Full wwPDB X-ray Structure Validation Report (i)

Jul 5, 2022 – 03:02 PM JST

PDB ID	:	7Y8I
Title	:	Crystal structure of sDscam FNIII3 domain, isoform alpha7
Deposited on	:	2022-06-24
Resolution	:	1.90 Å(reported)

This wwPDB validation report is for manuscript review

This is a Full wwPDB X-ray Structure Validation Report.

This report is produced by the wwPDB biocuration pipeline after annotation of the structure.

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 (i) 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 (i)) 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	:	2.29
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)	:	2.29

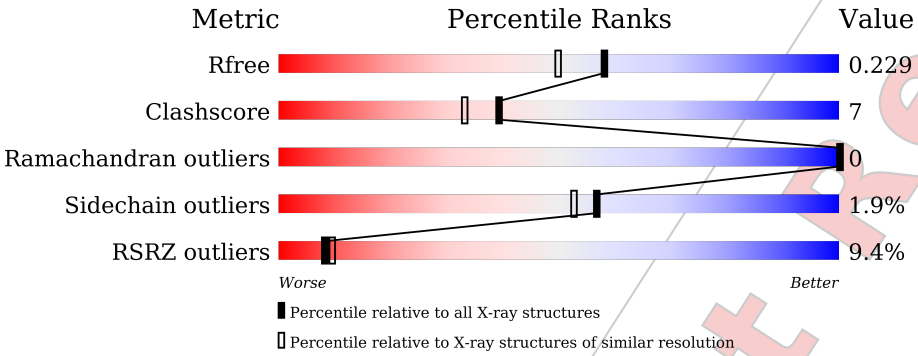
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1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

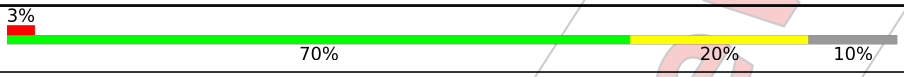
The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	97	4% 74% 10% 15%
1	B	97	16% 74% 12% • 12%
1	C	97	6% 73% 16% 10%
1	D	97	12% 78% 18% •
1	E	97	9% 85% 15%

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Mol	Chain	Length	Quality of chain
1	F	97	

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2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 8916 atoms, of which 4218 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 Dscam.

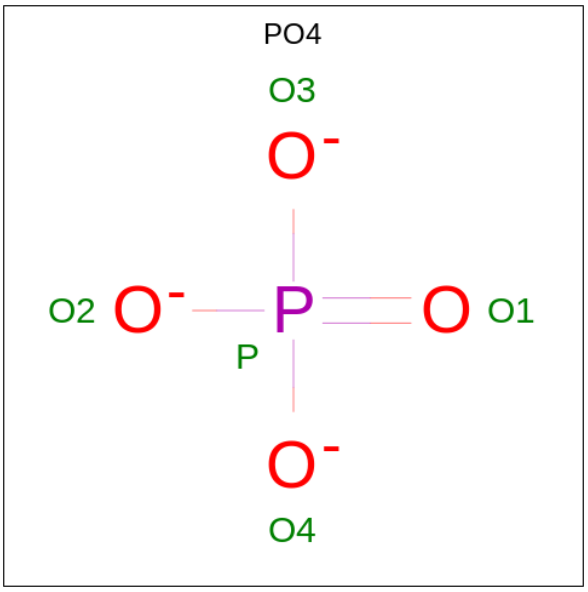
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	82	Total	C	H	N	O	0	0	0
			1313	425	648	121	119			
1	B	85	Total	C	H	N	O	0	0	0
			1351	437	662	125	127			
1	C	87	Total	C	H	N	O	0	2	0
			1397	452	688	129	128			
1	D	93	Total	C	H	N	O	0	4	0
			1503	486	742	137	138			
1	E	97	Total	C	H	N	O	0	1	0
			1551	494	770	143	144			
1	F	87	Total	C	H	N	O	0	0	0
			1382	443	682	130	127			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	493	ALA	ASP	conflict	UNP A0A159W468
A	494	SER	PRO	conflict	UNP A0A159W468
B	493	ALA	ASP	conflict	UNP A0A159W468
B	494	SER	PRO	conflict	UNP A0A159W468
C	493	ALA	ASP	conflict	UNP A0A159W468
C	494	SER	PRO	conflict	UNP A0A159W468
D	493	ALA	ASP	conflict	UNP A0A159W468
D	494	SER	PRO	conflict	UNP A0A159W468
E	493	ALA	ASP	conflict	UNP A0A159W468
E	494	SER	PRO	conflict	UNP A0A159W468
F	493	ALA	ASP	conflict	UNP A0A159W468
F	494	SER	PRO	conflict	UNP A0A159W468

- Molecule 2 is PHOSPHATE ION (three-letter code: PO4) (formula: O_4P) (labeled as "Ligand of Interest" by depositor).

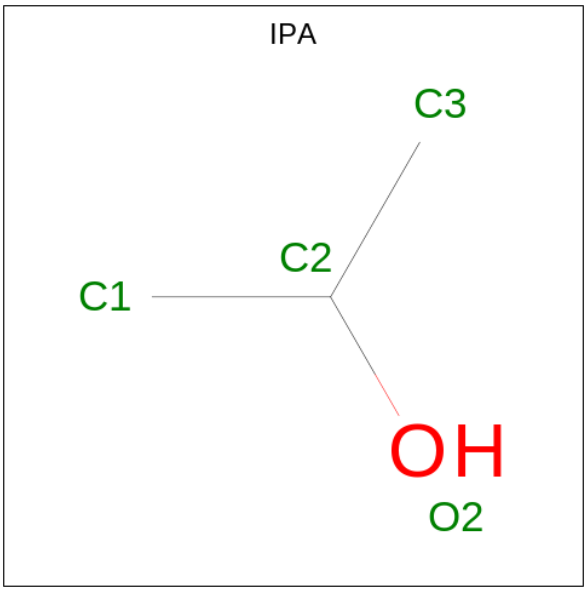
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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total O P 5 4 1	0	0
2	B	1	Total O P 5 4 1	0	0
2	B	1	Total O P 5 4 1	0	0
2	D	1	Total O P 5 4 1	0	0
2	E	1	Total O P 5 4 1	0	0
2	E	1	Total O P 5 4 1	0	0
2	F	1	Total O P 5 4 1	0	0
2	F	1	Total O P 5 4 1	0	0

- Molecule 3 is ISOPROPYL ALCOHOL (three-letter code: IPA) (formula: C_3H_8O) (labeled as "Ligand of Interest" by depositor).

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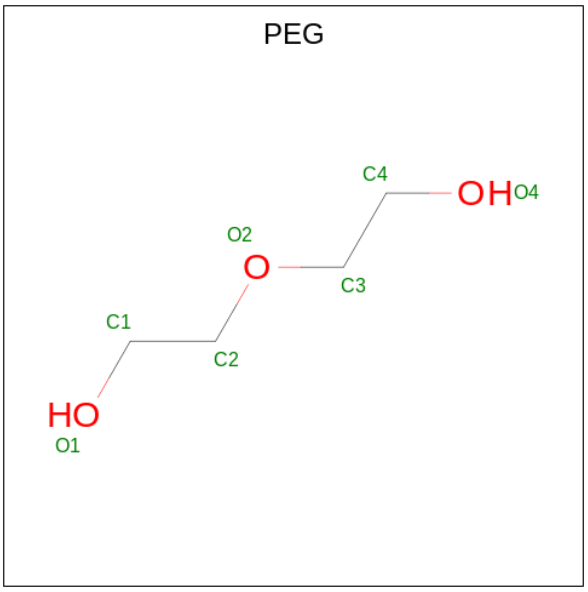
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total C H O 12 3 8 1	0	0
3	C	1	Total C H O 12 3 8 1	0	0

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	E	1	Total Cl 1 1	0	0

- Molecule 5 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: $C_4H_{10}O_3$) (labeled as "Ligand of Interest" by depositor).

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	F	1	Total C H O 17 4 10 3	0	0

- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	57	Total O 57 57	0	0
6	B	48	Total O 48 48	0	0
6	C	57	Total O 57 57	0	0
6	D	49	Total O 49 49	0	0
6	E	57	Total O 57 57	0	0
6	F	69	Total O 69 69	0	0

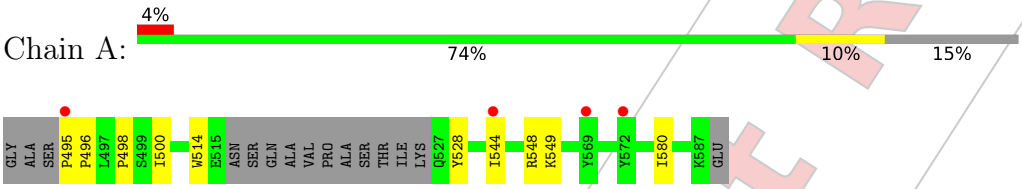
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3 Residue-property plots (i)

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.

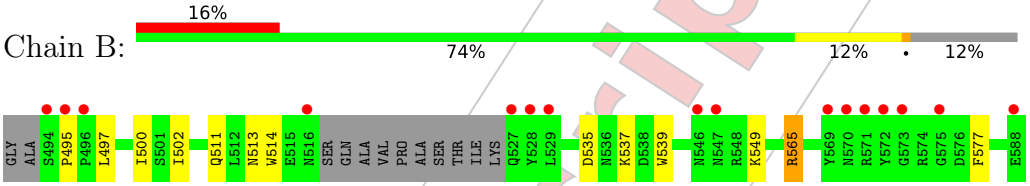
- Molecule 1: Dscam

Chain A:



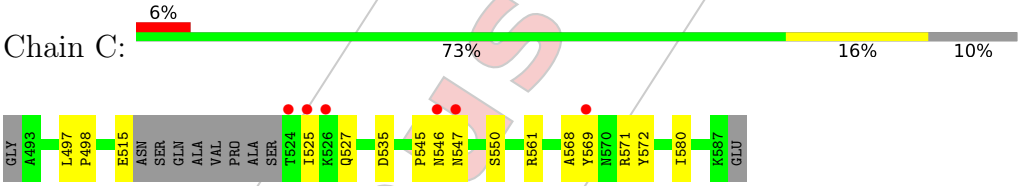
- Molecule 1: Dscam

Chain B:



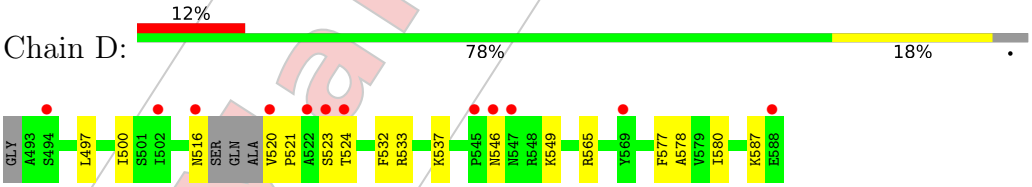
- Molecule 1: Dscam

Chain C:



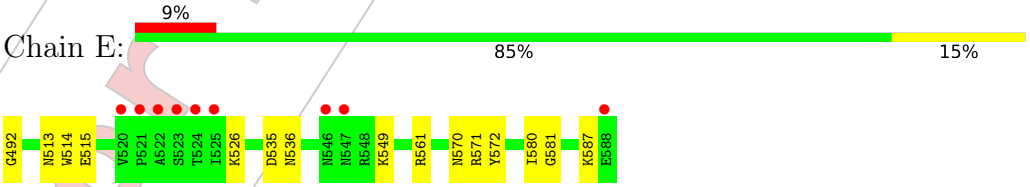
- Molecule 1: Dscam

Chain D:



- Molecule 1: Dscam

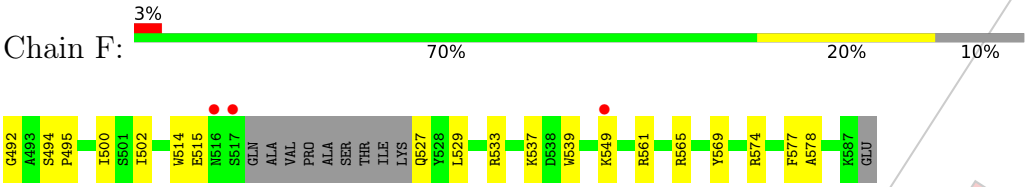
Chain E:



- Molecule 1: Dscam

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Chain F:



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4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	69.11Å 85.78Å 106.24Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.77 – 1.90 39.77 – 1.90	Depositor EDS
% Data completeness (in resolution range)	91.7 (39.77-1.90) 91.7 (39.77-1.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.00 (at 1.89Å)	Xtriage
Refinement program	PHENIX 1.15.2_3472	Depositor
R, R_{free}	0.198 , 0.229 0.198 , 0.229	Depositor DCC
R_{free} test set	911 reflections (1.96%)	wwPDB-VP
Wilson B-factor (Å^2)	17.6	Xtriage
Anisotropy	0.171	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.43 , 52.5	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	8916	wwPDB-VP
Average B, all atoms (Å^2)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 16.41% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

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5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CL, PO4, PEG, IPA

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	0.41	0/682	0.63	0/922
1	B	0.33	0/706	0.54	0/954
1	C	0.71	1/732 (0.1%)	0.58	0/989
1	D	0.35	0/789	0.54	0/1068
1	E	0.35	0/800	0.59	0/1083
1	F	0.40	0/717	0.62	0/969
All	All	0.44	1/4426 (0.0%)	0.58	0/5985

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	515	GLU	CD-OE2	-16.69	1.07	1.25

There are no bond angle outliers.

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	665	648	651	7	0
1	B	689	662	667	9	0
1	C	709	688	697	8	0
1	D	761	742	749	11	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	781	770	772	17	0
1	F	700	682	685	14	0
2	A	5	0	0	0	0
2	B	10	0	0	2	0
2	D	5	0	0	0	0
2	E	10	0	0	0	0
2	F	10	0	0	0	0
3	B	4	8	8	0	0
3	C	4	8	8	0	0
4	E	1	0	0	0	0
5	F	7	10	10	0	0
6	A	57	0	0	1	1
6	B	48	0	0	5	0
6	C	57	0	0	0	0
6	D	49	0	0	3	1
6	E	57	0	0	4	0
6	F	69	0	0	3	0
All	All	4698	4218	4247	64	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 7.

All (64) 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:516:ASN:ND2	6:D:702:HOH:O	2.21	0.71
1:D:546:ASN:O	6:D:701:HOH:O	2.09	0.70
2:B:703:PO4:O3	6:B:801:HOH:O	2.11	0.68
1:A:498:PRO:O	6:A:701:HOH:O	2.11	0.67
1:E:515:GLU:HG3	1:E:549:LYS:HB2	1.77	0.66
1:E:515:GLU:N	6:E:701:HOH:O	2.16	0.66
1:E:515:GLU:OE2	6:E:701:HOH:O	2.14	0.65
2:B:702:PO4:O1	1:E:492:GLY:N	2.31	0.62
1:B:495:PRO:O	6:B:802:HOH:O	2.16	0.61
1:F:565:ARG:HG3	1:F:577:PHE:CD1	2.36	0.61
1:E:561:ARG:HB2	1:E:561:ARG:HH21	1.66	0.60
1:E:571:ARG:HD3	1:E:572:TYR:CZ	2.37	0.60
1:F:515:GLU:HA	1:F:515:GLU:OE1	2.02	0.59
1:D:497:LEU:HD11	1:D:577:PHE:O	2.03	0.58
1:E:514:TRP:N	6:E:701:HOH:O	2.37	0.57
1:D:500:ILE:HD11	1:D:580[B]:ILE:HG23	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:521:PRO:O	1:D:524:THR:HG22	2.08	0.54
1:C:571:ARG:HD3	1:C:572:TYR:CZ	2.44	0.53
1:B:513:ASN:ND2	6:B:803:HOH:O	2.28	0.53
1:E:526:LYS:N	1:E:526:LYS:HD2	2.22	0.53
1:F:561:ARG:NH1	6:F:703:HOH:O	2.41	0.53
1:A:528:TYR:HB2	1:A:544:ILE:HD11	1.91	0.52
1:B:511:GLN:NE2	6:B:805:HOH:O	2.44	0.51
1:E:535:ASP:OD2	1:E:561:ARG:NH1	2.44	0.51
1:E:513:ASN:HB3	6:E:701:HOH:O	2.12	0.50
1:B:502:ILE:HD11	1:F:502:ILE:HD11	1.93	0.50
1:D:500:ILE:HD12	1:D:580[B]:ILE:HG12	1.94	0.50
1:E:515:GLU:HG3	1:E:549:LYS:CB	2.41	0.49
1:D:546:ASN:N	1:D:546:ASN:OD1	2.46	0.49
1:A:548:ARG:C	1:A:549:LYS:HD3	2.33	0.48
1:B:565:ARG:HD3	1:B:577:PHE:CD1	2.49	0.48
1:B:497:LEU:HD11	1:B:577:PHE:O	2.15	0.47
1:C:497:LEU:HD12	1:C:498:PRO:O	2.14	0.47
1:F:527:GLN:N	1:F:527:GLN:OE1	2.48	0.46
1:F:529:LEU:HD12	6:F:707:HOH:O	2.15	0.46
1:B:537:LYS:NZ	6:B:807:HOH:O	2.49	0.46
1:E:526:LYS:HE3	1:E:570:ASN:HD22	1.81	0.46
1:D:565:ARG:HA	1:D:578:ALA:O	2.16	0.45
1:D:533:ARG:HD2	1:D:537:LYS:O	2.16	0.45
1:E:561:ARG:HB2	1:E:561:ARG:NH2	2.31	0.45
1:C:545:PRO:HB2	1:C:547:ASN:OD1	2.17	0.45
1:B:539:TRP:CZ3	1:B:565:ARG:HG2	2.51	0.45
1:C:561:ARG:HD3	1:E:571:ARG:O	2.16	0.45
1:A:548:ARG:O	1:A:549:LYS:HD3	2.18	0.44
1:F:539:TRP:CE3	1:F:565:ARG:HD2	2.53	0.44
1:A:500:ILE:HG22	1:A:514:TRP:HB3	1.99	0.44
1:F:500:ILE:HG22	1:F:514:TRP:HB3	2.00	0.43
1:F:492:GLY:N	1:F:574:ARG:O	2.51	0.43
1:A:495:PRO:HA	1:A:496:PRO:HD3	1.87	0.43
1:C:527:GLN:HG3	1:C:569:TYR:CD2	2.54	0.43
1:F:565:ARG:HA	1:F:578:ALA:O	2.19	0.42
1:E:535:ASP:O	1:E:536:ASN:HB2	2.19	0.42
1:F:494:SER:HA	1:F:495:PRO:C	2.39	0.42
1:F:533:ARG:HD2	1:F:537:LYS:O	2.20	0.42
1:A:580:ILE:HG23	1:C:580:ILE:HD11	2.00	0.42
1:D:549:LYS:NZ	6:D:707:HOH:O	2.54	0.41
1:D:520:VAL:HB	1:D:524:THR:HG21	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:549:LYS:NZ	6:F:710:HOH:O	2.53	0.41
1:B:500:ILE:HG22	1:B:514:TRP:HB3	2.01	0.41
1:C:525:ILE:HD12	1:C:568:ALA:HB1	2.02	0.41
1:C:545:PRO:CB	1:C:547:ASN:OD1	2.69	0.41
1:E:580:ILE:HD12	1:E:581:GLY:H	1.85	0.41
1:F:527:GLN:HB2	1:F:569:TYR:CZ	2.57	0.40
1:E:587:LYS:HD3	1:E:587:LYS:HA	1.85	0.40

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:A:731:HOH:O	6:D:746:HOH:O[4_555]	2.00	0.20

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	78/97 (80%)	76 (97%)	2 (3%)	0	100	100
1	B	81/97 (84%)	80 (99%)	1 (1%)	0	100	100
1	C	85/97 (88%)	84 (99%)	1 (1%)	0	100	100
1	D	93/97 (96%)	92 (99%)	1 (1%)	0	100	100
1	E	96/97 (99%)	94 (98%)	2 (2%)	0	100	100
1	F	83/97 (86%)	81 (98%)	2 (2%)	0	100	100
All	All	516/582 (89%)	507 (98%)	9 (2%)	0	100	100

There are no Ramachandran outliers to report.

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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	72/84 (86%)	72 (100%)	0	100	100
1	B	75/84 (89%)	72 (96%)	3 (4%)	31	22
1	C	78/84 (93%)	75 (96%)	3 (4%)	33	24
1	D	84/84 (100%)	81 (96%)	3 (4%)	35	26
1	E	85/84 (101%)	85 (100%)	0	100	100
1	F	76/84 (90%)	76 (100%)	0	100	100
All	All	470/504 (93%)	461 (98%)	9 (2%)	57	53

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	535	ASP
1	B	549	LYS
1	B	565	ARG
1	C	535	ASP
1	C	546	ASN
1	C	550	SER
1	D	523	SER
1	D	532	PHE
1	D	587	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	547	ASN
1	E	543	HIS
1	F	527	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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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

Of 12 ligands modelled in this entry, 1 is monoatomic - leaving 11 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	F	601	-	6,6,6	0.45	0	5,5,5	0.77	0
2	PO4	B	703	-	4,4,4	0.88	0	6,6,6	0.83	0
2	PO4	A	601	-	4,4,4	0.73	0	6,6,6	0.53	0
2	PO4	E	602	-	4,4,4	0.85	0	6,6,6	0.65	0
2	PO4	B	702	-	4,4,4	1.05	0	6,6,6	0.89	0
3	IPA	C	601	-	3,3,3	0.59	0	3,3,3	0.11	0
2	PO4	F	602	-	4,4,4	0.82	0	6,6,6	0.58	0
2	PO4	D	601	-	4,4,4	1.05	0	6,6,6	0.60	0
2	PO4	F	603	-	4,4,4	1.07	0	6,6,6	0.43	0
2	PO4	E	601	-	4,4,4	0.85	0	6,6,6	0.70	0
3	IPA	B	701	-	3,3,3	0.53	0	3,3,3	0.15	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	F	601	-	-	1/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

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There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	F	601	PEG	O1-C1-C2-O2

There are no ring outliers.

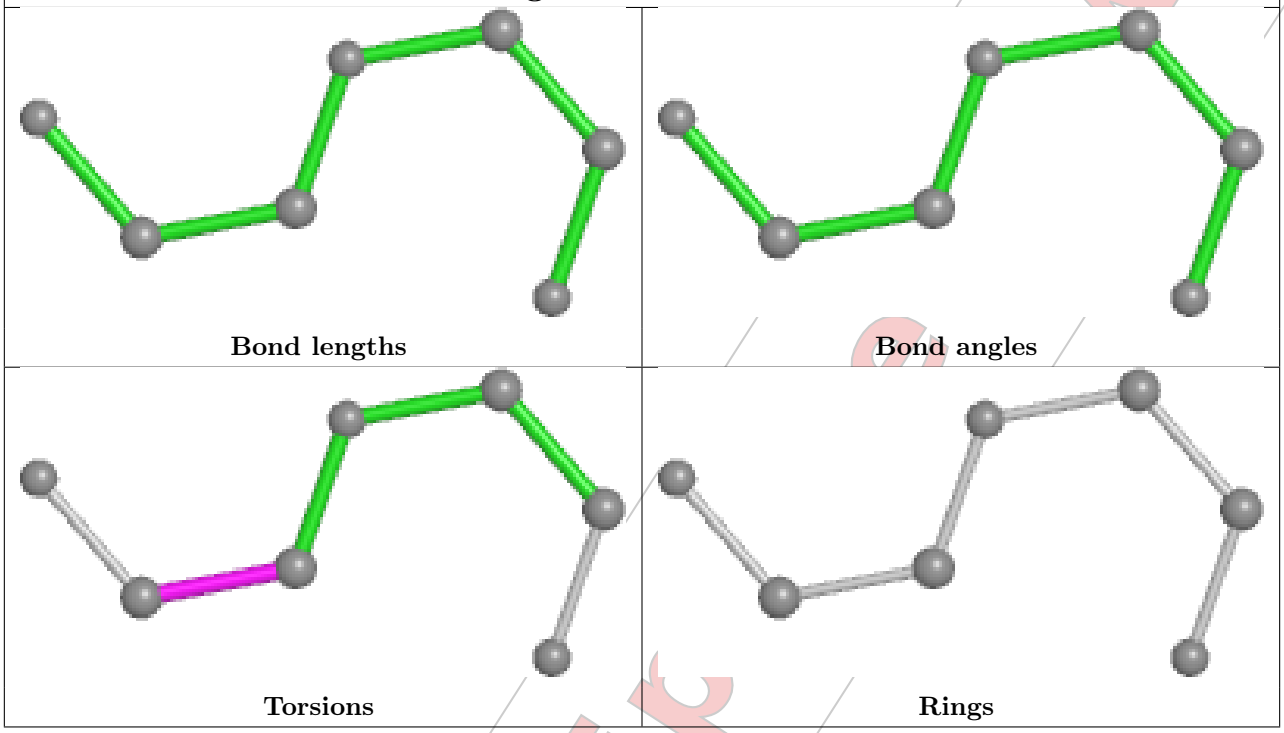
2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	703	PO4	1	0
2	B	702	PO4	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 then 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.

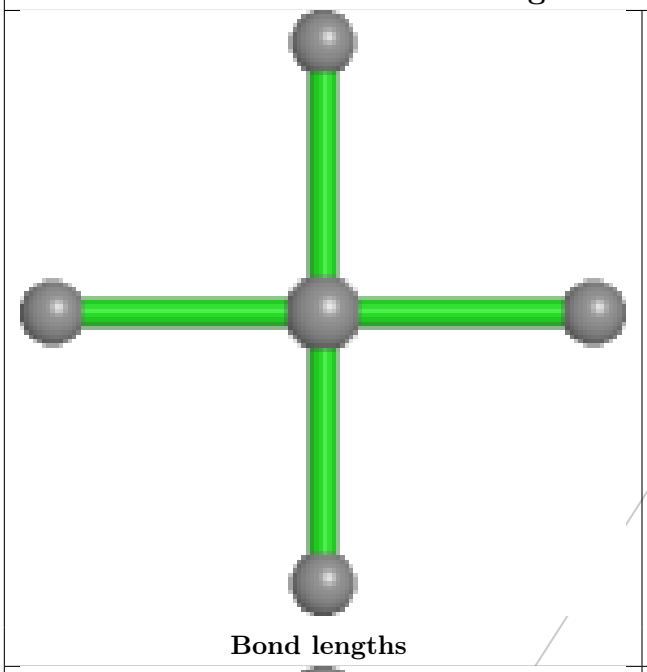
For Manuscript Review

Ligand PEG F 601

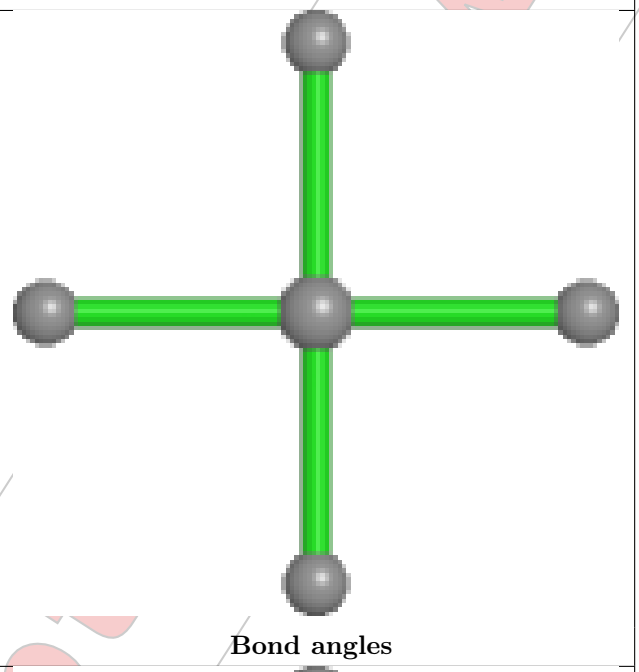


For Manuscript Review

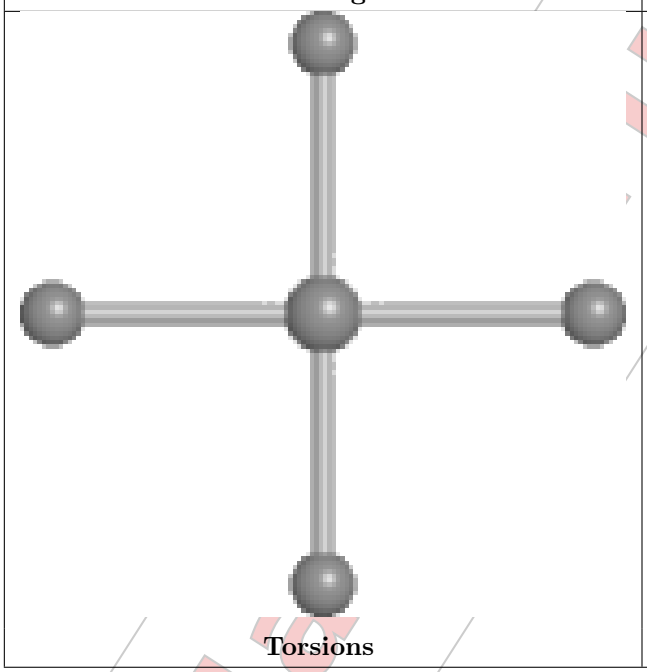
Ligand PO4 B 703



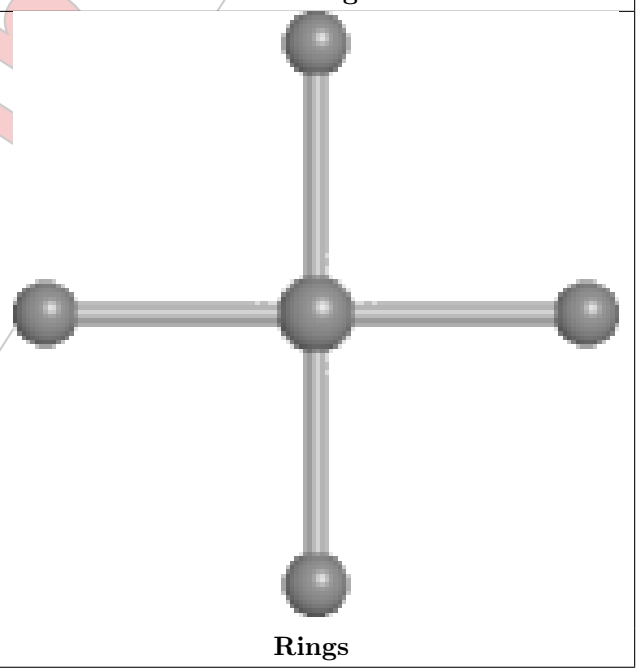
Bond lengths



Bond angles



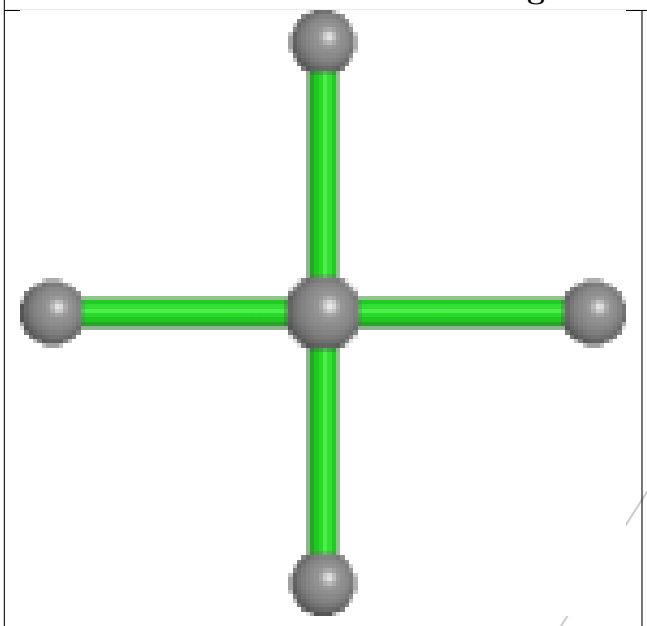
Torsions



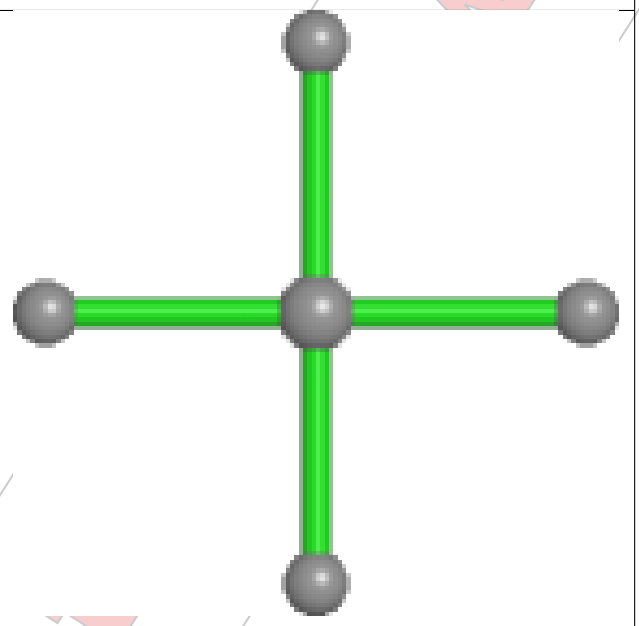
Rings

For Manuscript Review

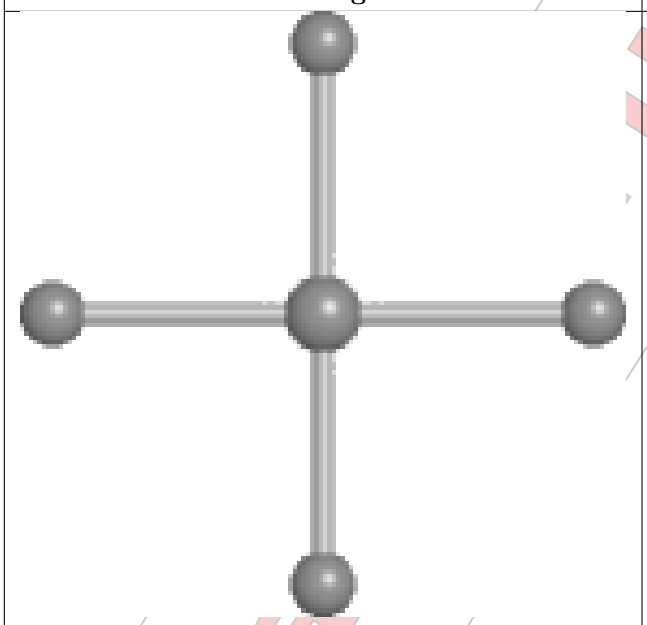
Ligand PO4 A 601



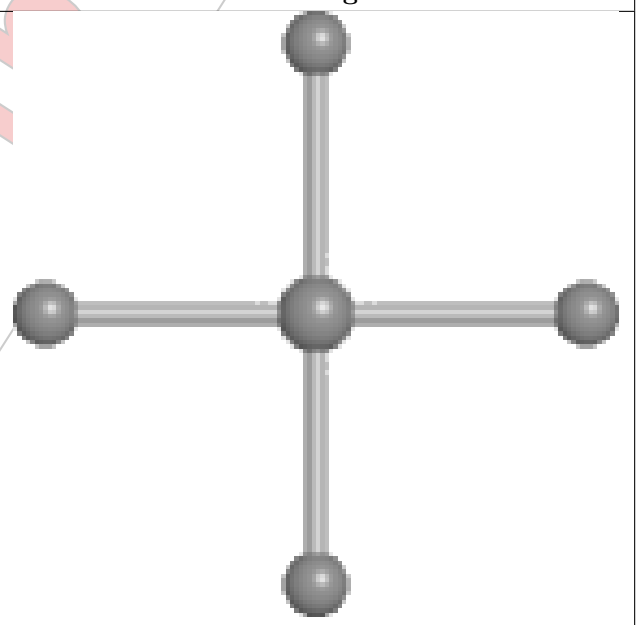
Bond lengths



Bond angles



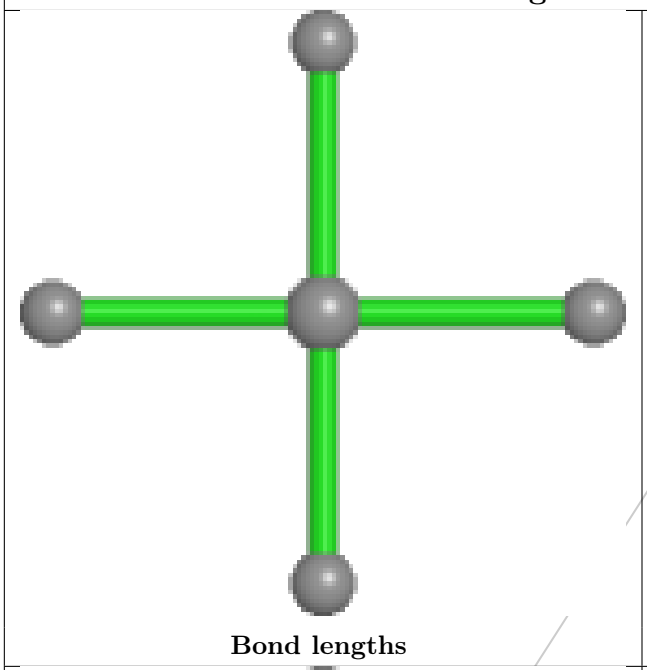
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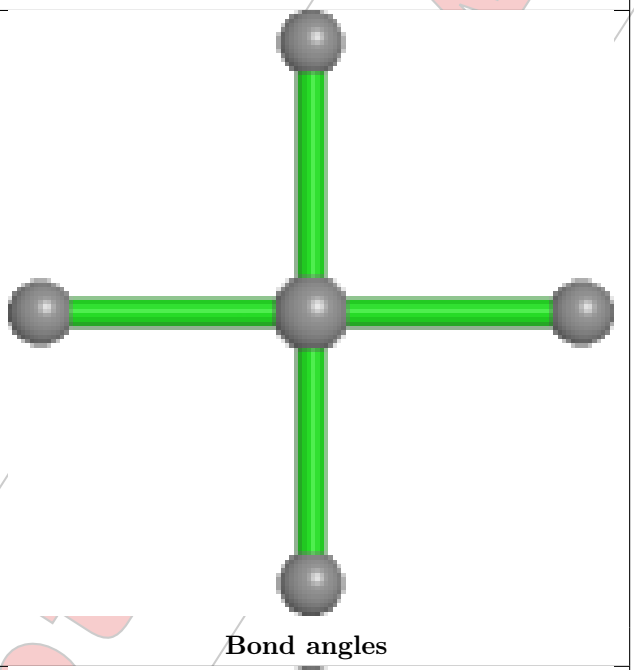
Rings

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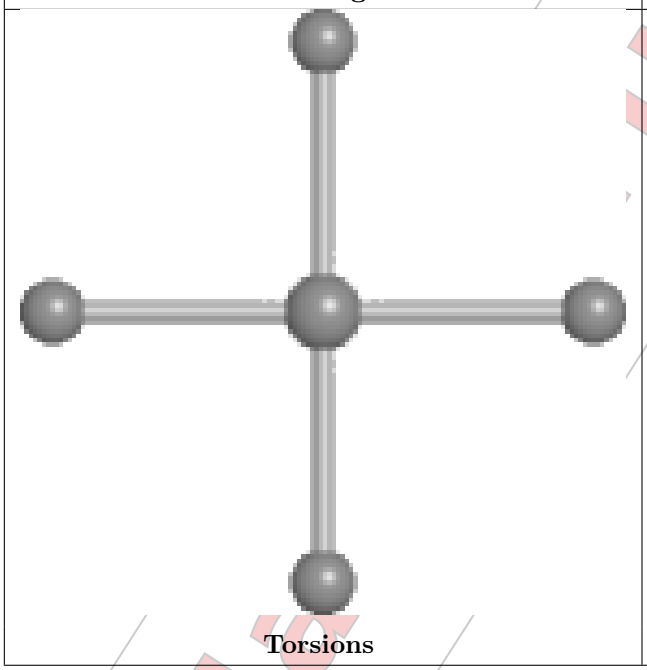
Ligand PO4 E 602



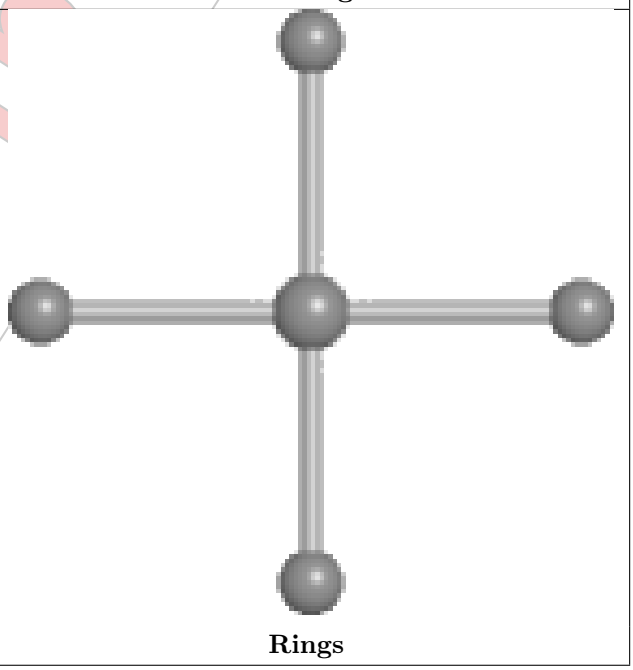
Bond lengths



Bond angles



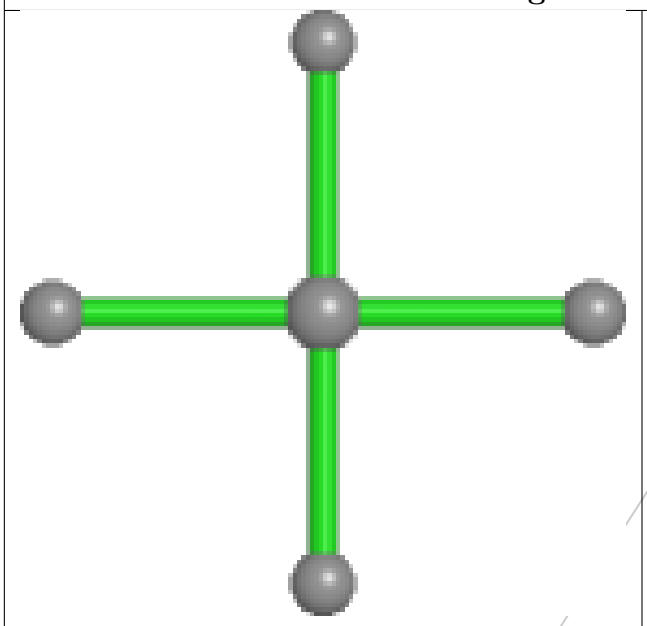
Torsions



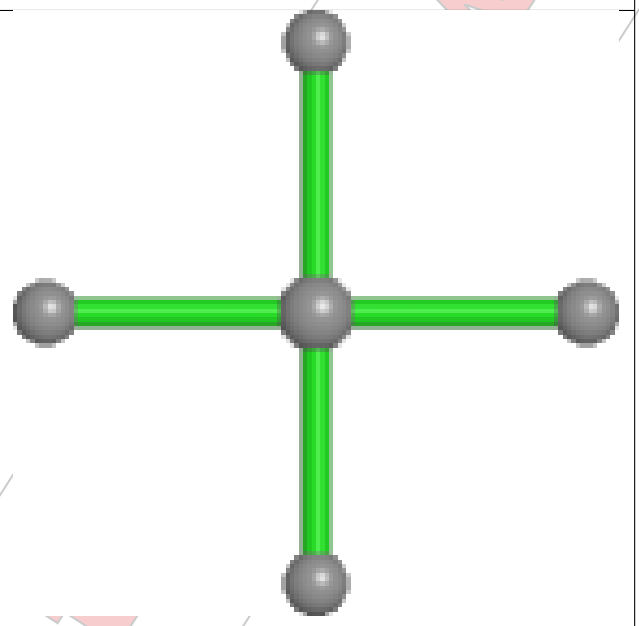
Rings

For Manuscript Review

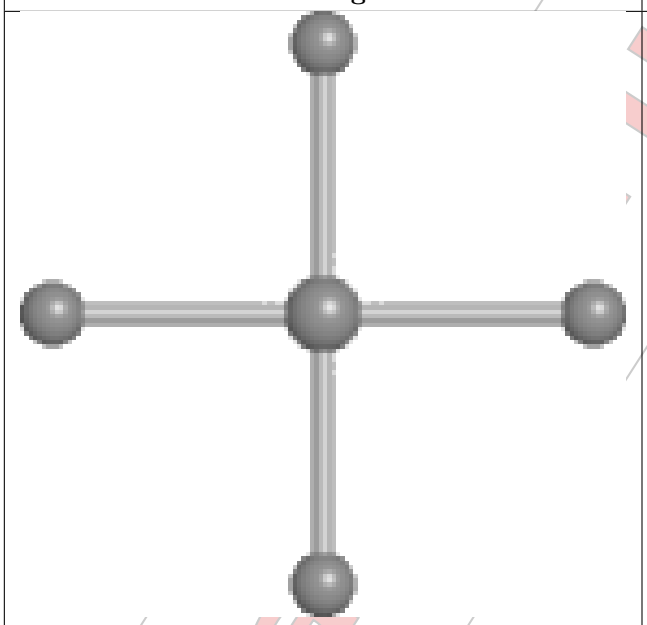
Ligand PO4 B 702



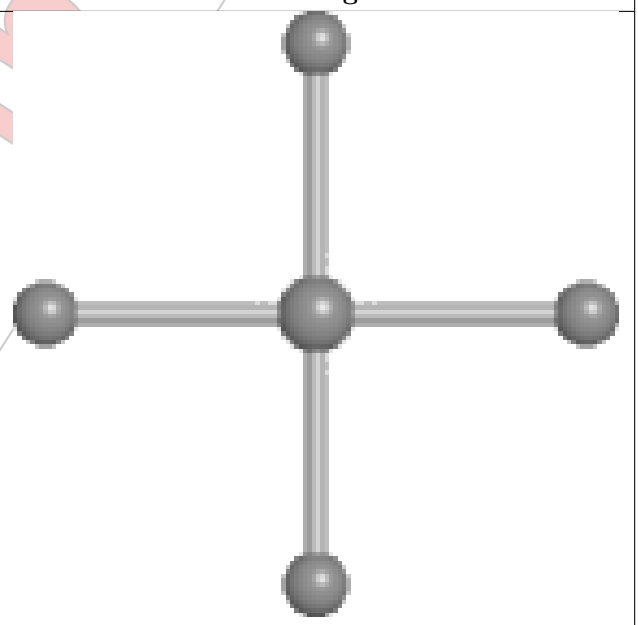
Bond lengths



Bond angles



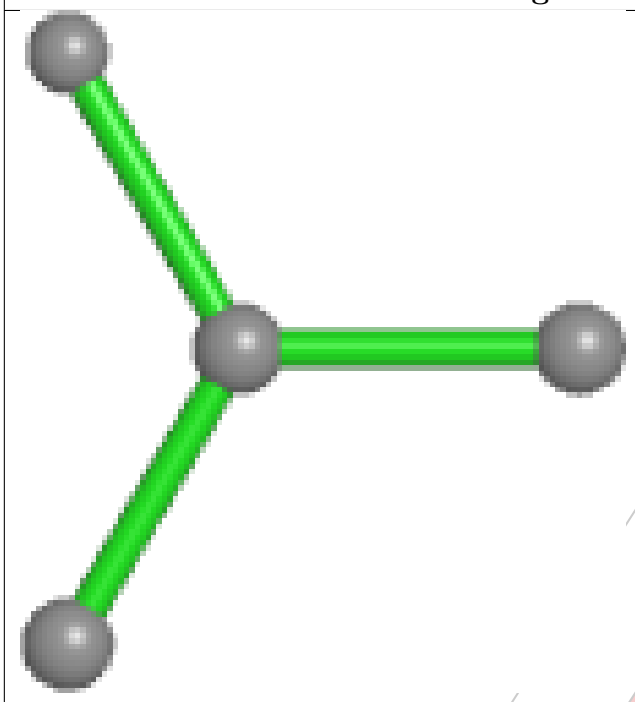
Torsions



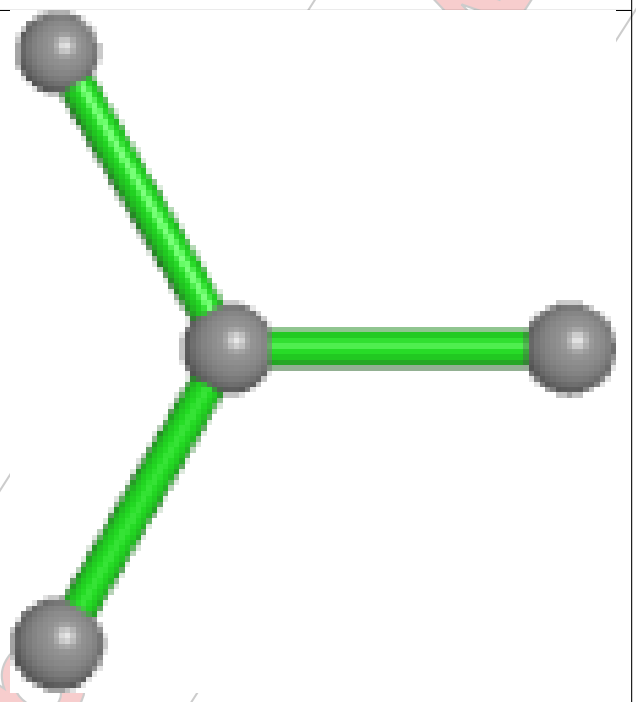
Rings

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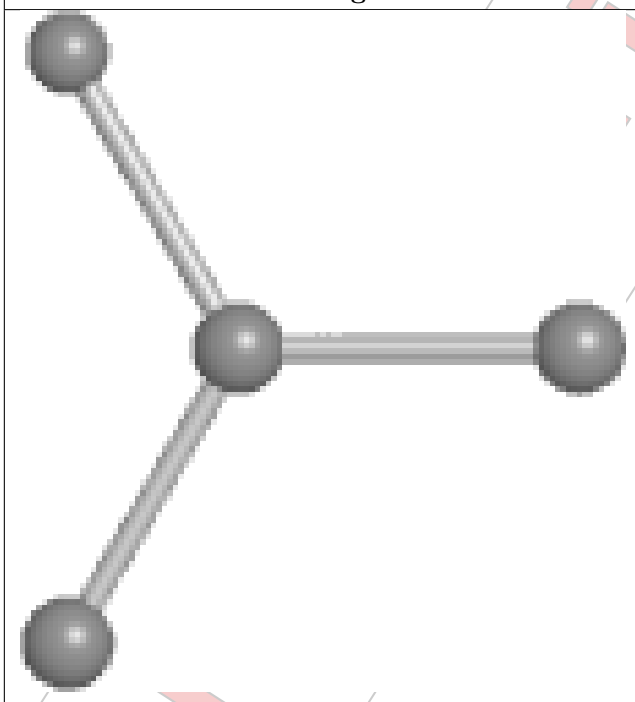
Ligand IPA C 601



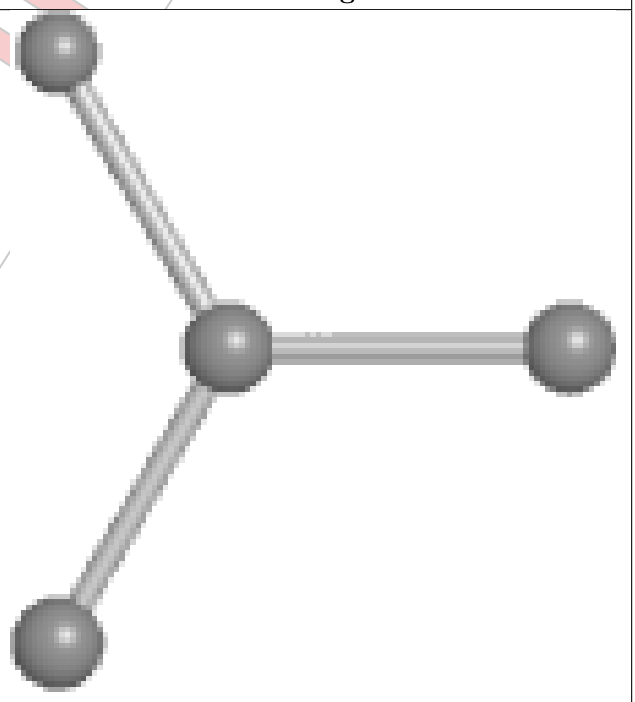
Bond lengths



Bond angles



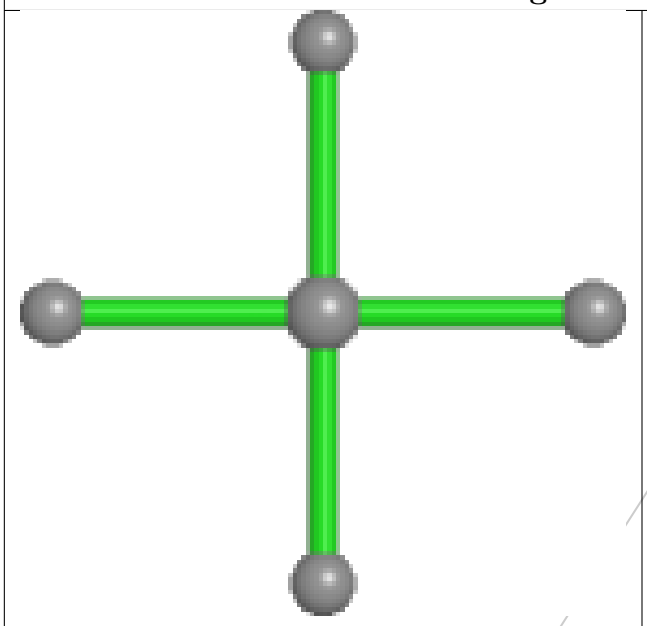
Torsions



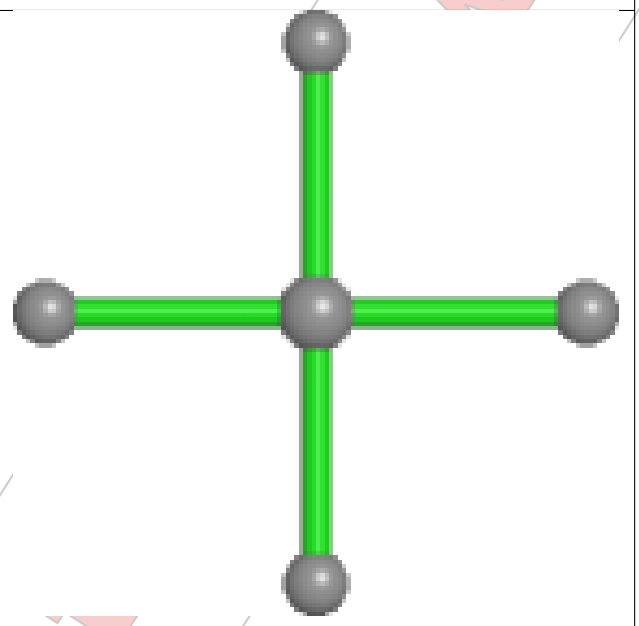
Rings

For Review Only

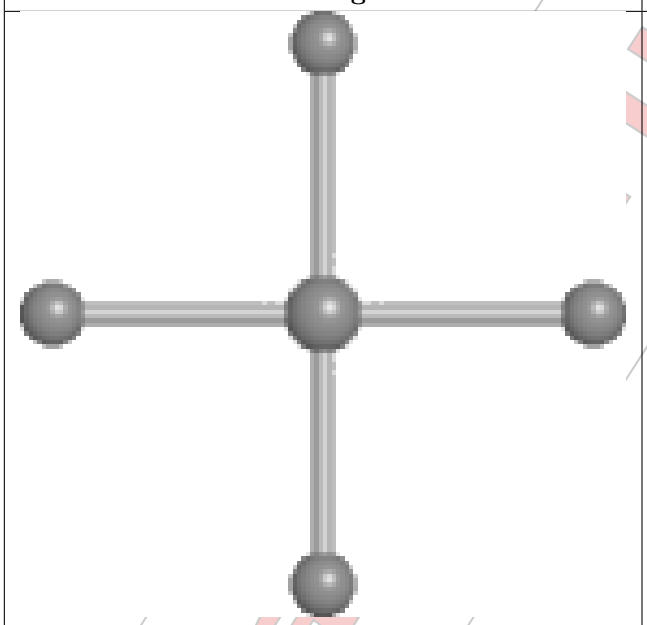
Ligand PO4 F 602



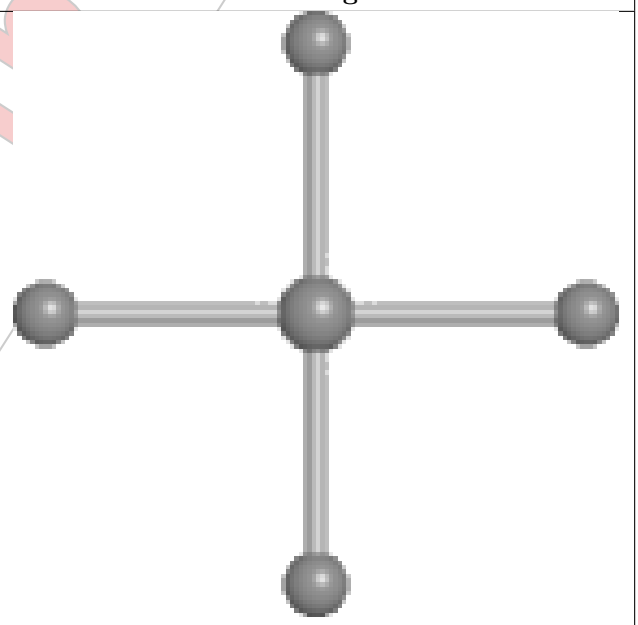
Bond lengths



Bond angles



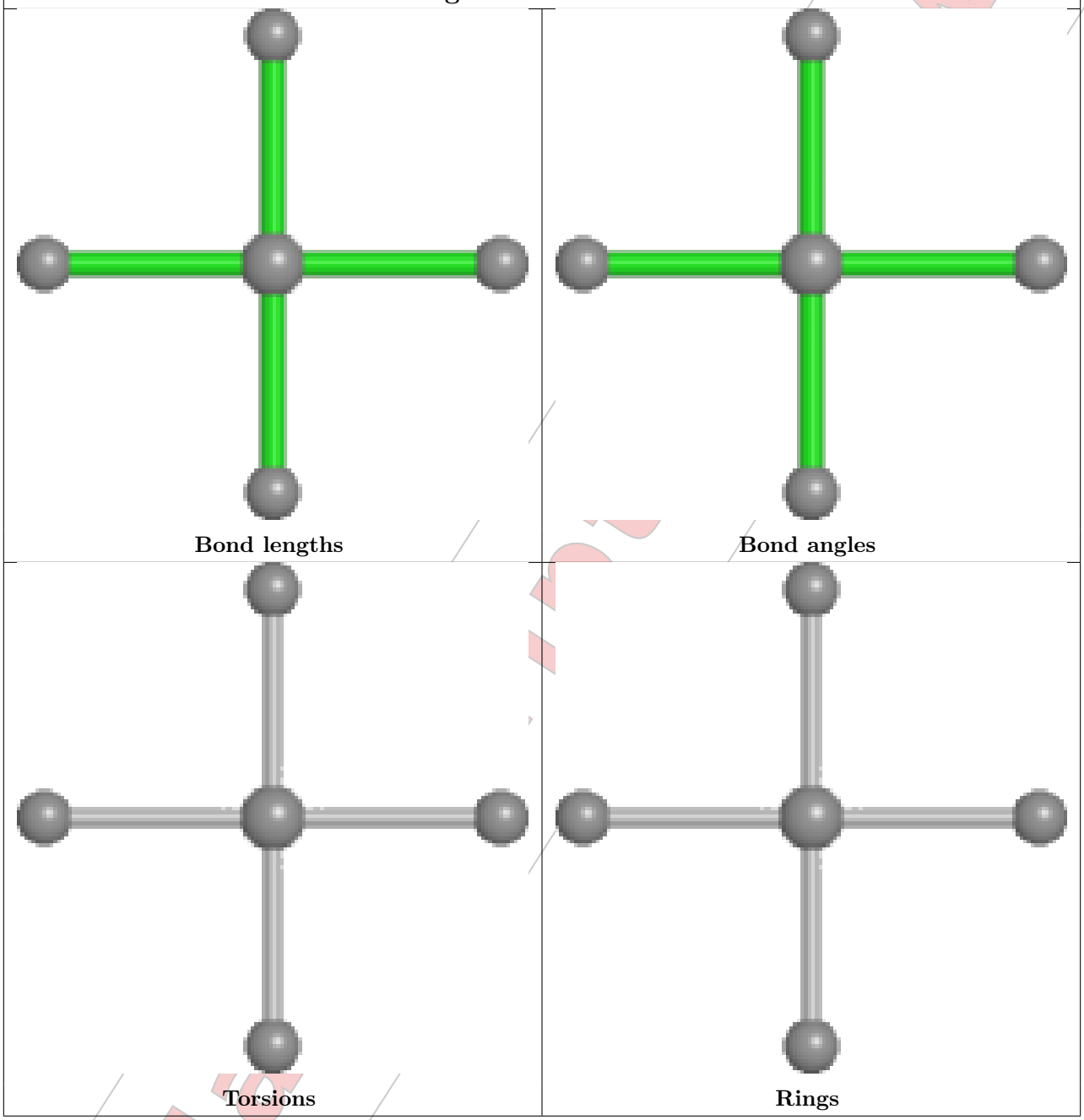
Torsions



Rings

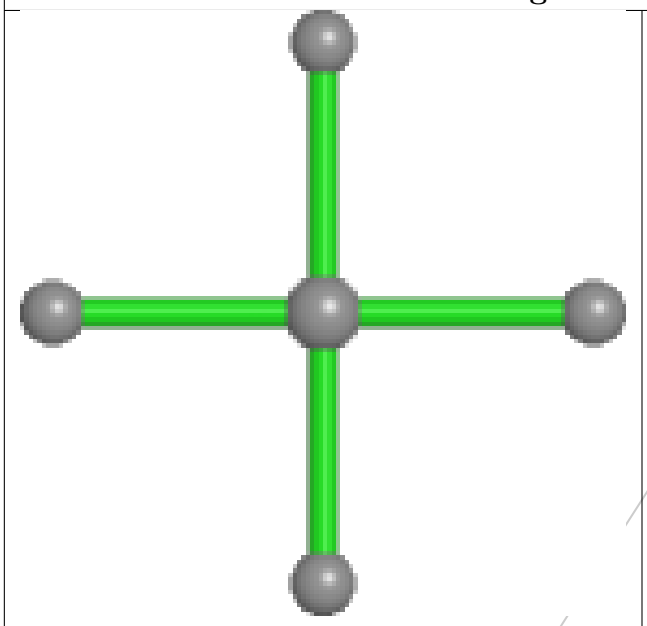
For Manuscript Review

Ligand PO4 D 601

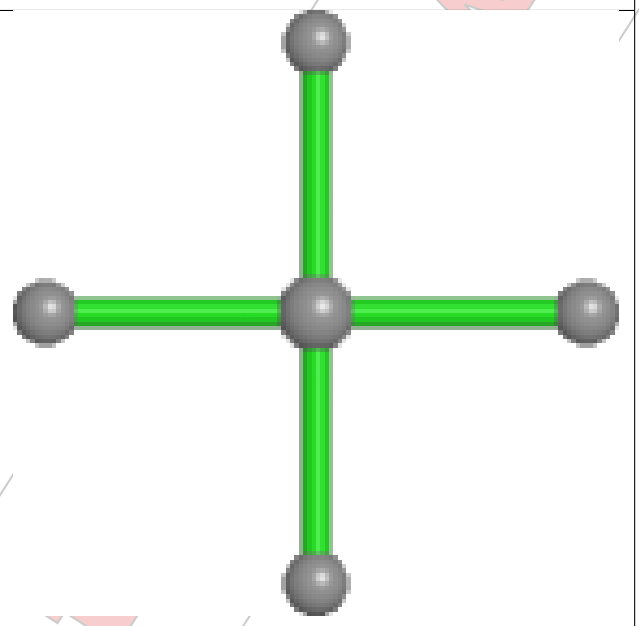


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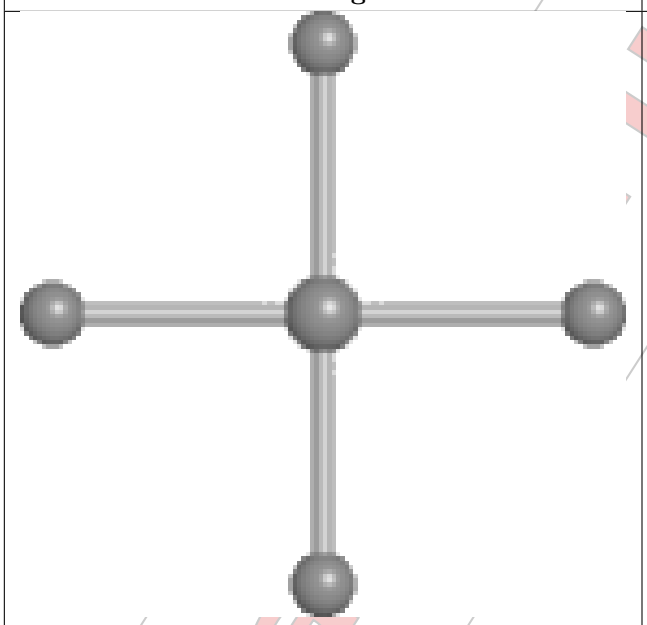
Ligand PO4 F 603



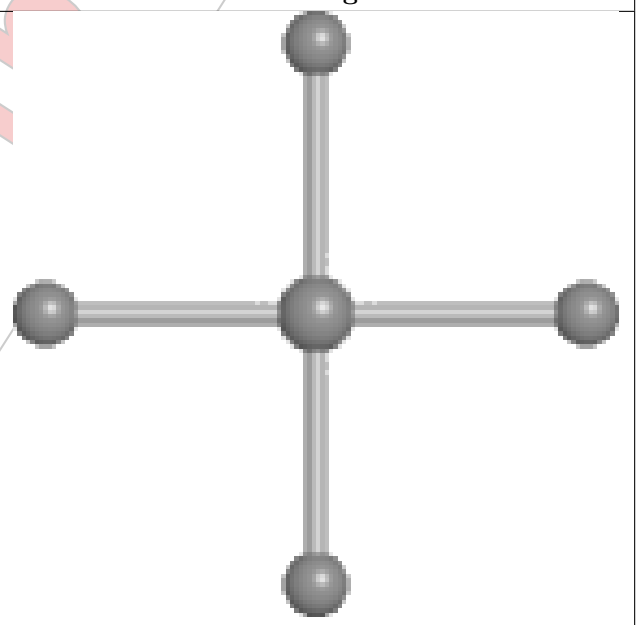
Bond lengths



Bond angles



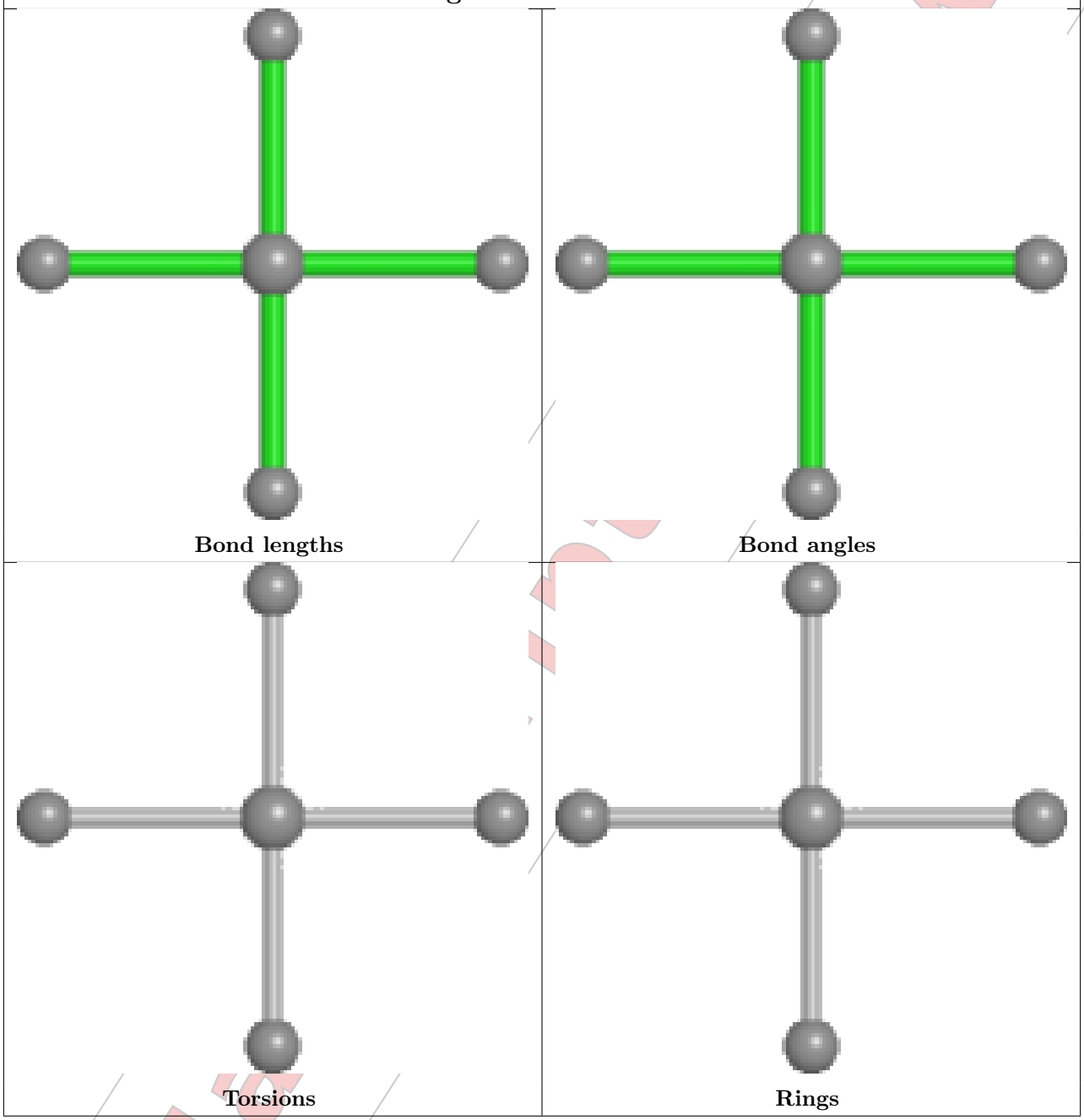
Torsions



Rings

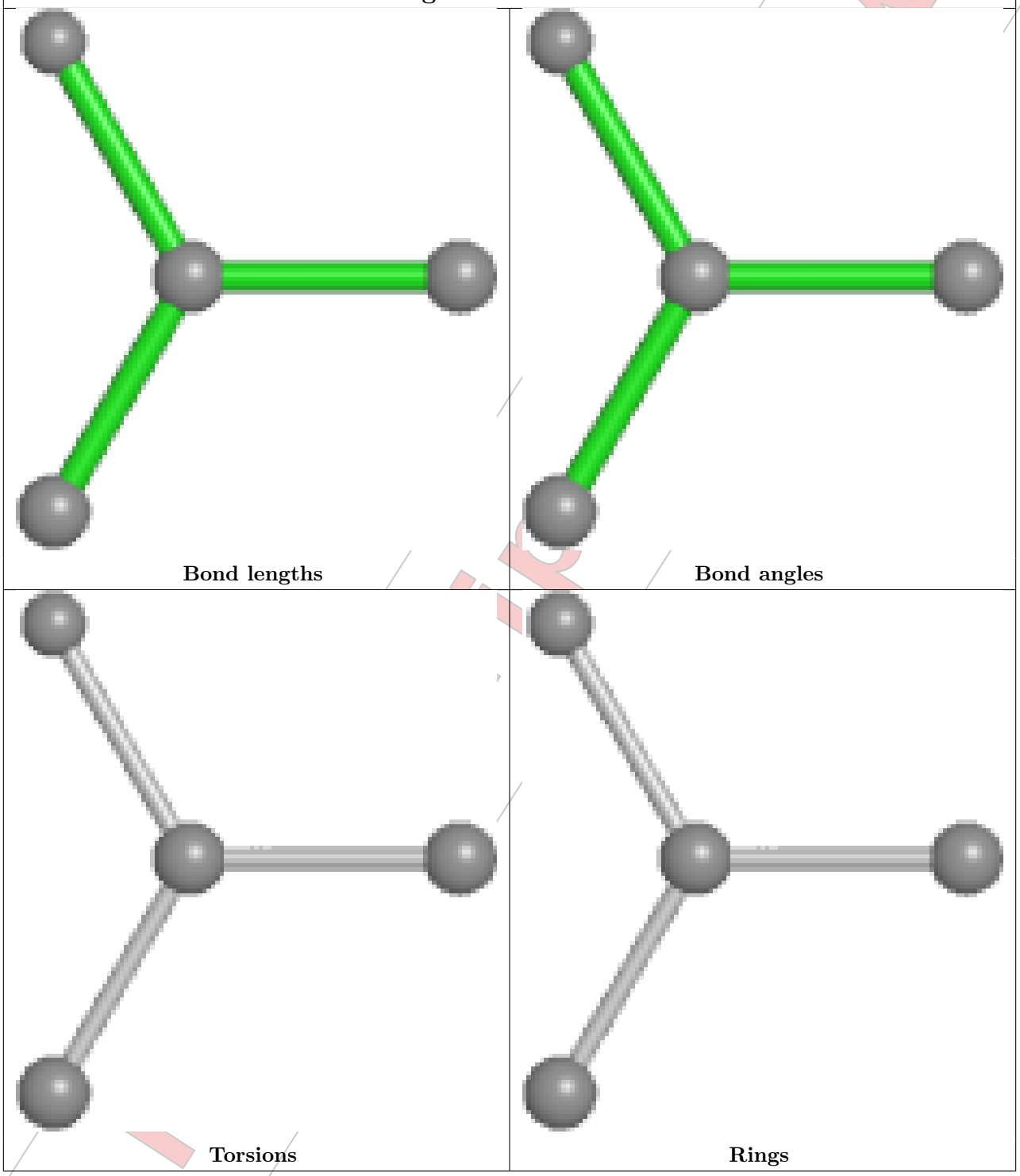
For Manuscript Review

Ligand PO4 E 601



For Manuscript Review

Ligand IPA B 701



5.7 Other polymers (i)

There are no such residues in this entry.

For Review Only

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

For Manuscript Review

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^{th} 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	82/97 (84%)	0.39	4 (4%) 29 33	9, 17, 52, 55	0
1	B	85/97 (87%)	0.96	16 (18%) 1 1	9, 27, 61, 70	0
1	C	87/97 (89%)	0.55	6 (6%) 16 19	11, 20, 51, 70	0
1	D	93/97 (95%)	0.70	12 (12%) 3 3	12, 23, 63, 68	0
1	E	97/97 (100%)	0.54	9 (9%) 8 10	7, 20, 60, 64	0
1	F	87/97 (89%)	0.26	3 (3%) 45 48	7, 17, 51, 62	0
All	All	531/582 (91%)	0.57	50 (9%) 8 9	7, 21, 59, 70	0

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	569	TYR	8.7
1	B	572	TYR	7.5
1	D	520	VAL	7.5
1	C	546	ASN	5.5
1	B	516	ASN	5.2
1	C	524	THR	5.2
1	C	525	ILE	4.8
1	E	523	SER	4.7
1	A	495	PRO	4.6
1	D	523	SER	4.6
1	D	588	GLU	4.4
1	F	517	SER	4.2
1	B	570	ASN	4.2
1	D	547	ASN	4.2
1	D	546	ASN	4.1
1	A	569	TYR	4.1
1	B	495	PRO	4.0
1	D	522	ALA	4.0
1	B	494	SER	3.9

Continued on next page...

For Manuscript Review

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	E	520	VAL	3.9
1	E	522	ALA	3.9
1	E	521	PRO	3.7
1	E	524	THR	3.6
1	B	573	GLY	3.6
1	F	516	ASN	3.4
1	B	571	ARG	3.3
1	C	526	LYS	3.3
1	B	529	LEU	3.3
1	C	547	ASN	3.2
1	E	588	GLU	3.2
1	D	569	TYR	3.1
1	B	496	PRO	3.0
1	D	516	ASN	2.9
1	E	546	ASN	2.6
1	B	575	GLY	2.5
1	A	572	TYR	2.5
1	F	549	LYS	2.4
1	E	525	ILE	2.3
1	C	569	TYR	2.3
1	B	546	ASN	2.2
1	B	547	ASN	2.2
1	D	494	SER	2.2
1	E	547	ASN	2.1
1	D	545	PRO	2.1
1	B	527	GLN	2.1
1	D	502[A]	ILE	2.1
1	A	544	ILE	2.0
1	B	588	GLU	2.0
1	D	524	THR	2.0
1	B	528	TYR	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.

For Manuscript Review

6.4 Ligands

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

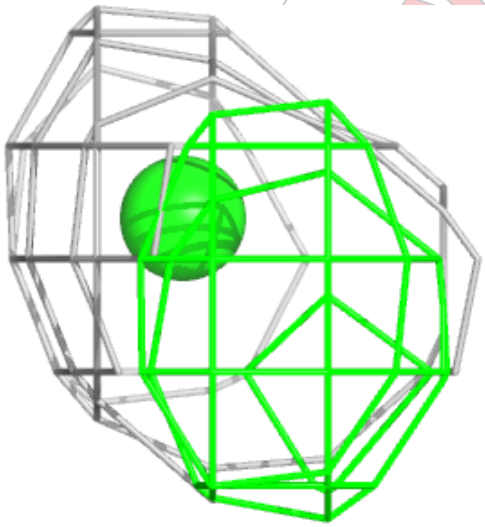
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
4	CL	E	603	1/1	0.74	0.11	48,48,48,48	0
5	PEG	F	601	7/7	0.77	0.25	25,52,61,67	0
2	PO4	B	703	5/5	0.89	0.14	43,46,52,64	0
3	IPA	B	701	4/4	0.89	0.17	18,25,37,42	0
2	PO4	E	602	5/5	0.92	0.18	27,33,50,53	0
3	IPA	C	601	4/4	0.93	0.15	19,29,48,48	0
2	PO4	D	601	5/5	0.97	0.15	31,35,38,41	0
2	PO4	A	601	5/5	0.97	0.07	28,29,39,44	0
2	PO4	E	601	5/5	0.98	0.12	24,30,36,38	0
2	PO4	F	602	5/5	0.98	0.07	16,17,22,25	0
2	PO4	B	702	5/5	0.99	0.08	18,22,23,24	0
2	PO4	F	603	5/5	0.99	0.10	22,23,26,28	0

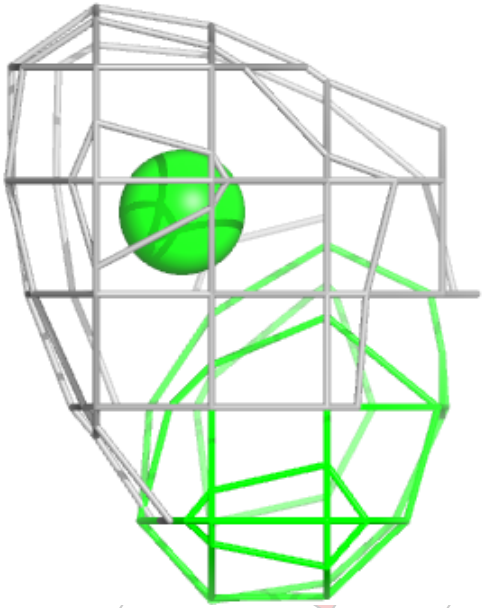
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

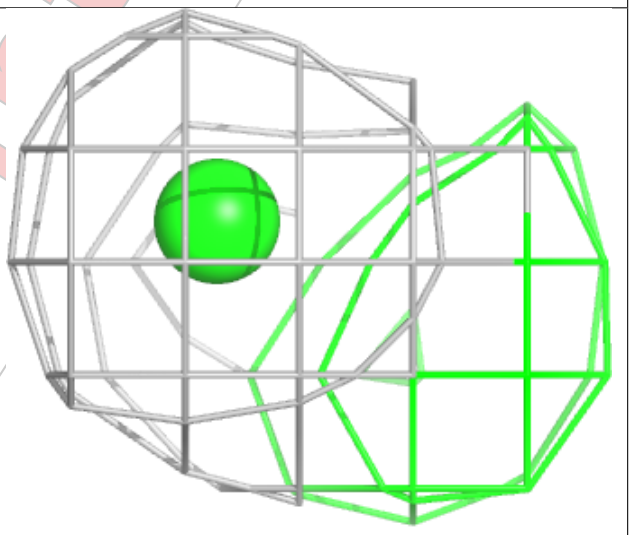
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Electron density around CL E 603:

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



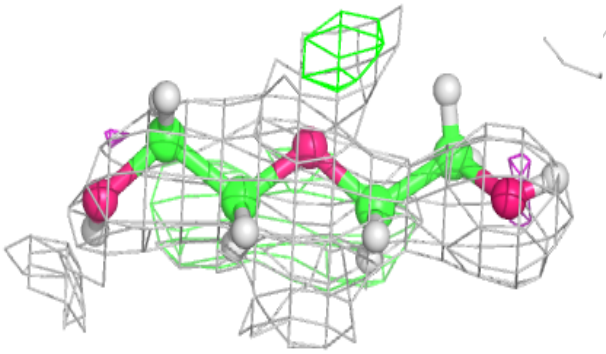


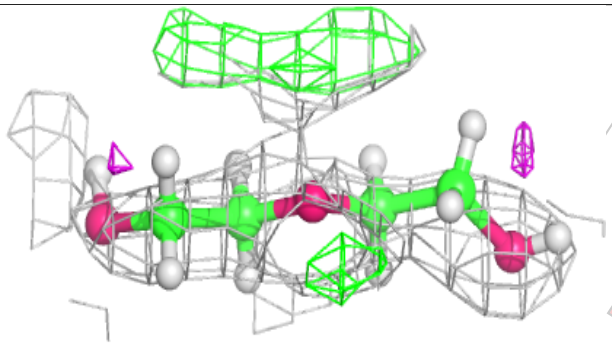


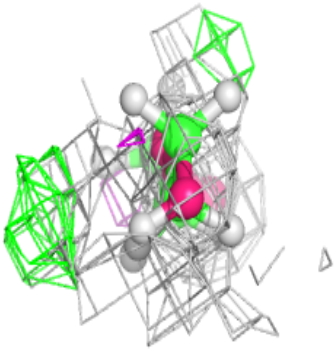
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Electron density around PEG F 601:

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



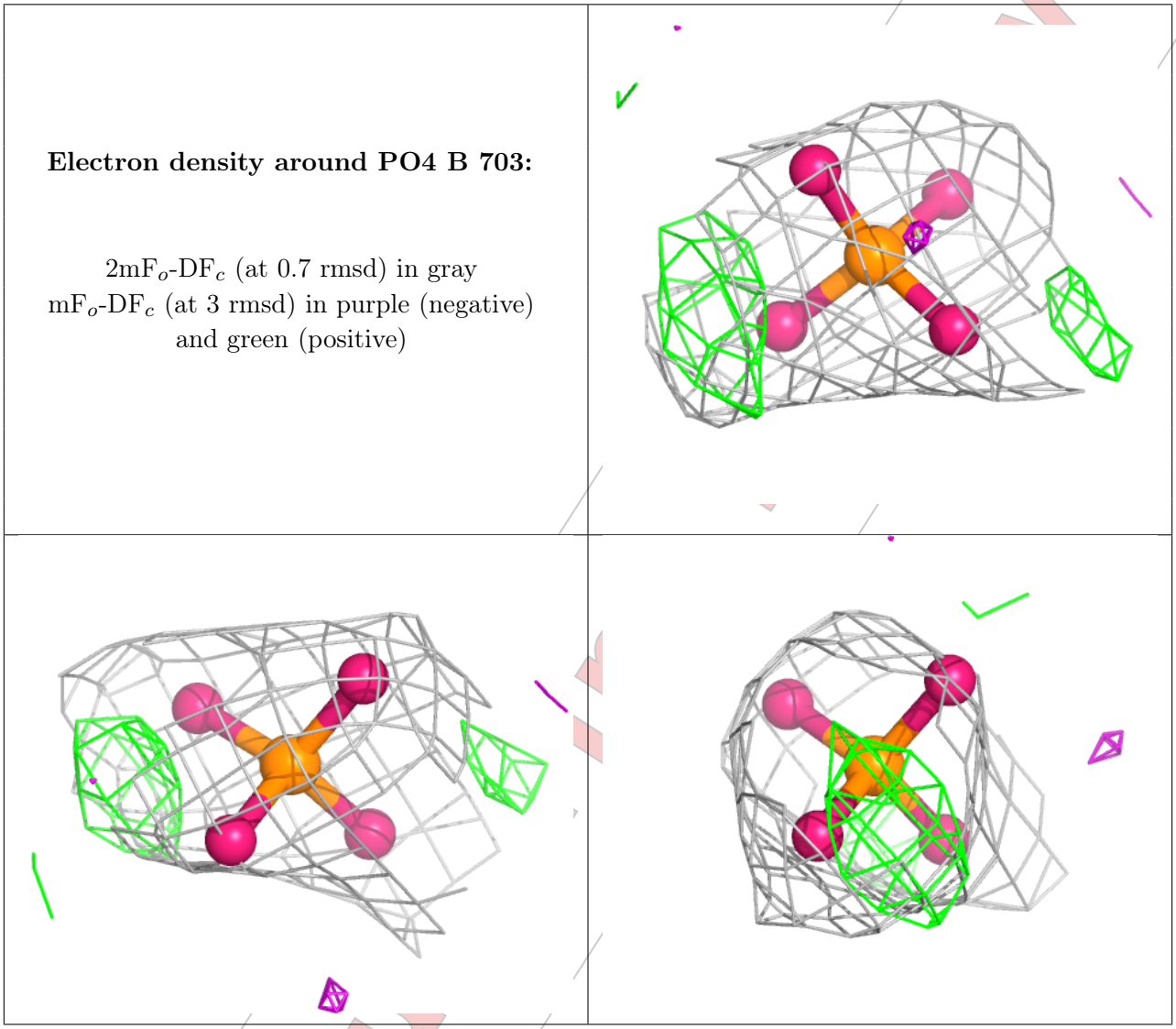




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Electron density around PO4 B 703:

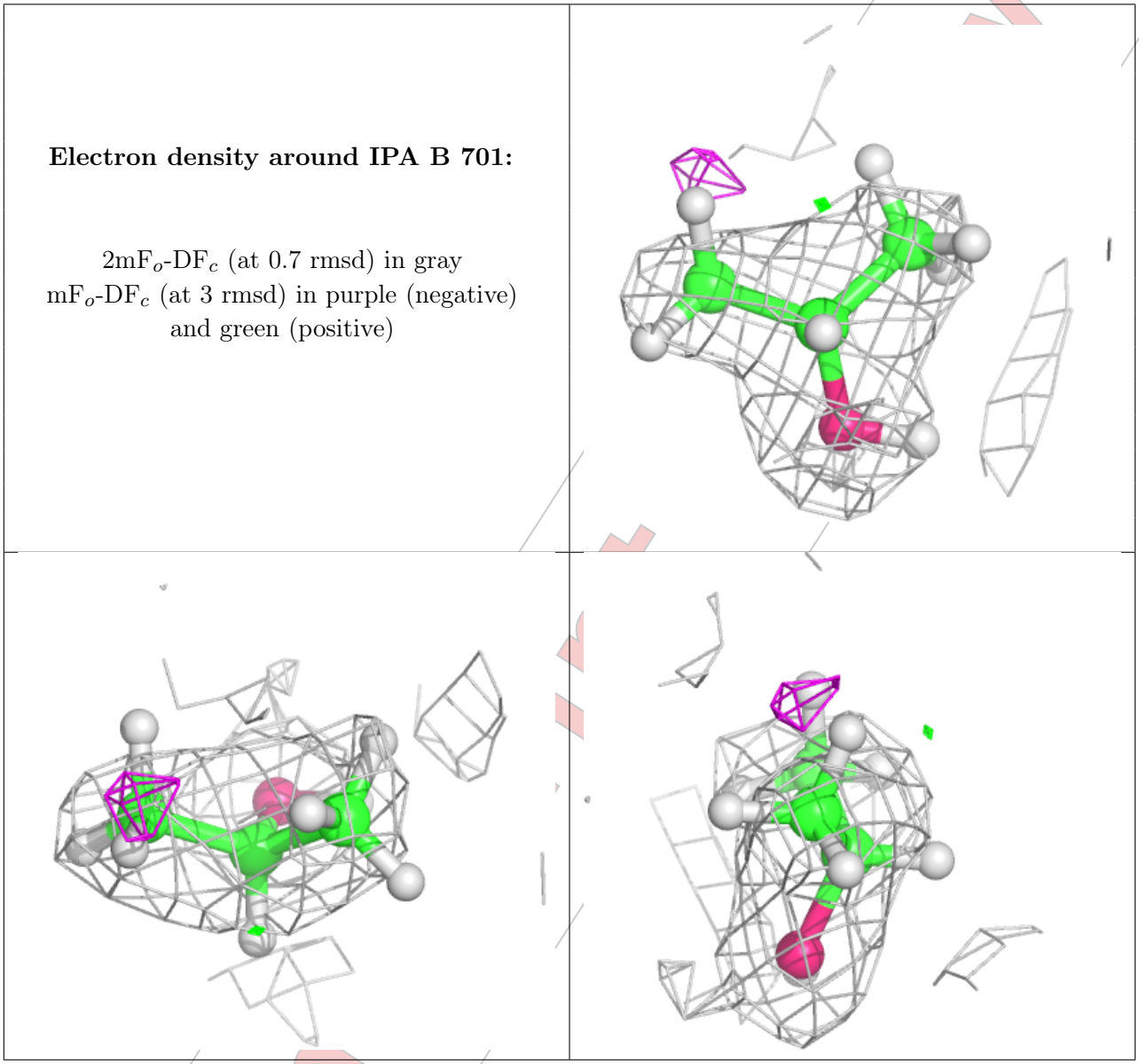
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



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Electron density around IPA B 701:

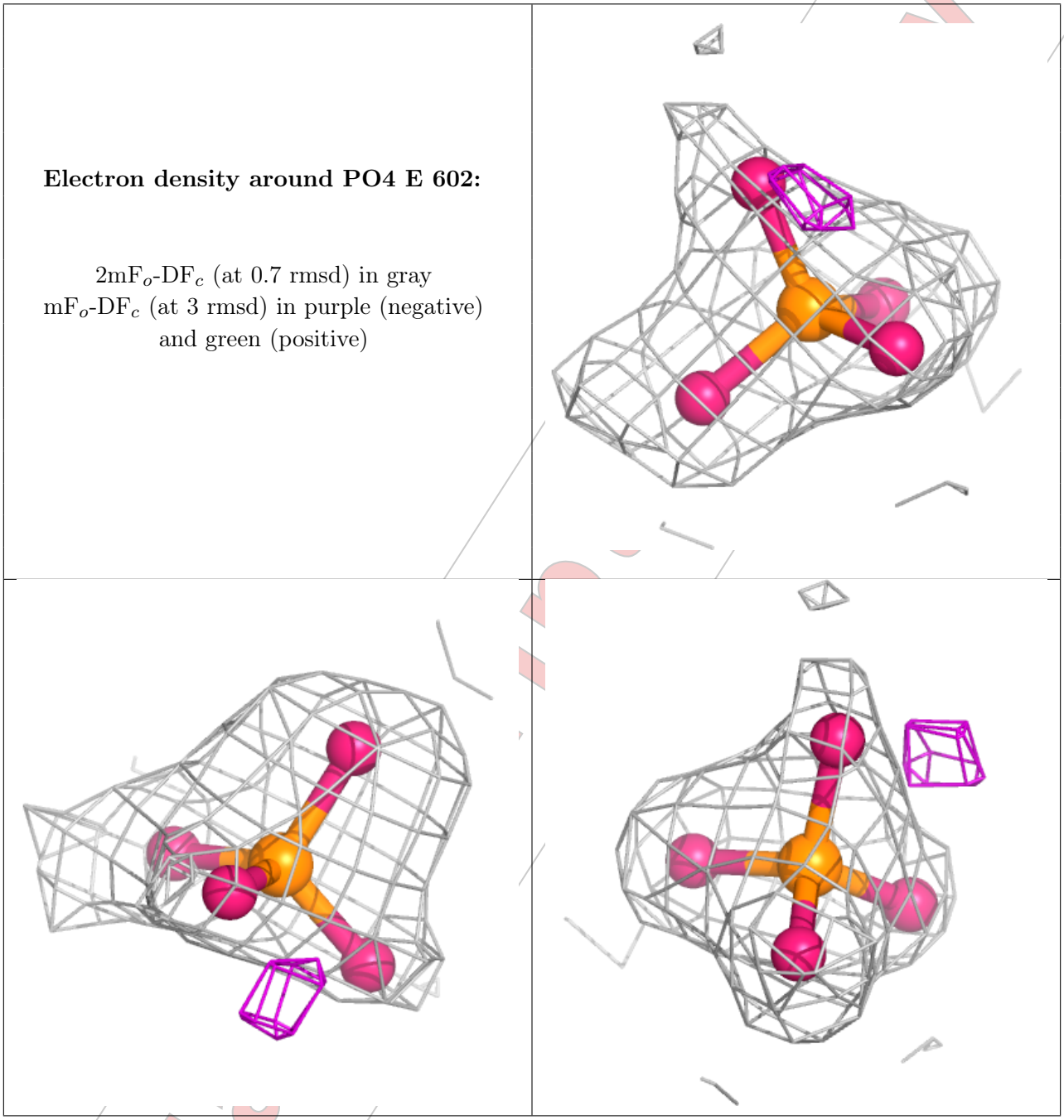
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



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Electron density around PO4 E 602:

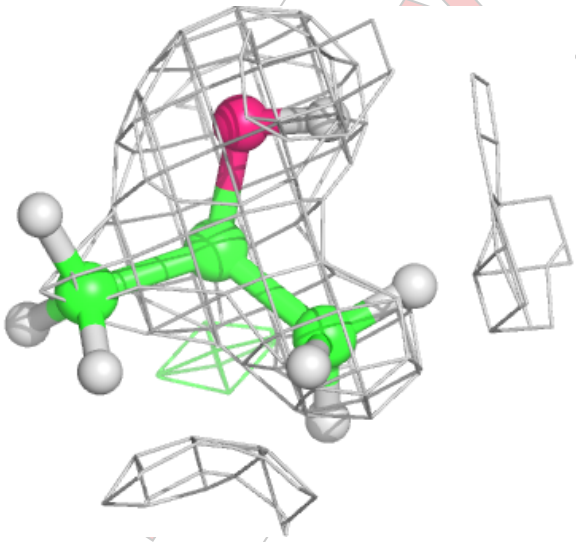
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

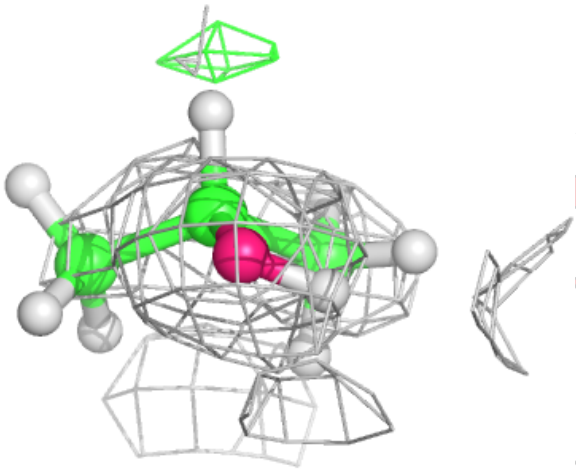


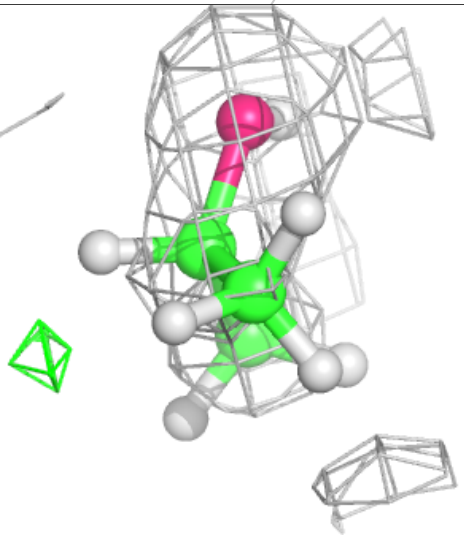
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Electron density around IPA C 601:

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



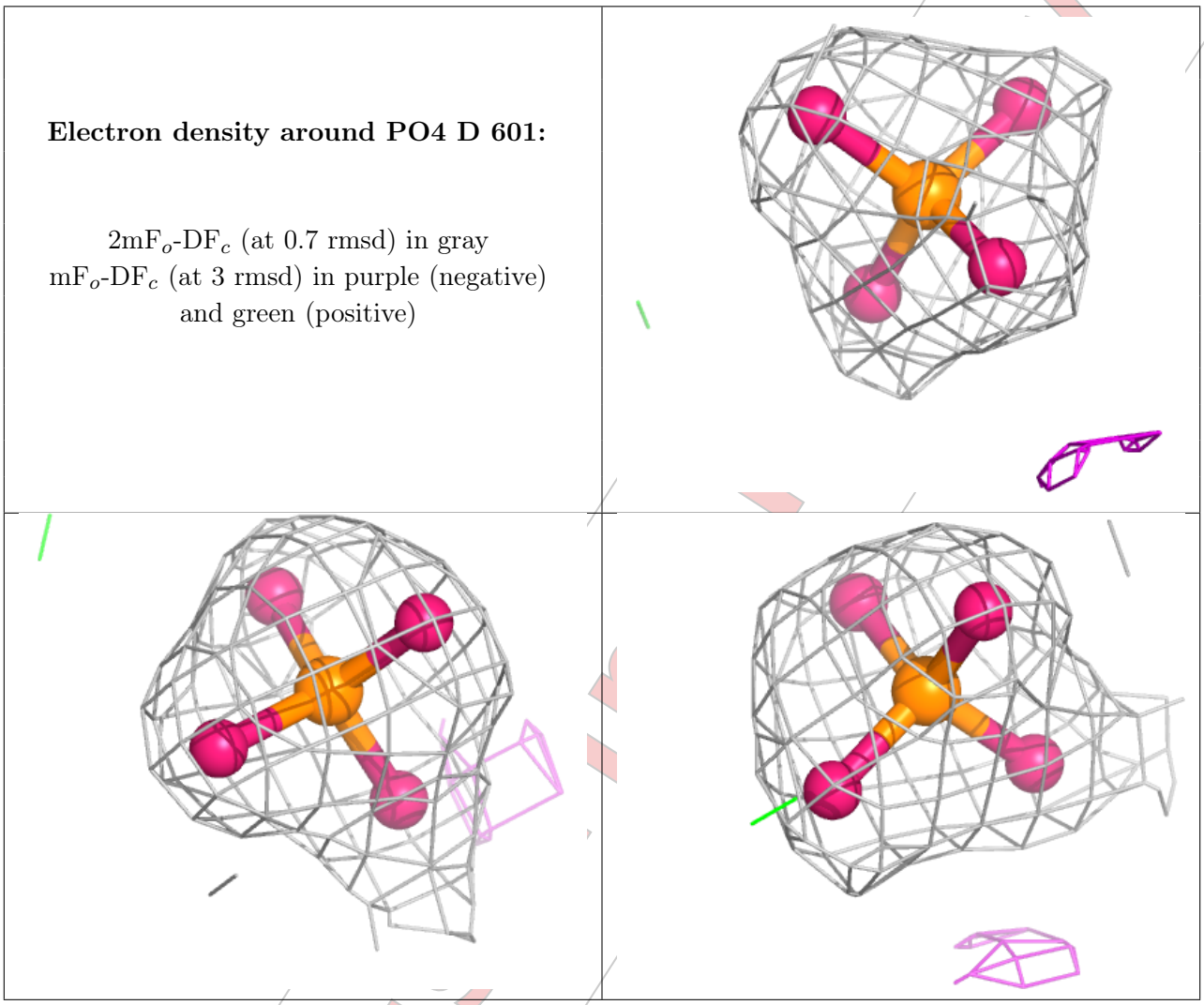




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Electron density around PO4 D 601:

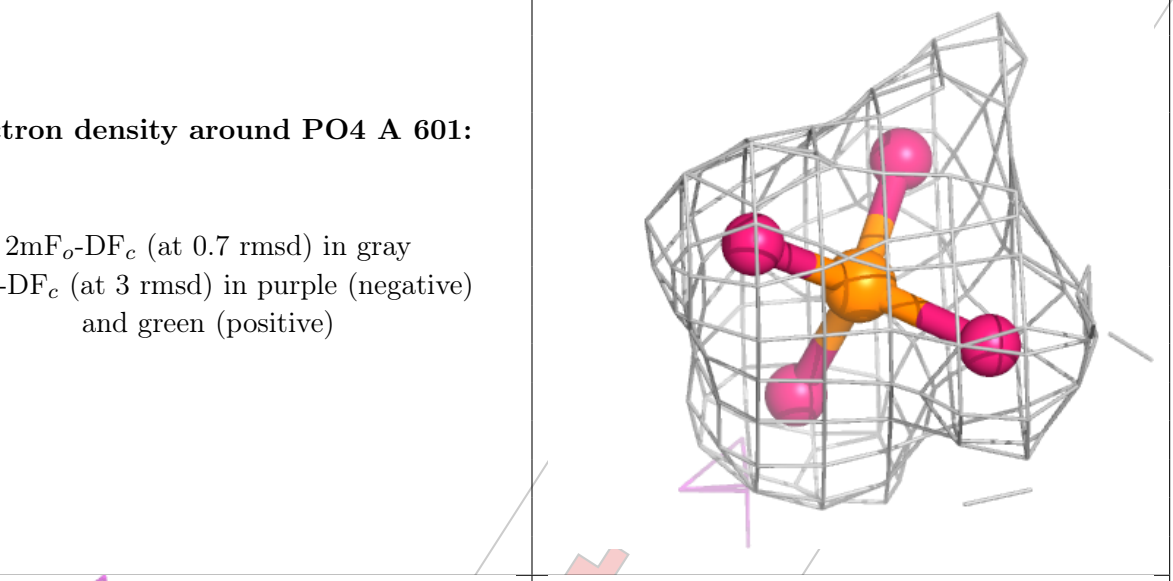
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

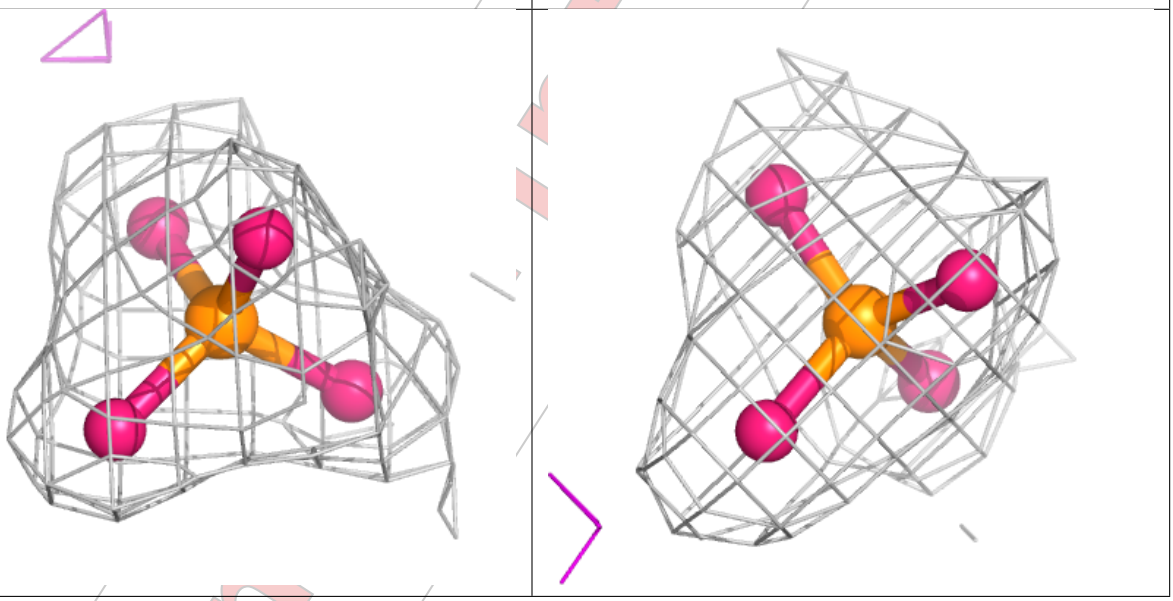


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Electron density around PO4 A 601:

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

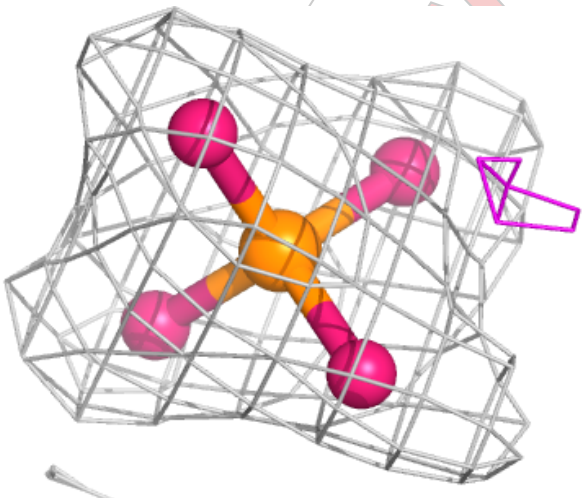


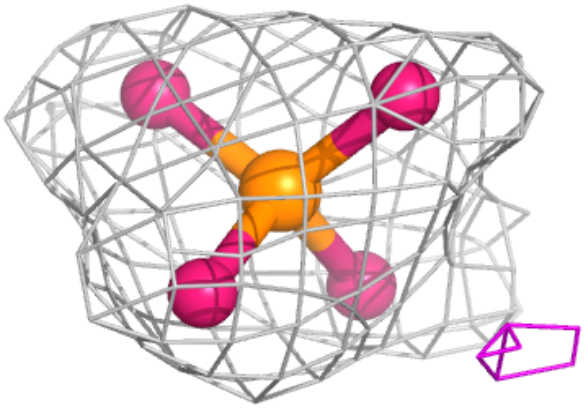


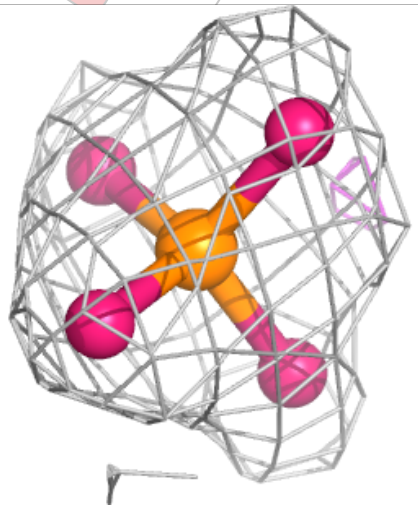
For Manuscript Review

Electron density around PO4 E 601:

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



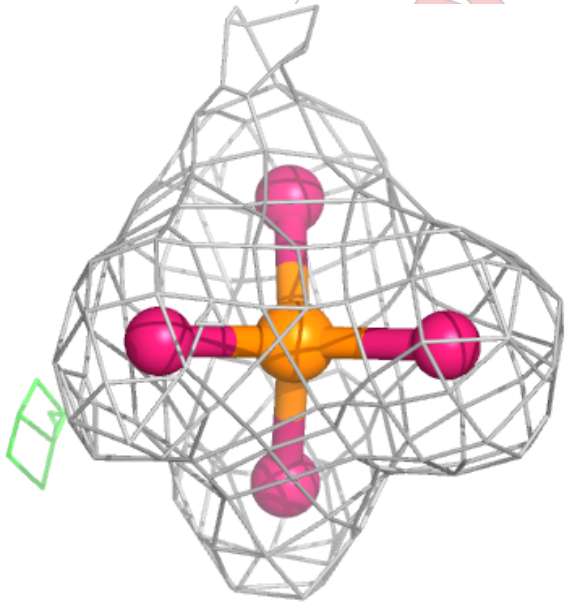


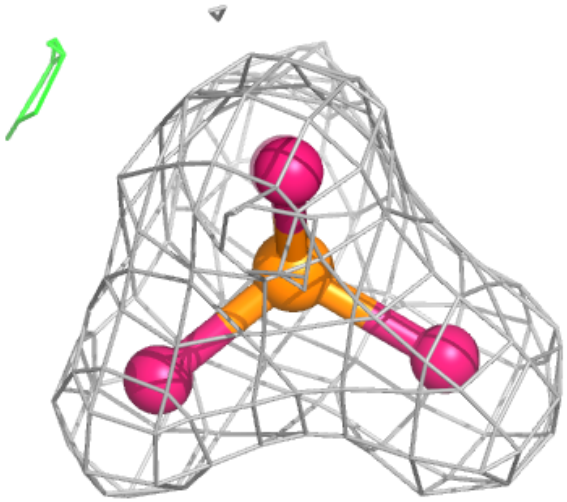


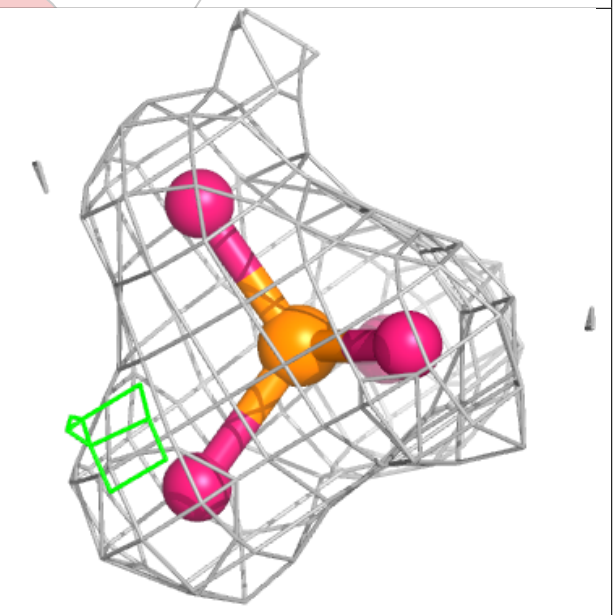
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Electron density around PO4 F 602:

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



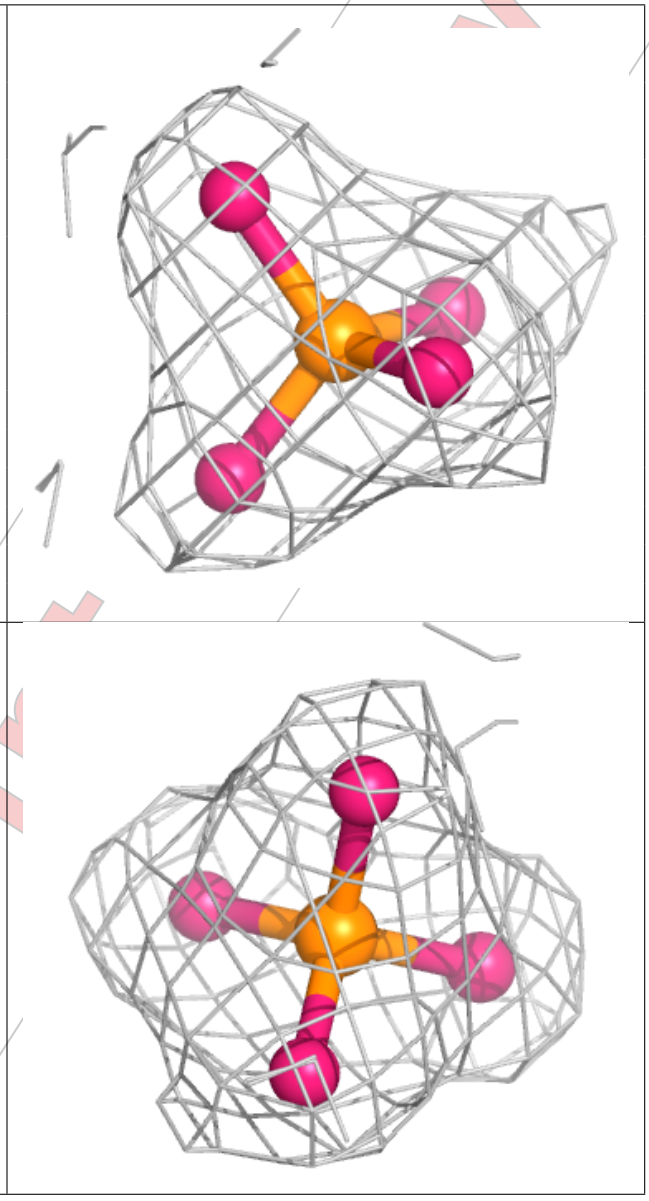




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Electron density around PO4 B 702:

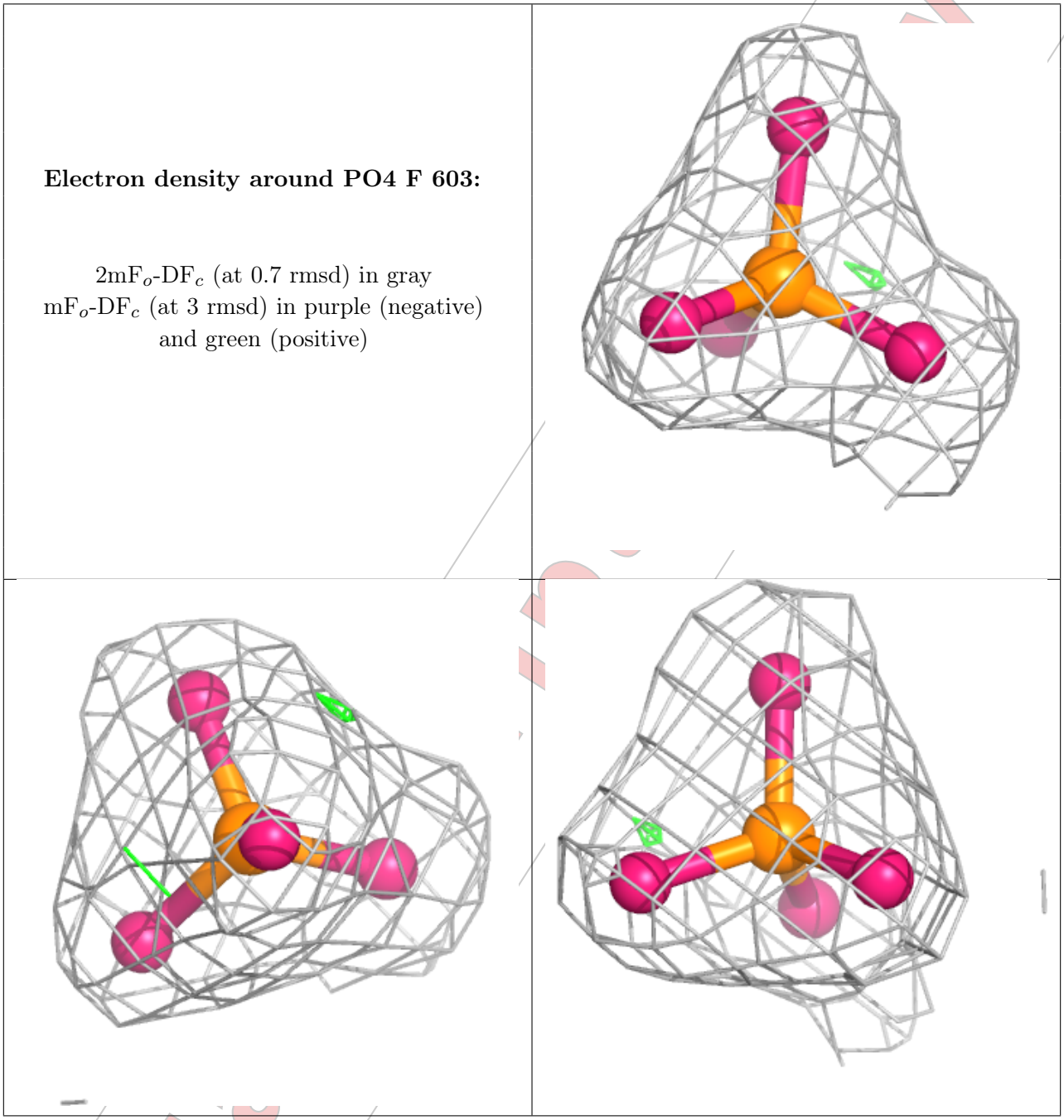
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



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Electron density around PO4 F 603:

$2mF_o$-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.

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