



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 28, 2021 – 12:16 PM JST

PDB ID : 7C5M  
Title : Crystal Structure of C150A+H177A mutant of Glyceraldehyde-3-phosphate-dehydrogenase1 from Escherichia coli complexed with G3P at 1.8 Angstrom resolution  
Authors : Zhang, L.; Liu, M.R.; Bao, L.Y.; Yao, Y.C.; Bostrom, I.K.; Wang, Y.D.; Chen, A.Q.; Li, J.X.; Gu, S.H.; Ji, C.N.  
Deposited on : 2020-05-20  
Resolution : 1.80 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](mailto: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.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.23.2  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.23.2

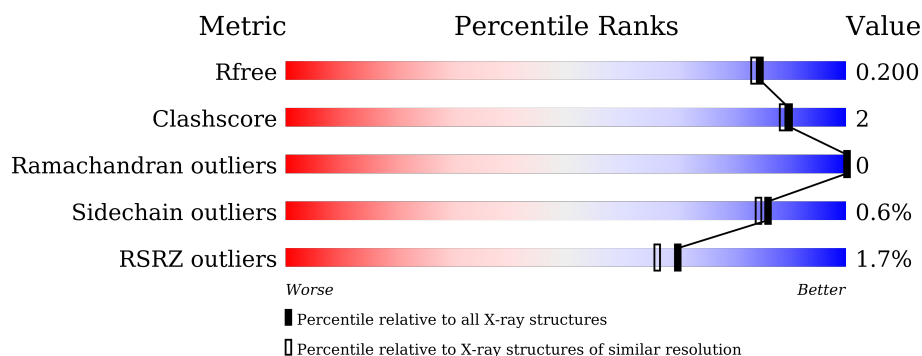
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.80 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))
$R_{free}$	130704	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

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  $\geq 3$ , 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	O	352	<div><div></div><div>91%</div><div>5%</div></div>
1	P	352	<div><div></div><div>89%</div><div>5%</div><div>5%</div></div>
1	Q	352	<div><div></div><div>91%</div><div>5%</div></div>
1	R	352	<div><div>6%</div><div>88%</div><div>6%</div><div>5%</div></div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	G3H	P	402	-	X	-	-
5	G3H	Q	402	-	X	-	-
5	G3H	R	402	-	X	-	-

## 2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 11850 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 Glyceraldehyde-3-phosphate dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	O	334	Total	C	N	O	S	0	2	0
			2531	1599	425	501	6			
1	P	334	Total	C	N	O	S	0	2	1
			2528	1598	424	500	6			
1	Q	336	Total	C	N	O	S	0	2	1
			2538	1605	426	501	6			
1	R	333	Total	C	N	O	S	0	2	0
			2528	1599	426	497	6			

There are 84 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
O	-18	HIS	-	expression tag	UNP A0A140NCK4
O	-17	HIS	-	expression tag	UNP A0A140NCK4
O	-16	HIS	-	expression tag	UNP A0A140NCK4
O	-15	HIS	-	expression tag	UNP A0A140NCK4
O	-14	HIS	-	expression tag	UNP A0A140NCK4
O	-13	HIS	-	expression tag	UNP A0A140NCK4
O	-12	SER	-	expression tag	UNP A0A140NCK4
O	-11	SER	-	expression tag	UNP A0A140NCK4
O	-10	GLY	-	expression tag	UNP A0A140NCK4
O	-9	LEU	-	expression tag	UNP A0A140NCK4
O	-8	VAL	-	expression tag	UNP A0A140NCK4
O	-7	PRO	-	expression tag	UNP A0A140NCK4
O	-6	ARG	-	expression tag	UNP A0A140NCK4
O	-5	GLY	-	expression tag	UNP A0A140NCK4
O	-4	SER	-	expression tag	UNP A0A140NCK4
O	-3	HIS	-	expression tag	UNP A0A140NCK4
O	-2	MET	-	expression tag	UNP A0A140NCK4
O	-1	ALA	-	expression tag	UNP A0A140NCK4
O	0	SER	-	expression tag	UNP A0A140NCK4
O	150	ALA	CYS	engineered mutation	UNP A0A140NCK4
O	177	ALA	HIS	engineered mutation	UNP A0A140NCK4

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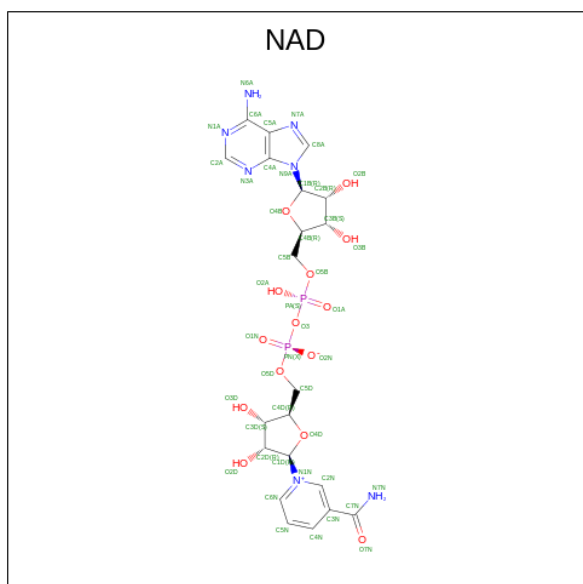
Chain	Residue	Modelled	Actual	Comment	Reference
P	-18	HIS	-	expression tag	UNP A0A140NCK4
P	-17	HIS	-	expression tag	UNP A0A140NCK4
P	-16	HIS	-	expression tag	UNP A0A140NCK4
P	-15	HIS	-	expression tag	UNP A0A140NCK4
P	-14	HIS	-	expression tag	UNP A0A140NCK4
P	-13	HIS	-	expression tag	UNP A0A140NCK4
P	-12	SER	-	expression tag	UNP A0A140NCK4
P	-11	SER	-	expression tag	UNP A0A140NCK4
P	-10	GLY	-	expression tag	UNP A0A140NCK4
P	-9	LEU	-	expression tag	UNP A0A140NCK4
P	-8	VAL	-	expression tag	UNP A0A140NCK4
P	-7	PRO	-	expression tag	UNP A0A140NCK4
P	-6	ARG	-	expression tag	UNP A0A140NCK4
P	-5	GLY	-	expression tag	UNP A0A140NCK4
P	-4	SER	-	expression tag	UNP A0A140NCK4
P	-3	HIS	-	expression tag	UNP A0A140NCK4
P	-2	MET	-	expression tag	UNP A0A140NCK4
P	-1	ALA	-	expression tag	UNP A0A140NCK4
P	0	SER	-	expression tag	UNP A0A140NCK4
P	150	ALA	CYS	engineered mutation	UNP A0A140NCK4
P	177	ALA	HIS	engineered mutation	UNP A0A140NCK4
Q	-18	HIS	-	expression tag	UNP A0A140NCK4
Q	-17	HIS	-	expression tag	UNP A0A140NCK4
Q	-16	HIS	-	expression tag	UNP A0A140NCK4
Q	-15	HIS	-	expression tag	UNP A0A140NCK4
Q	-14	HIS	-	expression tag	UNP A0A140NCK4
Q	-13	HIS	-	expression tag	UNP A0A140NCK4
Q	-12	SER	-	expression tag	UNP A0A140NCK4
Q	-11	SER	-	expression tag	UNP A0A140NCK4
Q	-10	GLY	-	expression tag	UNP A0A140NCK4
Q	-9	LEU	-	expression tag	UNP A0A140NCK4
Q	-8	VAL	-	expression tag	UNP A0A140NCK4
Q	-7	PRO	-	expression tag	UNP A0A140NCK4
Q	-6	ARG	-	expression tag	UNP A0A140NCK4
Q	-5	GLY	-	expression tag	UNP A0A140NCK4
Q	-4	SER	-	expression tag	UNP A0A140NCK4
Q	-3	HIS	-	expression tag	UNP A0A140NCK4
Q	-2	MET	-	expression tag	UNP A0A140NCK4
Q	-1	ALA	-	expression tag	UNP A0A140NCK4
Q	0	SER	-	expression tag	UNP A0A140NCK4
Q	150	ALA	CYS	engineered mutation	UNP A0A140NCK4
Q	177	ALA	HIS	engineered mutation	UNP A0A140NCK4

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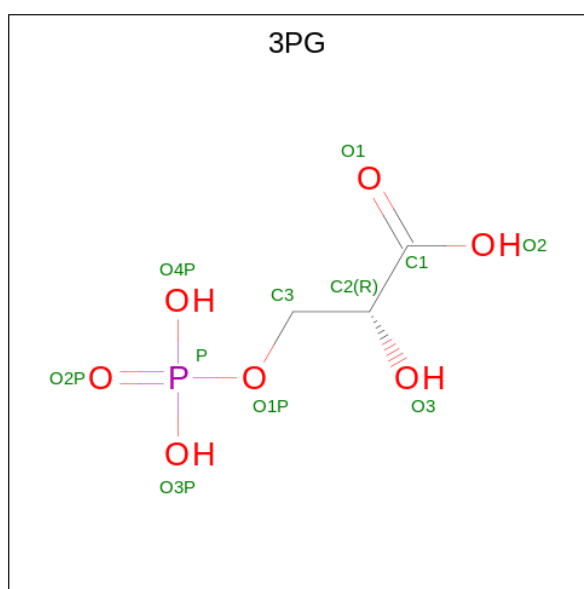
Chain	Residue	Modelled	Actual	Comment	Reference
R	-18	HIS	-	expression tag	UNP A0A140NCK4
R	-17	HIS	-	expression tag	UNP A0A140NCK4
R	-16	HIS	-	expression tag	UNP A0A140NCK4
R	-15	HIS	-	expression tag	UNP A0A140NCK4
R	-14	HIS	-	expression tag	UNP A0A140NCK4
R	-13	HIS	-	expression tag	UNP A0A140NCK4
R	-12	SER	-	expression tag	UNP A0A140NCK4
R	-11	SER	-	expression tag	UNP A0A140NCK4
R	-10	GLY	-	expression tag	UNP A0A140NCK4
R	-9	LEU	-	expression tag	UNP A0A140NCK4
R	-8	VAL	-	expression tag	UNP A0A140NCK4
R	-7	PRO	-	expression tag	UNP A0A140NCK4
R	-6	ARG	-	expression tag	UNP A0A140NCK4
R	-5	GLY	-	expression tag	UNP A0A140NCK4
R	-4	SER	-	expression tag	UNP A0A140NCK4
R	-3	HIS	-	expression tag	UNP A0A140NCK4
R	-2	MET	-	expression tag	UNP A0A140NCK4
R	-1	ALA	-	expression tag	UNP A0A140NCK4
R	0	SER	-	expression tag	UNP A0A140NCK4
R	150	ALA	CYS	engineered mutation	UNP A0A140NCK4
R	177	ALA	HIS	engineered mutation	UNP A0A140NCK4

- Molecule 2 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula:  $C_{21}H_{27}N_7O_{14}P_2$ ) (labeled as "Ligand of Interest" by depositor).



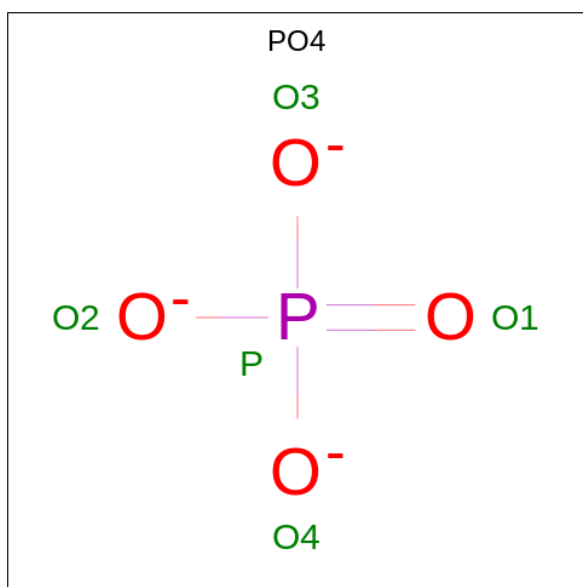
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	O	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	P	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	Q	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	R	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

- Molecule 3 is 3-PHOSPHOGLYCERIC ACID (three-letter code: 3PG) (formula:  $C_3H_7O_7P$ ) (labeled as "Ligand of Interest" by depositor).



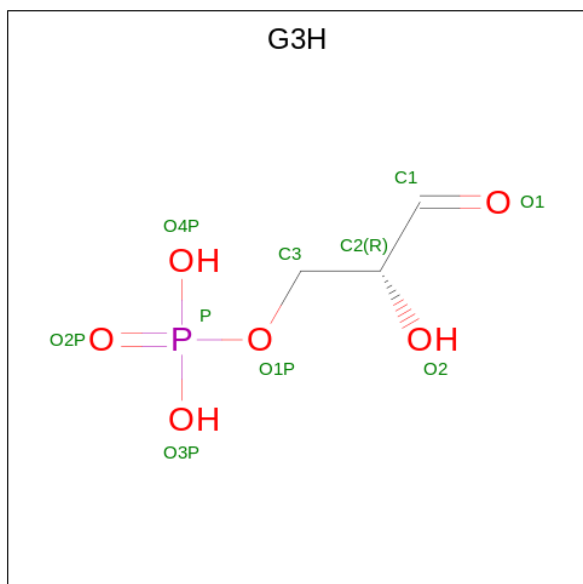
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	O	1	Total	C	O	P	0	0
			11	3	7	1		

- Molecule 4 is PHOSPHATE ION (three-letter code: PO4) (formula:  $O_4P$ ) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	O	1	Total	O	P	0	0
			5	4	1		
4	O	1	Total	O	P	0	0
			5	4	1		

- Molecule 5 is GLYCERALDEHYDE-3-PHOSPHATE (three-letter code: G3H) (formula:  $C_3H_7O_6P$ ) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	P	1	Total	C	O	P	0	0
			10	3	6	1		

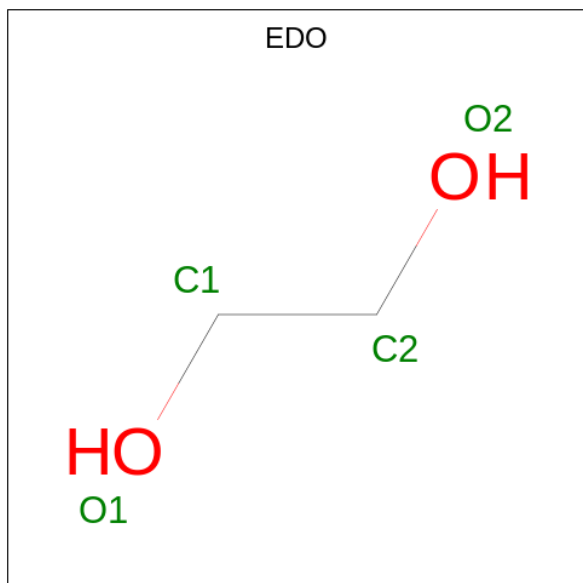
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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	Q	1	Total	C	O	P	0	0
			10	3	6	1		
5	R	1	Total	C	O	P	0	0
			10	3	6	1		

- Molecule 6 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula:  $C_2H_6O_2$ ) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	P	1	Total	C	O	0	1
			8	4	4		
6	Q	1	Total	C	O	0	0
			4	2	2		
6	Q	1	Total	C	O	0	1
			8	4	4		
6	R	1	Total	C	O	0	0
			4	2	2		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	O	419	Total	O	0	0
			419	419		
7	P	388	Total	O	0	0
			388	388		
7	Q	375	Total	O	0	0
			375	375		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	R	292	Total 292	O 292	0	0

### 3 Residue-property plots [i](#)


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

- Molecule 1: Glyceraldehyde-3-phosphate dehydrogenase

Chain O:  91% 5%

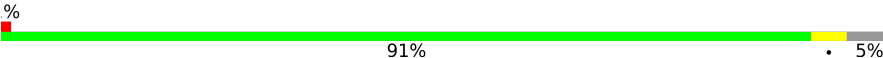


- Molecule 1: Glyceraldehyde-3-phosphate dehydrogenase

Chain P:  89% 5% 5%




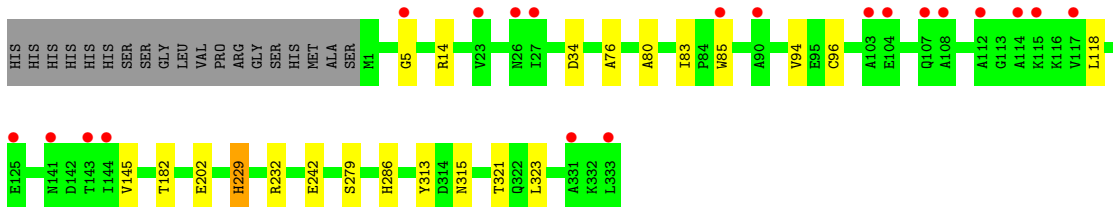
- Molecule 1: Glyceraldehyde-3-phosphate dehydrogenase

Chain Q:  91% 5%



- Molecule 1: Glyceraldehyde-3-phosphate dehydrogenase

Chain R:  88% 6% 5%



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	89.92Å 89.92Å 342.38Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.00 – 1.80 44.96 – 1.80	Depositor EDS
% Data completeness (in resolution range)	99.8 (45.00-1.80) 99.8 (44.96-1.80)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.55 (at 1.79Å)	Xtriage
Refinement program	REFMAC 5.8.0131	Depositor
R, $R_{free}$	0.145 , 0.189 0.159 , 0.200	Depositor DCC
$R_{free}$ test set	6457 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	22.7	Xtriage
Anisotropy	0.027	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 47.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	11850	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.17% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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, PO4, 3PG, NAD, G3H

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	O	0.97	0/2571	1.01	12/3489 (0.3%)
1	P	0.95	2/2568 (0.1%)	1.01	10/3484 (0.3%)
1	Q	0.94	1/2578 (0.0%)	1.00	7/3498 (0.2%)
1	R	0.93	1/2569 (0.0%)	0.95	3/3486 (0.1%)
All	All	0.95	4/10286 (0.0%)	0.99	32/13957 (0.2%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	P	202	GLU	CD-OE2	8.03	1.34	1.25
1	R	202	GLU	CD-OE2	7.21	1.33	1.25
1	Q	202	GLU	CD-OE2	6.84	1.33	1.25
1	P	179	TYR	CE1-CZ	5.87	1.46	1.38

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	P	193	ASP	CB-CG-OD1	12.06	129.15	118.30
1	Q	18	ARG	NE-CZ-NH1	10.57	125.58	120.30
1	P	18	ARG	NE-CZ-NH2	10.21	125.41	120.30
1	O	193	ASP	CB-CG-OD1	8.36	125.83	118.30
1	P	18	ARG	NE-CZ-NH1	-7.87	116.36	120.30
1	R	14	ARG	NE-CZ-NH2	7.75	124.18	120.30
1	Q	187	ASP	CB-CG-OD1	7.73	125.26	118.30
1	Q	18	ARG	NE-CZ-NH2	-7.47	116.56	120.30
1	P	232	ARG	NE-CZ-NH1	-7.30	116.65	120.30
1	O	139	ASP	CB-CG-OD1	7.24	124.81	118.30
1	O	193	ASP	CB-CG-OD2	-7.06	111.94	118.30
1	O	195	ARG	NE-CZ-NH1	-6.80	116.90	120.30
1	P	187	ASP	CB-CG-OD1	6.70	124.33	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	P	193	ASP	CB-CG-OD2	-6.64	112.33	118.30
1	P	187	ASP	CB-CG-OD2	-6.58	112.38	118.30
1	O	232	ARG	NE-CZ-NH1	-6.54	117.03	120.30
1	P	14	ARG	NE-CZ-NH1	-6.49	117.06	120.30
1	R	232	ARG	NE-CZ-NH2	6.27	123.43	120.30
1	Q	187	ASP	CB-CG-OD2	-5.82	113.06	118.30
1	Q	14	ARG	NE-CZ-NH2	5.79	123.19	120.30
1	O	59	ASP	CB-CG-OD2	-5.75	113.13	118.30
1	O	274	ASP	CB-CG-OD1	5.62	123.36	118.30
1	O	232	ARG	NE-CZ-NH2	5.56	123.08	120.30
1	O	14	ARG	NE-CZ-NH2	5.53	123.06	120.30
1	Q	63	ASP	CB-CG-OD1	5.45	123.20	118.30
1	Q	193	ASP	CB-CG-OD1	5.21	122.99	118.30
1	P	19	ARG	NE-CZ-NH1	-5.20	117.70	120.30
1	P	59	ASP	CB-CG-OD2	-5.15	113.67	118.30
1	O	325	ARG	NE-CZ-NH2	5.11	122.86	120.30
1	O	274	ASP	CB-CG-OD2	-5.10	113.71	118.30
1	R	14	ARG	NE-CZ-NH1	-5.06	117.77	120.30
1	O	270	PHE	CB-CG-CD1	5.01	124.31	120.80

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	O	2531	0	2556	3	0
1	P	2528	0	2551	8	0
1	Q	2538	0	2566	9	0
1	R	2528	0	2556	14	0
2	O	44	0	26	2	0
2	P	44	0	26	2	0
2	Q	44	0	26	1	0
2	R	44	0	26	1	0
3	O	11	0	4	1	0
4	O	10	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	P	10	0	5	1	0
5	Q	10	0	5	2	0
5	R	10	0	5	2	0
6	P	8	0	12	3	0
6	Q	12	0	18	1	0
6	R	4	0	6	0	0
7	O	419	0	0	1	1
7	P	388	0	0	4	0
7	Q	375	0	0	3	0
7	R	292	0	0	1	2
All	All	11850	0	10388	37	2

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 (37) 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)
5:Q:402:G3H:H31	7:Q:557:HOH:O	1.53	1.04
1:P:182:THR:OG1	5:P:402:G3H:H32	1.95	0.66
1:Q:229:HIS:HD2	7:Q:814:HOH:O	1.87	0.58
1:Q:315:ASN:HD22	2:Q:401:NAD:H72N	1.56	0.54
1:Q:229:HIS:HE1	1:R:242:GLU:OE2	1.90	0.54
1:R:323:LEU:C	1:R:323:LEU:HD23	2.28	0.53
1:P:118:LEU:HD12	1:P:145:VAL:HG23	1.91	0.53
1:P:118:LEU:CD1	1:P:145:VAL:HG23	2.39	0.52
1:P:323:LEU:C	1:P:323:LEU:HD23	2.29	0.52
1:R:94:VAL:HG12	1:R:96:CYS:SG	2.50	0.51
1:P:18:ARG:HD3	7:P:766:HOH:O	2.11	0.50
1:Q:242:GLU:OE2	1:R:229[B]:HIS:HE1	1.94	0.49
5:Q:402:G3H:H32	7:Q:763:HOH:O	2.13	0.49
1:R:5:GLY:HA3	1:R:85:TRP:CZ3	2.48	0.49
6:P:403[B]:EDO:C1	7:P:805:HOH:O	2.60	0.48
1:R:315:ASN:O	2:R:401:NAD:H4N	2.16	0.46
2:O:401:NAD:C5N	7:O:768:HOH:O	2.64	0.46
1:P:286:HIS:HA	1:P:321:THR:HG21	1.98	0.46
1:O:182:THR:OG1	3:O:402:3PG:H32	2.16	0.45
1:R:229[B]:HIS:HD2	7:R:760:HOH:O	1.98	0.45
1:P:120:SER:O	2:P:401:NAD:H6N	2.17	0.45
1:R:286:HIS:HA	1:R:321:THR:HG21	2.00	0.44
1:Q:242:GLU:OE2	1:R:229[A]:HIS:HE1	2.01	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:P:403[B]:EDO:H11	7:P:805:HOH:O	2.18	0.43
1:Q:204:ILE:HB	1:R:279:SER:HB3	2.01	0.43
1:R:118:LEU:CD1	1:R:145:VAL:HG23	2.48	0.43
1:Q:19:ARG:HH12	6:Q:404[A]:EDO:H21	1.84	0.42
1:R:80:ALA:HA	1:R:83:ILE:CD1	2.49	0.42
1:O:286:HIS:HA	1:O:321:THR:HG21	2.02	0.41
1:P:315:ASN:O	2:P:401:NAD:H4N	2.20	0.41
1:Q:286:HIS:HA	1:Q:321:THR:HG21	2.03	0.41
1:Q:323:LEU:C	1:Q:323:LEU:HD23	2.41	0.41
1:R:182:THR:OG1	5:R:402:G3H:H32	2.20	0.41
1:O:315:ASN:O	2:O:401:NAD:H4N	2.21	0.41
6:P:403[B]:EDO:H12	7:P:805:HOH:O	2.21	0.41
1:R:34:ASP:O	1:R:76:ALA:HA	2.22	0.40
5:R:402:G3H:H2	5:R:402:G3H:O4P	2.21	0.40

All (2) 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)
7:O:803:HOH:O	7:R:629:HOH:O[1_655]	1.96	0.24
7:R:532:HOH:O	7:R:532:HOH:O[7_465]	2.17	0.03

## 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	O	334/352 (95%)	322 (96%)	12 (4%)	0	100	100
1	P	334/352 (95%)	322 (96%)	12 (4%)	0	100	100
1	Q	336/352 (96%)	325 (97%)	11 (3%)	0	100	100
1	R	333/352 (95%)	320 (96%)	13 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	1337/1408 (95%)	1289 (96%)	48 (4%)	0	100	100

There are no Ramachandran outliers to report.

### 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	O	274/287 (96%)	272 (99%)	2 (1%)	84	81
1	P	273/287 (95%)	271 (99%)	2 (1%)	84	81
1	Q	274/287 (96%)	273 (100%)	1 (0%)	91	89
1	R	273/287 (95%)	270 (99%)	3 (1%)	73	68
All	All	1094/1148 (95%)	1086 (99%)	8 (1%)	86	81

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	O	209	THR
1	O	229	HIS
1	P	209	THR
1	P	229	HIS
1	Q	229	HIS
1	R	229[A]	HIS
1	R	229[B]	HIS
1	R	313	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	Q	229	HIS

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

## 5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

## 5.6 Ligand geometry ⓘ

16 ligands are modelled in this entry.

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	G3H	Q	402	-	8,9,9	2.28	3 (37%)	10,12,12	3.36	7 (70%)
4	PO4	O	403	-	4,4,4	0.87	0	6,6,6	2.21	2 (33%)
5	G3H	P	402	-	8,9,9	2.32	2 (25%)	10,12,12	3.62	4 (40%)
6	EDO	P	403[B]	-	3,3,3	0.48	0	2,2,2	0.34	0
6	EDO	Q	404[B]	-	3,3,3	0.43	0	2,2,2	0.71	0
2	NAD	R	401	-	42,48,48	1.45	5 (11%)	50,73,73	1.71	12 (24%)
4	PO4	O	404	-	4,4,4	0.71	0	6,6,6	0.97	0
2	NAD	P	401	-	42,48,48	1.18	6 (14%)	50,73,73	1.43	8 (16%)
2	NAD	Q	401	-	42,48,48	1.30	7 (16%)	50,73,73	1.66	10 (20%)
6	EDO	R	403	-	3,3,3	0.48	0	2,2,2	0.82	0
6	EDO	Q	403	-	3,3,3	0.46	0	2,2,2	0.47	0
2	NAD	O	401	-	42,48,48	1.29	5 (11%)	50,73,73	1.54	9 (18%)
6	EDO	P	403[A]	-	3,3,3	0.55	0	2,2,2	0.49	0
6	EDO	Q	404[A]	-	3,3,3	0.46	0	2,2,2	0.44	0
3	3PG	O	402	-	7,10,10	1.47	1 (14%)	10,14,14	2.37	4 (40%)
5	G3H	R	402	-	8,9,9	1.88	2 (25%)	10,12,12	2.36	5 (50%)

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	G3H	Q	402	-	-	4/7/8/8	-
5	G3H	P	402	-	-	7/7/8/8	-
6	EDO	P	403[B]	-	-	1/1/1/1	-
6	EDO	Q	404[B]	-	-	1/1/1/1	-
2	NAD	R	401	-	-	5/26/62/62	0/5/5/5
2	NAD	P	401	-	-	5/26/62/62	0/5/5/5
2	NAD	Q	401	-	-	5/26/62/62	0/5/5/5
6	EDO	R	403	-	-	1/1/1/1	-
6	EDO	Q	403	-	-	1/1/1/1	-
2	NAD	O	401	-	-	5/26/62/62	0/5/5/5
6	EDO	P	403[A]	-	-	0/1/1/1	-
6	EDO	Q	404[A]	-	-	1/1/1/1	-
3	3PG	O	402	-	-	6/6/10/10	-
5	G3H	R	402	-	-	6/7/8/8	-

All (31) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	P	402	G3H	C3-C2	5.28	1.59	1.51
2	R	401	NAD	C2A-N3A	4.89	1.40	1.32
5	Q	402	G3H	P-O1P	-4.30	1.46	1.60
5	R	402	G3H	C3-C2	4.24	1.57	1.51
2	O	401	NAD	C2N-C3N	3.79	1.44	1.39
2	Q	401	NAD	C2A-N3A	3.35	1.37	1.32
2	Q	401	NAD	C7N-N7N	-3.23	1.26	1.33
5	Q	402	G3H	O1P-C3	-2.90	1.33	1.44
2	Q	401	NAD	C2N-C3N	2.83	1.43	1.39
2	R	401	NAD	O4B-C1B	2.83	1.45	1.41
2	Q	401	NAD	C8A-N7A	2.82	1.39	1.34
2	P	401	NAD	C2A-N3A	2.80	1.36	1.32
2	P	401	NAD	C5A-C4A	2.68	1.48	1.40
2	O	401	NAD	C2N-N1N	2.67	1.38	1.35
5	P	402	G3H	O1P-C3	2.64	1.54	1.44
2	R	401	NAD	C2D-C1D	-2.63	1.49	1.53
5	R	402	G3H	P-O2P	2.47	1.58	1.50
2	R	401	NAD	C5A-C4A	2.44	1.47	1.40
5	Q	402	G3H	C3-C2	2.39	1.54	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	O	401	NAD	O4B-C1B	2.35	1.44	1.41
2	Q	401	NAD	C5A-C4A	2.31	1.47	1.40
2	R	401	NAD	C2N-C3N	2.29	1.42	1.39
2	P	401	NAD	O4B-C4B	-2.22	1.40	1.45
2	P	401	NAD	C2A-N1A	2.22	1.38	1.33
2	P	401	NAD	O4D-C1D	2.12	1.44	1.41
2	P	401	NAD	PA-O2A	-2.11	1.45	1.55
2	Q	401	NAD	C5N-C4N	2.09	1.43	1.38
3	O	402	3PG	O1P-C3	2.07	1.52	1.44
2	O	401	NAD	C5N-C4N	2.06	1.43	1.38
2	O	401	NAD	C2A-N3A	2.03	1.35	1.32
2	Q	401	NAD	C2A-N1A	2.02	1.37	1.33

All (61) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	P	402	G3H	O1P-C3-C2	10.04	134.96	108.33
2	Q	401	NAD	N3A-C2A-N1A	-6.20	118.98	128.68
5	Q	402	G3H	O2-C2-C1	5.63	120.26	109.17
5	Q	402	G3H	P-O1P-C3	5.61	133.74	118.30
5	Q	402	G3H	O2-C2-C3	4.66	124.14	110.17
2	P	401	NAD	N3A-C2A-N1A	-4.53	121.60	128.68
5	R	402	G3H	O1P-C3-C2	4.45	120.12	108.33
2	R	401	NAD	N6A-C6A-N1A	4.35	127.61	118.57
3	O	402	3PG	P-O1P-C3	4.29	130.11	118.30
2	R	401	NAD	C1B-N9A-C4A	-4.28	119.12	126.64
3	O	402	3PG	O3-C2-C3	4.27	124.12	110.40
2	Q	401	NAD	C1B-N9A-C4A	-4.17	119.31	126.64
2	O	401	NAD	N3A-C2A-N1A	-3.98	122.46	128.68
2	O	401	NAD	PN-O3-PA	-3.75	119.95	132.83
4	O	403	PO4	O3-P-O2	3.72	119.90	107.97
2	R	401	NAD	N3A-C2A-N1A	-3.49	123.23	128.68
2	R	401	NAD	PN-O3-PA	-3.46	120.96	132.83
2	R	401	NAD	O2A-PA-O1A	3.30	128.57	112.24
2	O	401	NAD	O7N-C7N-C3N	-3.25	115.74	119.63
5	R	402	G3H	P-O1P-C3	3.24	127.22	118.30
2	P	401	NAD	C6N-N1N-C2N	-3.20	119.06	121.97
2	P	401	NAD	PN-O3-PA	-3.18	121.93	132.83
2	Q	401	NAD	O7N-C7N-C3N	3.14	123.39	119.63
2	R	401	NAD	C2A-N1A-C6A	3.12	124.09	118.75
5	P	402	G3H	O2-C2-C1	-3.06	103.14	109.17
5	R	402	G3H	O3P-P-O1P	3.04	114.82	106.73

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	O	401	NAD	C1B-N9A-C4A	-3.01	121.35	126.64
5	Q	402	G3H	O1P-P-O2P	-3.01	98.03	106.47
3	O	402	3PG	O1P-C3-C2	2.90	116.30	107.94
2	O	401	NAD	N6A-C6A-N1A	2.79	124.37	118.57
2	Q	401	NAD	C2A-N1A-C6A	2.77	123.50	118.75
4	O	403	PO4	O3-P-O1	-2.77	100.77	110.89
5	P	402	G3H	P-O1P-C3	2.65	125.60	118.30
5	P	402	G3H	O2-C2-C3	2.64	118.09	110.17
2	O	401	NAD	C4A-C5A-N7A	2.57	112.08	109.40
2	R	401	NAD	C5A-C6A-N6A	-2.53	116.50	120.35
2	Q	401	NAD	C3N-C7N-N7N	-2.51	114.74	117.75
2	Q	401	NAD	N6A-C6A-N1A	2.50	123.77	118.57
2	O	401	NAD	C3N-C7N-N7N	2.48	120.73	117.75
2	Q	401	NAD	C5A-C6A-N6A	-2.48	116.59	120.35
2	R	401	NAD	C6N-N1N-C2N	-2.47	119.72	121.97
3	O	402	3PG	O3P-P-O1P	2.42	113.17	106.73
2	P	401	NAD	C1B-N9A-C4A	-2.41	122.40	126.64
5	Q	402	G3H	O4P-P-O3P	2.40	116.81	107.64
2	R	401	NAD	O5D-PN-O1N	-2.35	99.89	109.07
2	P	401	NAD	O2A-PA-O1A	2.31	123.66	112.24
5	R	402	G3H	O4P-P-O1P	-2.27	100.70	106.73
2	P	401	NAD	N6A-C6A-N1A	2.26	123.26	118.57
5	Q	402	G3H	O4P-P-O1P	-2.25	100.73	106.73
2	Q	401	NAD	PN-O3-PA	-2.25	125.10	132.83
5	Q	402	G3H	O3P-P-O1P	2.23	112.67	106.73
2	P	401	NAD	C3N-C2N-N1N	2.22	122.60	120.43
2	O	401	NAD	O2A-PA-O1A	2.21	123.18	112.24
2	Q	401	NAD	O5D-PN-O1N	-2.18	100.56	109.07
2	O	401	NAD	C5A-C6A-N6A	-2.16	117.07	120.35
2	P	401	NAD	O2N-PN-O1N	2.12	122.70	112.24
5	R	402	G3H	O2-C2-C3	2.09	116.44	110.17
2	R	401	NAD	O4D-C1D-C2D	2.05	109.93	106.93
2	R	401	NAD	C2N-N1N-C1D	2.04	123.69	119.14
2	Q	401	NAD	O2N-PN-O1N	2.03	122.25	112.24
2	R	401	NAD	O4D-C4D-C3D	2.00	109.07	105.11

There are no chirality outliers.

All (48) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	O	401	NAD	O4D-C1D-N1N-C2N
2	O	401	NAD	O4D-C1D-N1N-C6N

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Mol	Chain	Res	Type	Atoms
2	O	401	NAD	C2D-C1D-N1N-C2N
2	O	401	NAD	C2D-C1D-N1N-C6N
2	P	401	NAD	O4D-C1D-N1N-C2N
2	P	401	NAD	O4D-C1D-N1N-C6N
2	P	401	NAD	C2D-C1D-N1N-C2N
2	P	401	NAD	C2D-C1D-N1N-C6N
2	Q	401	NAD	O4D-C1D-N1N-C2N
2	Q	401	NAD	O4D-C1D-N1N-C6N
2	Q	401	NAD	C2D-C1D-N1N-C2N
2	Q	401	NAD	C2D-C1D-N1N-C6N
2	R	401	NAD	O4D-C1D-N1N-C2N
2	R	401	NAD	O4D-C1D-N1N-C6N
2	R	401	NAD	C2D-C1D-N1N-C2N
2	R	401	NAD	C2D-C1D-N1N-C6N
3	O	402	3PG	C1-C2-C3-O1P
3	O	402	3PG	O3-C2-C3-O1P
3	O	402	3PG	C2-C3-O1P-P
3	O	402	3PG	C3-O1P-P-O3P
5	P	402	G3H	O1-C1-C2-C3
5	P	402	G3H	C1-C2-C3-O1P
5	P	402	G3H	C2-C3-O1P-P
5	P	402	G3H	C3-O1P-P-O3P
5	P	402	G3H	C3-O1P-P-O4P
5	Q	402	G3H	O1-C1-C2-C3
5	Q	402	G3H	C3-O1P-P-O3P
5	Q	402	G3H	C3-O1P-P-O4P
5	R	402	G3H	C1-C2-C3-O1P
5	R	402	G3H	O2-C2-C3-O1P
5	R	402	G3H	C2-C3-O1P-P
5	R	402	G3H	C3-O1P-P-O3P
6	P	403[B]	EDO	O1-C1-C2-O2
6	Q	404[A]	EDO	O1-C1-C2-O2
6	Q	404[B]	EDO	O1-C1-C2-O2
3	O	402	3PG	C3-O1P-P-O2P
5	Q	402	G3H	C3-O1P-P-O2P
5	R	402	G3H	C3-O1P-P-O2P
5	P	402	G3H	O2-C2-C3-O1P
2	Q	401	NAD	O4B-C4B-C5B-O5B
6	R	403	EDO	O1-C1-C2-O2
2	R	401	NAD	O4B-C4B-C5B-O5B
3	O	402	3PG	C3-O1P-P-O4P
5	R	402	G3H	C3-O1P-P-O4P

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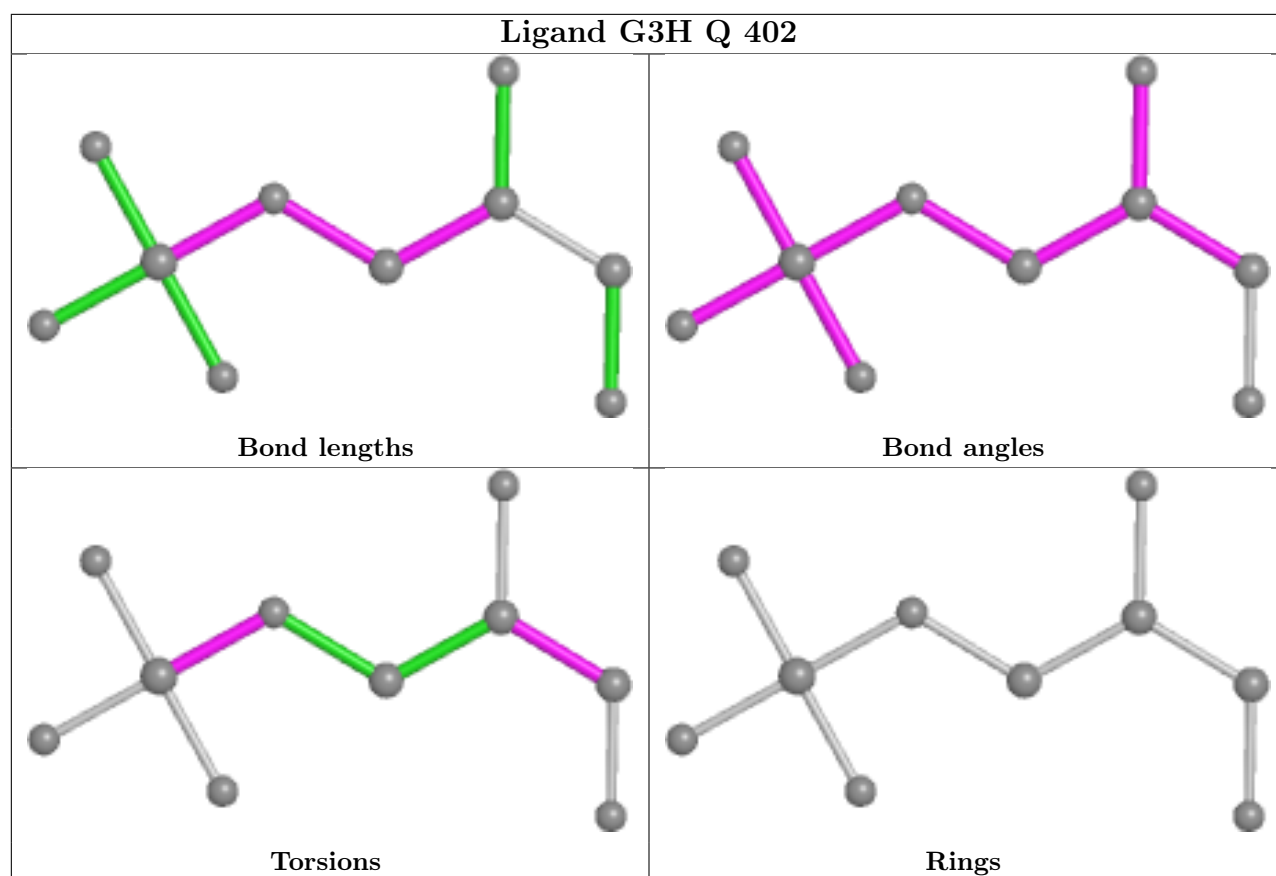
Mol	Chain	Res	Type	Atoms
5	P	402	G3H	C3-O1P-P-O2P
6	Q	403	EDO	O1-C1-C2-O2
2	O	401	NAD	O4B-C4B-C5B-O5B
2	P	401	NAD	O4B-C4B-C5B-O5B

There are no ring outliers.

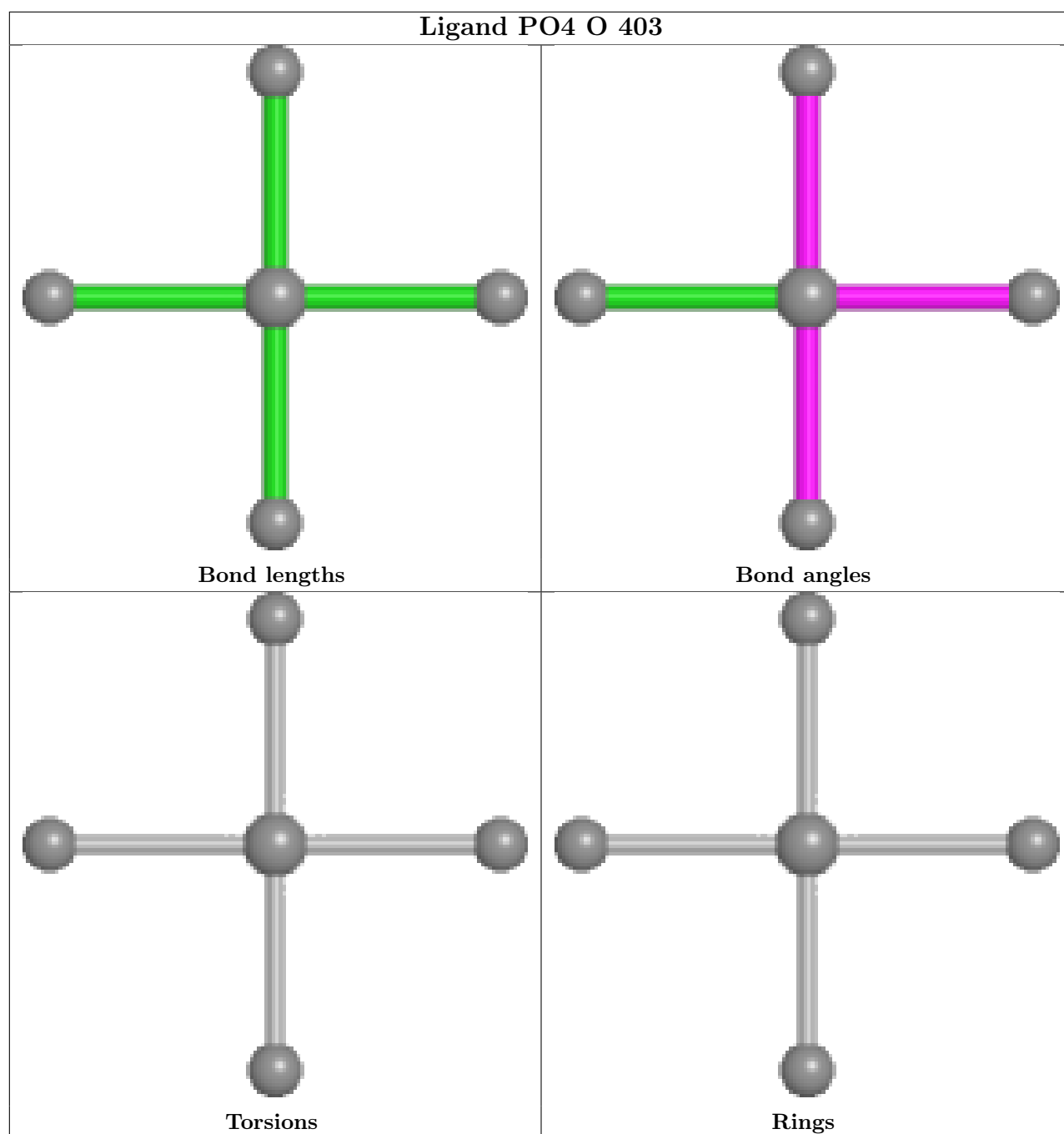
10 monomers are involved in 16 short contacts:

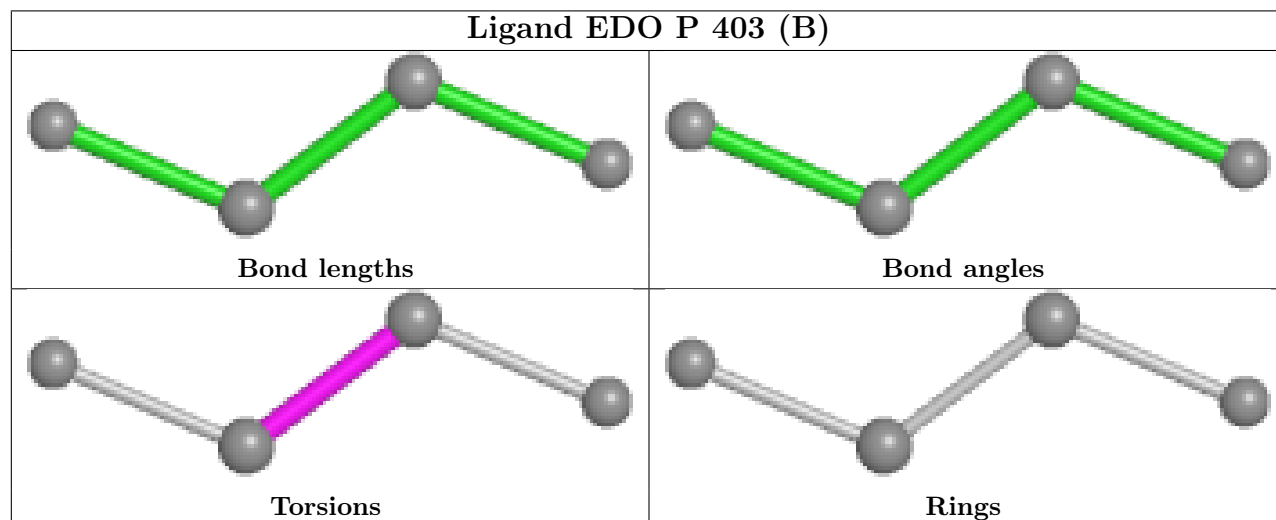
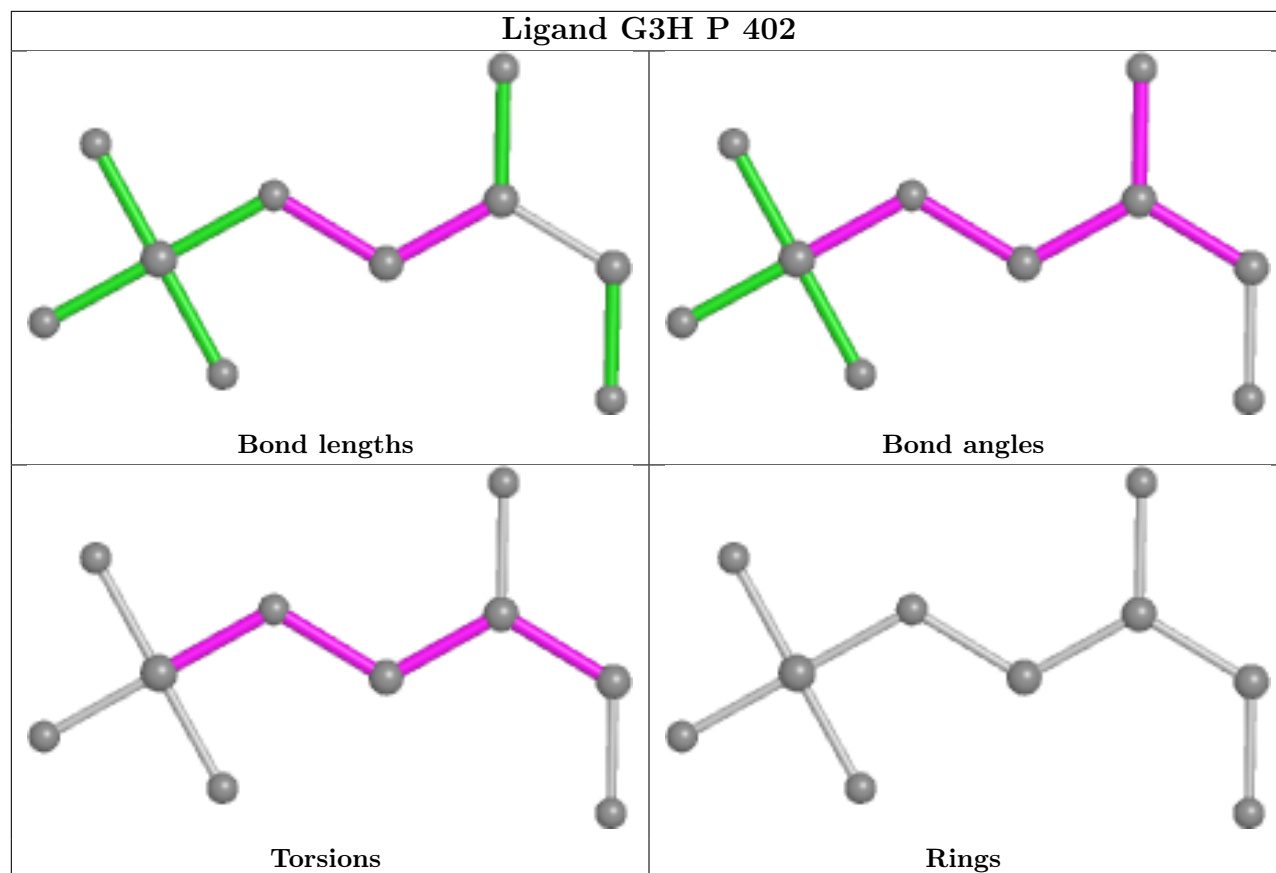
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	Q	402	G3H	2	0
5	P	402	G3H	1	0
6	P	403[B]	EDO	3	0
2	R	401	NAD	1	0
2	P	401	NAD	2	0
2	Q	401	NAD	1	0
2	O	401	NAD	2	0
6	Q	404[A]	EDO	1	0
3	O	402	3PG	1	0
5	R	402	G3H	2	0

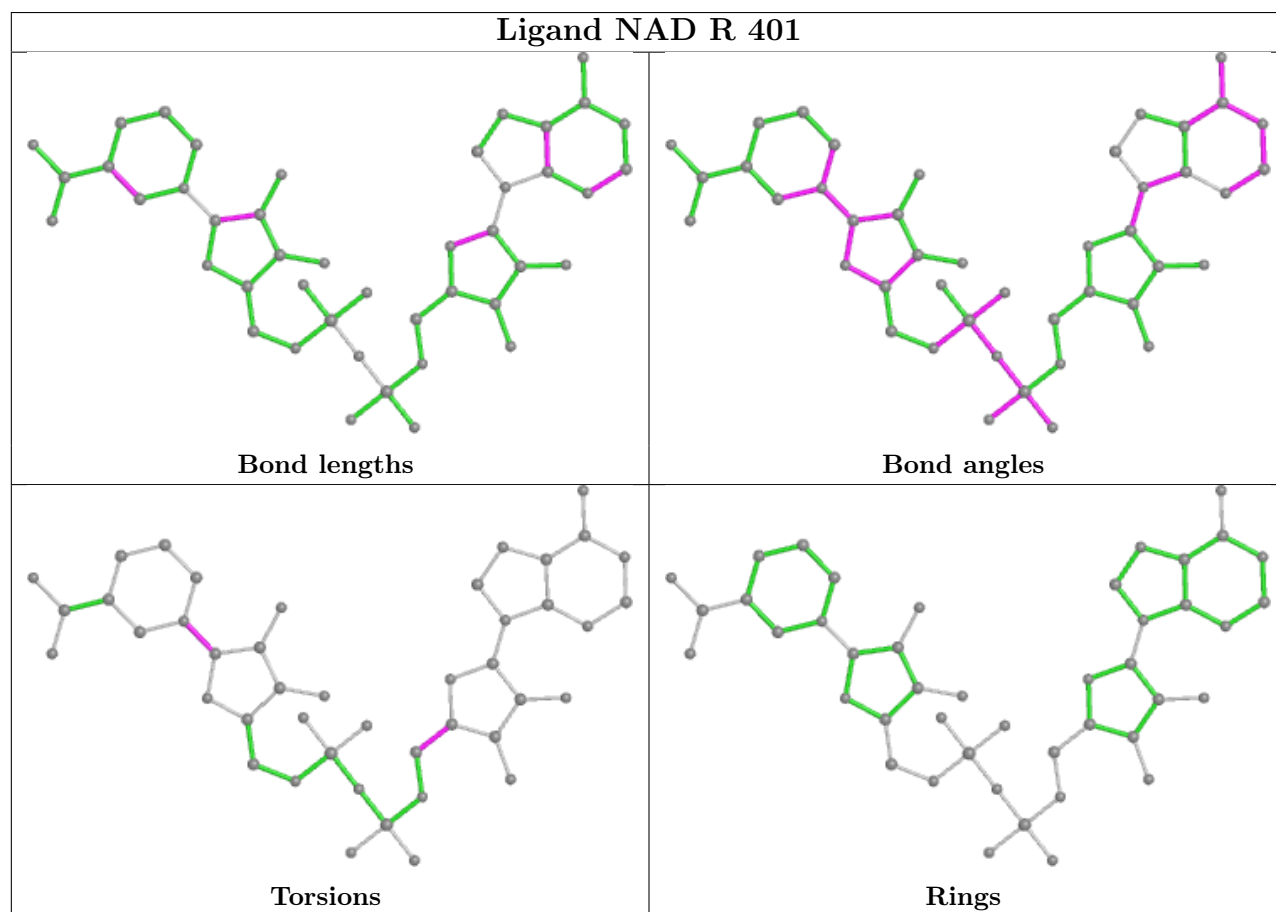
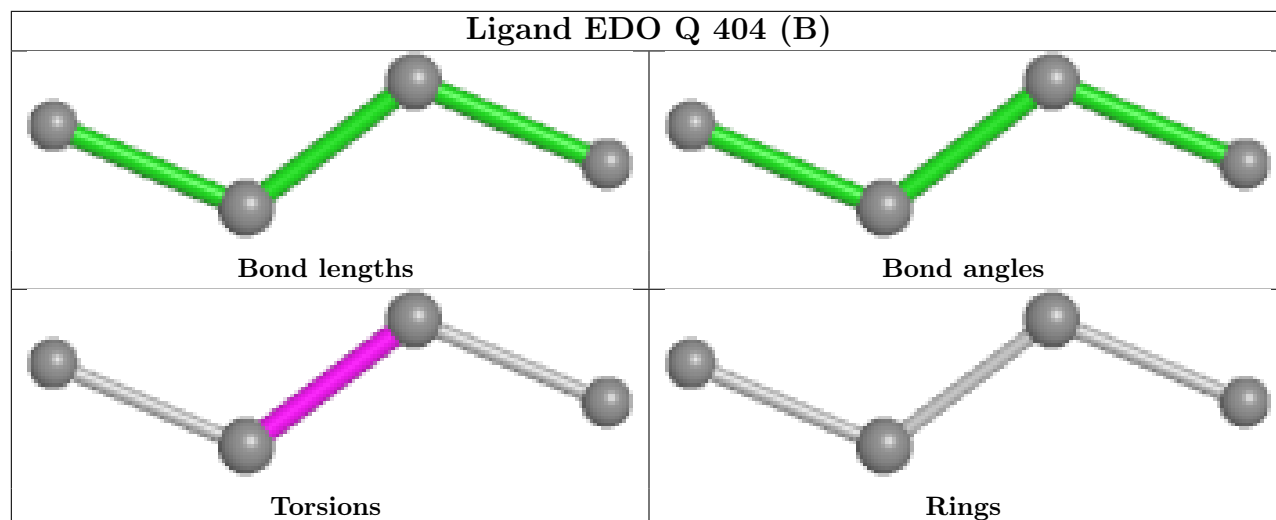
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

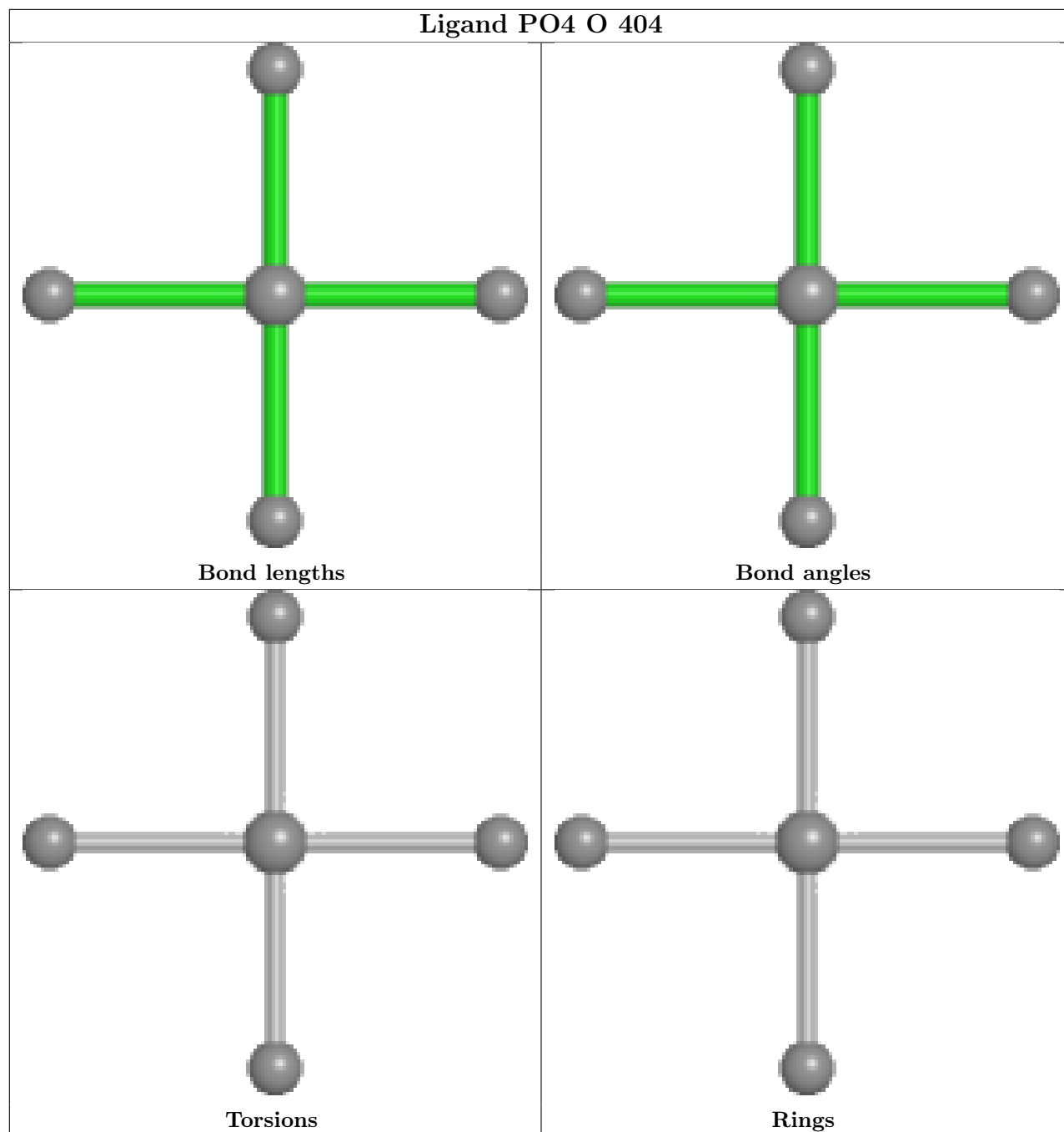


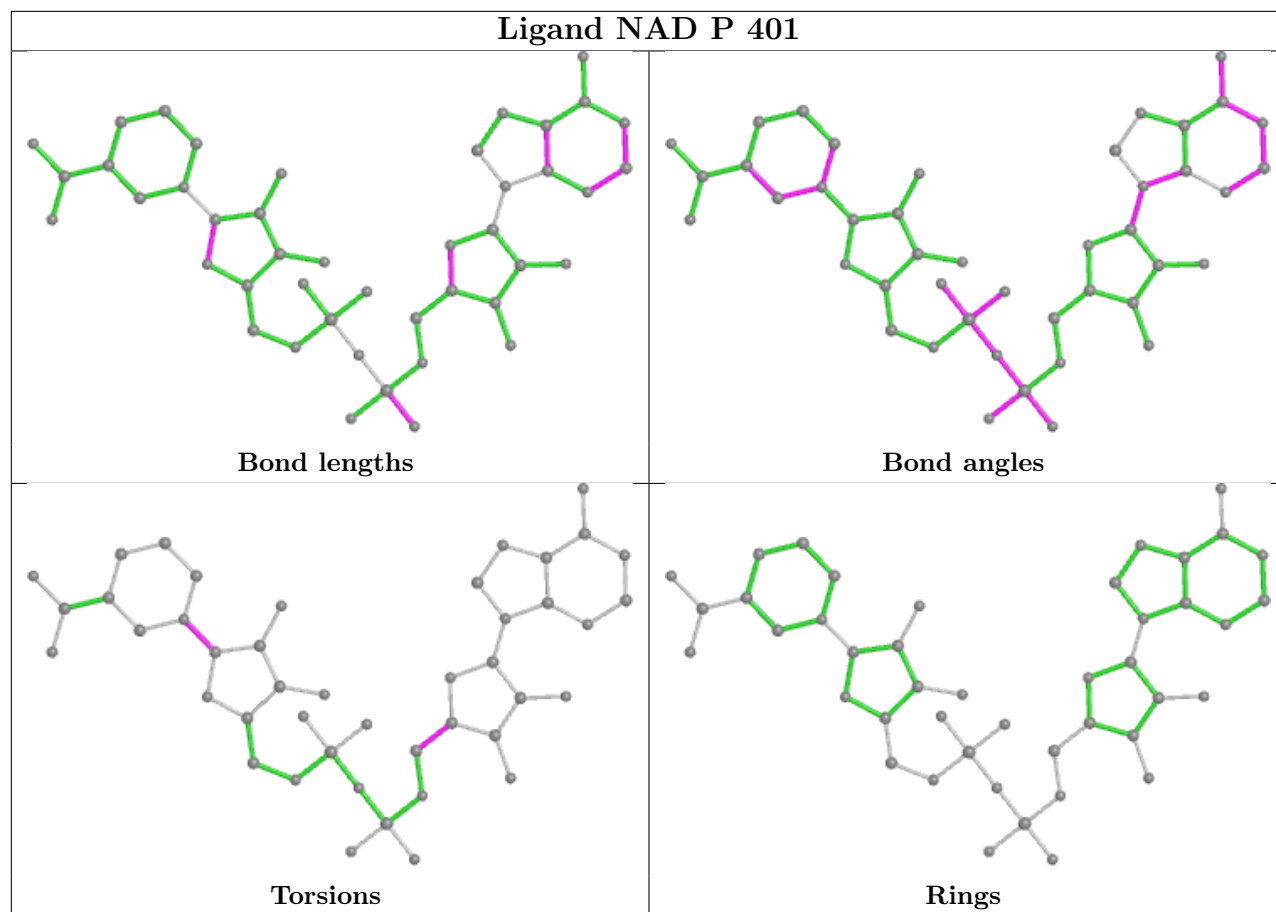


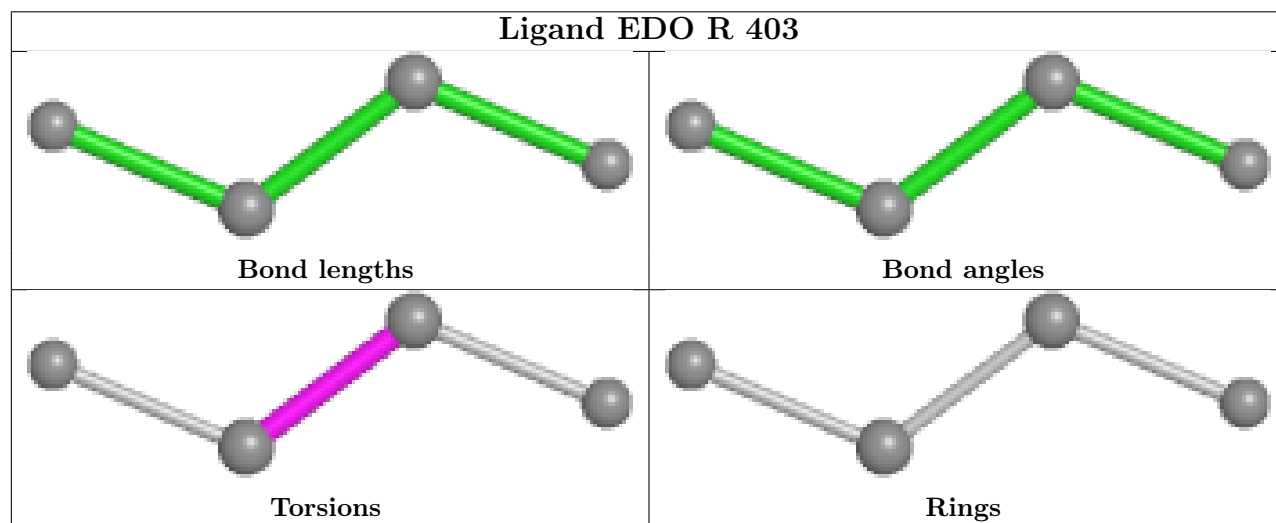
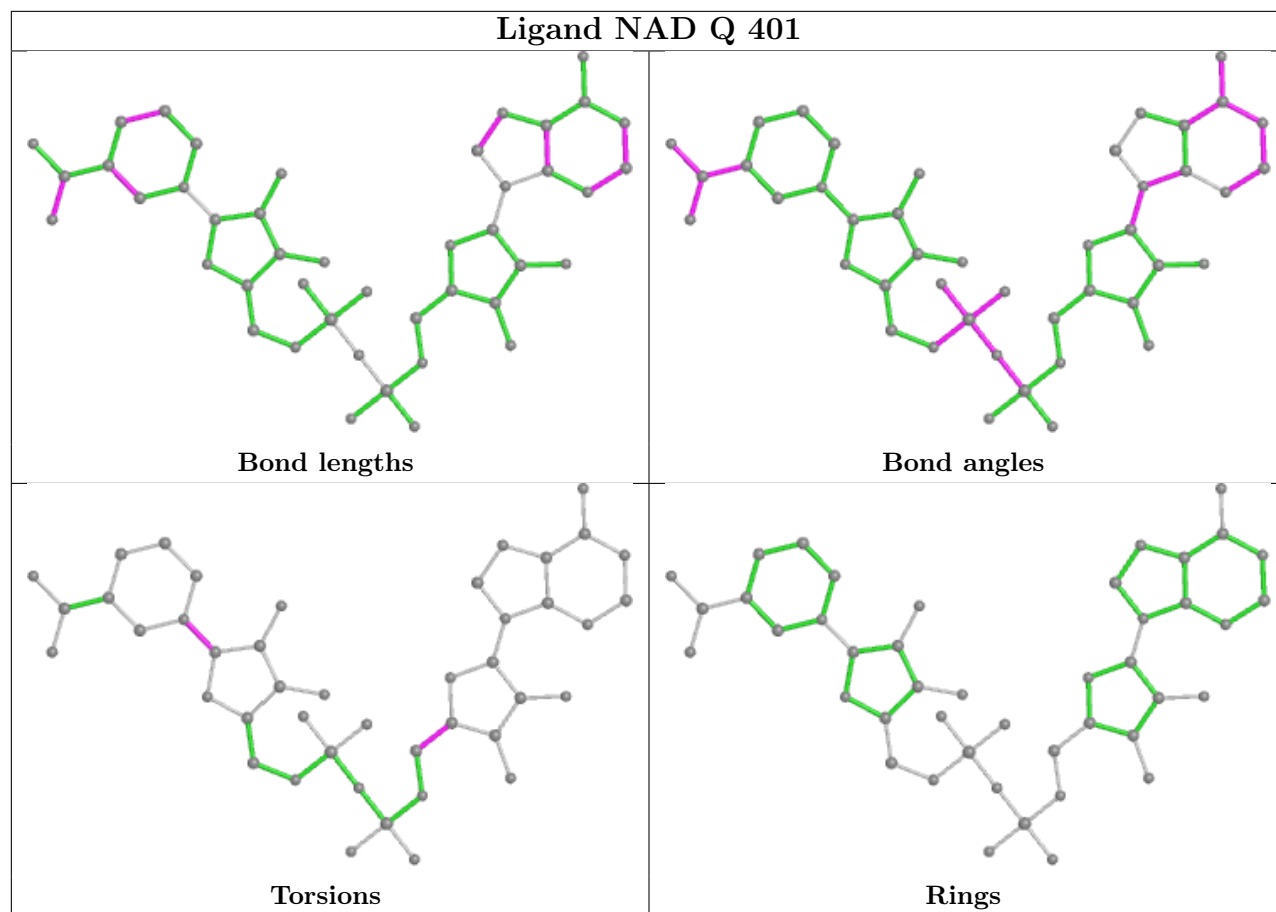


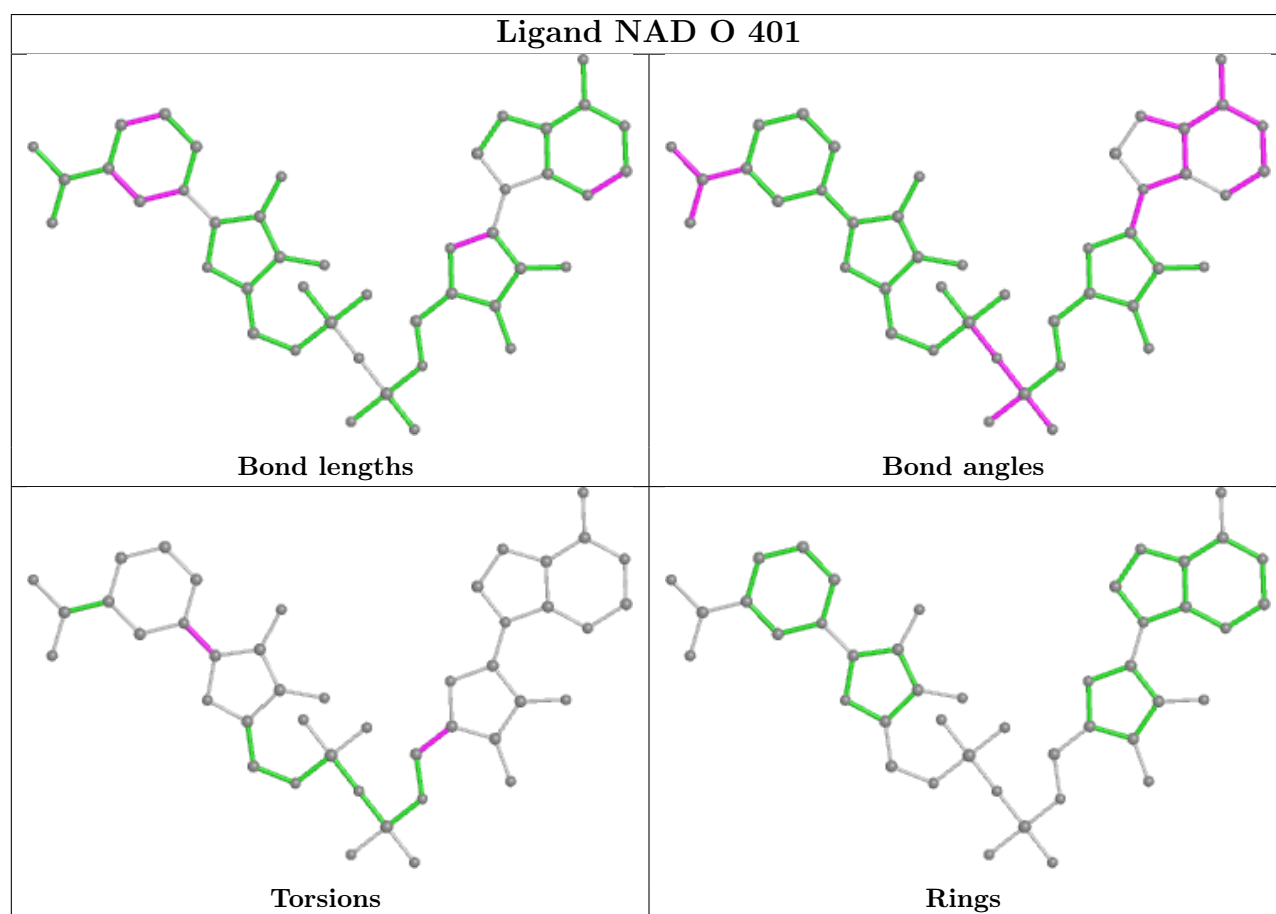
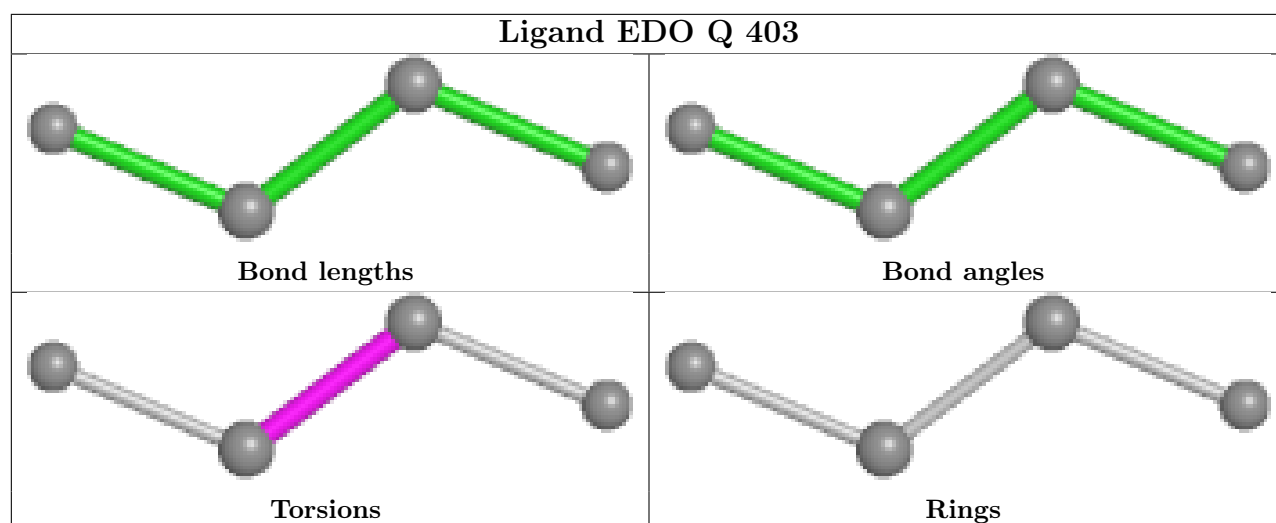


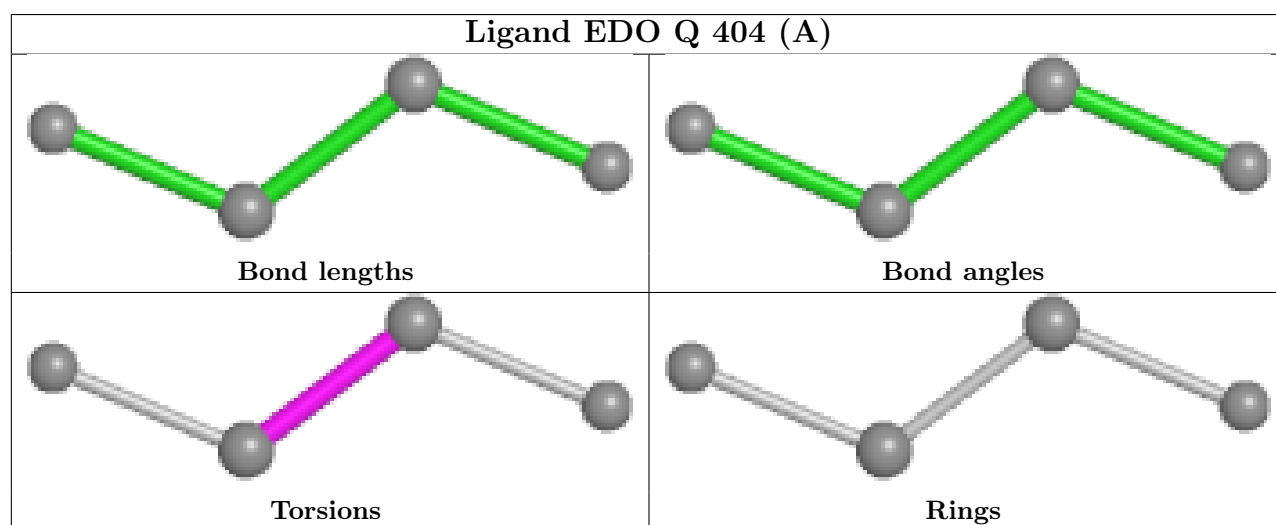
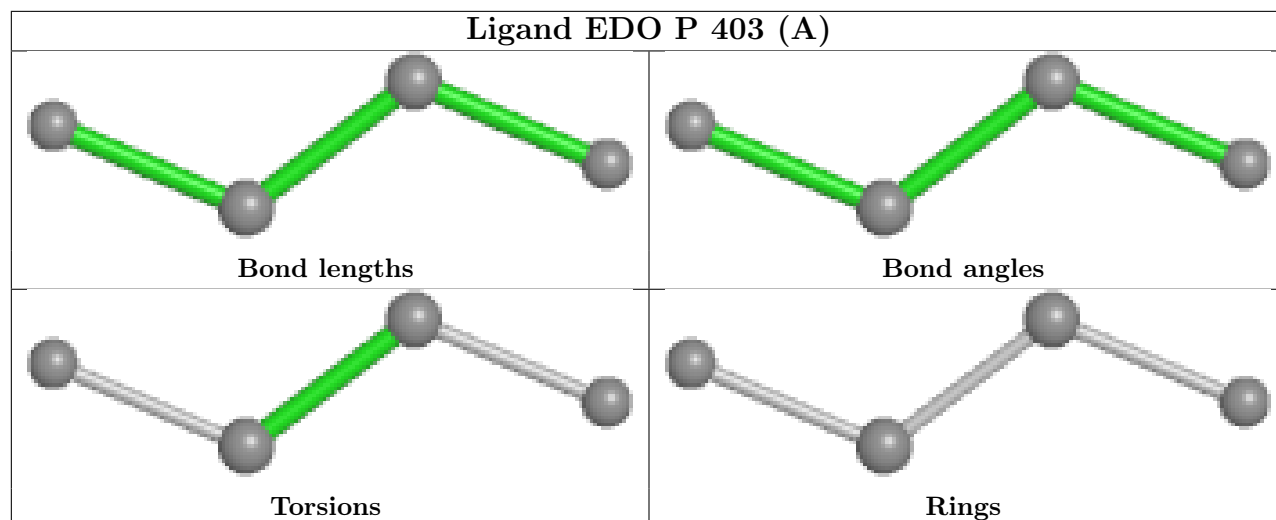






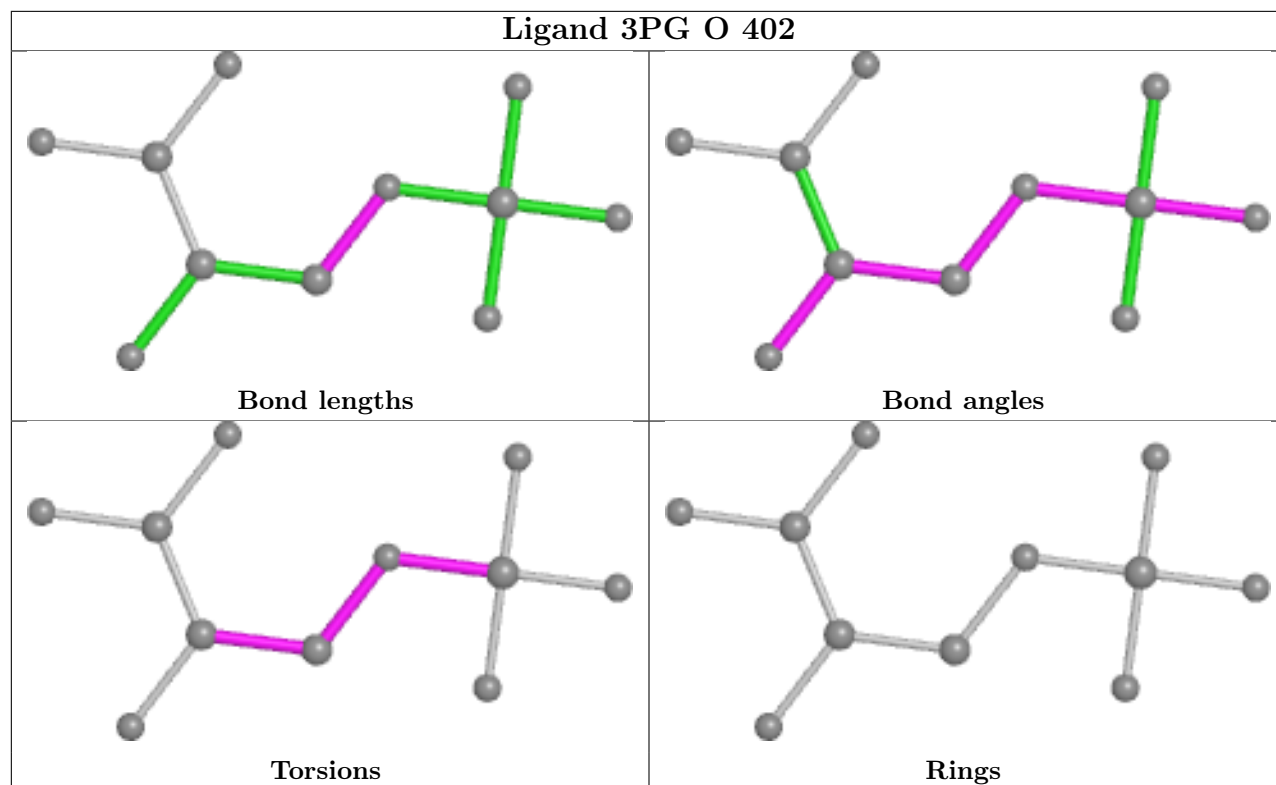




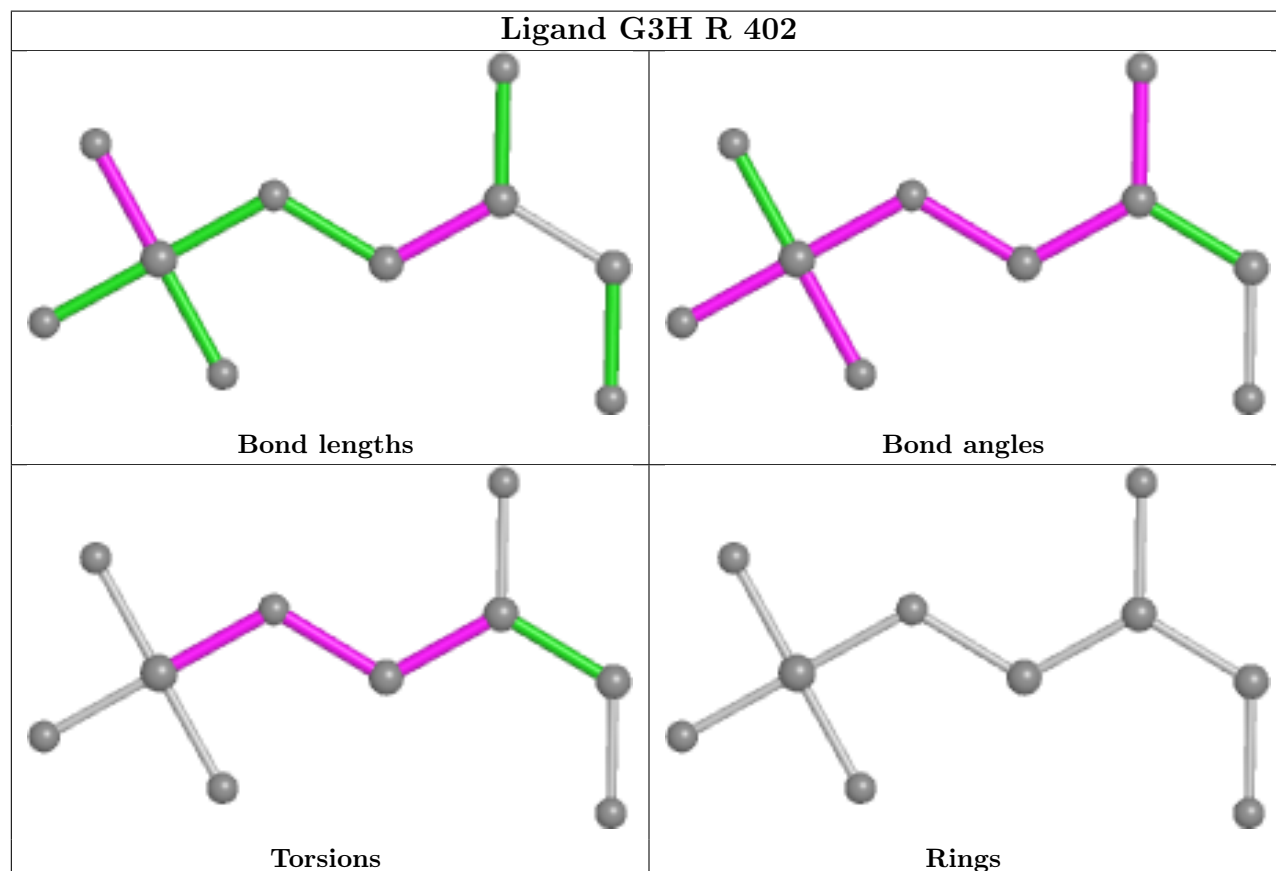




## Ligand 3PG O 402



## Ligand G3H R 402



## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	O	334/352 (94%)	-0.66	0  	15, 20, 32, 47	0
1	P	334/352 (94%)	-0.58	0  	15, 21, 35, 60	0
1	Q	336/352 (95%)	-0.39	3 (0%)  	16, 22, 39, 56	0
1	R	333/352 (94%)	0.01	20 (6%)  	17, 29, 50, 62	0
All	All	1337/1408 (94%)	-0.40	23 (1%)  	15, 22, 43, 62	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	R	333	LEU	3.5
1	R	117	VAL	3.3
1	R	90	ALA	3.2
1	R	331	ALA	3.1
1	Q	-1	ALA	3.1
1	Q	-2	MET	3.0
1	R	103	ALA	2.7
1	R	143	THR	2.7
1	R	23	VAL	2.6
1	R	141	ASN	2.5
1	R	144	ILE	2.5
1	R	112	ALA	2.5
1	R	108	ALA	2.5
1	R	115	LYS	2.3
1	R	5	GLY	2.3
1	R	26	ASN	2.1
1	Q	1	MET	2.1
1	R	27	ILE	2.1
1	R	125	GLU	2.1
1	R	107	GLN	2.0
1	R	85	TRP	2.0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	R	114	ALA	2.0
1	R	104	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 monosaccharides in this entry.

## 6.4 Ligands [i](#)

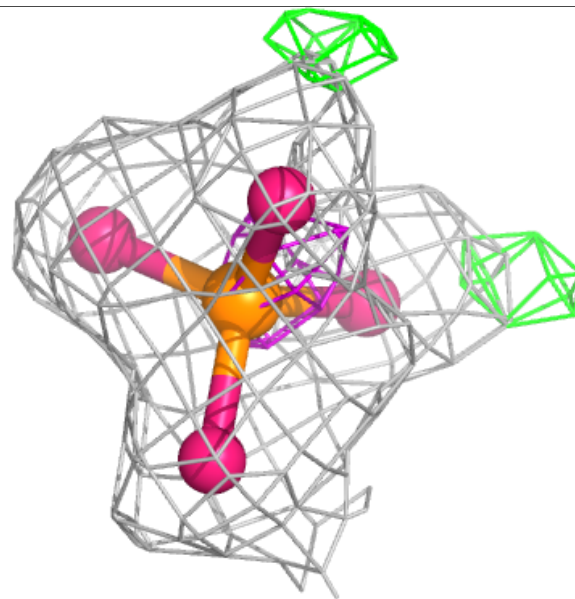
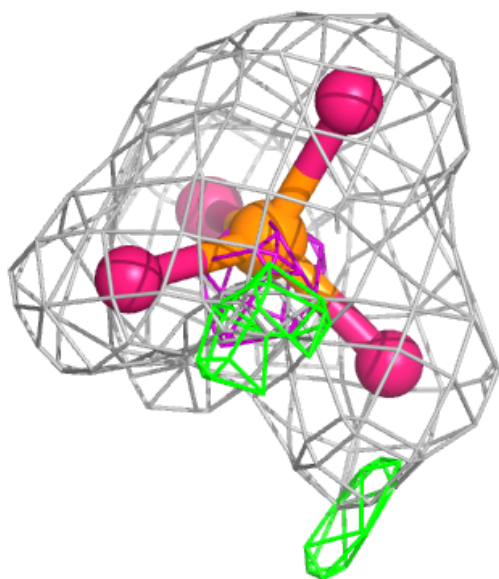
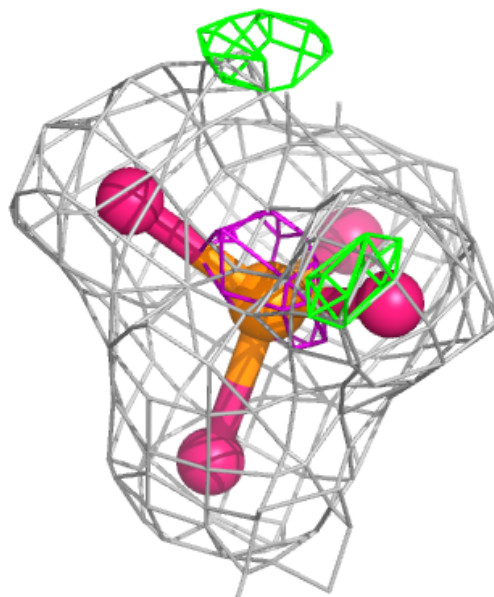
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
4	PO4	O	404	5/5	0.70	0.23	58,64,65,98	0
6	EDO	Q	403	4/4	0.76	0.27	54,56,56,59	0
6	EDO	R	403	4/4	0.86	0.13	45,46,52,53	0
6	EDO	Q	404[B]	4/4	0.88	0.19	24,28,29,32	4
6	EDO	Q	404[A]	4/4	0.88	0.19	26,28,28,28	4
6	EDO	P	403[A]	4/4	0.91	0.15	25,32,32,36	4
6	EDO	P	403[B]	4/4	0.91	0.15	23,24,25,26	4
5	G3H	R	402	10/10	0.93	0.14	34,42,53,59	0
2	NAD	R	401	44/44	0.95	0.11	25,31,34,36	0
3	3PG	O	402	11/11	0.95	0.12	28,33,57,66	0
5	G3H	P	402	10/10	0.96	0.10	28,33,47,49	0
4	PO4	O	403	5/5	0.96	0.18	32,35,37,41	0
2	NAD	O	401	44/44	0.97	0.08	18,20,23,27	0
2	NAD	P	401	44/44	0.97	0.10	18,22,25,27	0
5	G3H	Q	402	10/10	0.98	0.09	21,27,31,33	0
2	NAD	Q	401	44/44	0.98	0.08	16,21,23,25	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.

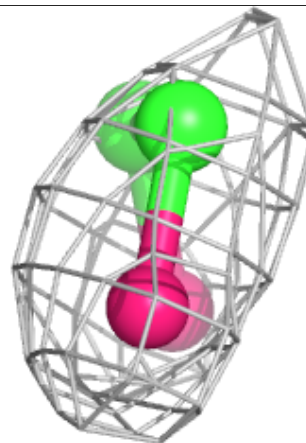
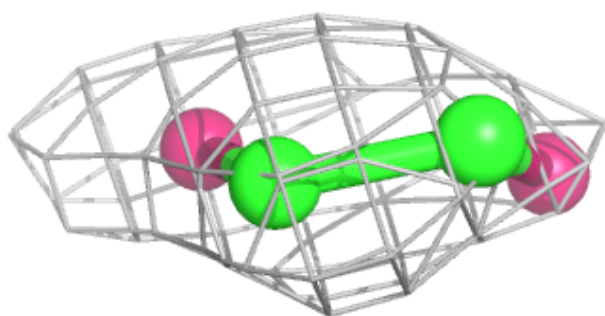
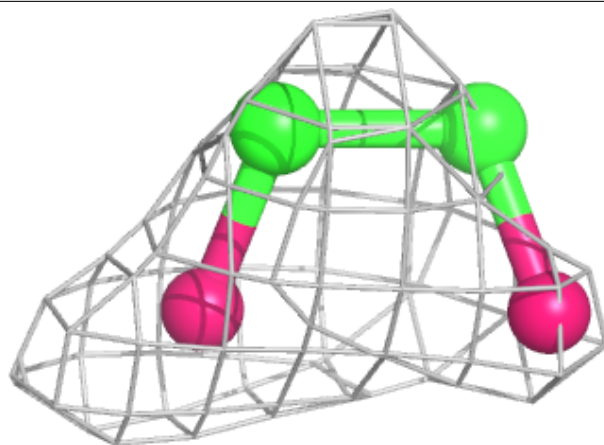
**Electron density around PO4 O 404:**

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



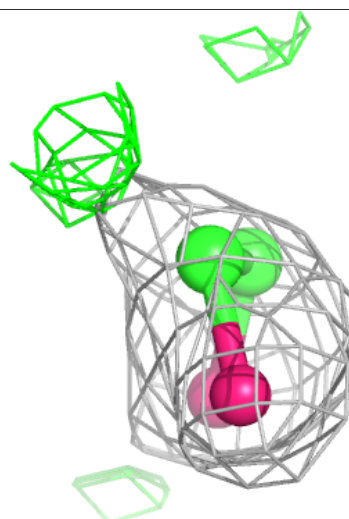
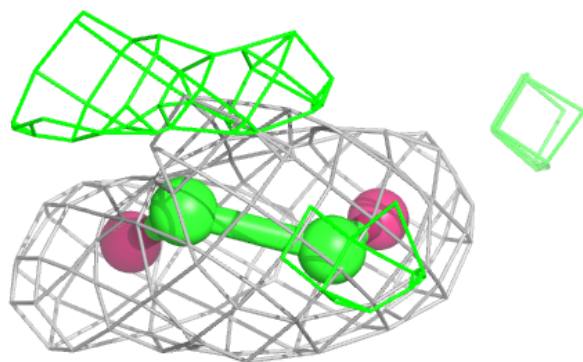
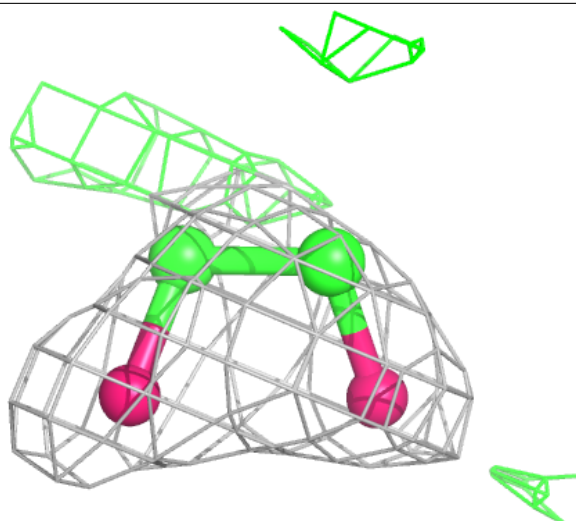
**Electron density around EDO Q 403:**

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



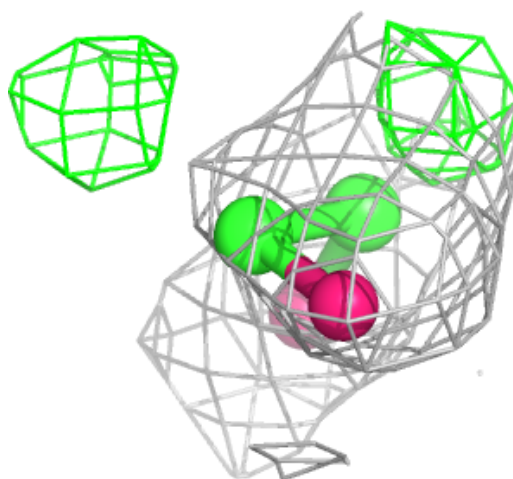
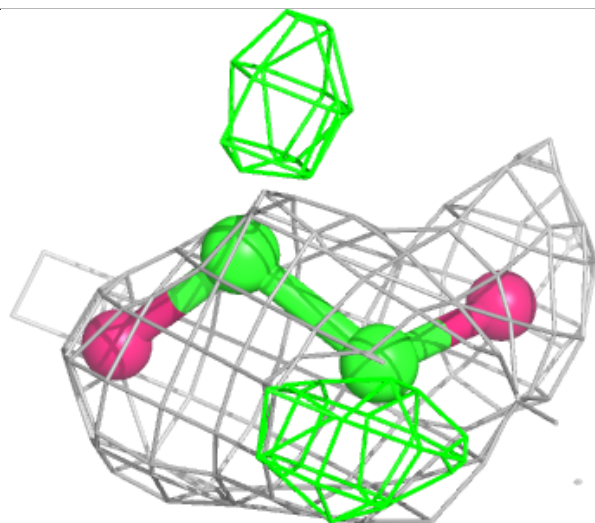
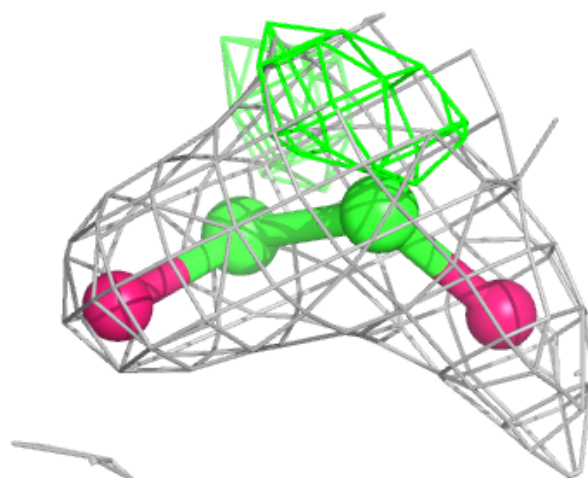
**Electron density around EDO R 403:**

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



**Electron density around EDO Q 404 (B):**

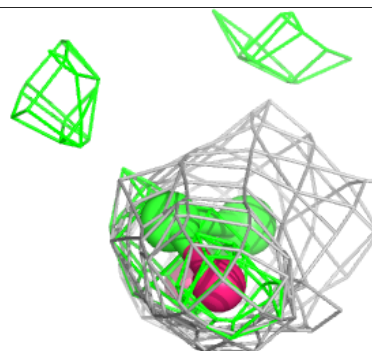
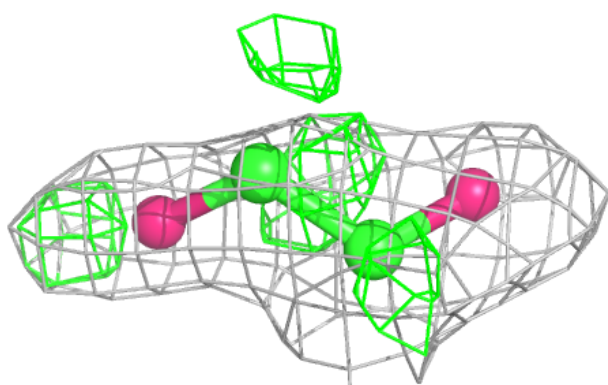
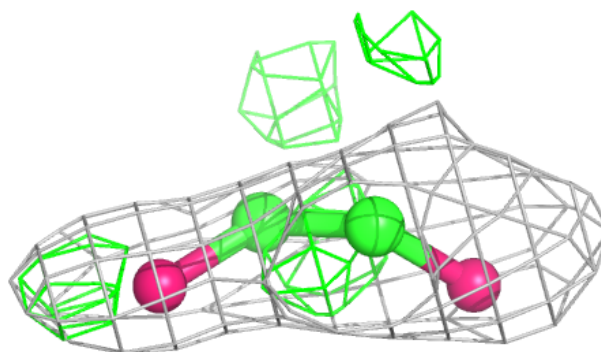
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





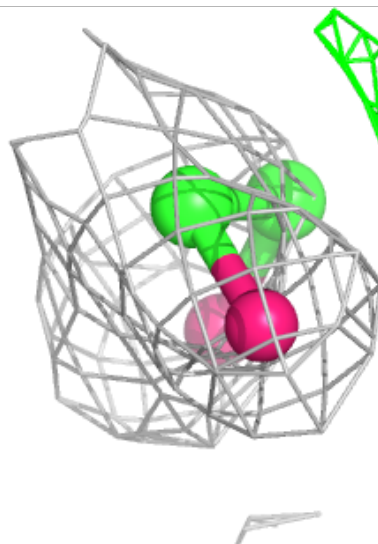
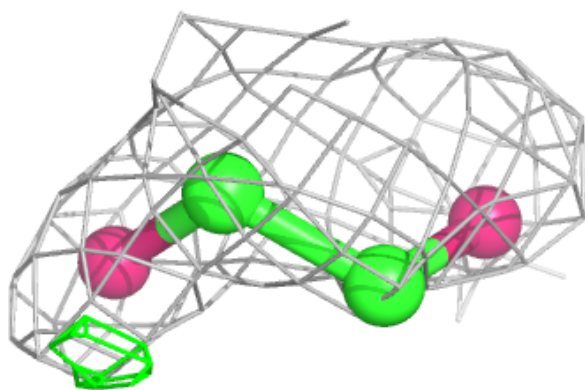
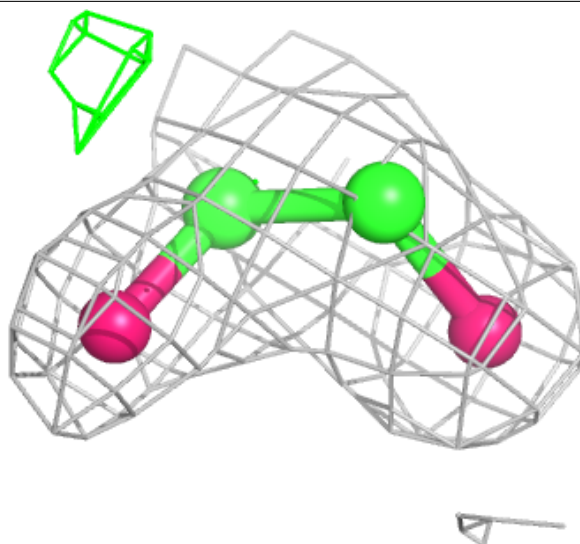
**Electron density around EDO Q 404 (A):**

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



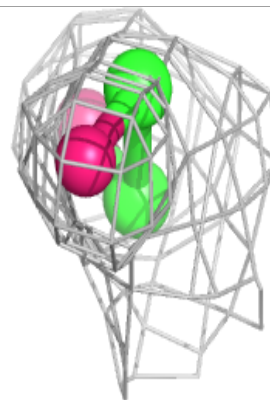
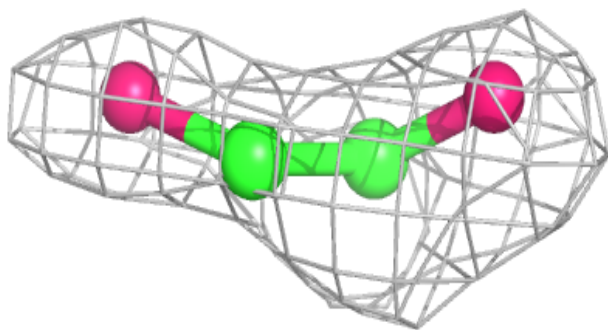
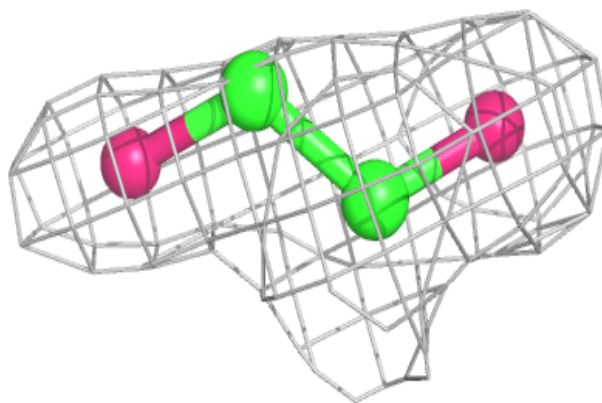
**Electron density around EDO P 403 (A):**

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



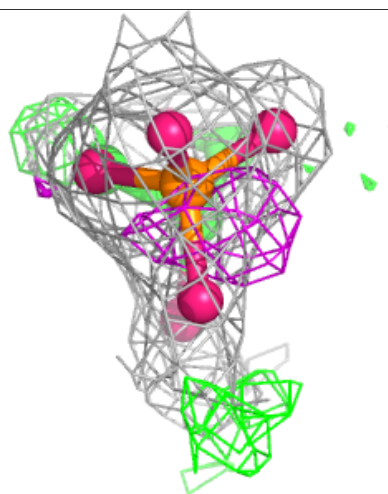
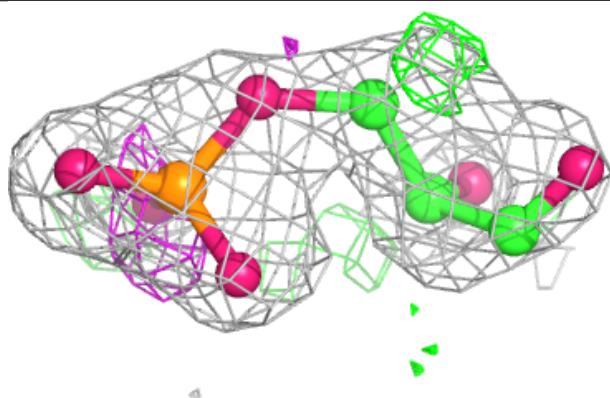
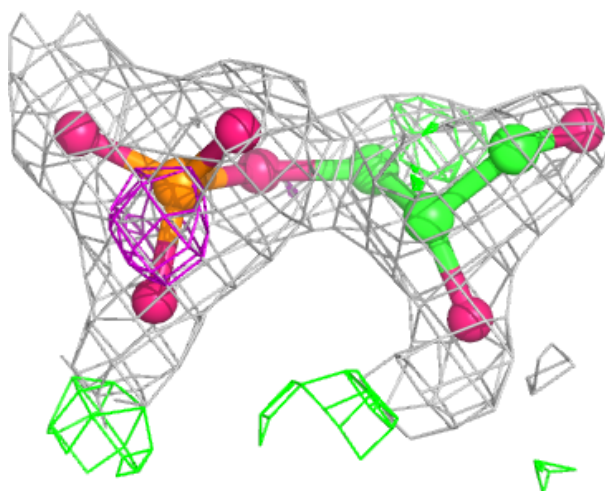
**Electron density around EDO P 403 (B):**

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



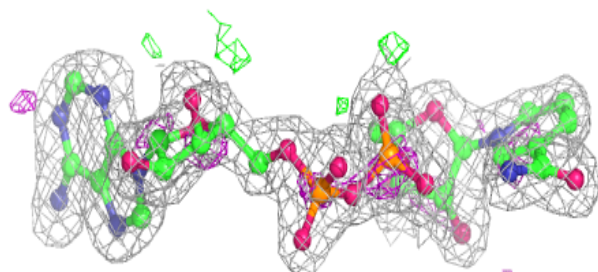
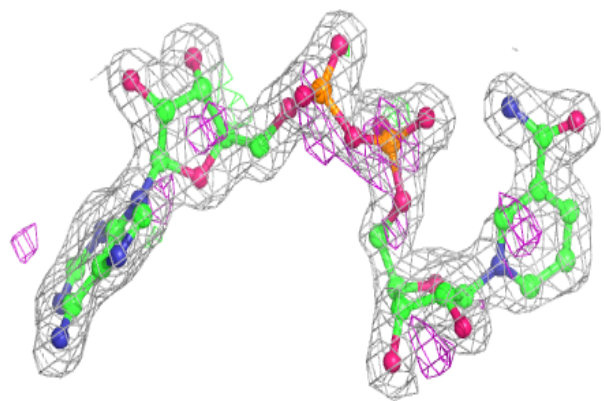
**Electron density around G3H R 402:**

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



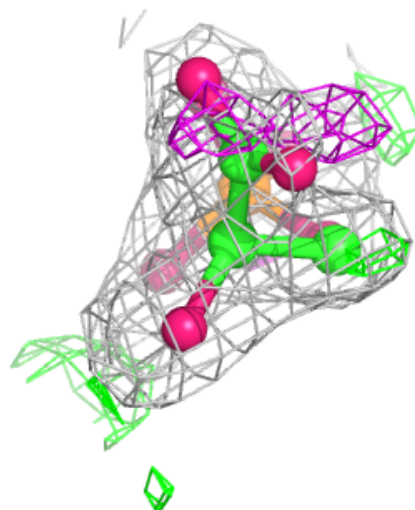
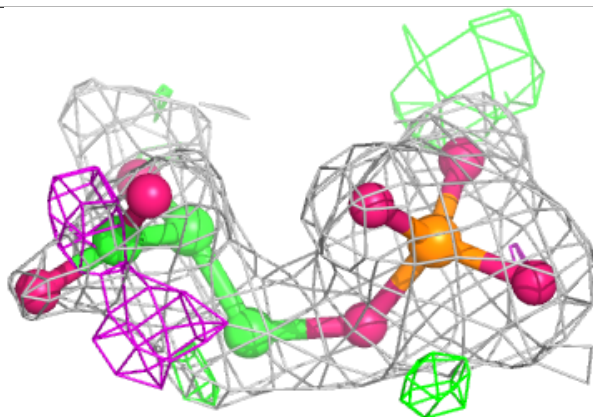
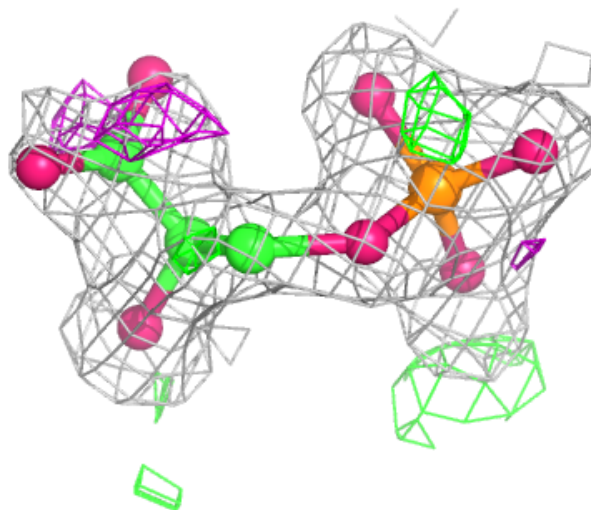
**Electron density around NAD R 401:**

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



**Electron density around 3PG O 402:**

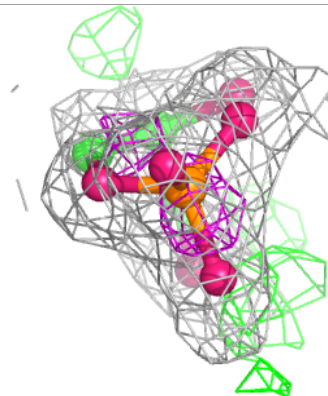
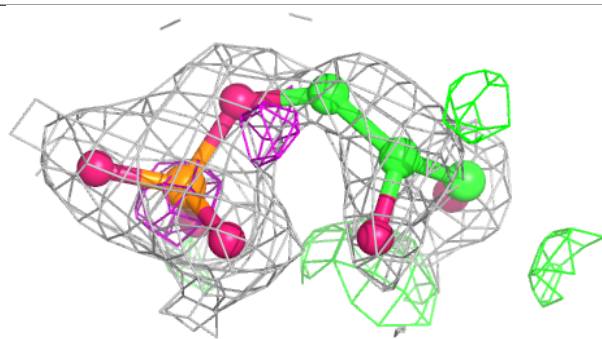
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





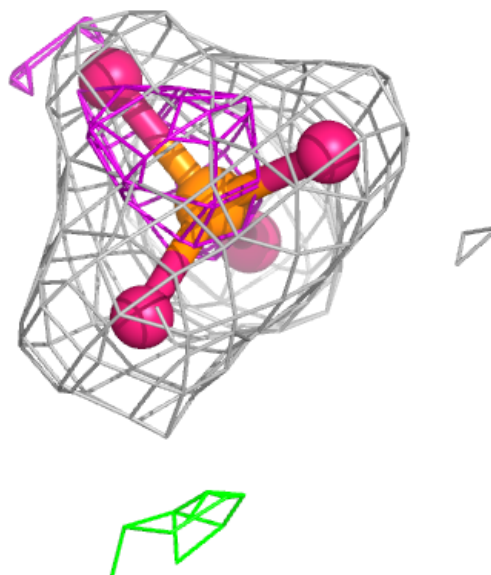
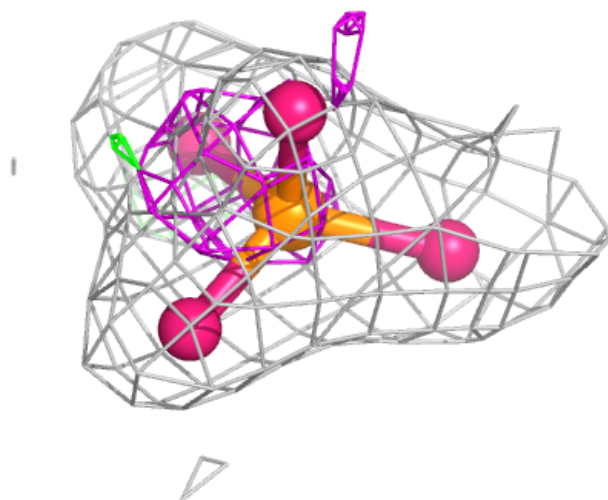
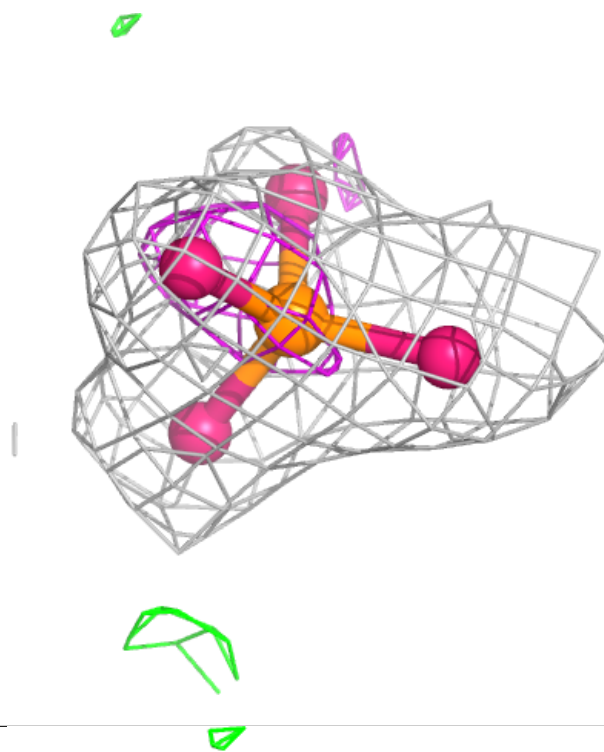
**Electron density around G3H P 402:**

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



**Electron density around PO4 O 403:**

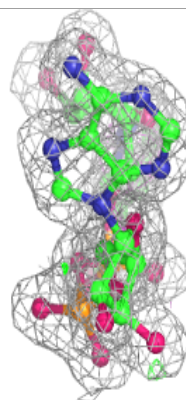
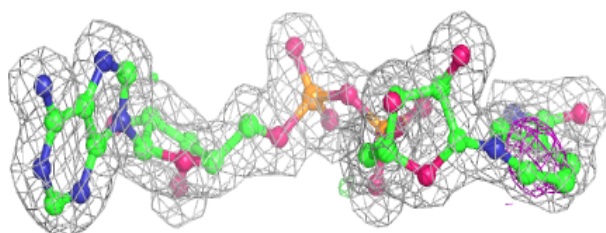
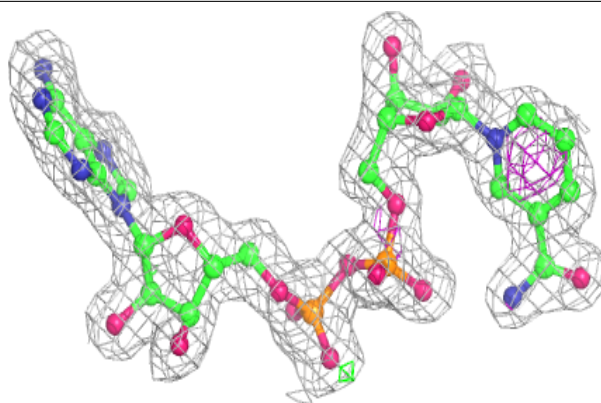
$2mF_o - DF_c$  (at 0.7 rmsd) in gray  
 $mF_o - DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



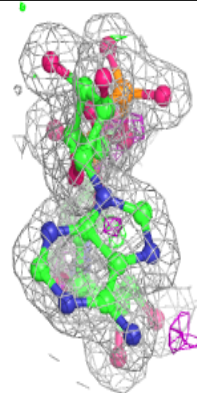
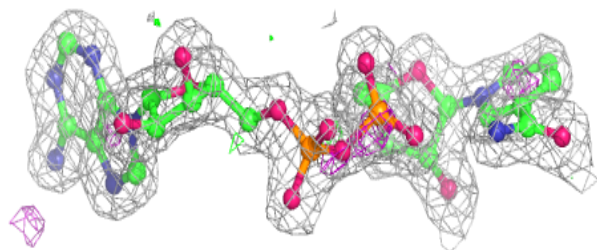
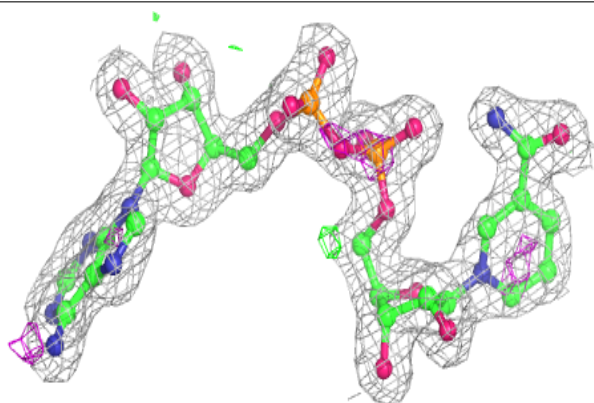


**Electron density around NAD O 401:**

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

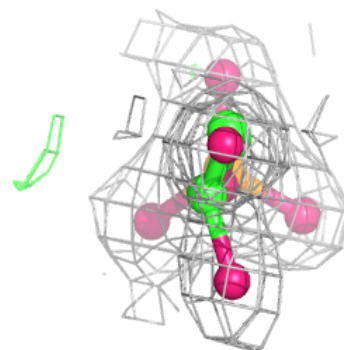
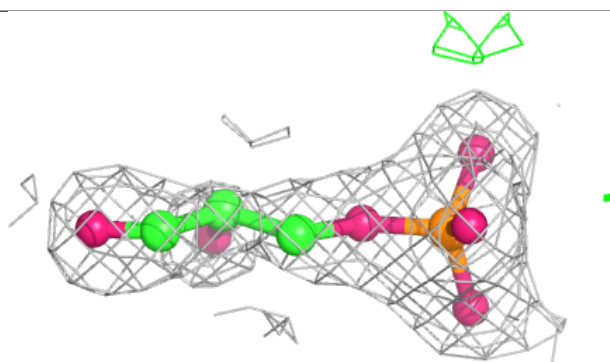
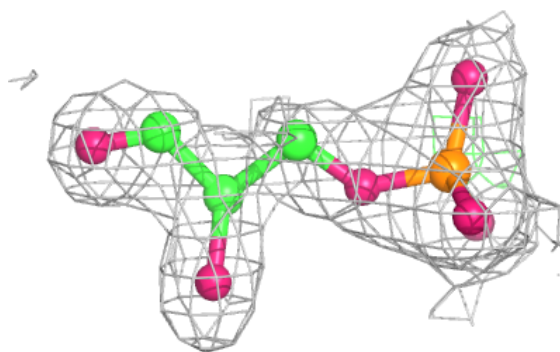
**Electron density around NAD P 401:**

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

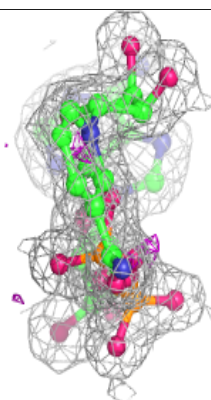
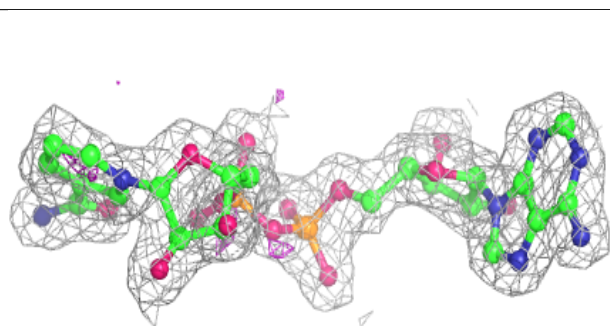
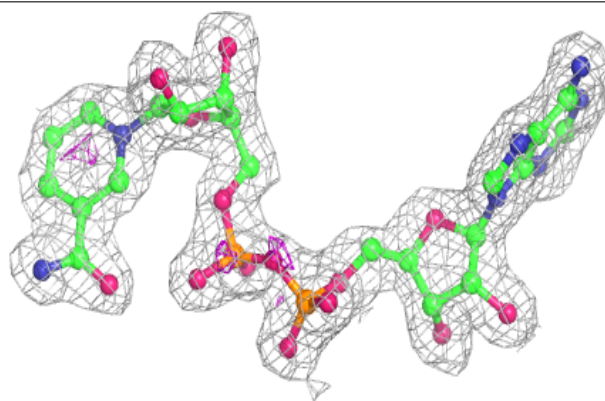


**Electron density around G3H Q 402:**

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

**Electron density around NAD Q 401:**

$2mF_o-DF_c$  (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.