



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 06:17 PM UTC

PDB ID : 5O0X / pdb_00005o0x
Title : Crystal structure of dehydrogenase domain of *Cylindrospermum stagnale* NADPH-Oxidase 5 (NOX5)
Authors : Magnani, F.; Nenci, S.; Mattevi, A.
Deposited on : 2017-05-17
Resolution : 2.20 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

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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

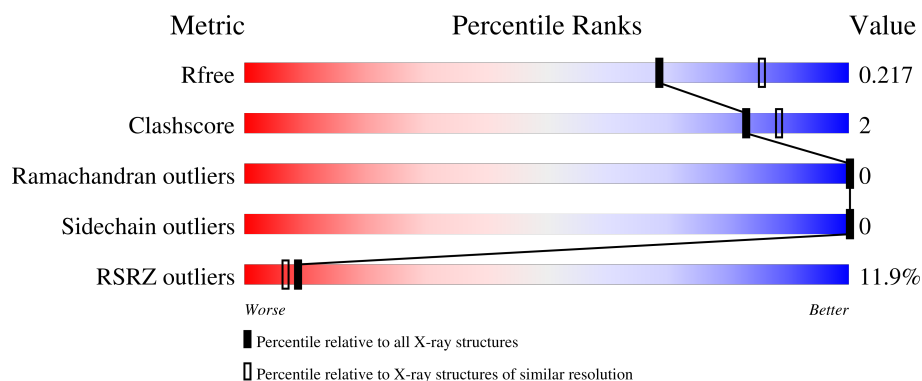
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	291	11% 85% 10%

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 2378 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.

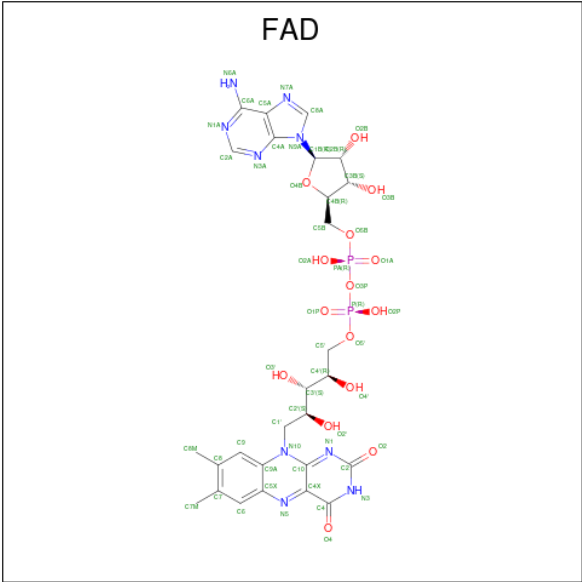
- Molecule 1 is a protein called Putative ferric reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	261	Total	C	N	O	S	0	0	0
			2066	1348	343	369	6			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	411	GLY	-	expression tag	UNP K9WT99
A	412	SER	-	expression tag	UNP K9WT99
A	694	PRO	-	expression tag	UNP K9WT99
A	695	TRP	-	expression tag	UNP K9WT99
A	696	LEU	-	expression tag	UNP K9WT99
A	697	GLU	-	expression tag	UNP K9WT99
A	698	LEU	-	expression tag	UNP K9WT99
A	699	ALA	-	expression tag	UNP K9WT99
A	700	ALA	-	expression tag	UNP K9WT99
A	701	ALA	-	expression tag	UNP K9WT99

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).

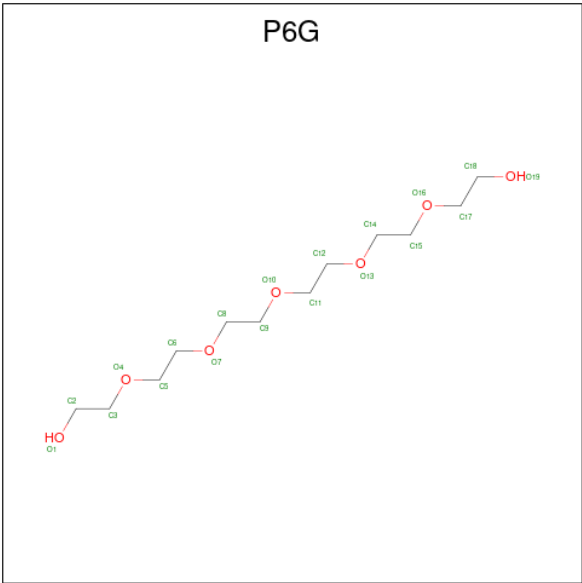


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

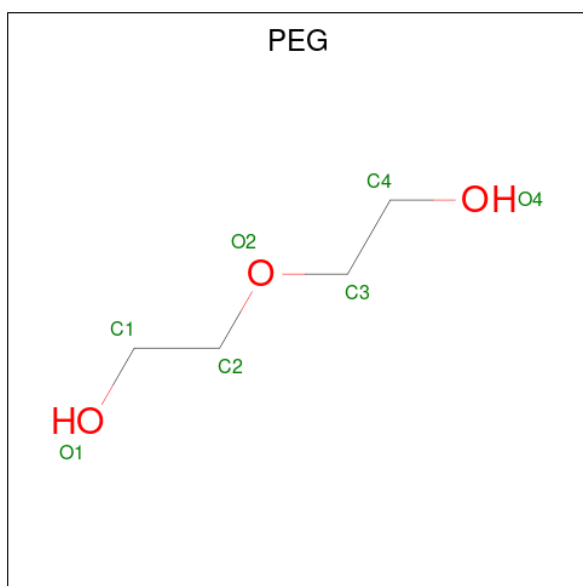
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	7	Total	Cl	0	0
			7	7		

- Molecule 4 is HEXAETHYLENE GLYCOL (CCD ID: P6G) (formula: C₁₂H₂₆O₇).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			9	6	3		
4	A	1	Total	C	O	0	0
			9	6	3		
4	A	1	Total	C	O	0	0
			10	6	4		
4	A	1	Total	C	O	0	0
			13	8	5		
4	A	1	Total	C	O	0	0
			13	8	5		

- Molecule 5 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



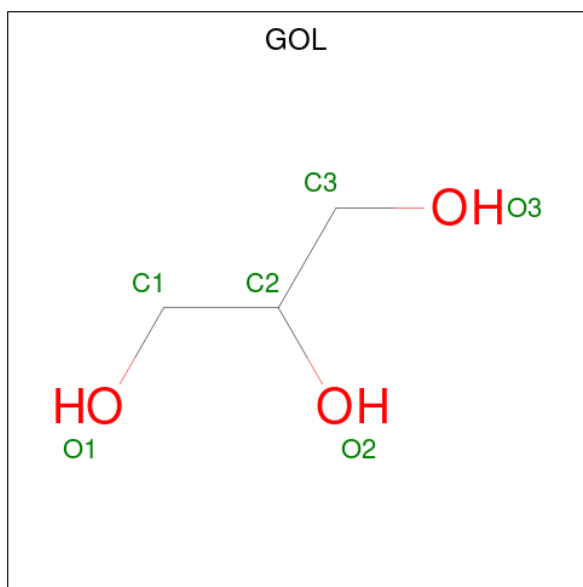
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			7	4	3		
5	A	1	Total	C	O	0	0
			7	4	3		
5	A	1	Total	C	O	0	0
			7	4	3		
5	A	1	Total	C	O	0	0
			7	4	3		
5	A	1	Total	C	O	0	0
			7	4	3		
5	A	1	Total	C	O	0	0
			7	4	3		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			7	4	3		
5	A	1	Total	C	O	0	0
			7	4	3		

- Molecule 6 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

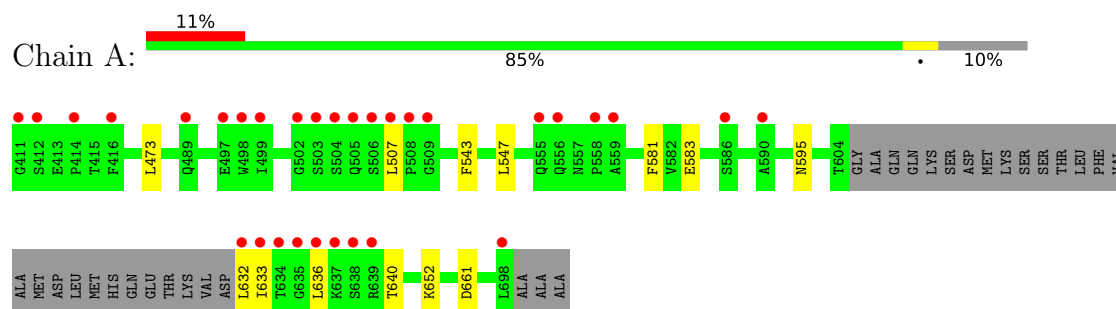
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	99	Total	O	0	0
			99	99		

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: Putative ferric reductase



4 Data and refinement statistics

Property	Value	Source
Space group	P 3 2 1	Depositor
Cell constants a, b, c, α , β , γ	128.28Å 128.28Å 72.52Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	111.09 – 2.20 111.09 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.0 (111.09-2.20) 99.0 (111.09-2.20)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.73 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.8.0155	Depositor
R, R_{free}	0.184 , 0.209 0.191 , 0.217	Depositor DCC
R_{free} test set	1660 reflections (4.72%)	wwPDB-VP
Wilson B-factor (Å ²)	39.1	Xtriage
Anisotropy	0.029	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 43.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.032 for -h,-k,l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	2378	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.73% 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: FAD, PEG, GOL, P6G, CL

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	0.87	0/2130	0.91	1/2902 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	473	LEU	N-CA-C	-5.16	101.75	109.95

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	2066	0	1991	10	0
2	A	53	0	31	1	0
3	A	7	0	0	0	0
4	A	54	0	67	1	1
5	A	63	0	90	4	0
6	A	36	0	48	1	0
7	A	99	0	0	1	0
All	All	2378	0	2227	11	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 (11) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:583:GLU:HB2	5:A:816:PEG:H32	1.63	0.80
1:A:632:LEU:CB	1:A:636:LEU:O	2.36	0.74
1:A:652:LYS:HG3	4:A:813:P6G:H172	1.87	0.56
1:A:633:ILE:CB	1:A:640:THR:HB	2.38	0.54
1:A:581:PHE:HA	5:A:816:PEG:H31	1.93	0.51
2:A:801:FAD:H1'2	2:A:801:FAD:O4'	2.15	0.45
1:A:661:ASP:O	6:A:822:GOL:O2	2.33	0.45
1:A:507:LEU:HD23	7:A:917:HOH:O	2.17	0.44
1:A:595:ASN:OD1	5:A:818:PEG:H21	2.19	0.43
1:A:595:ASN:OD1	5:A:818:PEG:C2	2.67	0.42
1:A:543:PHE:O	1:A:547:LEU:HG	2.20	0.41

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)
4:A:813:P6G:O7	4:A:813:P6G:O7[2_655]	2.10	0.10

5.3 Torsion angles

5.3.1 Protein backbone

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	257/291 (88%)	254 (99%)	3 (1%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	217/255 (85%)	217 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 28 ligands modelled in this entry, 7 are monoatomic - leaving 21 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	$\# Z > 2$
5	PEG	A	814	-	6,6,6	0.56	0	5,5,5	0.48	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	PEG	A	828	-	6,6,6	0.53	0	5,5,5	0.54	0
5	PEG	A	821	-	6,6,6	0.62	0	5,5,5	0.47	0
6	GOL	A	822	-	5,5,5	0.41	0	5,5,5	0.33	0
6	GOL	A	823	-	5,5,5	0.38	0	5,5,5	1.33	1 (20%)
4	P6G	A	813	-	12,12,18	0.58	0	11,11,17	0.30	0
5	PEG	A	815	-	6,6,6	0.57	0	5,5,5	0.53	0
6	GOL	A	826	-	5,5,5	0.26	0	5,5,5	0.52	0
5	PEG	A	816	-	6,6,6	0.49	0	5,5,5	0.93	0
5	PEG	A	820	-	6,6,6	0.57	0	5,5,5	0.38	0
4	P6G	A	810	-	8,8,18	0.58	0	7,7,17	0.40	0
6	GOL	A	824	-	5,5,5	0.18	0	5,5,5	0.24	0
6	GOL	A	825	-	5,5,5	0.36	0	5,5,5	0.50	0
5	PEG	A	817	-	6,6,6	0.46	0	5,5,5	0.47	0
6	GOL	A	827	-	5,5,5	0.43	0	5,5,5	0.82	0
5	PEG	A	818	-	6,6,6	0.67	0	5,5,5	0.88	0
5	PEG	A	819	-	6,6,6	0.47	0	5,5,5	0.47	0
4	P6G	A	811	-	9,9,18	0.29	0	8,8,17	0.73	0
4	P6G	A	812	-	12,12,18	0.54	0	11,11,17	0.38	0
2	FAD	A	801	-	58,58,58	2.19	21 (36%)	85,89,89	2.22	26 (30%)
4	P6G	A	809	-	8,8,18	0.62	0	7,7,17	0.57	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
5	PEG	A	814	-	-	2/4/4/4	-
5	PEG	A	828	-	-	1/4/4/4	-
5	PEG	A	821	-	-	3/4/4/4	-
6	GOL	A	822	-	-	3/4/4/4	-
6	GOL	A	823	-	-	4/4/4/4	-
4	P6G	A	813	-	-	4/10/10/16	-
5	PEG	A	815	-	-	3/4/4/4	-
6	GOL	A	826	-	-	0/4/4/4	-
5	PEG	A	816	-	-	1/4/4/4	-
5	PEG	A	820	-	-	1/4/4/4	-
4	P6G	A	810	-	-	1/6/6/16	-
6	GOL	A	824	-	-	2/4/4/4	-
6	GOL	A	825	-	-	2/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	PEG	A	817	-	-	0/4/4/4	-
6	GOL	A	827	-	-	0/4/4/4	-
5	PEG	A	818	-	-	2/4/4/4	-
5	PEG	A	819	-	-	2/4/4/4	-
4	P6G	A	811	-	-	1/7/7/16	-
4	P6G	A	812	-	-	3/10/10/16	-
2	FAD	A	801	-	-	8/34/50/50	0/6/6/6
4	P6G	A	809	-	-	4/6/6/16	-

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	801	FAD	C9A-C5X	4.95	1.49	1.41
2	A	801	FAD	C4-N3	-4.92	1.29	1.38
2	A	801	FAD	P-O3P	-4.39	1.54	1.59
2	A	801	FAD	C4'-C3'	-3.98	1.46	1.53
2	A	801	FAD	C5A-C4A	3.87	1.46	1.39
2	A	801	FAD	C5X-N5	-3.79	1.32	1.39
2	A	801	FAD	C5A-N7A	-3.66	1.32	1.39
2	A	801	FAD	C2-N3	-3.66	1.30	1.39
2	A	801	FAD	C4A-N9A	-3.52	1.30	1.37
2	A	801	FAD	P-O2P	-3.04	1.41	1.55
2	A	801	FAD	PA-O3P	-3.00	1.56	1.59
2	A	801	FAD	C8A-N9A	-2.93	1.32	1.37
2	A	801	FAD	P-O1P	-2.74	1.41	1.50
2	A	801	FAD	C1'-C2'	2.67	1.56	1.52
2	A	801	FAD	C2'-C3'	2.57	1.57	1.53
2	A	801	FAD	PA-O1A	-2.53	1.42	1.50
2	A	801	FAD	C8-C7	2.46	1.46	1.40
2	A	801	FAD	PA-O2A	-2.34	1.44	1.55
2	A	801	FAD	C5'-C4'	2.09	1.54	1.51
2	A	801	FAD	O2'-C2'	-2.08	1.39	1.43
2	A	801	FAD	C6-C5X	-2.02	1.36	1.40

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	801	FAD	C1'-C2'-C3'	6.98	128.59	109.66
2	A	801	FAD	O5'-C5'-C4'	5.67	124.49	109.36
2	A	801	FAD	C5A-C4A-N3A	-5.25	119.48	126.72
2	A	801	FAD	O3'-C3'-C2'	4.92	120.12	108.93

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	801	FAD	O3'-C3'-C4'	-4.58	98.52	108.93
2	A	801	FAD	C2B-C1B-N9A	-4.37	102.45	113.30
2	A	801	FAD	N3A-C4A-N9A	4.25	134.39	127.17
2	A	801	FAD	C4X-C10-N10	4.18	122.47	116.48
2	A	801	FAD	C9A-N10-C10	-3.98	114.69	120.75
2	A	801	FAD	C4'-C3'-C2'	3.85	119.98	113.57
2	A	801	FAD	O4'-C4'-C3'	-3.70	100.58	109.25
2	A	801	FAD	C4X-C10-N1	-3.59	115.78	124.59
2	A	801	FAD	C2A-N3A-C4A	3.14	119.49	111.83
2	A	801	FAD	C10-N1-C2	3.03	123.42	116.85
2	A	801	FAD	N3A-C2A-N1A	-2.93	124.14	128.58
2	A	801	FAD	C4A-C5A-N7A	-2.86	107.31	110.58
2	A	801	FAD	O4'-C4'-C5'	2.75	116.06	109.99
2	A	801	FAD	C4-C4X-N5	2.54	121.72	118.21
2	A	801	FAD	C5X-C9A-N10	2.54	120.26	117.97
2	A	801	FAD	C5'-C4'-C3'	-2.47	107.56	112.22
2	A	801	FAD	O2P-P-O3P	-2.42	100.74	107.27
2	A	801	FAD	C4-N3-C2	-2.35	121.47	125.64
2	A	801	FAD	C4A-N9A-C8A	2.28	108.13	105.74
2	A	801	FAD	C2A-N1A-C6A	2.15	122.26	118.73
2	A	801	FAD	C1'-N10-C9A	2.07	124.66	120.63
2	A	801	FAD	C4X-C4-N3	2.06	118.50	113.25
6	A	823	GOL	O3-C3-C2	-2.05	101.16	110.38

There are no chirality outliers.

All (47) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	801	FAD	C1'-C2'-C3'-O3'
2	A	801	FAD	C1'-C2'-C3'-C4'
2	A	801	FAD	O2'-C2'-C3'-O3'
2	A	801	FAD	O2'-C2'-C3'-C4'
6	A	822	GOL	C1-C2-C3-O3
6	A	823	GOL	O1-C1-C2-C3
5	A	816	PEG	O2-C3-C4-O4
5	A	819	PEG	O2-C3-C4-O4
5	A	820	PEG	O1-C1-C2-O2
5	A	828	PEG	O2-C3-C4-O4
6	A	823	GOL	C1-C2-C3-O3
6	A	824	GOL	O1-C1-C2-C3
6	A	825	GOL	C1-C2-C3-O3
6	A	822	GOL	O2-C2-C3-O3

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Mol	Chain	Res	Type	Atoms
2	A	801	FAD	C2'-C3'-C4'-O4'
4	A	812	P6G	O16-C17-C18-O19
4	A	809	P6G	O10-C11-C12-O13
4	A	811	P6G	O16-C17-C18-O19
5	A	821	PEG	O1-C1-C2-O2
6	A	823	GOL	O2-C2-C3-O3
6	A	824	GOL	O1-C1-C2-O2
2	A	801	FAD	O4B-C4B-C5B-O5B
5	A	814	PEG	O2-C3-C4-O4
2	A	801	FAD	C3B-C4B-C5B-O5B
4	A	813	P6G	O10-C11-C12-O13
6	A	823	GOL	O1-C1-C2-O2
5	A	821	PEG	C1-C2-O2-C3
5	A	819	PEG	C1-C2-O2-C3
4	A	809	P6G	C5-C6-O7-C8
4	A	812	P6G	C8-C9-O10-C11
5	A	814	PEG	C1-C2-O2-C3
4	A	813	P6G	O16-C17-C18-O19
5	A	818	PEG	C1-C2-O2-C3
4	A	812	P6G	C11-C12-O13-C14
5	A	821	PEG	C4-C3-O2-C2
4	A	813	P6G	C18-C17-O16-C15
2	A	801	FAD	C2'-C3'-C4'-C5'
5	A	818	PEG	C4-C3-O2-C2
5	A	815	PEG	C1-C2-O2-C3
6	A	825	GOL	O2-C2-C3-O3
4	A	813	P6G	O13-C14-C15-O16
4	A	809	P6G	O7-C8-C9-O10
6	A	822	GOL	O1-C1-C2-C3
5	A	815	PEG	C4-C3-O2-C2
4	A	810	P6G	O7-C8-C9-O10
5	A	815	PEG	O1-C1-C2-O2
4	A	809	P6G	C12-C11-O10-C9

There are no ring outliers.

5 monomers are involved in 8 short contacts:

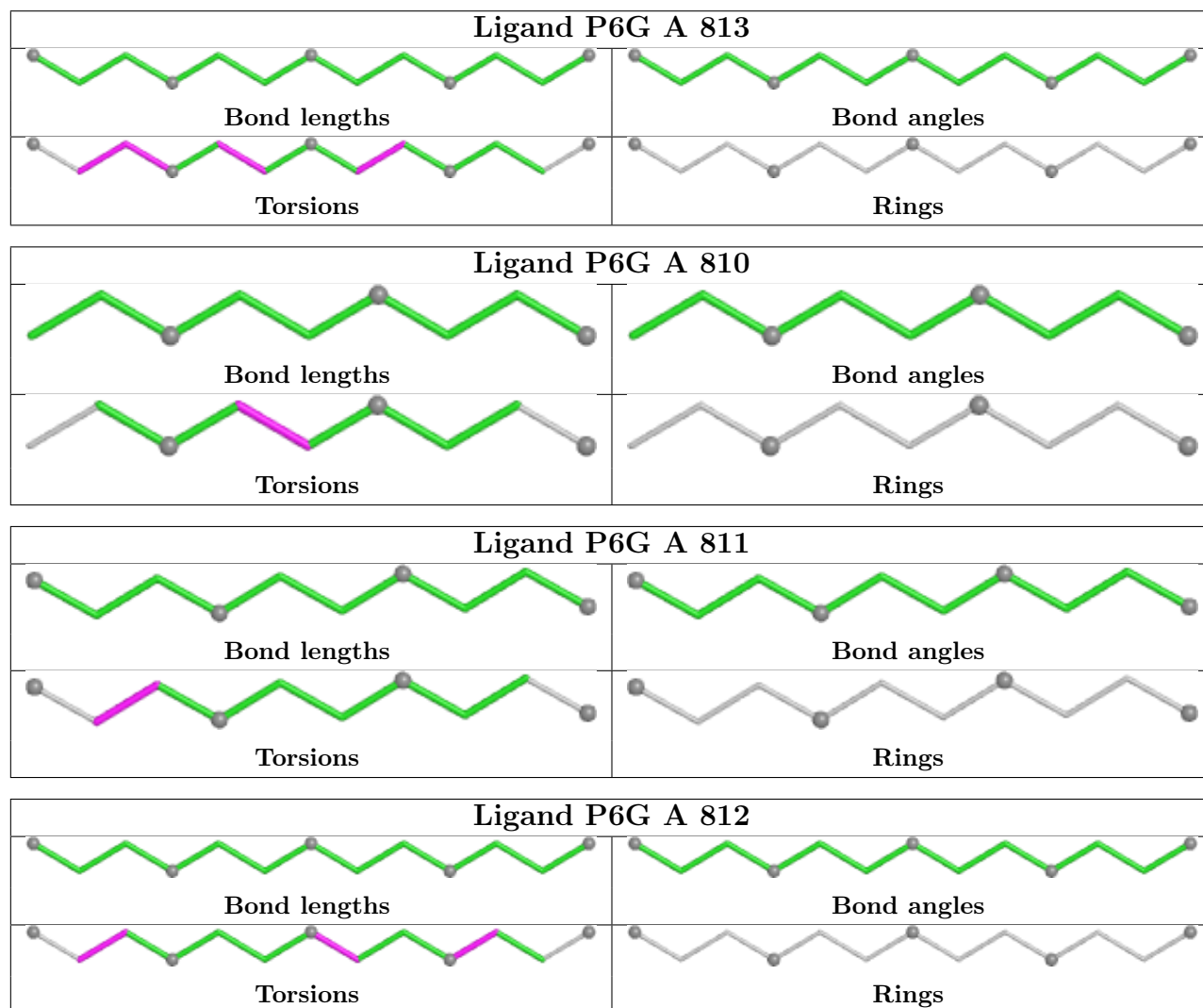
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	822	GOL	1	0
4	A	813	P6G	1	1
5	A	816	PEG	2	0
5	A	818	PEG	2	0

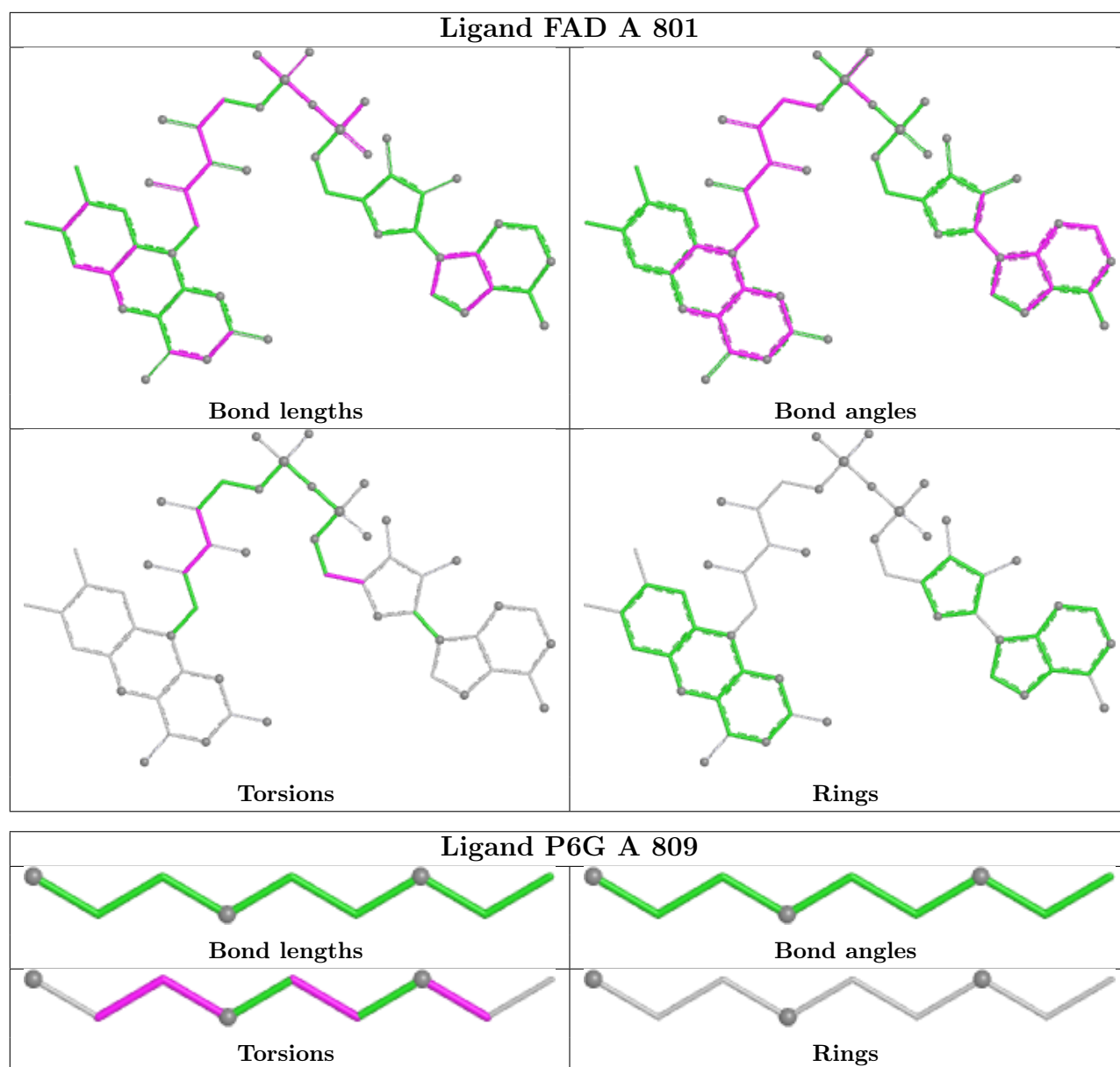
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	801	FAD	1	0

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





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

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	261/291 (89%)	0.40	31 (11%) 9 7	25, 40, 86, 104	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	634	THR	7.0
1	A	504	SER	6.5
1	A	506	SER	6.5
1	A	411	GLY	6.3
1	A	507	LEU	5.9
1	A	503	SER	5.8
1	A	509	GLY	5.2
1	A	508	PRO	5.1
1	A	633	ILE	5.1
1	A	635	GLY	4.4
1	A	505	GLN	4.2
1	A	632	LEU	4.2
1	A	556	GLN	3.9
1	A	698	LEU	3.9
1	A	559	ALA	3.7
1	A	499	ILE	3.6
1	A	636	LEU	3.5
1	A	414	PRO	3.1
1	A	590	ALA	2.9
1	A	638	SER	2.7
1	A	498	TRP	2.7
1	A	555	GLN	2.5
1	A	497	GLU	2.5
1	A	558	PRO	2.5
1	A	502	GLY	2.4
1	A	416	PHE	2.4
1	A	639	ARG	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	489	GLN	2.3
1	A	637	LYS	2.2
1	A	412	SER	2.1
1	A	586	SER	2.1

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	PEG	A	828	7/7	0.72	0.31	80,83,91,91	0
5	PEG	A	816	7/7	0.79	0.26	61,66,69,70	0
5	PEG	A	814	7/7	0.79	0.24	69,77,81,83	0
4	P6G	A	813	13/19	0.84	0.27	60,85,99,99	0
5	PEG	A	818	7/7	0.85	0.27	70,75,80,83	0
6	GOL	A	826	6/6	0.87	0.20	82,86,86,98	0
4	P6G	A	809	9/19	0.88	0.20	56,61,68,69	0
6	GOL	A	823	6/6	0.89	0.16	51,55,58,62	0
6	GOL	A	827	6/6	0.89	0.17	73,75,79,80	0
6	GOL	A	825	6/6	0.90	0.21	75,80,86,92	0
5	PEG	A	820	7/7	0.90	0.23	70,77,87,89	0
5	PEG	A	821	7/7	0.90	0.18	52,72,88,91	0
4	P6G	A	810	9/19	0.91	0.17	49,57,66,70	0
5	PEG	A	815	7/7	0.91	0.17	66,73,78,78	0
5	PEG	A	819	7/7	0.91	0.16	74,76,80,81	0
6	GOL	A	822	6/6	0.91	0.17	65,72,77,79	0
3	CL	A	807	1/1	0.92	0.12	97,97,97,97	0
3	CL	A	804	1/1	0.92	0.13	88,88,88,88	0
3	CL	A	806	1/1	0.92	0.16	101,101,101,101	0

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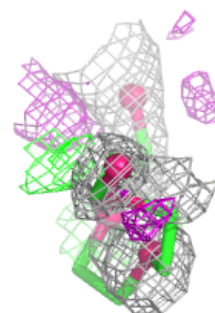
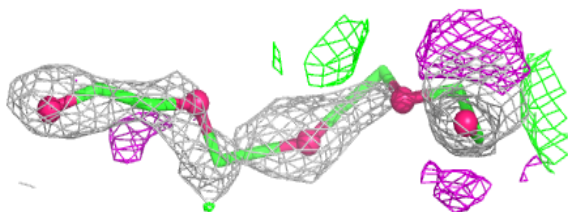
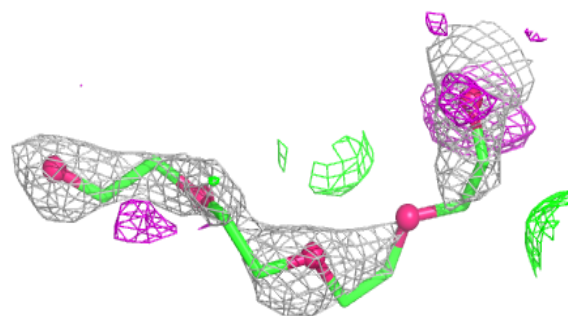
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	P6G	A	812	13/19	0.92	0.16	54,64,71,73	0
5	PEG	A	817	7/7	0.92	0.15	61,65,77,78	0
3	CL	A	808	1/1	0.93	0.20	82,82,82,82	0
3	CL	A	803	1/1	0.93	0.15	87,87,87,87	0
6	GOL	A	824	6/6	0.93	0.18	75,79,85,86	0
4	P6G	A	811	10/19	0.94	0.15	42,68,73,74	0
3	CL	A	805	1/1	0.94	0.11	83,83,83,83	0
3	CL	A	802	1/1	0.95	0.11	82,82,82,82	0
2	FAD	A	801	53/53	0.96	0.09	25,32,74,85	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.

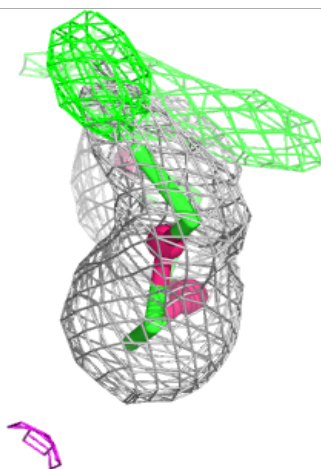
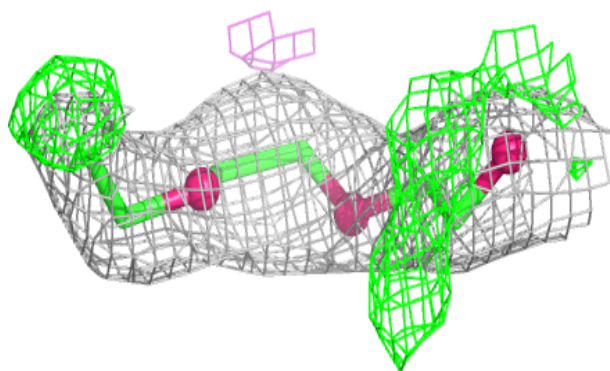
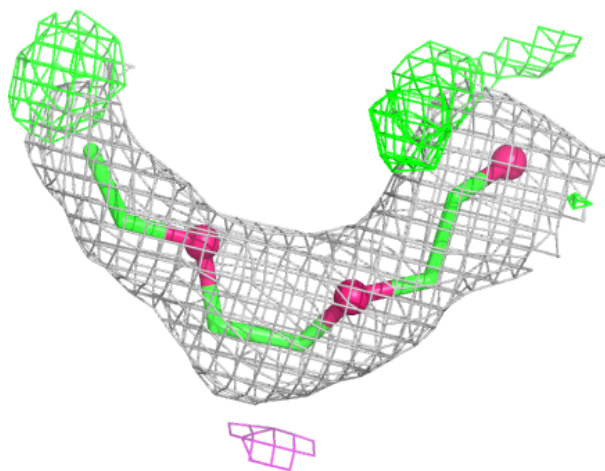
Electron density around P6G A 813:

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



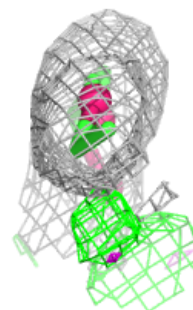
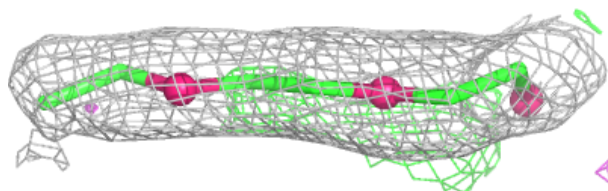
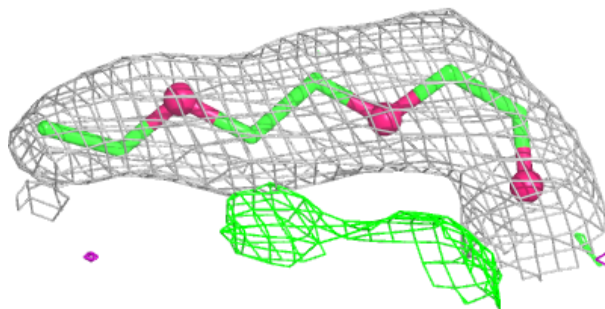
Electron density around P6G A 809:

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

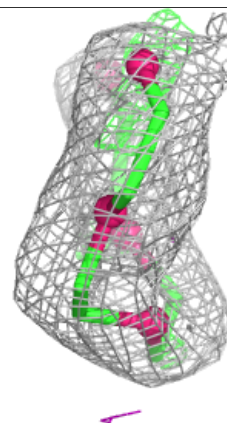
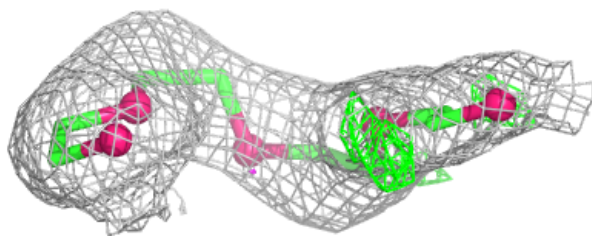
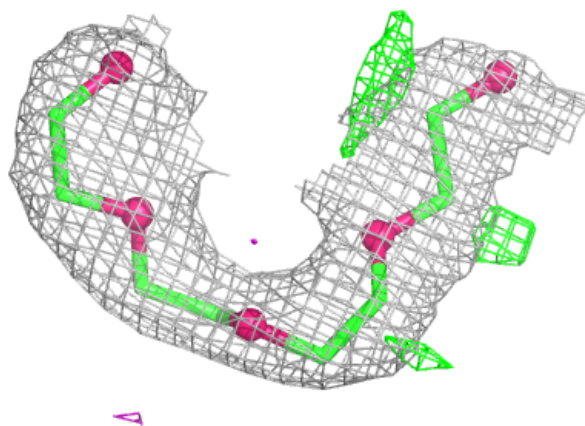


Electron density around P6G A 810:

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

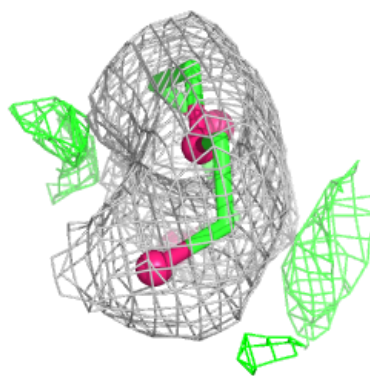
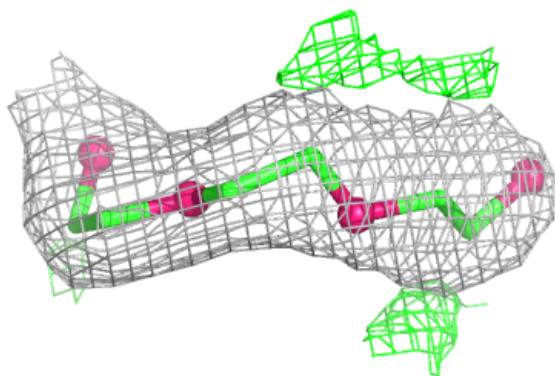
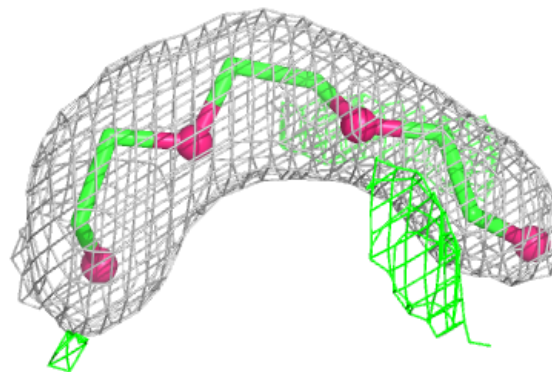
**Electron density around P6G A 812:**

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

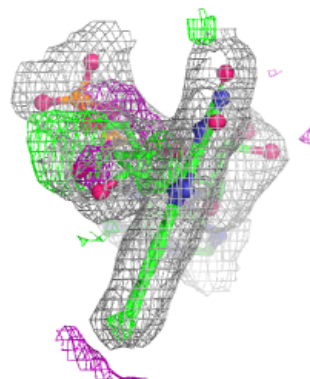
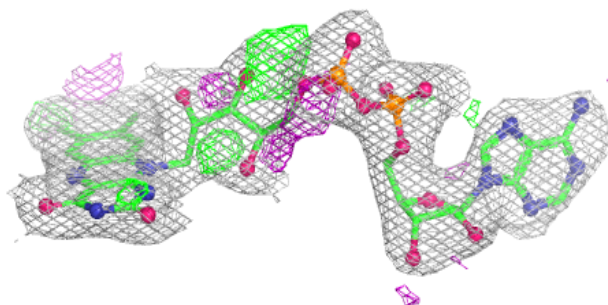
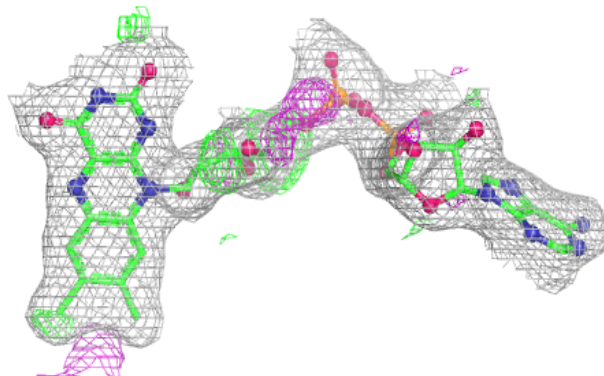


Electron density around P6G A 811:

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

**Electron density around FAD A 801:**

$2mF_o-DF_c$ (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.