



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 08:37 PM UTC

PDB ID : 8E0V / pdb_00008e0v
Title : DAHP (3-deoxy-D-arabinoheptulosonate-7-phosphate) Synthase complexed with Mn(II), PEP, and Pi in unbound:(bound)2:other Conformations
Authors : Berti, P.J.; Junop, M.S.; Grainger, R.
Deposited on : 2022-08-09
Resolution : 2.30 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

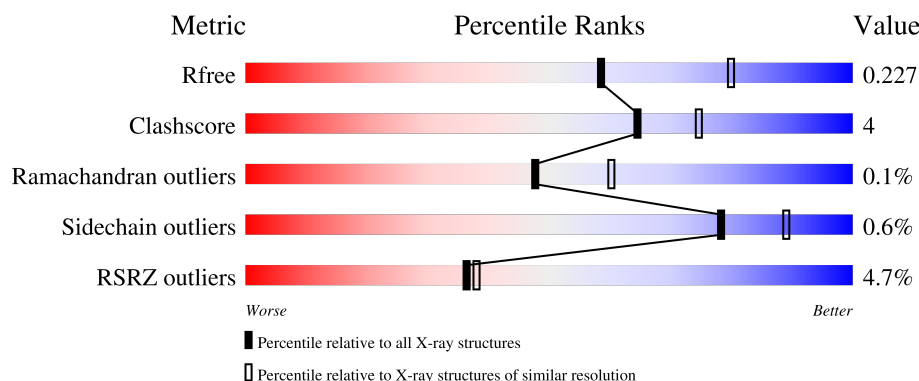
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	351	7% 88% 9% ..
1	B	351	2% 88% 8% ..
1	C	351	% 93% 5% .
1	D	351	9% 83% 14% ..

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 11265 atoms, of which 36 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 Phospho-2-dehydro-3-deoxyheptonate aldolase, Phe-sensitive.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	344	Total	C	N	O	S	0	1	0
			2572	1614	458	486	14			
1	B	340	Total	C	N	O	S	0	0	0
			2547	1601	451	481	14			
1	C	344	Total	C	N	O	S	0	0	0
			2572	1615	455	488	14			
1	D	343	Total	C	N	O	S	0	0	0
			2543	1600	451	478	14			

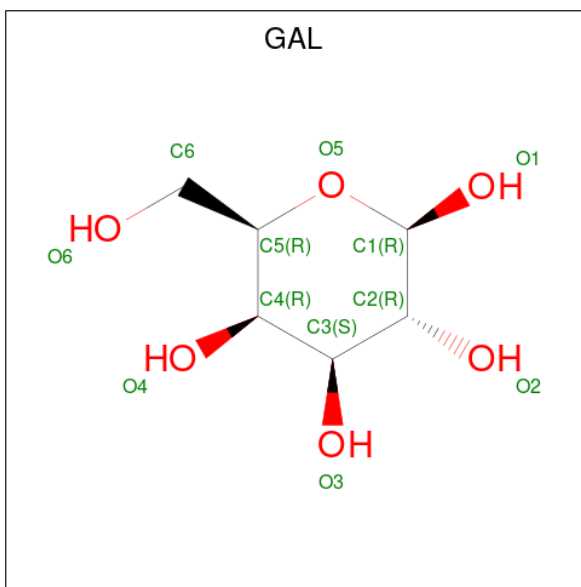
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	expression tag	UNP P0AB91
B	0	GLY	-	expression tag	UNP P0AB91
C	0	GLY	-	expression tag	UNP P0AB91
D	0	GLY	-	expression tag	UNP P0AB91

- Molecule 2 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn) (labeled as "Ligand of Interest" by depositor).

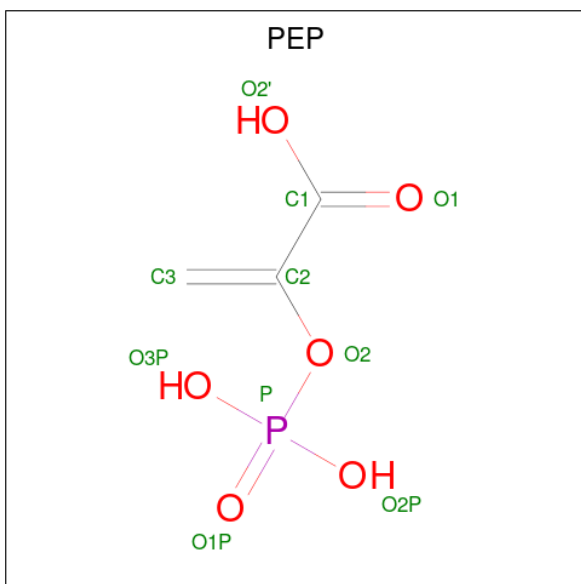
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mn	0	0
			1	1		
2	B	1	Total	Mn	0	0
			1	1		
2	C	1	Total	Mn	0	0
			1	1		
2	D	1	Total	Mn	0	0
			1	1		

- Molecule 3 is beta-D-galactopyranose (CCD ID: GAL) (formula: C₆H₁₂O₆).



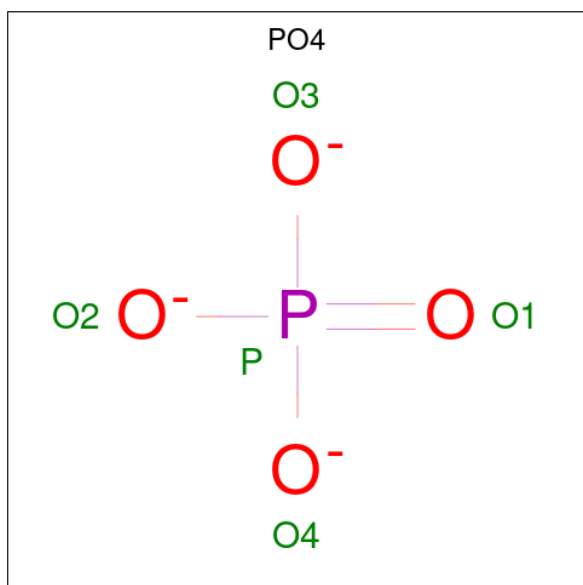
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	H	O	0	0
			24	6	12	6		
3	A	1	Total	C	H	O	0	0
			24	6	12	6		
3	C	1	Total	C	H	O	0	0
			24	6	12	6		

- Molecule 4 is PHOSPHOENOLPYRUVATE (CCD ID: PEP) (formula: $C_3H_5O_6P$) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 5 is PHOSPHATE ION (CCD ID: PO4) (formula: O_4P) (labeled as "Ligand of Interest" by depositor).



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

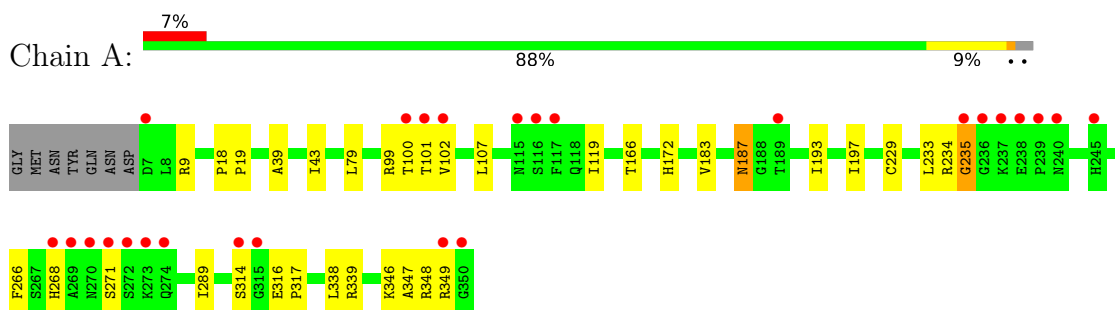
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	231	Total	O	0	0
			231	231		
6	B	165	Total	O	0	0
			165	165		
6	C	329	Total	O	0	0
			329	329		
6	D	200	Total	O	0	0
			200	200		

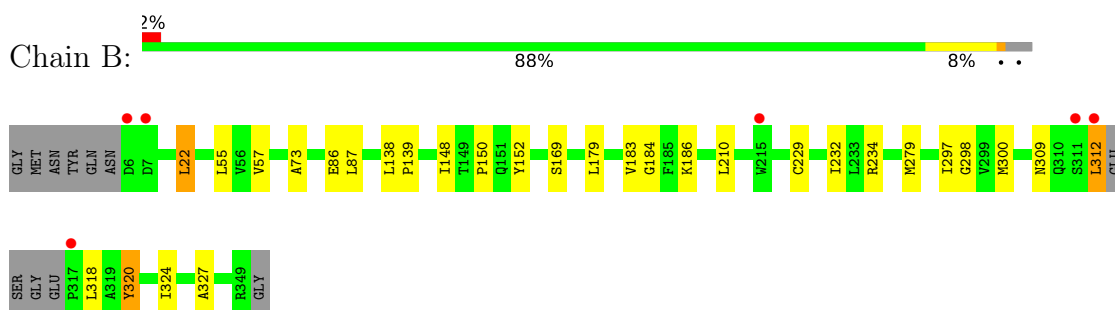
3 Residue-property plots [i](#)

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.

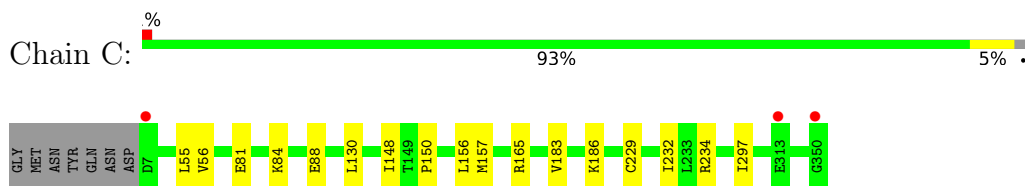
- Molecule 1: Phospho-2-dehydro-3-deoxyheptonate aldolase, Phe-sensitive



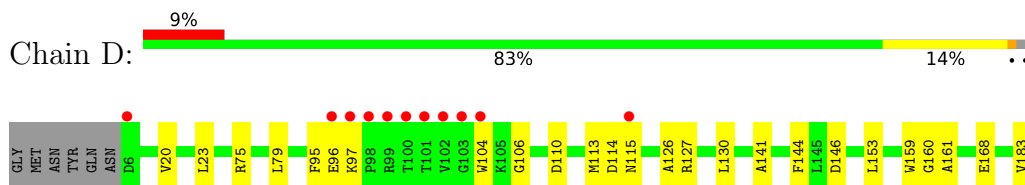
- Molecule 1: Phospho-2-dehydro-3-deoxyheptonate aldolase, Phe-sensitive

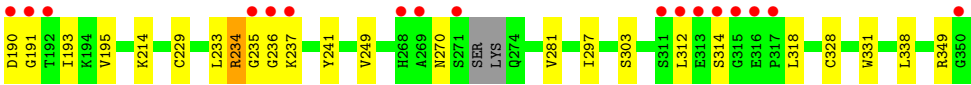


- Molecule 1: Phospho-2-dehydro-3-deoxyheptonate aldolase, Phe-sensitive



- Molecule 1: Phospho-2-dehydro-3-deoxyheptonate aldolase, Phe-sensitive





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	210.19Å 52.91Å 150.48Å 90.00° 115.71° 90.00°	Depositor
Resolution (Å)	49.65 – 2.30 49.65 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.7 (49.65-2.30) 99.6 (49.65-2.30)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.58 (at 2.29Å)	Xtriage
Refinement program	PHENIX 1.19.2_4158	Depositor
R, R_{free}	0.194 , 0.228 0.194 , 0.227	Depositor DCC
R_{free} test set	1998 reflections (2.97%)	wwPDB-VP
Wilson B-factor (Å ²)	23.5	Xtriage
Anisotropy	0.738	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 37.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	11265	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.61% 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: PEP, PO4, GAL, MN

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	0.66	0/2620	0.87	3/3554 (0.1%)
1	B	0.66	0/2591	0.91	2/3514 (0.1%)
1	C	0.62	0/2617	0.77	2/3549 (0.1%)
1	D	0.76	0/2587	1.10	6/3510 (0.2%)
All	All	0.68	0/10415	0.92	13/14127 (0.1%)

There are no bond length outliers.

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	190	ASP	N-CA-C	-12.40	97.79	113.16
1	D	270	ASN	N-CA-C	-7.51	104.00	113.01
1	A	100	THR	N-CA-C	-6.98	102.75	112.45
1	D	237	LYS	N-CA-C	-6.15	105.63	113.01
1	D	146	ASP	CA-C-N	-6.03	113.01	122.65
1	D	146	ASP	C-N-CA	-6.03	113.01	122.65
1	B	279	MET	CA-C-N	-5.83	112.47	120.28
1	B	279	MET	C-N-CA	-5.83	112.47	120.28
1	A	268	HIS	CB-CA-C	-5.80	109.88	116.54
1	C	88	GLU	CA-C-N	-5.67	113.43	122.50
1	C	88	GLU	C-N-CA	-5.67	113.43	122.50
1	A	235	GLY	CA-C-O	-5.25	118.61	122.23
1	D	314	SER	N-CA-C	-5.03	107.81	114.04

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	2572	0	2540	22	0
1	B	2547	0	2516	21	0
1	C	2572	0	2535	11	0
1	D	2543	0	2490	30	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	24	24	24	0	0
3	C	12	12	12	0	0
4	B	10	0	2	0	0
4	C	10	0	2	0	0
5	B	5	0	0	0	0
5	C	5	0	0	0	0
6	A	231	0	0	1	0
6	B	165	0	0	0	0
6	C	329	0	0	0	0
6	D	200	0	0	2	0
All	All	11229	36	10121	80	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

All (80) 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:D:20:VAL:HG12	1:D:214:LYS:HE2	1.80	0.64
1:B:312:LEU:HD23	1:B:324:ILE:HD12	1.80	0.63
1:A:107:LEU:C	1:A:107:LEU:HD23	2.28	0.59
1:B:312:LEU:HD22	1:B:318:LEU:HD11	1.85	0.57
1:C:186:LYS:HB3	1:C:234:ARG:HG2	1.85	0.57
1:A:193:ILE:HD12	1:A:197:ILE:HD11	1.86	0.56
1:B:57:VAL:O	1:B:300:MET:HA	2.05	0.56
1:D:234:ARG:HD2	1:D:235:GLY:O	2.06	0.56
1:A:9:ARG:HD2	1:B:179:LEU:O	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:148:ILE:C	1:C:150:PRO:HD2	2.32	0.54
1:B:86:GLU:C	1:B:87:LEU:HD23	2.33	0.54
1:C:130:LEU:HD13	1:C:156:LEU:HB2	1.89	0.54
1:A:187:ASN:OD1	1:A:233:LEU:HA	2.09	0.53
1:D:96:GLU:HG2	1:D:97:LYS:N	2.24	0.53
1:D:193:ILE:HG21	1:D:249:VAL:HG22	1.90	0.53
1:B:312:LEU:CD2	1:B:318:LEU:HD11	2.40	0.52
1:D:110:ASP:OD2	1:D:113:MET:HA	2.10	0.51
1:D:95:PHE:CZ	1:D:126:ALA:HB1	2.46	0.50
1:A:183:VAL:O	1:A:229:CYS:HA	2.12	0.49
1:D:186:LYS:HD2	6:D:616:HOH:O	2.12	0.49
1:A:79:LEU:HD12	1:A:338:LEU:HD12	1.94	0.49
1:C:297:ILE:HD12	1:C:297:ILE:C	2.38	0.49
1:B:312:LEU:HD23	1:B:324:ILE:CD1	2.43	0.48
1:D:195:VAL:HG22	6:D:501:HOH:O	2.14	0.48
1:D:189:THR:O	1:D:236:GLY:HA3	2.14	0.48
1:A:101:THR:HG22	1:A:102:VAL:N	2.30	0.47
1:A:107:LEU:HD23	1:A:107:LEU:O	2.14	0.47
1:B:138:LEU:HD12	1:B:139:PRO:HD2	1.96	0.47
1:B:186:LYS:HB3	1:B:234:ARG:HG2	1.96	0.47
1:C:183:VAL:O	1:C:229:CYS:HA	2.15	0.47
1:B:22:LEU:HD13	1:B:152:TYR:HD2	1.79	0.47
1:A:266:PHE:O	1:A:271:SER:HB3	2.15	0.46
1:B:184:GLY:HA3	1:B:232:ILE:HD11	1.97	0.46
1:B:148:ILE:C	1:B:150:PRO:HD2	2.41	0.46
1:A:39:ALA:O	1:A:43:ILE:HG13	2.15	0.46
1:C:130:LEU:HD11	1:C:157:MET:HE2	1.97	0.46
1:A:314:SER:HA	6:A:529:HOH:O	2.15	0.45
1:D:144:PHE:HE2	1:D:160:GLY:HA3	1.80	0.45
1:C:186:LYS:HD3	1:C:232:ILE:HB	1.99	0.45
1:D:297:ILE:C	1:D:297:ILE:HD12	2.41	0.45
1:A:234:ARG:HD2	1:A:235:GLY:O	2.17	0.45
1:D:20:VAL:CG1	1:D:214:LYS:HE2	2.46	0.45
1:D:144:PHE:CE2	1:D:160:GLY:HA3	2.52	0.45
1:D:159:TRP:CH2	1:D:161:ALA:HB2	2.51	0.45
1:D:141:ALA:HB1	1:D:159:TRP:HE3	1.82	0.44
1:B:22:LEU:HD13	1:B:152:TYR:CD2	2.51	0.44
1:A:166:THR:HG21	1:A:172:HIS:CE1	2.53	0.44
1:B:320:TYR:CD1	1:B:320:TYR:C	2.96	0.43
1:D:75:ARG:HB3	1:D:331:TRP:CZ2	2.53	0.43
1:D:96:GLU:CG	1:D:97:LYS:N	2.81	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:130:LEU:HD12	1:D:153:LEU:HD22	2.00	0.43
1:D:79:LEU:HD23	1:D:338:LEU:HD12	2.00	0.43
1:A:166:THR:HG22	1:B:169:SER:HB2	2.01	0.43
1:B:183:VAL:O	1:B:229:CYS:HA	2.19	0.43
1:A:289:ILE:O	1:A:348:ARG:HB2	2.19	0.43
1:A:346:LYS:O	1:A:347:ALA:C	2.60	0.43
1:B:73:ALA:HB1	1:B:138:LEU:HD23	2.01	0.42
1:D:113:MET:HB3	1:D:312:LEU:HD21	2.00	0.42
1:A:349:ARG:O	1:A:349:ARG:HG3	2.20	0.42
1:C:148:ILE:C	1:C:150:PRO:CD	2.92	0.42
1:C:165:ARG:HG2	1:D:168:GLU:HG2	2.02	0.42
1:D:312:LEU:HD13	1:D:318:LEU:HD11	2.01	0.42
1:D:104:TRP:CH2	1:D:106:GLY:HA2	2.55	0.42
1:D:303:SER:HA	1:D:328:CYS:HB3	2.01	0.42
1:D:187:ASN:OD1	1:D:233:LEU:HA	2.20	0.41
1:A:18:PRO:HA	1:A:19:PRO:HD3	1.89	0.41
1:A:316:GLU:HA	1:A:317:PRO:HD3	1.93	0.41
1:A:339:ARG:HH11	1:A:339:ARG:HG2	1.85	0.41
1:A:187:ASN:C	1:A:187:ASN:HD22	2.29	0.41
1:B:297:ILE:C	1:B:297:ILE:HD12	2.45	0.41
1:A:119:ILE:HG21	1:B:210:LEU:HD21	2.02	0.41
1:D:189:THR:C	1:D:191:GLY:H	2.25	0.41
1:B:55:LEU:O	1:B:298:GLY:HA2	2.21	0.41
1:B:309:ASN:HB3	1:B:327:ALA:HA	2.03	0.41
1:D:241:TYR:CE2	1:D:281:VAL:HG22	2.56	0.41
1:C:81:GLU:O	1:C:84:LYS:HD3	2.22	0.40
1:D:23:LEU:HD23	1:D:127:ARG:CZ	2.51	0.40
1:C:55:LEU:HD12	1:C:56:VAL:N	2.36	0.40
1:D:114:ASP:O	1:D:115:ASN:HB2	2.20	0.40
1:D:183:VAL:O	1:D:229:CYS:HA	2.22	0.40

There are no symmetry-related clashes.

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	343/351 (98%)	336 (98%)	7 (2%)	0	100	100
1	B	336/351 (96%)	328 (98%)	8 (2%)	0	100	100
1	C	342/351 (97%)	334 (98%)	8 (2%)	0	100	100
1	D	339/351 (97%)	320 (94%)	17 (5%)	2 (1%)	21	27
All	All	1360/1404 (97%)	1318 (97%)	40 (3%)	2 (0%)	48	60

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	189	THR
1	D	349	ARG

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	264/283 (93%)	262 (99%)	2 (1%)	73	86
1	B	262/283 (93%)	259 (99%)	3 (1%)	65	81
1	C	263/283 (93%)	263 (100%)	0	100	100
1	D	255/283 (90%)	254 (100%)	1 (0%)	84	92
All	All	1044/1132 (92%)	1038 (99%)	6 (1%)	78	89

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	99	ARG
1	A	187	ASN
1	B	22	LEU
1	B	312	LEU
1	B	320	TYR
1	D	234	ARG

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

Mol	Chain	Res	Type
1	A	112	HIS
1	A	118	GLN
1	A	134	ASN
1	B	34	ASN
1	B	112	HIS
1	B	217	HIS
1	C	112	HIS
1	C	217	HIS
1	C	288	GLN
1	D	134	ASN
1	D	288	GLN
1	D	340	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 11 ligands modelled in this entry, 4 are monoatomic - leaving 7 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	$\# Z > 2$
3	GAL	A	402	-	12,12,12	0.51	0	17,17,17	0.60	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	PEP	B	402	-	9,9,9	3.20	3 (33%)	11,13,13	1.91	4 (36%)
5	PO4	B	403	-	4,4,4	0.94	0	6,6,6	0.27	0
3	GAL	C	404	-	12,12,12	0.45	0	17,17,17	0.67	0
4	PEP	C	402	-	9,9,9	3.01	3 (33%)	11,13,13	1.85	3 (27%)
3	GAL	A	403	-	12,12,12	1.24	0	17,17,17	1.50	5 (29%)
5	PO4	C	403	-	4,4,4	0.69	0	6,6,6	1.12	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
3	GAL	A	402	-	-	0/2/22/22	0/1/1/1
4	PEP	B	402	-	-	0/9/9/9	-
3	GAL	C	404	-	-	2/2/22/22	0/1/1/1
4	PEP	C	402	-	-	0/9/9/9	-
3	GAL	A	403	-	-	1/2/22/22	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	402	PEP	C2-C1	-8.40	1.41	1.49
4	C	402	PEP	C2-C1	-7.79	1.41	1.49
4	B	402	PEP	C3-C2	3.31	1.40	1.31
4	C	402	PEP	C3-C2	3.26	1.40	1.31
4	C	402	PEP	O2'-C1	-2.60	1.23	1.30
4	B	402	PEP	O2'-C1	-2.59	1.23	1.30

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	402	PEP	O2-C2-C3	-4.59	116.08	124.85
4	B	402	PEP	O2-C2-C3	-4.12	116.98	124.85
3	A	403	GAL	C3-C4-C5	3.14	115.92	110.23
4	B	402	PEP	O1-C1-C2	-2.75	117.64	121.79
4	B	402	PEP	O2'-C1-C2	2.68	118.48	113.91
4	B	402	PEP	C3-C2-C1	2.45	127.26	122.73
3	A	403	GAL	O3-C3-C4	-2.27	105.02	110.38
4	C	402	PEP	C3-C2-C1	2.15	126.70	122.73
3	A	403	GAL	O5-C5-C4	-2.11	105.90	109.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	403	GAL	O2-C2-C3	-2.11	105.41	110.38
4	C	402	PEP	O1-C1-C2	-2.06	118.69	121.79
3	A	403	GAL	O4-C4-C5	-2.01	104.37	109.32

There are no chirality outliers.

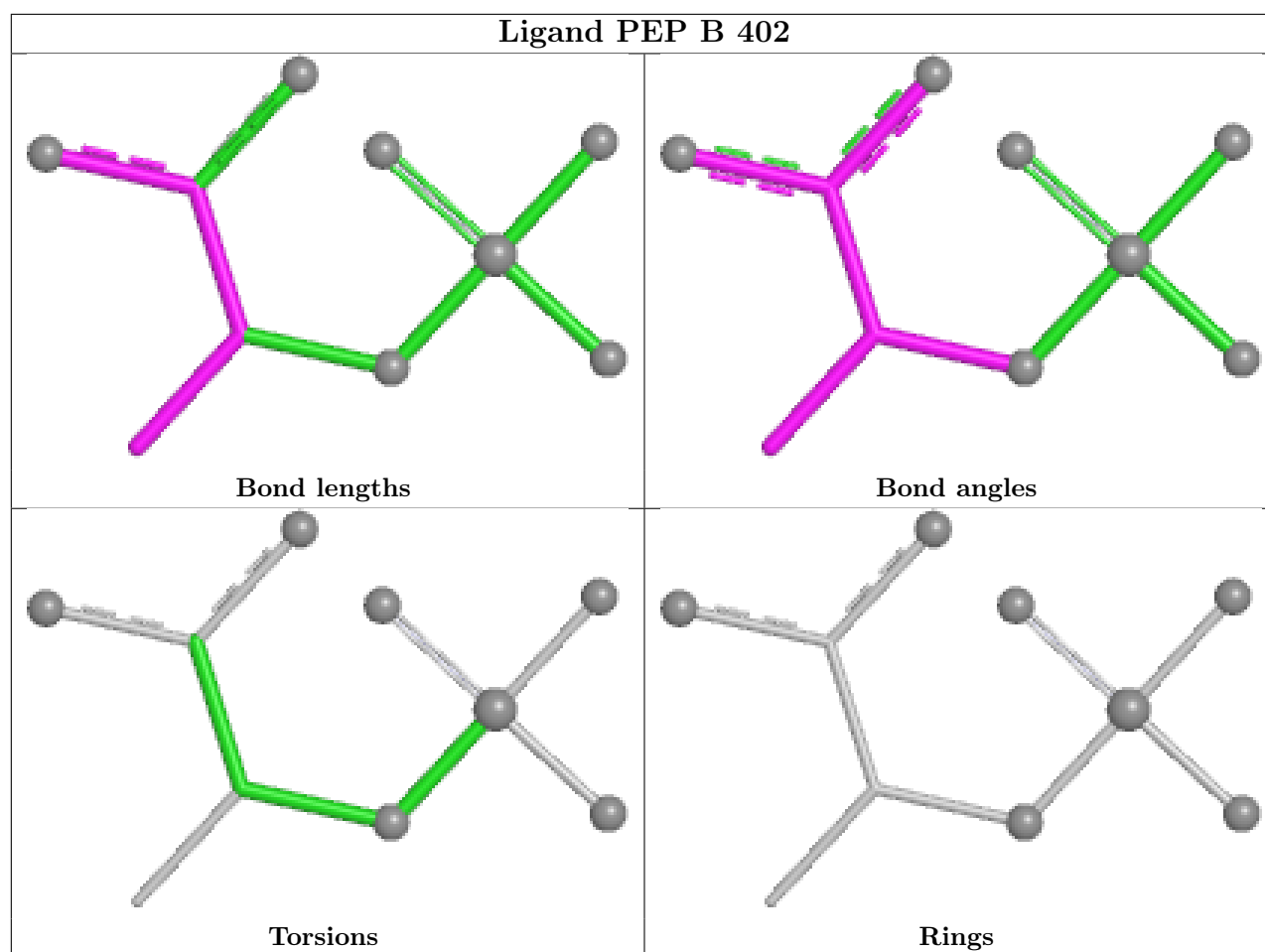
All (3) torsion outliers are listed below:

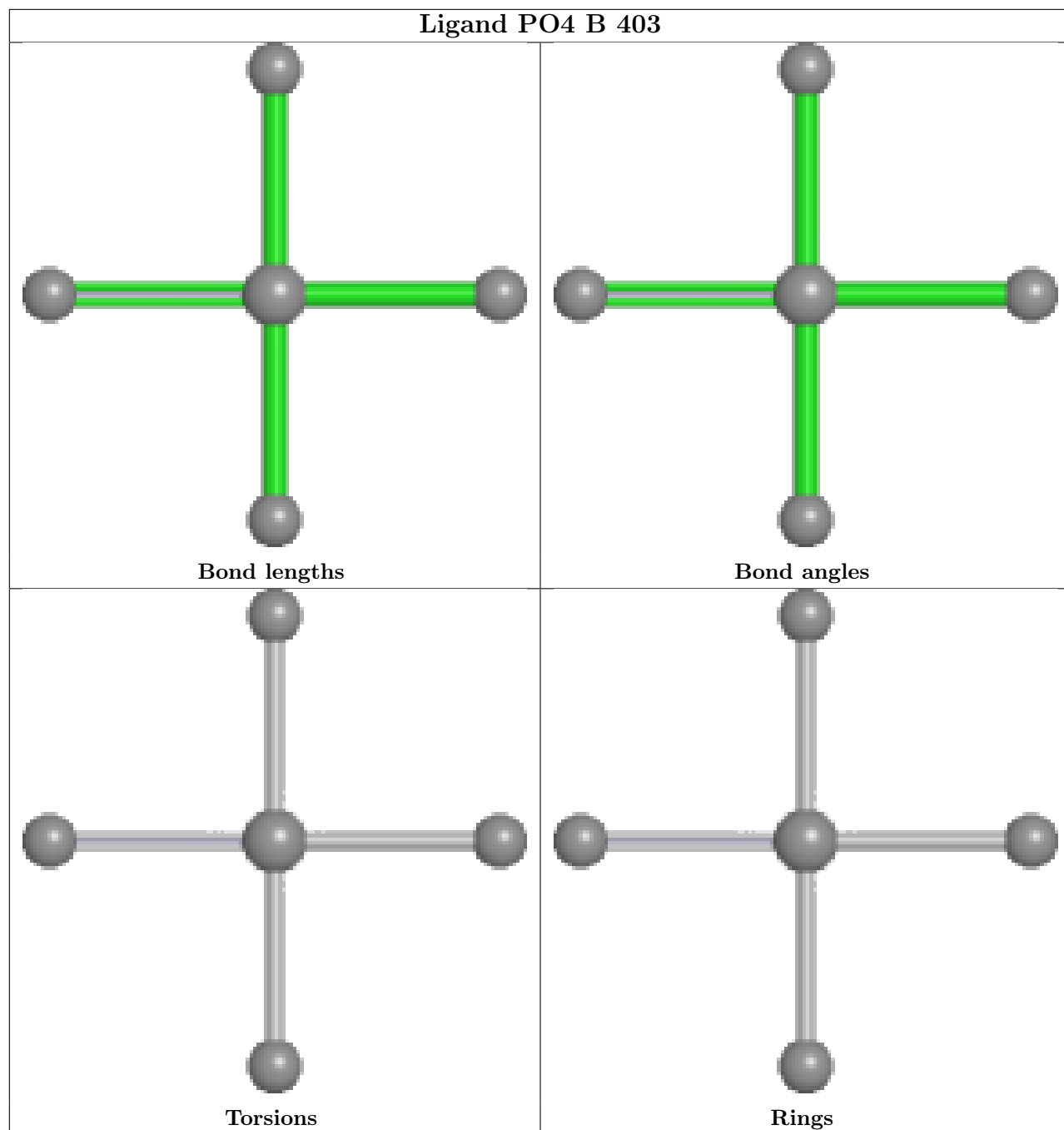
Mol	Chain	Res	Type	Atoms
3	C	404	GAL	C4-C5-C6-O6
3	C	404	GAL	O5-C5-C6-O6
3	A	403	GAL	O5-C5-C6-O6

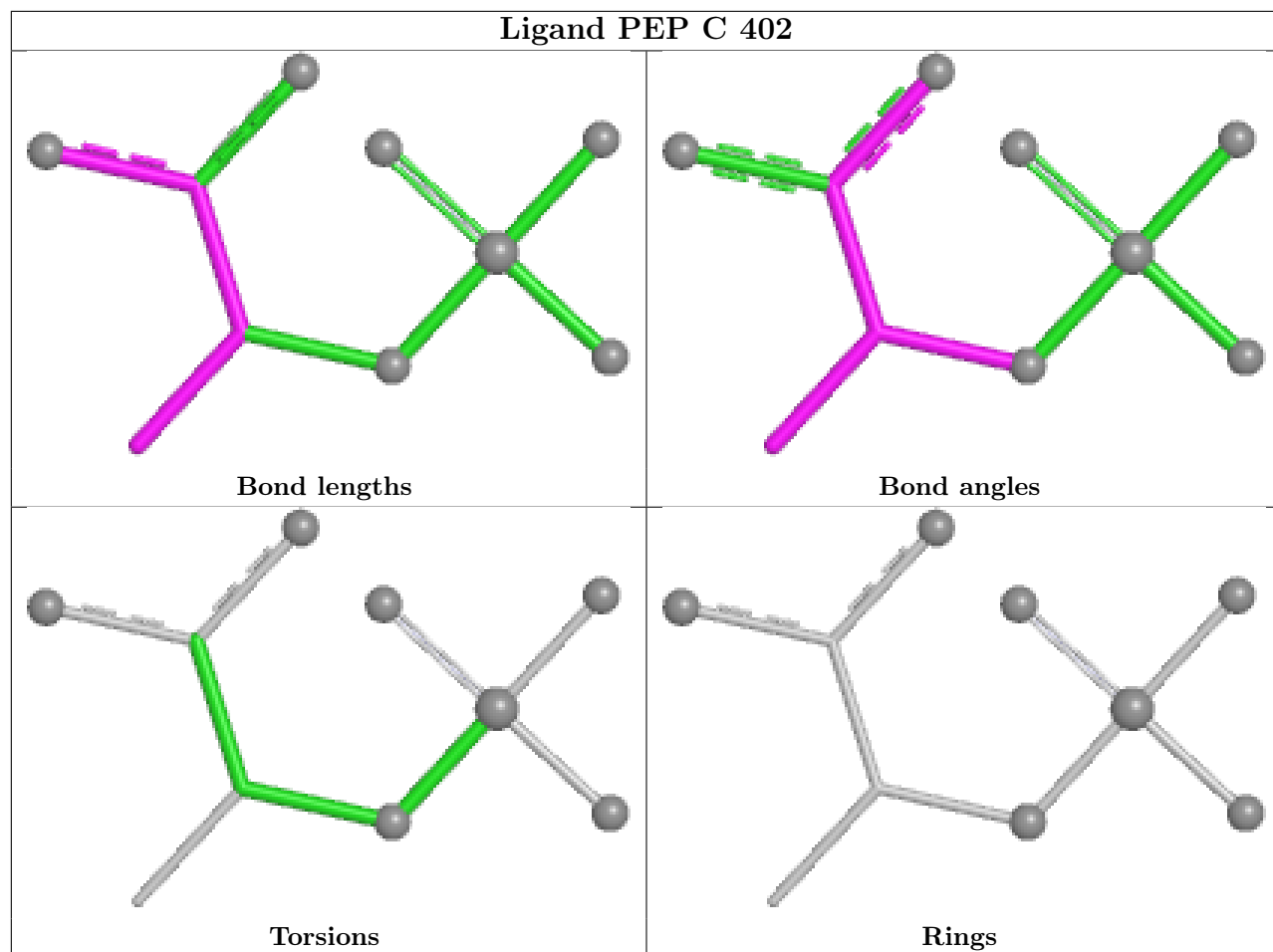
There are no ring outliers.

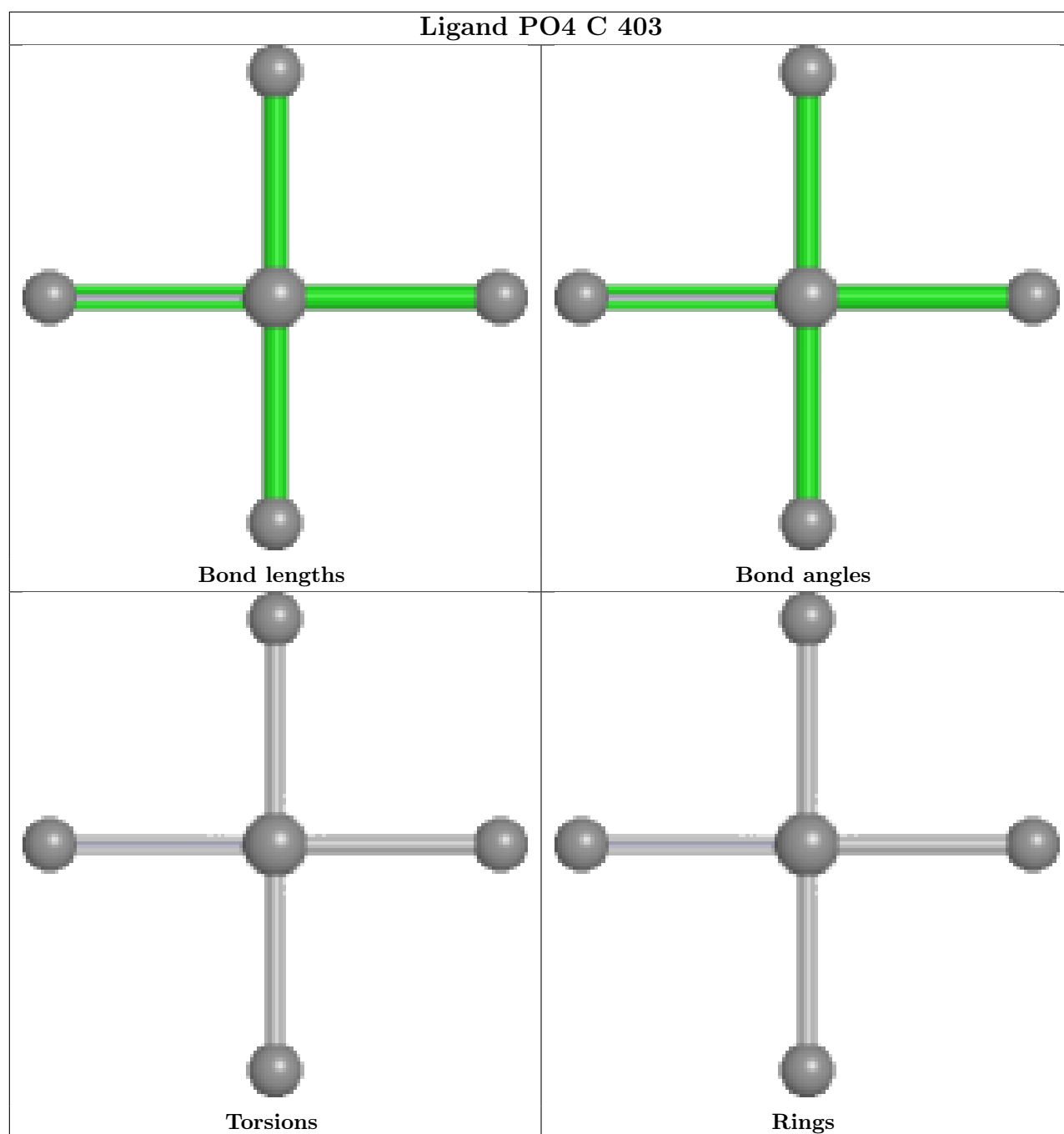
No monomer is involved in short contacts.

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	344/351 (98%)	0.27	26 (7%) 20 21	19, 28, 38, 44	1 (0%)
1	B	340/351 (96%)	0.09	6 (1%) 67 69	24, 32, 38, 43	0
1	C	344/351 (98%)	-0.27	3 (0%) 81 82	16, 22, 27, 34	0
1	D	343/351 (97%)	0.43	30 (8%) 16 17	20, 30, 37, 42	0
All	All	1371/1404 (97%)	0.13	65 (4%) 36 38	16, 28, 37, 44	1 (0%)

All (65) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	315	GLY	6.0
1	D	103	GLY	5.6
1	C	350	GLY	5.6
1	A	236	GLY	5.4
1	D	100	THR	5.2
1	D	189	THR	5.1
1	A	315	GLY	4.9
1	D	314	SER	4.7
1	D	6	ASP	4.7
1	D	102	VAL	4.6
1	A	271	SER	4.6
1	B	6	ASP	4.6
1	A	314	SER	4.5
1	D	317	PRO	4.5
1	D	350	GLY	4.5
1	D	101	THR	4.5
1	D	236	GLY	4.3
1	D	271	SER	3.8
1	D	237	LYS	3.8
1	A	269	ALA	3.7
1	A	268	HIS	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	350	GLY	3.4
1	D	269	ALA	3.3
1	D	99	ARG	3.3
1	B	312	LEU	3.2
1	A	235	GLY	3.2
1	D	235	GLY	3.1
1	D	311	SER	3.1
1	A	101	THR	2.9
1	A	272	SER	2.9
1	D	313	GLU	2.9
1	A	100	THR	2.9
1	A	237	LYS	2.8
1	D	98	PRO	2.8
1	B	317	PRO	2.8
1	A	239	PRO	2.7
1	B	7	ASP	2.7
1	D	104	TRP	2.6
1	A	238	GLU	2.6
1	D	192	THR	2.6
1	D	268	HIS	2.5
1	D	191	GLY	2.5
1	C	313	GLU	2.4
1	D	312	LEU	2.4
1	A	102	VAL	2.4
1	A	7	ASP	2.4
1	A	115	ASN	2.3
1	A	273	LYS	2.3
1	A	270	ASN	2.3
1	D	188	GLY	2.3
1	A	189	THR	2.3
1	D	316	GLU	2.3
1	A	274	GLN	2.2
1	D	190	ASP	2.1
1	A	245	HIS	2.1
1	A	116	SER	2.1
1	A	117	PHE	2.1
1	A	349	ARG	2.1
1	A	240	ASN	2.1
1	B	215	TRP	2.1
1	C	7	ASP	2.0
1	D	115	ASN	2.0
1	D	97	LYS	2.0

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Mol	Chain	Res	Type	RSRZ
1	B	311	SER	2.0
1	D	96	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](#)

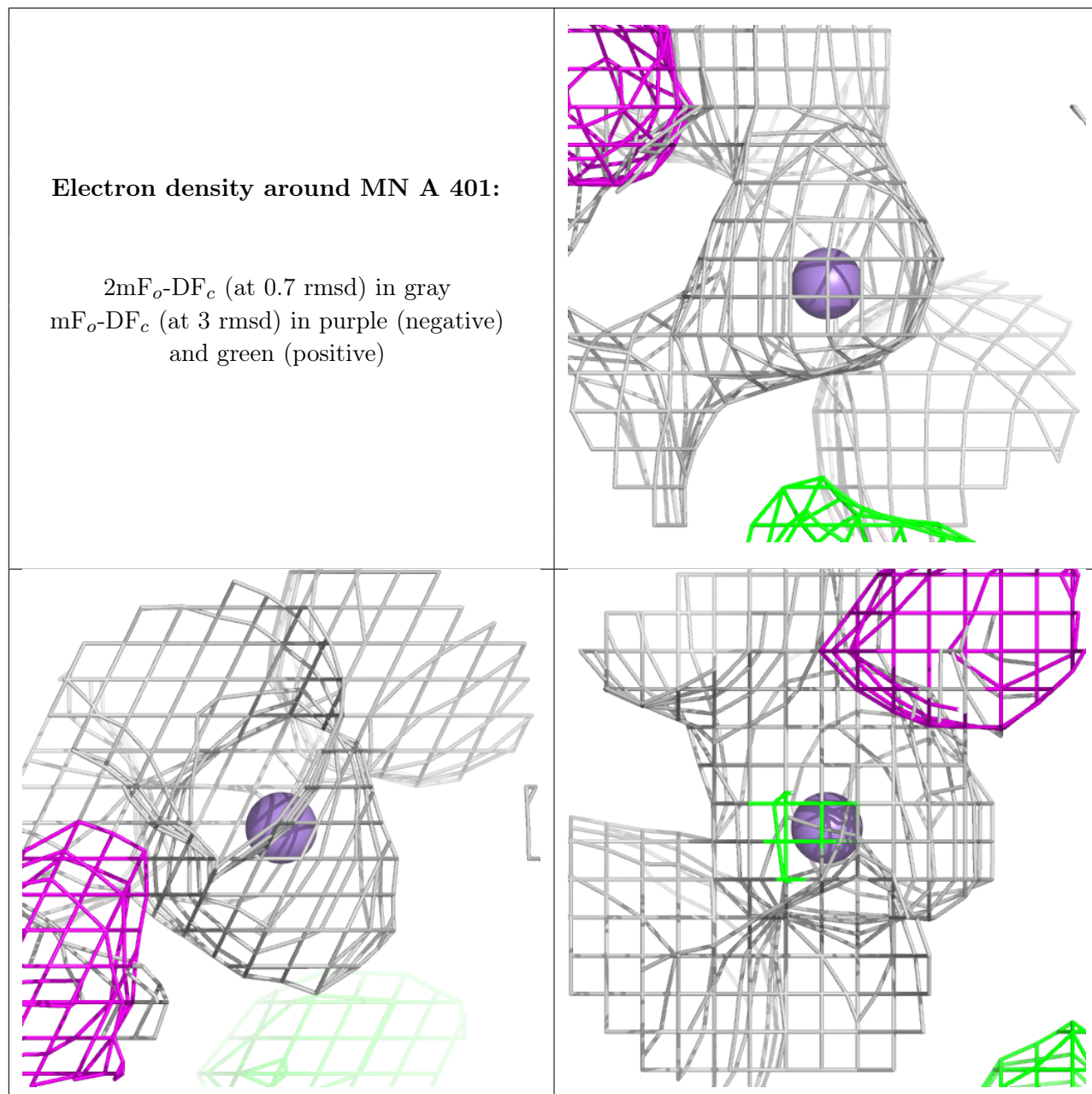
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	GAL	C	404	12/12	0.83	0.13	24,28,37,42	24
2	MN	A	401	1/1	0.87	0.11	31,31,31,31	1
5	PO4	B	403	5/5	0.87	0.13	31,34,35,36	5
3	GAL	A	402	12/12	0.93	0.08	25,28,32,38	24
4	PEP	B	402	10/10	0.95	0.09	30,31,32,32	0
4	PEP	C	402	10/10	0.95	0.08	19,20,21,21	0
3	GAL	A	403	12/12	0.95	0.08	22,25,27,28	0
5	PO4	C	403	5/5	0.95	0.09	23,25,26,26	5
2	MN	D	401	1/1	0.96	0.27	38,38,38,38	0
2	MN	B	401	1/1	0.98	0.03	32,32,32,32	1
2	MN	C	401	1/1	0.99	0.03	20,20,20,20	1

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.

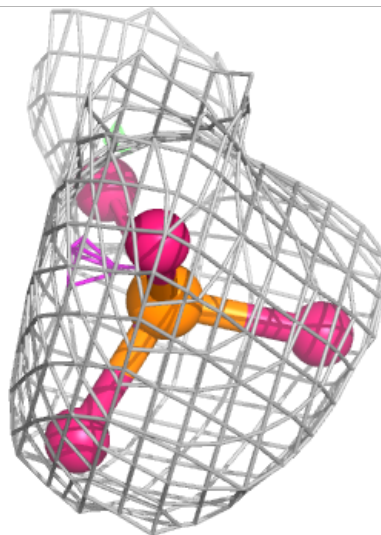
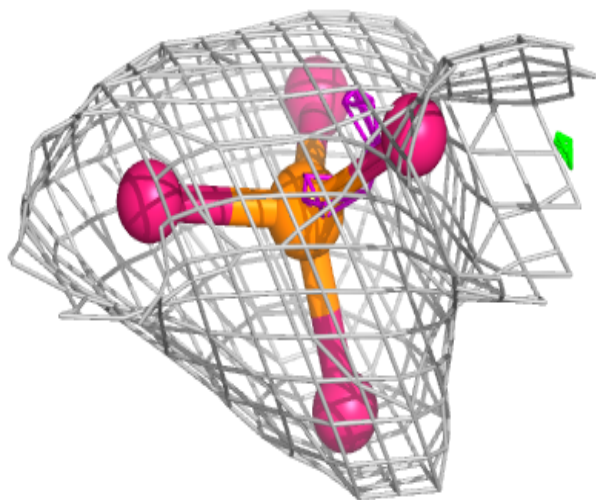
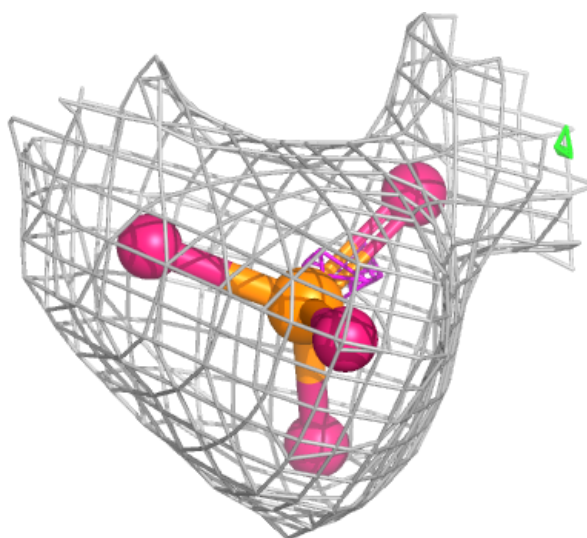
Electron density around MN A 401:

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



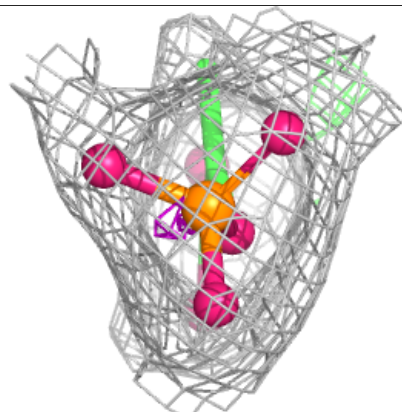
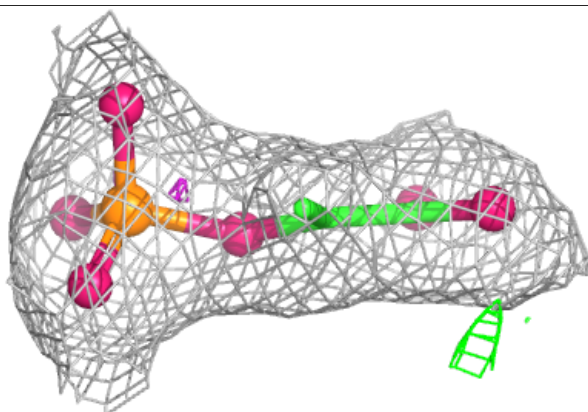
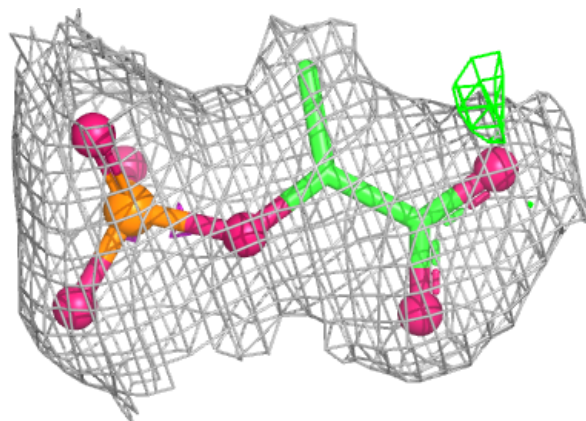
Electron density around PO4 B 403:

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



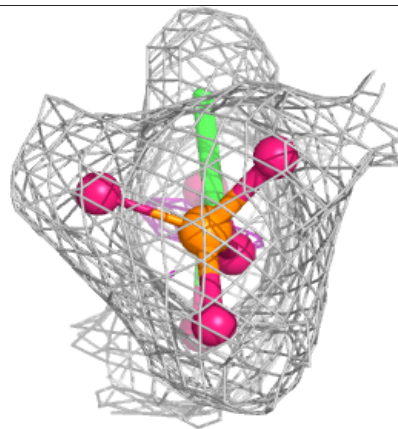
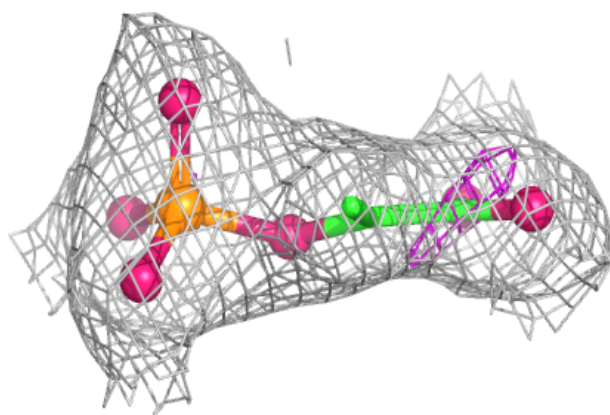
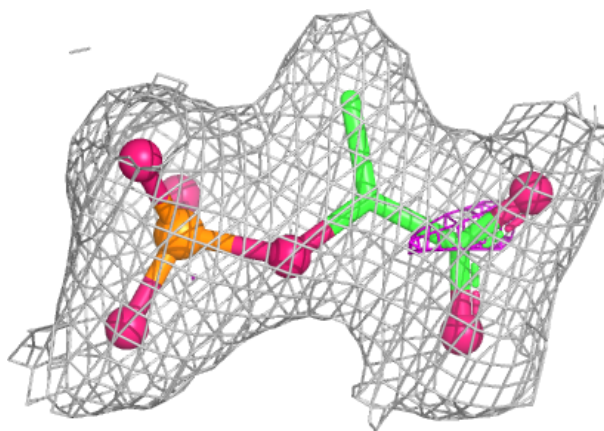
Electron density around PEP B 402:

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



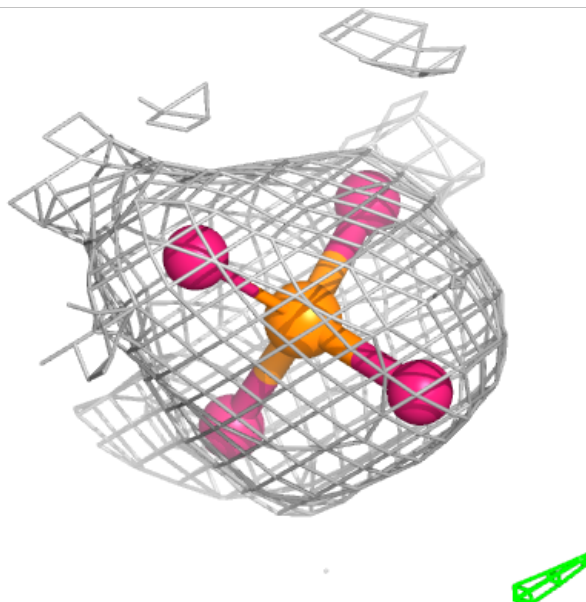
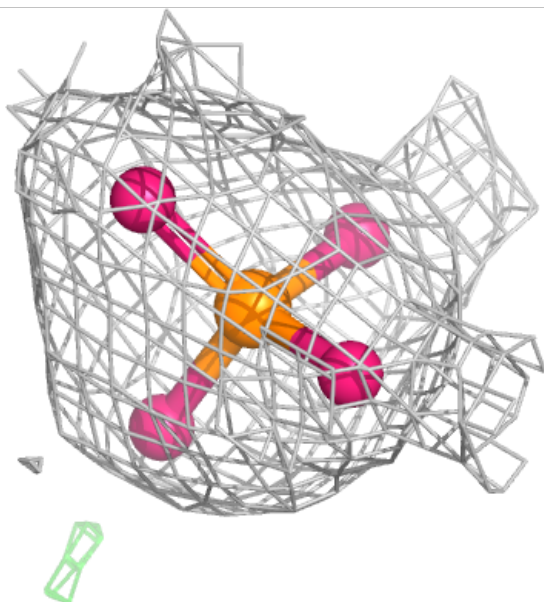
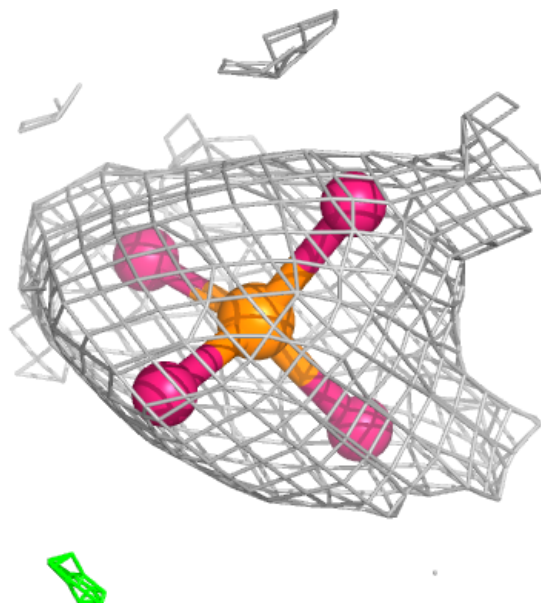
Electron density around PEP C 402:

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



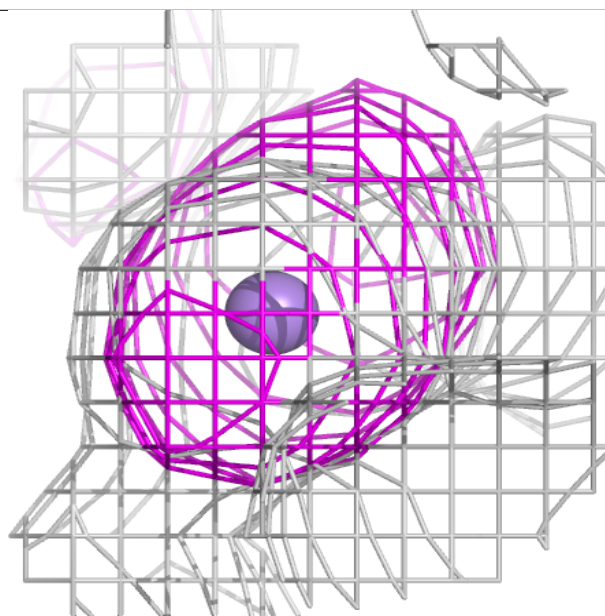
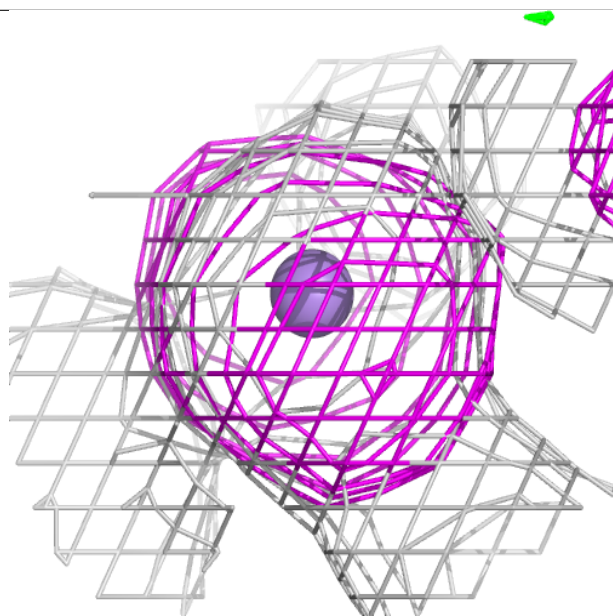
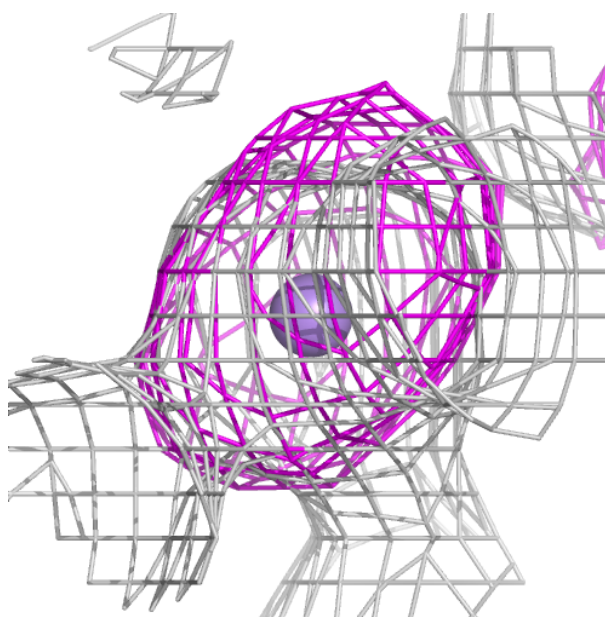
Electron density around PO4 C 403:

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



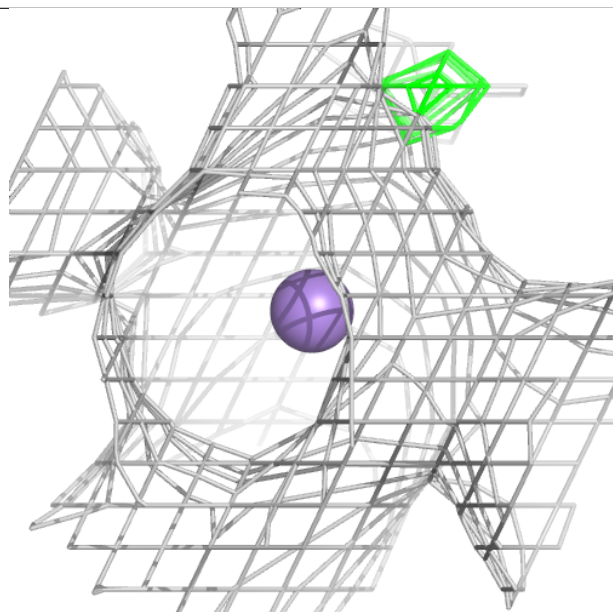
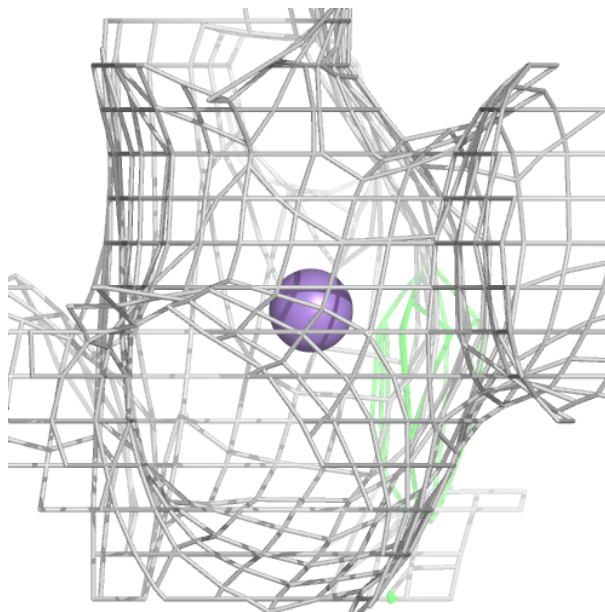
Electron density around MN D 401:

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



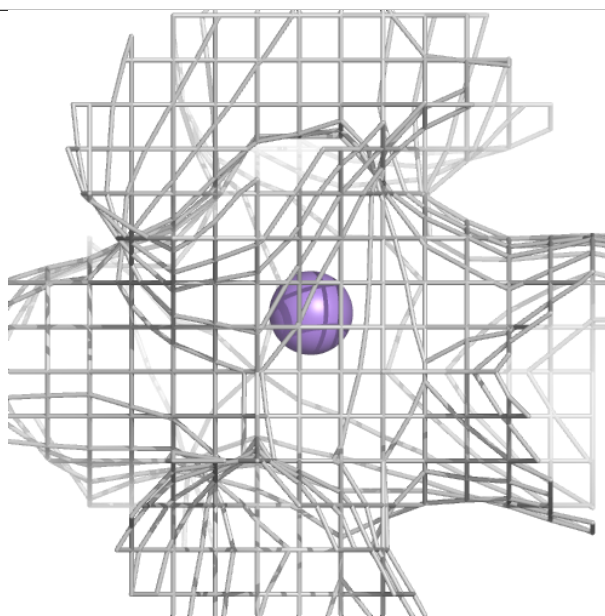
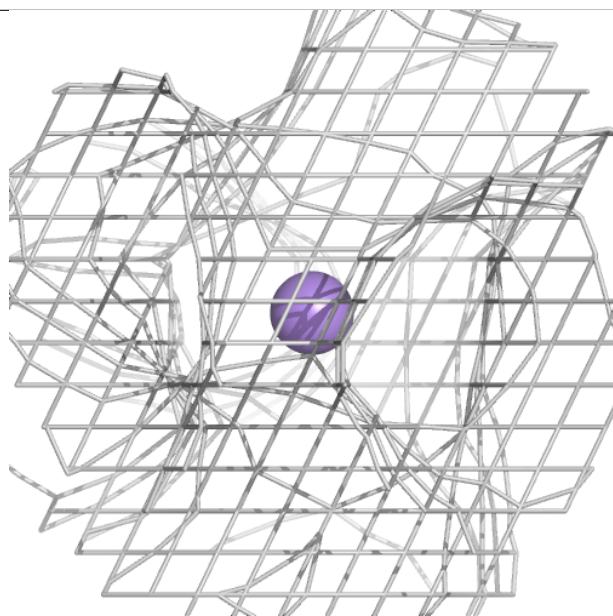
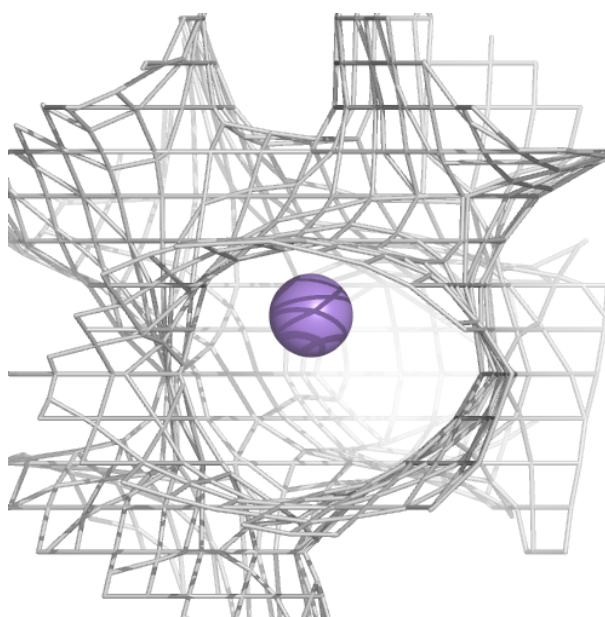
Electron density around MN B 401:

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



Electron density around MN C 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.