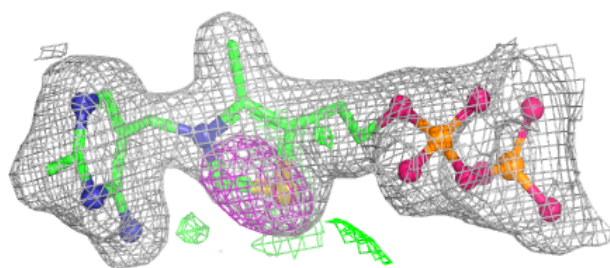
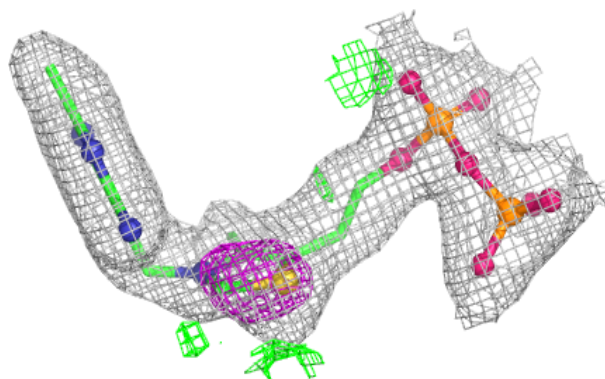
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	TPP	C	702	26/26	0.96	0.06	22,29,34,37	0
3	TPP	D	702	26/26	0.96	0.07	31,42,44,44	0
2	FAD	A	701	53/53	0.97	0.05	24,27,30,31	0
3	TPP	B	702	26/26	0.97	0.06	22,29,34,37	0
2	FAD	C	701	53/53	0.97	0.05	24,27,30,32	0
2	FAD	D	701	53/53	0.97	0.05	29,32,35,38	0
4	MG	C	703	1/1	0.97	0.04	30,30,30,30	0
2	FAD	B	701	53/53	0.98	0.04	22,24,26,28	0
4	MG	B	703	1/1	0.99	0.01	28,28,28,28	0
4	MG	D	703	1/1	0.99	0.03	49,49,49,49	0
4	MG	A	703	1/1	1.00	0.01	27,27,27,27	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.

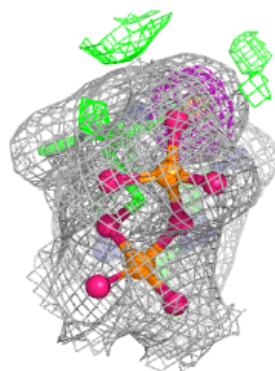
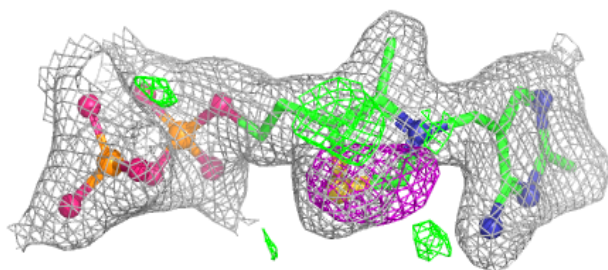
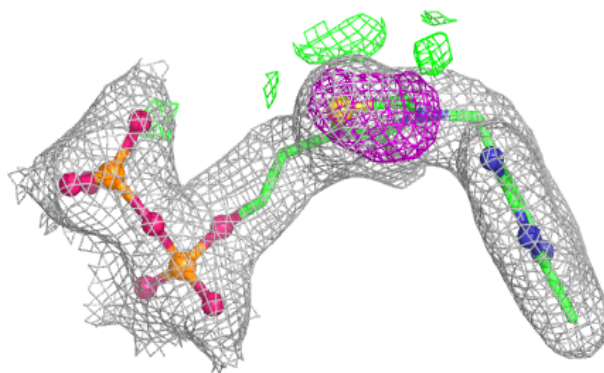
Electron density around TPP A 702:

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

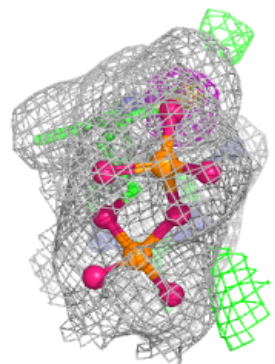
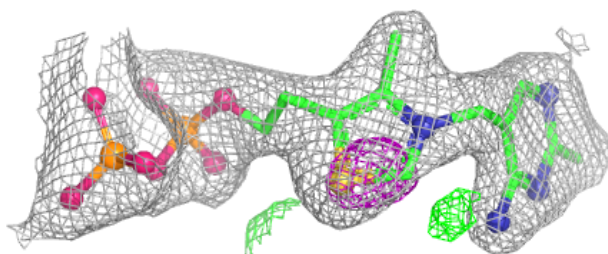
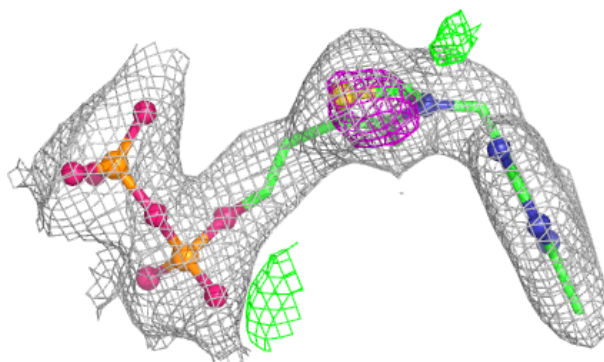


Electron density around TPP C 702:

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

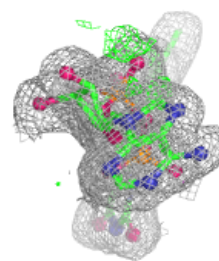
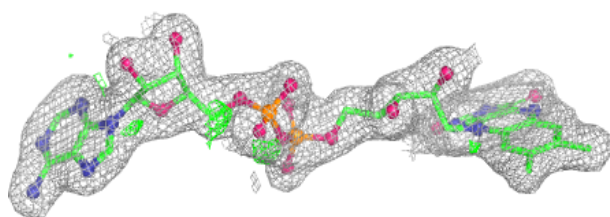
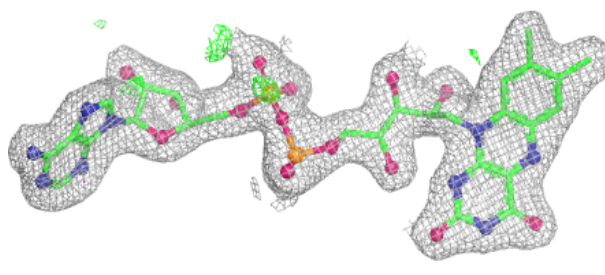
**Electron density around TPP D 702:**

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and green (positive)

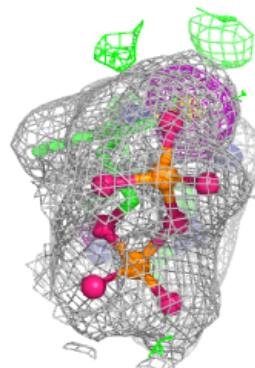
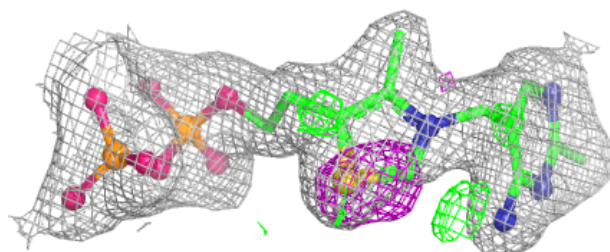
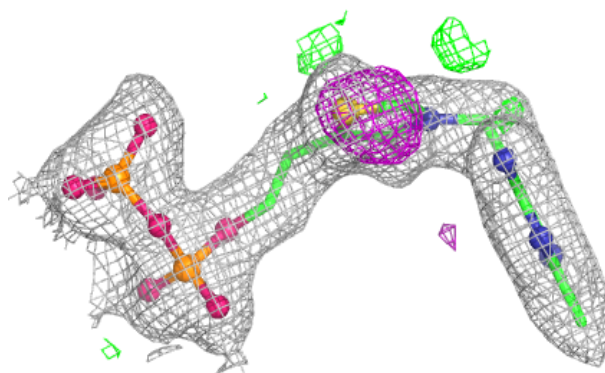


Electron density around FAD A 701:

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

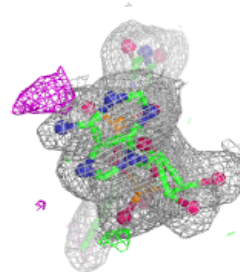
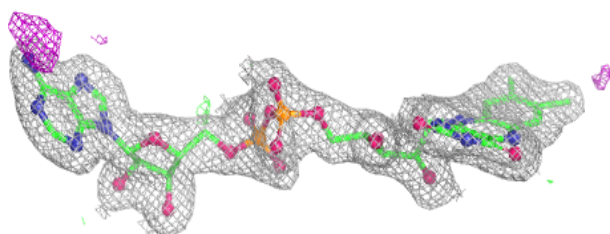
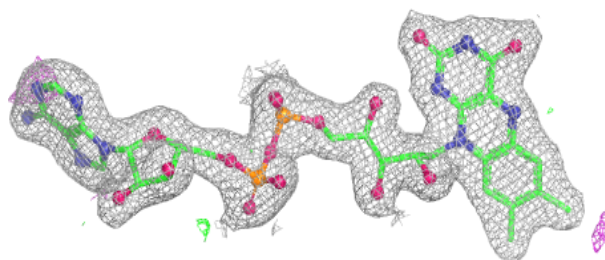
**Electron density around TPP B 702:**

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and green (positive)

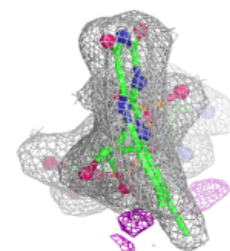
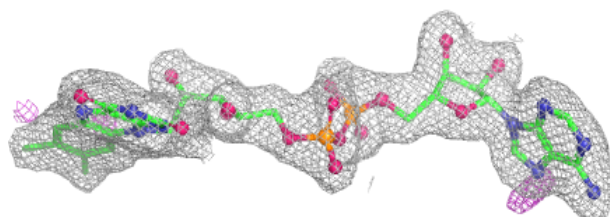
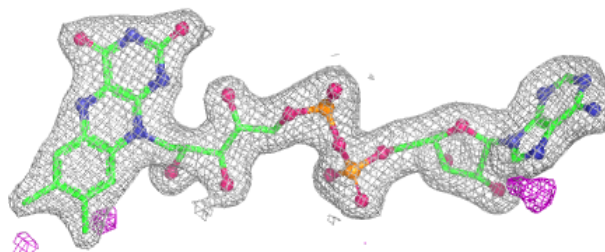


Electron density around FAD C 701:

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

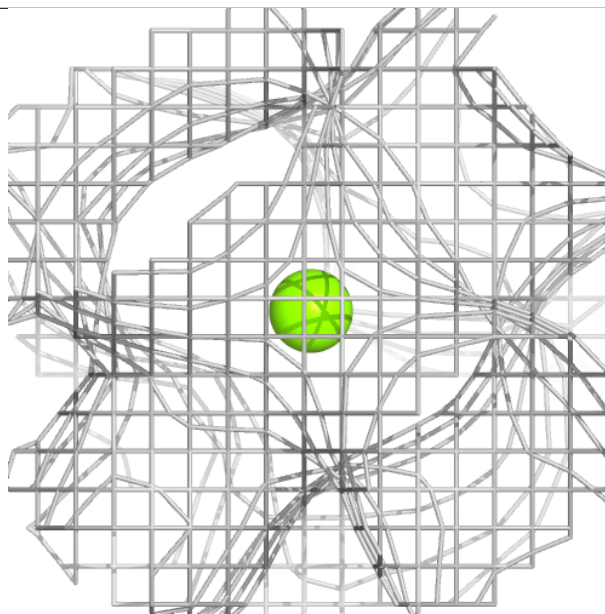
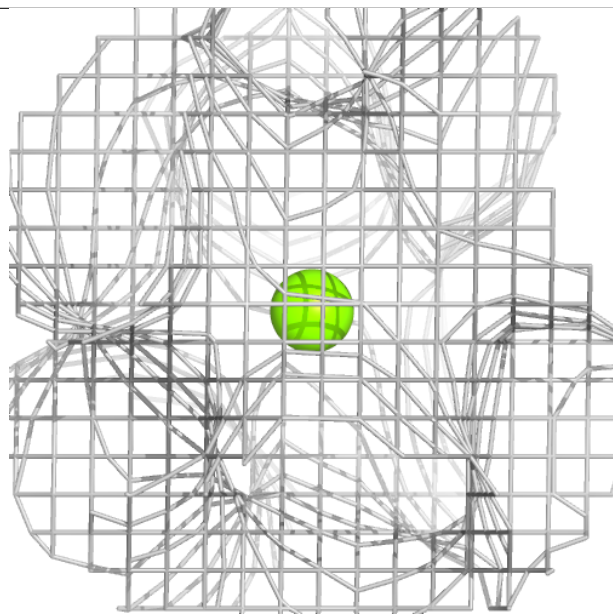
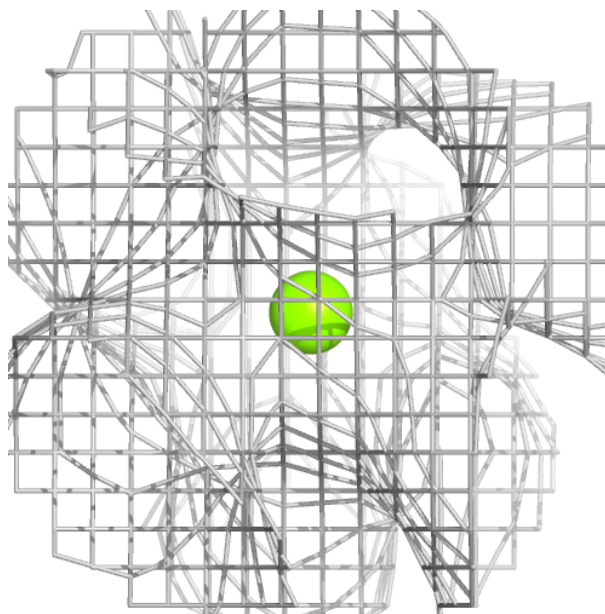
**Electron density around FAD D 701:**

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and green (positive)



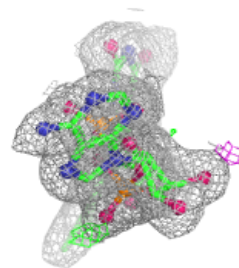
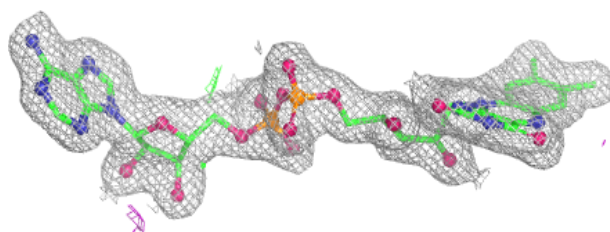
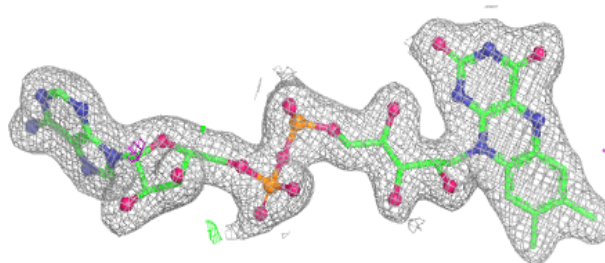
Electron density around MG C 703:

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and green (positive)



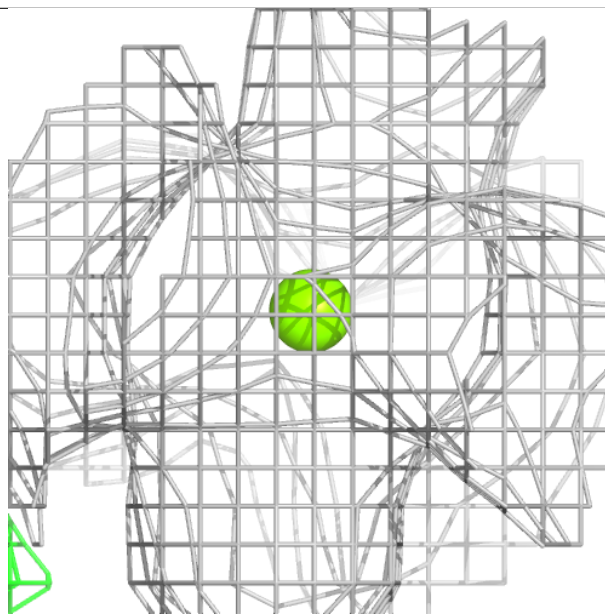
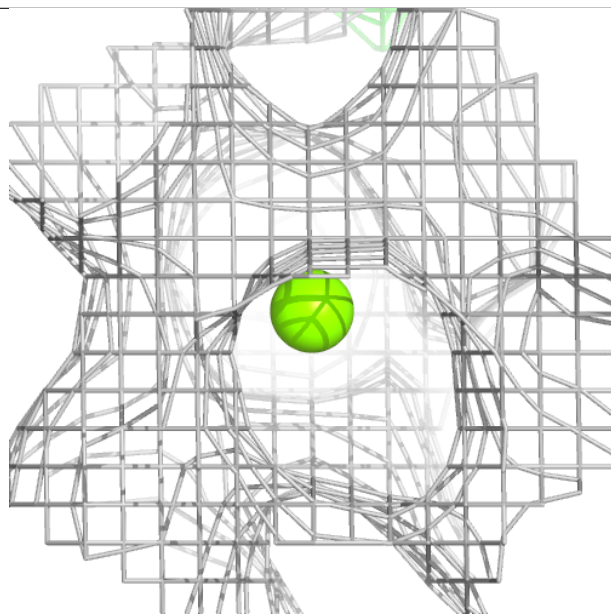
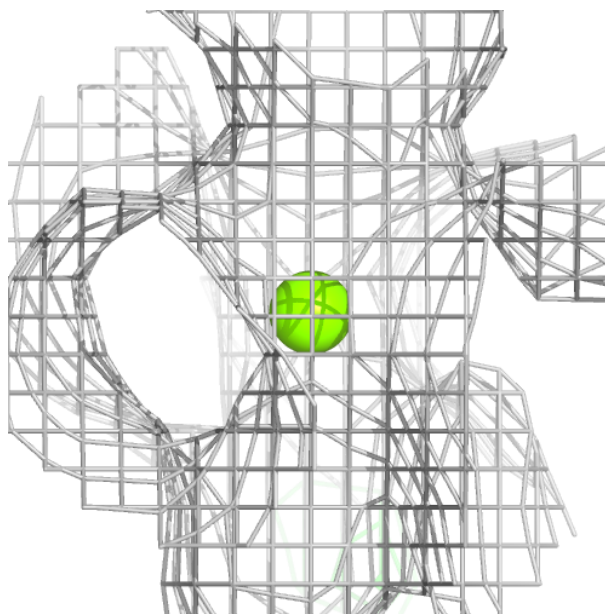
Electron density around FAD B 701:

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and green (positive)



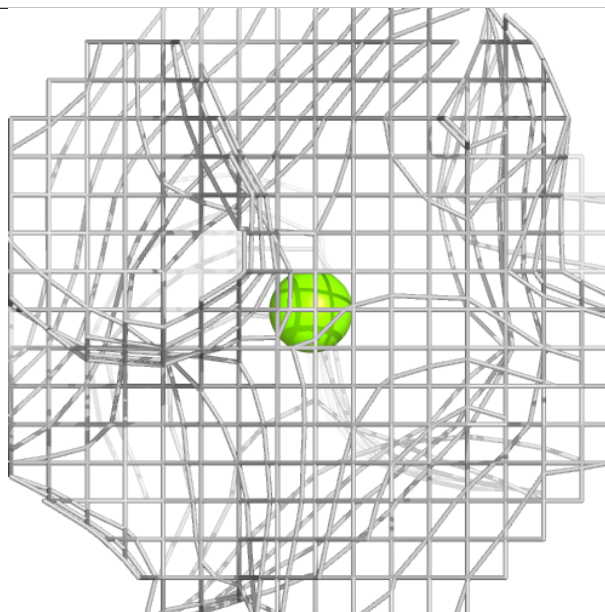
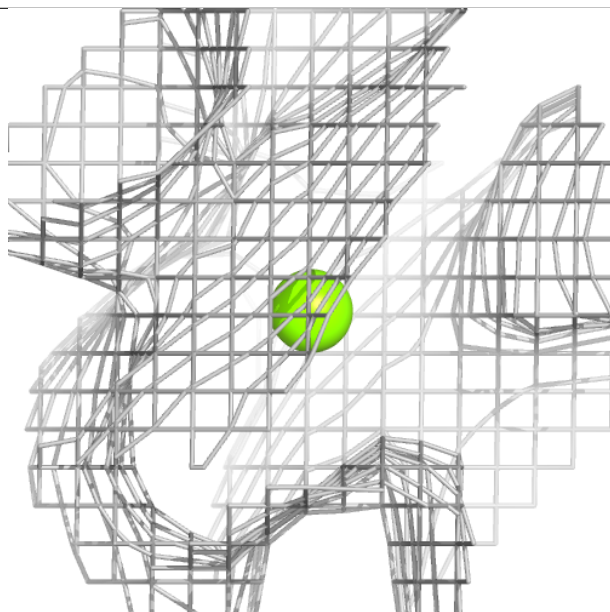
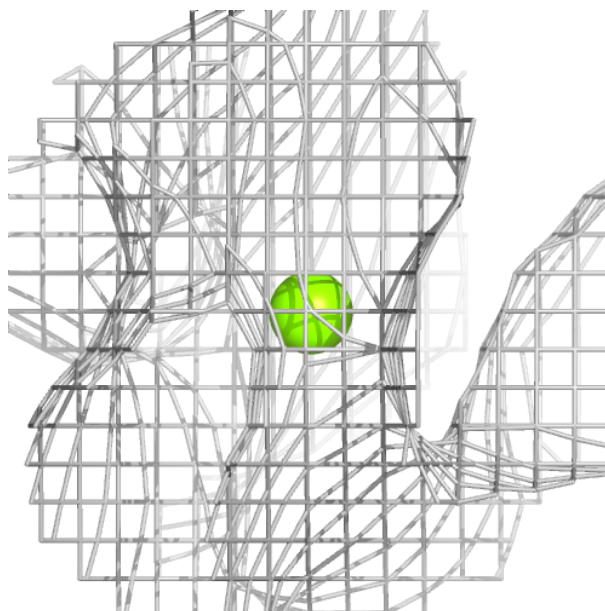
Electron density around MG B 703:

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



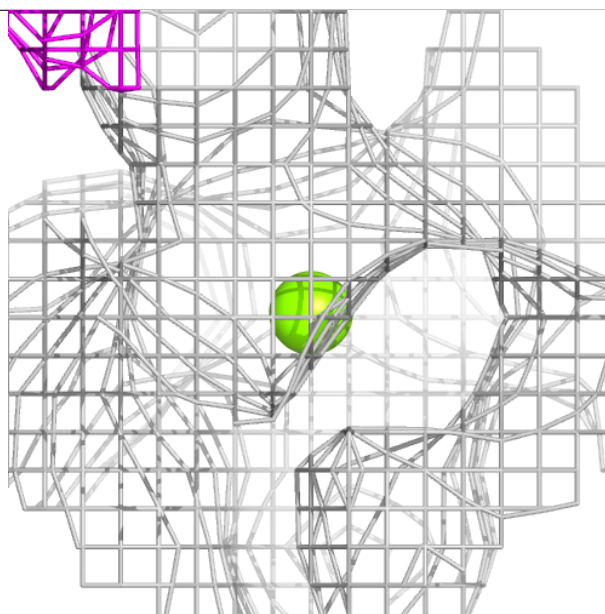
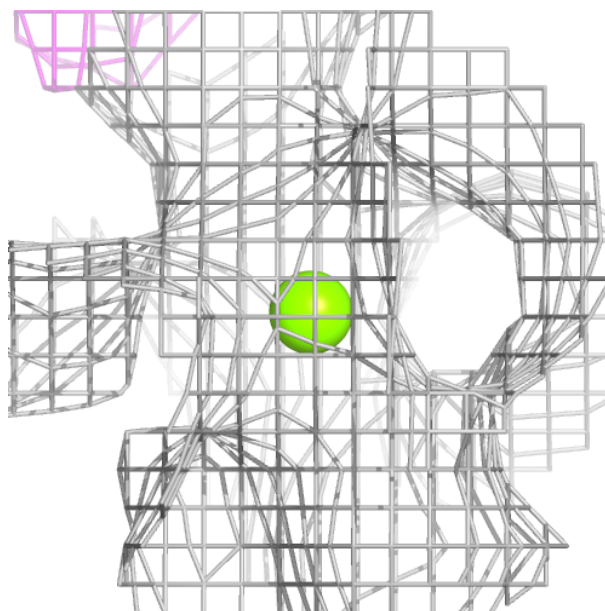
Electron density around MG D 703:

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



Electron density around MG A 703:

$2mF_o-DF_c$ (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.