wwPDB X-ray Structure Validation Summary Report ⓘ

May 22, 2020 – 12:34 pm BST

PDB ID	:	6BB4
Title	:	Fab/epitope complex of mouse monoclonal antibody C5.2 targeting a phospho-tau epitope.
Authors	:	Chukwu, J.E.; Kong, X.-P.
Deposited on	:	2017-10-16
Resolution	:	2.10 Å(reported)

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

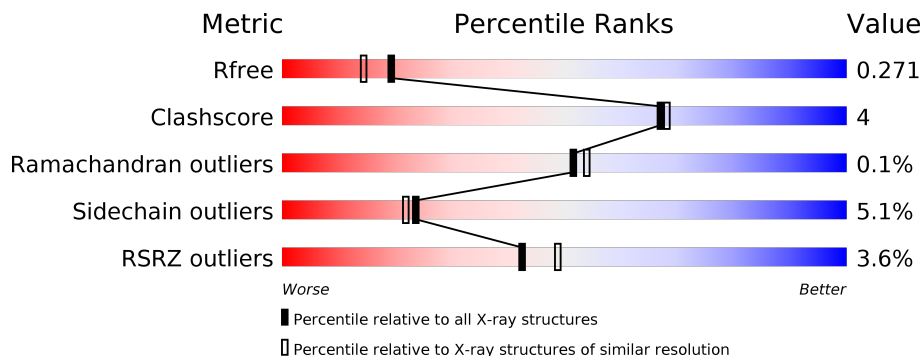|		
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

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 2.10 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	Percentile Ranks	Value
Rfree		0.271
Clashscore		4
Ramachandran outliers		0.1%
Sidechain outliers		5.1%
RSRZ outliers		3.6%

Worse ◄ ► *Better*

■ Percentile relative to all X-ray structures

▯ Percentile relative to X-ray structures of similar resolution

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R$_{free}$	130704	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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 on 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	L	213	3% / 89% / 9% •
1	M	213	5% / 83% / 15% ••
1	N	213	3% / 88% / 10% ••
2	H	220	4% / 85% / 10% ••
2	I	220	5% / 85% / 12% ••
2	J	220	2% / 88% / 10% •

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Mol	Chain	Length	Quality of chain
3	P	23	17% 9% • 70%
3	Q	23	26% • 70%
3	R	23	4% 17% 13% 70%

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 10894 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 Mouse monoclonal antibody C5.2 Fab light chain.

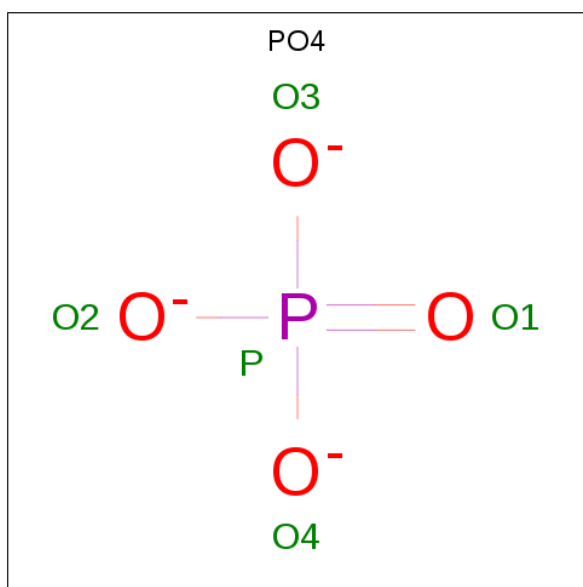| Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	211	Total	C	N	O	S	0	0	0
			1637	1023	270	336	8			
1	M	211	Total	C	N	O	S	0	0	0
			1637	1023	270	336	8			
1	N	211	Total	C	N	O	S	0	0	0
			1637	1023	270	336	8			

- Molecule 2 is a protein called Mouse monoclonal antibody C5.2 Fab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	216	Total	C	N	O	S	0	0	0
			1638	1030	271	329	8			
2	I	216	Total	C	N	O	S	0	0	0
			1638	1030	271	329	8			
2	J	216	Total	C	N	O	S	0	0	0
			1638	1030	271	329	8			

- Molecule 3 is a protein called Microtubule-associated protein tau.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	P	7	Total	C	N	O	P	0	0	0
			60	39	8	12	1			
3	Q	7	Total	C	N	O	P	0	0	0
			60	39	8	12	1			
3	R	7	Total	C	N	O	P	0	0	0
			60	39	8	12	1			

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

PO4

O3

O^-

O2 O^- —— P ==== O   O1

P

O^-

O4

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

- Molecule 5 is water.

| Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	L	167	Total O 167 167		0	0
5	H	172	Total O 172 172		0	0
5	P	10	Total O 10 10		0	0
5	M	115	Total O 115 115		0	0
5	I	138	Total O 138 138		0	0
5	Q	10	Total O 10 10		0	0
5	N	138	Total O 138 138		0	0
5	J	119	Total O 119 119		0	0
5	R	5	Total O 5 5		0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA 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.
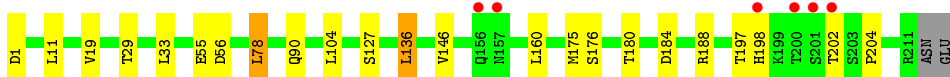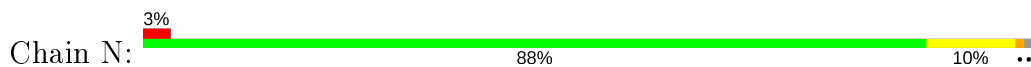
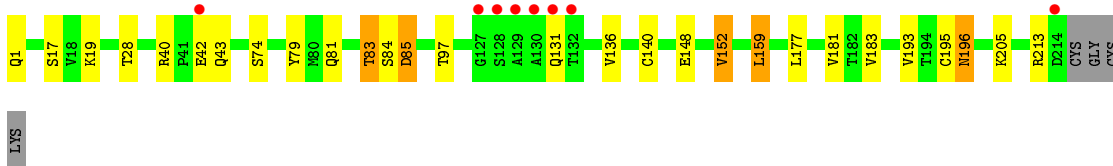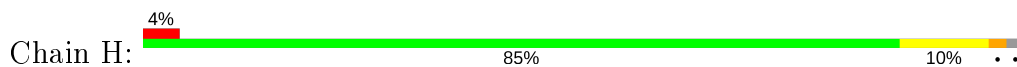- Molecule 1: Mouse monoclonal antibody C5.2 Fab light chain

Chain L:

- Molecule 1: Mouse monoclonal antibody C5.2 Fab light chain
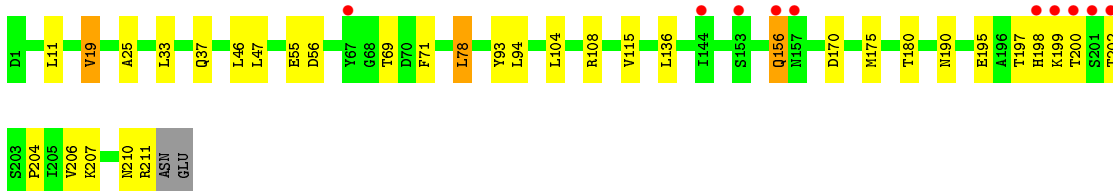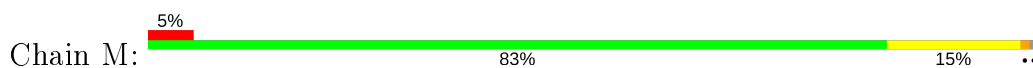
Chain M:

- Molecule 1: Mouse monoclonal antibody C5.2 Fab light chain
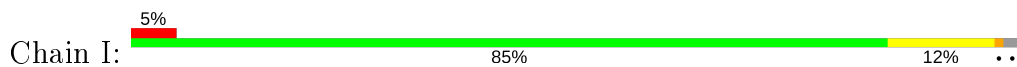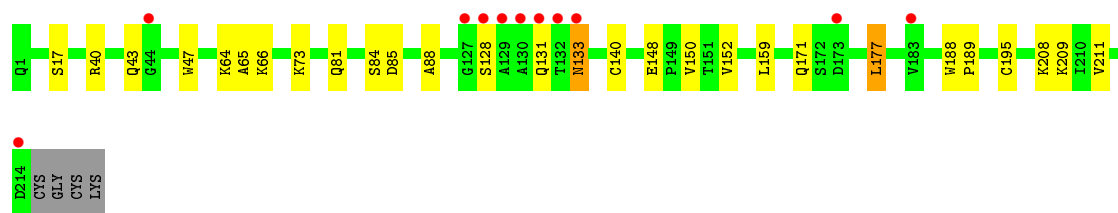
Chain N:

- Molecule 2: Mouse monoclonal antibody C5.2 Fab heavy chain

Chain H:

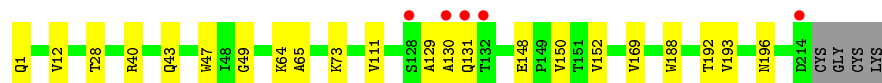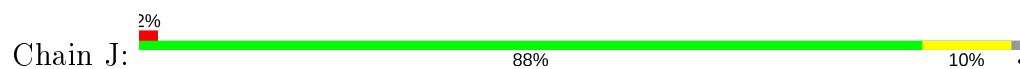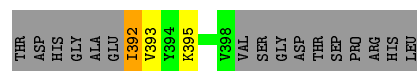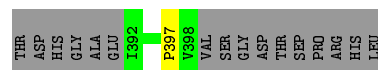- Molecule 2: Mouse monoclonal antibody C5.2 Fab heavy chain

Chain I:

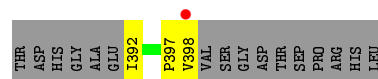● Molecule 2: Mouse monoclonal antibody C5.2 Fab heavy chain

● Molecule 3: Microtubule-associated protein tau

● Molecule 3: Microtubule-associated protein tau

● Molecule 3: Microtubule-associated protein tau

4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	161.16Å 40.53Å 228.06Å 90.00° 104.07° 90.00°	Depositor
Resolution (Å)	44.25 – 2.10 44.24 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.5 (44.25-2.10) 82.7 (44.24-2.10)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$<I/\sigma(I)>$ [1]	1.28 (at 2.10Å)	Xtriage
Refinement program	PHENIX 1.12_2829	Depositor
R, R_{free}	0.218 , 0.270 0.217 , 0.271	Depositor DCC
R_{free} test set	1910 reflections (2.36%)	wwPDB-VP
Wilson B-factor (Å²)	28.9	Xtriage
Anisotropy	0.439	Xtriage
Bulk solvent k_{sol}(e/Å³), B_{sol}(Å²)	0.34 , 47.0	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	0.96	EDS
Total number of atoms	10894	wwPDB-VP
Average B, all atoms (Å²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 73.03 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.9801e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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[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.

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: PO4, SEP

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	L	0.41	0/1673	0.58	1/2271 (0.0%)
1	M	0.39	0/1673	0.57	0/2271
1	N	0.41	0/1673	0.62	1/2271 (0.0%)
2	H	0.42	0/1679	0.59	1/2295 (0.0%)
2	I	0.40	0/1679	0.58	1/2295 (0.0%)
2	J	0.40	0/1679	0.59	0/2295
3	P	0.38	0/50	0.47	0/65
3	Q	0.30	0/50	0.36	0/65
3	R	0.33	0/50	0.44	0/65
All	All	0.40	0/10206	0.59	4/13893 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	L	11	LEU	CA-CB-CG	6.40	130.03	115.30
1	N	11	LEU	CA-CB-CG	6.39	130.00	115.30
2	I	177	LEU	CA-CB-CG	5.14	127.12	115.30
2	H	177	LEU	CA-CB-CG	5.01	126.83	115.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	L	1637	0	1571	9	0
1	M	1637	0	1571	19	0
1	N	1637	0	1571	11	0
2	H	1638	0	1596	10	0
2	I	1638	0	1596	12	0
2	J	1638	0	1596	11	0
3	P	60	0	61	3	0
3	Q	60	0	61	1	0
3	R	60	0	61	3	0
4	H	5	0	0	0	0
4	I	5	0	0	0	0
4	J	5	0	0	0	0
5	H	172	0	0	4	0
5	I	138	0	0	2	0
5	J	119	0	0	3	0
5	L	167	0	0	3	0
5	M	115	0	0	5	0
5	N	138	0	0	5	0
5	P	10	0	0	1	0
5	Q	10	0	0	1	0
5	R	5	0	0	3	0
All	All	10894	0	9684	77	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.

The worst 5 of 77 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)
3:R:392:ILE:N	5:R:501:HOH:O	1.97	0.97
1:L:45:LYS:NZ	5:L:301:HOH:O	1.90	0.96
3:P:392:ILE:N	5:P:501:HOH:O	2.11	0.81
1:M:170:ASP:OD2	5:M:301:HOH:O	2.04	0.75
2:I:148:GLU:OE1	5:I:401:HOH:O	2.04	0.75

There are no symmetry-related clashes.

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	L	209/213 (98%)	203 (97%)	6 (3%)	0	100	100
1	M	209/213 (98%)	202 (97%)	6 (3%)	1 (0%)	29	26
1	N	209/213 (98%)	201 (96%)	8 (4%)	0	100	100
2	H	214/220 (97%)	209 (98%)	5 (2%)	0	100	100
2	I	214/220 (97%)	211 (99%)	3 (1%)	0	100	100
2	J	214/220 (97%)	208 (97%)	6 (3%)	0	100	100
3	P	4/23 (17%)	4 (100%)	0	0	100	100
3	Q	4/23 (17%)	4 (100%)	0	0	100	100
3	R	4/23 (17%)	4 (100%)	0	0	100	100
All	All	1281/1368 (94%)	1246 (97%)	34 (3%)	1 (0%)	51	54

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	M	200	THR

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	L	189/191 (99%)	182 (96%)	7 (4%)	34	35
1	M	189/191 (99%)	181 (96%)	8 (4%)	30	30
1	N	189/191 (99%)	179 (95%)	10 (5%)	22	20

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	H	187/190 (98%)	171 (91%)	16 (9%)	10	7
2	I	187/190 (98%)	175 (94%)	12 (6%)	17	14
2	J	187/190 (98%)	183 (98%)	4 (2%)	53	59
3	P	6/18 (33%)	5 (83%)	1 (17%)	2	1
3	Q	6/18 (33%)	6 (100%)	0	100	100
3	R	6/18 (33%)	5 (83%)	1 (17%)	2	1
All	All	1146/1197 (96%)	1087 (95%)	59 (5%)	24	22

5 of 59 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	M	78	LEU
2	I	84	SER
2	J	1	GLN
1	M	156	GLN
1	M	198	HIS

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

5.3.3   RNA (i)

There are no RNA molecules in this entry.

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

3 non-standard protein/DNA/RNA residues 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	#$|Z| > 2$
3	SEP	P	396	3	8,9,10	1.47	1 (12%)	8,12,14	1.57	3 (37%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	#$|Z|$ > 2	Counts	RMSZ	#$|Z|$ > 2
3	SEP	Q	396	3	8,9,10	1.57	1 (12%)	8,12,14	1.83	2 (25%)
3	SEP	R	396	3	8,9,10	1.60	1 (12%)	8,12,14	1.29	1 (12%)

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	SEP	P	396	3	-	0/5/8/10	-
3	SEP	Q	396	3	-	2/5/8/10	-
3	SEP	R	396	3	-	0/5/8/10	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	R	396	SEP	P-O1P	3.52	1.61	1.50
3	Q	396	SEP	P-O1P	3.50	1.61	1.50
3	P	396	SEP	P-O1P	3.12	1.60	1.50

The worst 5 of 6 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	Q	396	SEP	OG-CB-CA	3.98	112.02	108.14
3	P	396	SEP	OG-CB-CA	2.95	111.01	108.14
3	R	396	SEP	P-OG-CB	-2.46	111.51	118.30
3	P	396	SEP	P-OG-CB	-2.32	111.89	118.30
3	P	396	SEP	O2P-P-OG	2.06	112.22	106.73

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	Q	396	SEP	CB-OG-P-O1P
3	Q	396	SEP	CA-CB-OG-P

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates (i)

There are no carbohydrates in this entry.

5.6 Ligand geometry (i)

3 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	#$|Z| > 2$
4	PO4	H	301	-	4,4,4	0.95	0	6,6,6	0.90	0
4	PO4	J	301	-	4,4,4	1.01	0	6,6,6	0.71	0
4	PO4	I	301	-	4,4,4	0.87	0	6,6,6	0.75	0

There are no bond length outliers.

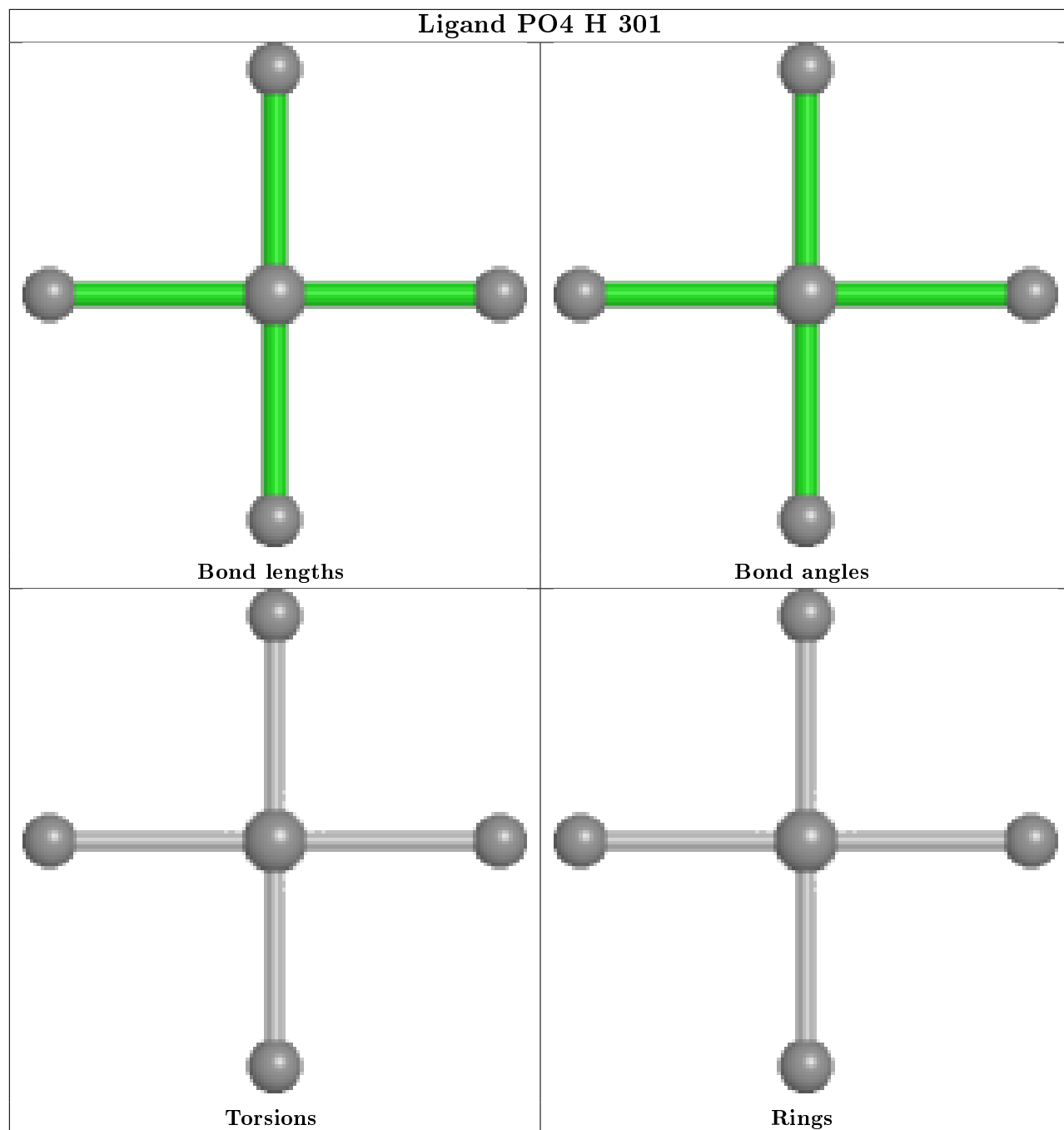There are no bond angle outliers.
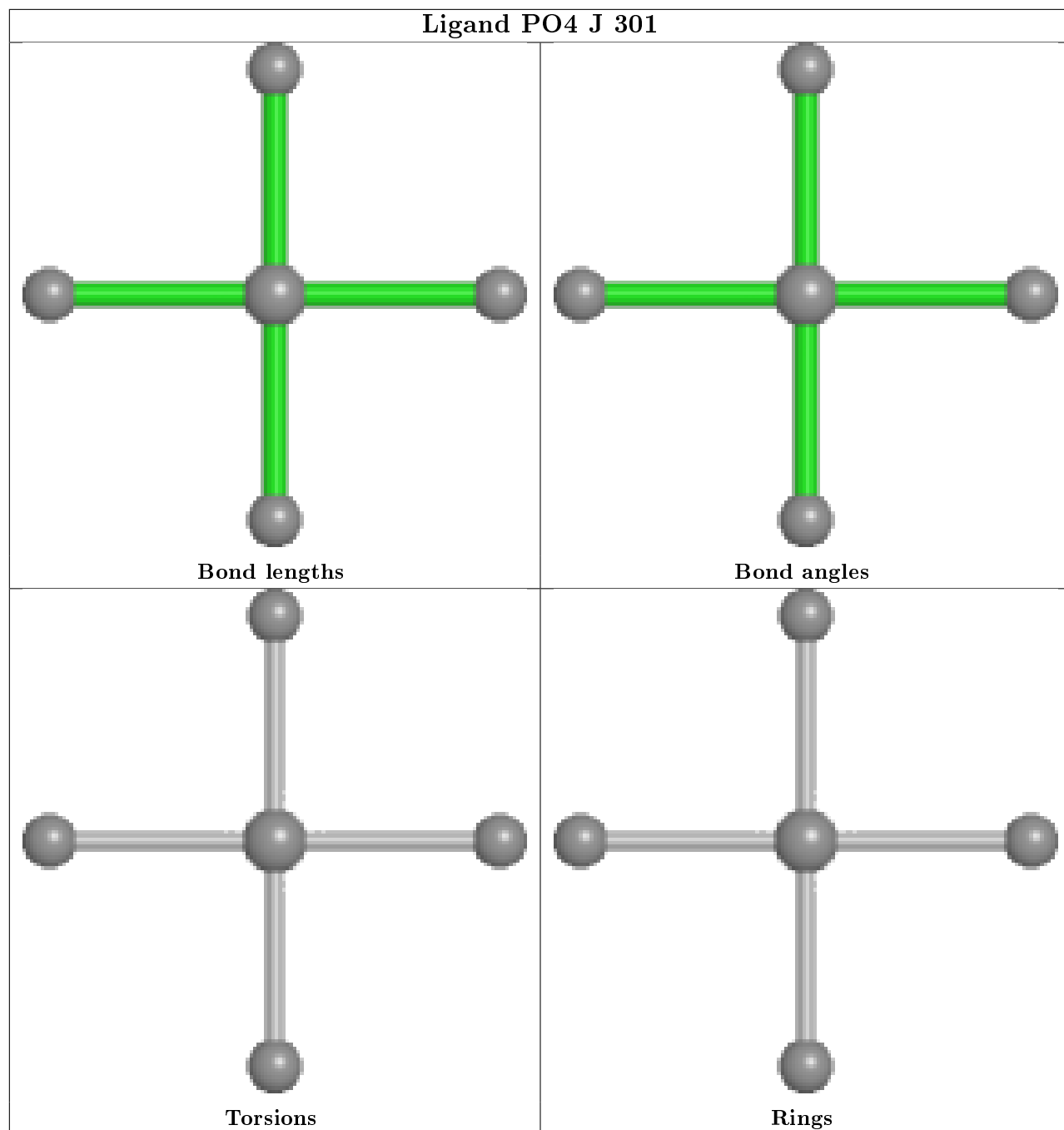
There are no chirality outliers.
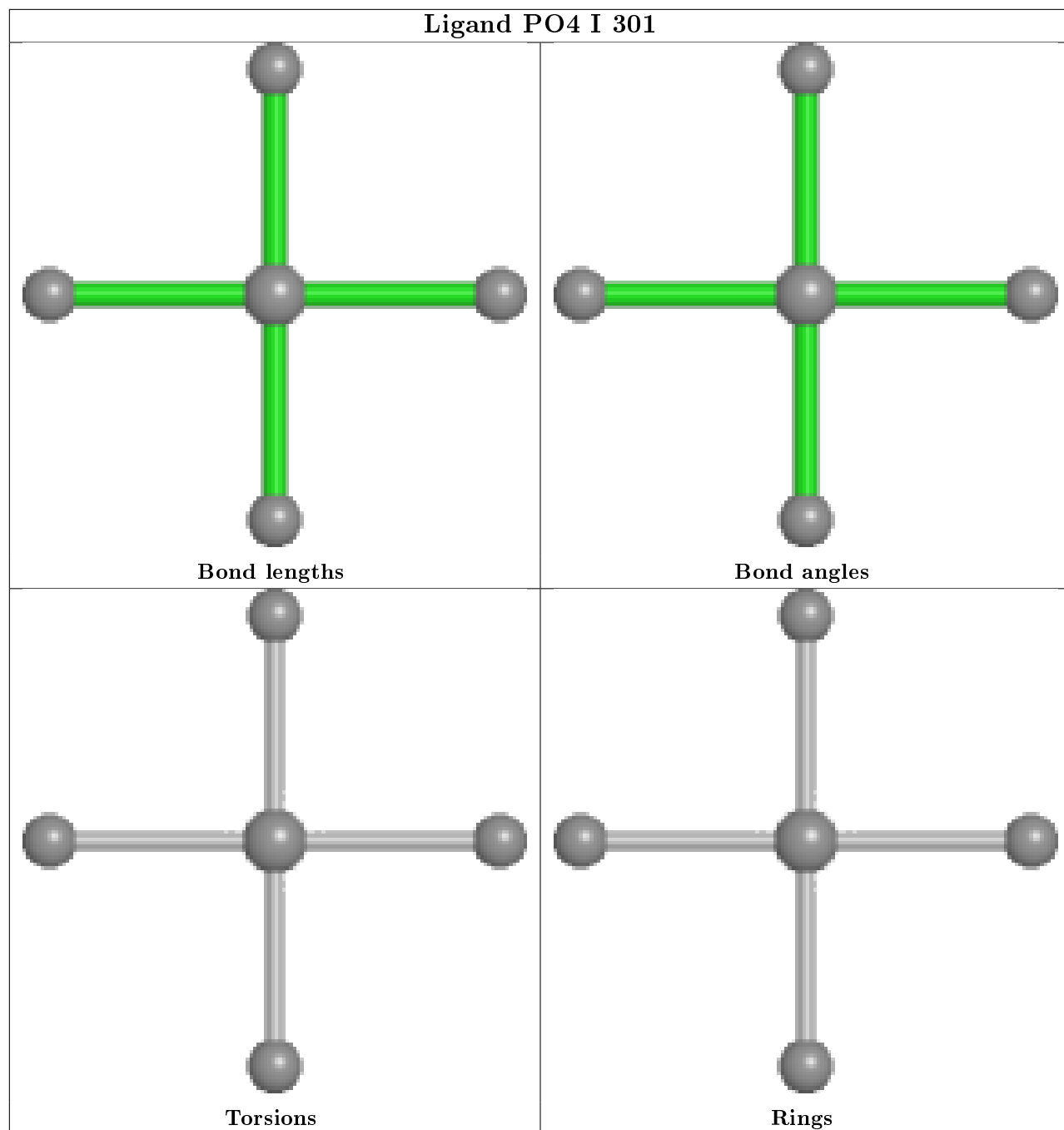
There are no torsion outliers.

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 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.

Ligand PO4 H 301

Bond lengths

Bond angles

Torsions

Rings

Ligand PO4 J 301

Bond lengths

Bond angles

Torsions

Rings

Ligand PO4 I 301

Bond lengths

Bond angles

Torsions

Rings

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

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	L	211/213 (99%)	0.06	6 (2%)	53	59	21, 32, 50, 80	0
1	M	211/213 (99%)	0.19	10 (4%)	31	37	27, 38, 61, 99	0
1	N	211/213 (99%)	0.00	6 (2%)	53	59	27, 36, 53, 81	0
2	H	216/220 (98%)	0.08	8 (3%)	41	48	19, 30, 46, 74	0
2	I	216/220 (98%)	0.19	11 (5%)	28	33	26, 35, 56, 88	0
2	J	216/220 (98%)	0.07	5 (2%)	60	65	27, 36, 54, 79	0
3	P	6/23 (26%)	0.36	0	100	100	36, 44, 51, 54	0
3	Q	6/23 (26%)	0.22	0	100	100	38, 45, 51, 52	0
3	R	6/23 (26%)	0.84	1 (16%)	1	2	38, 46, 53, 60	0
All	All	1299/1368 (94%)	0.10	47 (3%)	42	49	19, 35, 55, 99	0

The worst 5 of 47 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	I	131	GLN	8.4
2	J	130	ALA	7.7
1	L	200	THR	7.5
1	N	200	THR	7.3
1	M	198	HIS	7.0

## 6.2 Non-standard residues in protein, DNA, RNA chains (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(Å²)	Q<0.9
3	SEP	Q	396	10/11	0.97	0.11	31,36,40,42	0
3	SEP	R	396	10/11	0.97	0.10	34,37,46,47	0
3	SEP	P	396	10/11	0.98	0.10	27,35,42,44	0

6.3 Carbohydrates (i)

There are no carbohydrates 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(Å²)	Q<0.9
4	PO4	J	301	5/5	0.96	0.13	45,48,50,52	0
4	PO4	I	301	5/5	0.96	0.10	40,40,47,50	0
4	PO4	H	301	5/5	0.97	0.11	36,39,40,49	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.
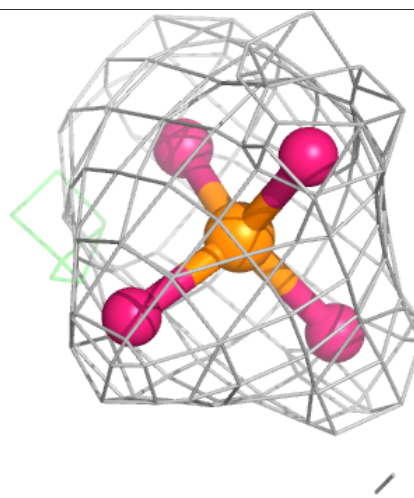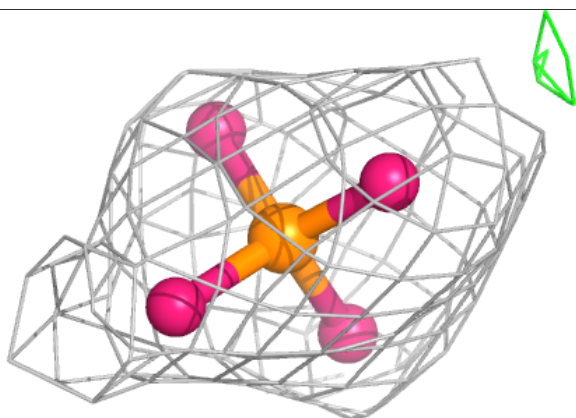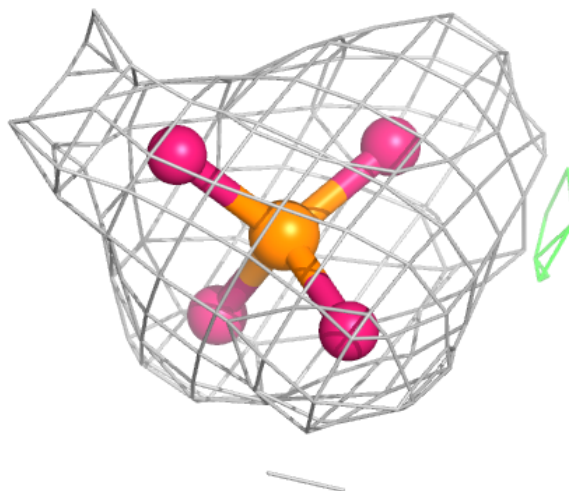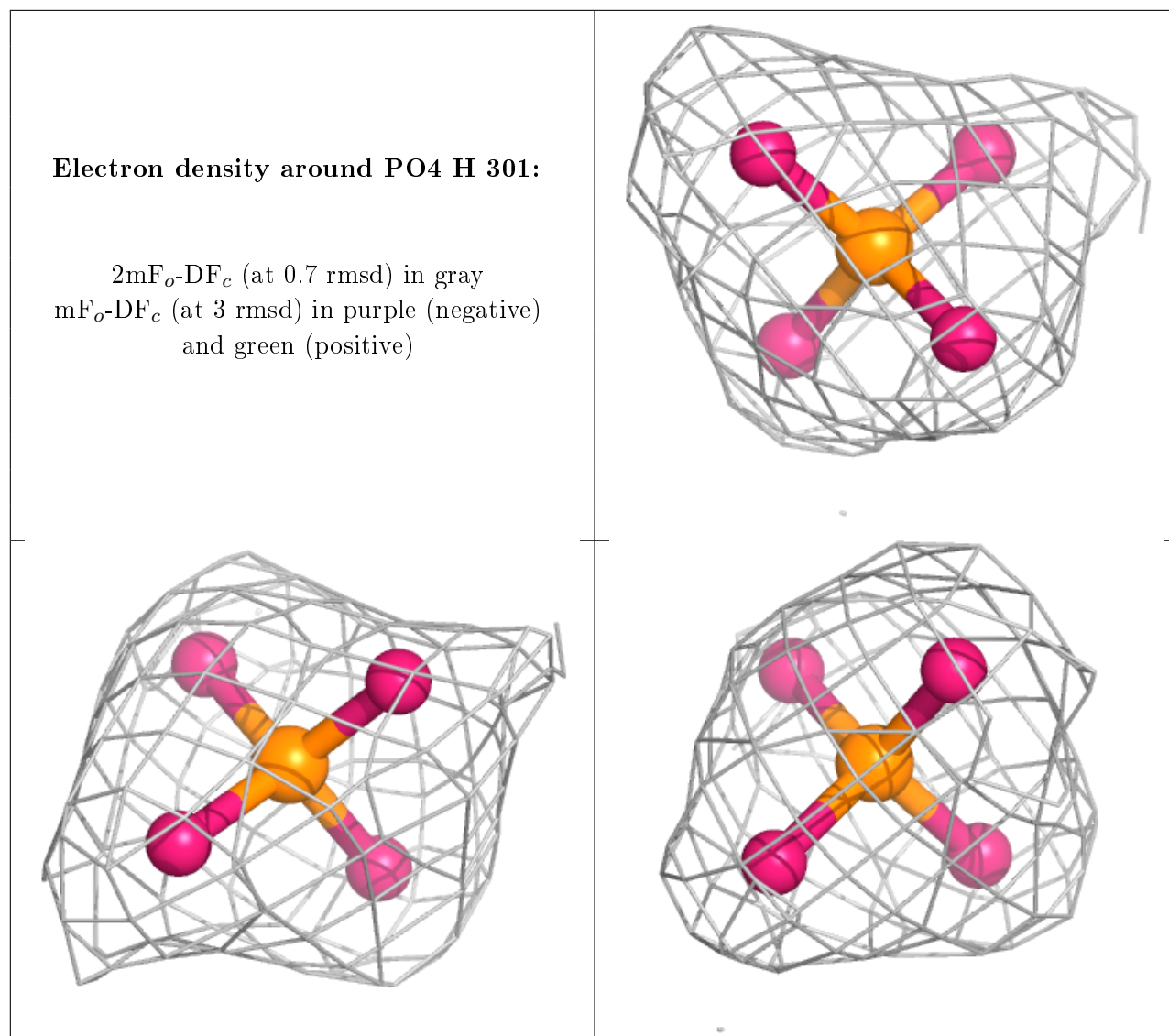
Electron density around PO4 J 301:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)
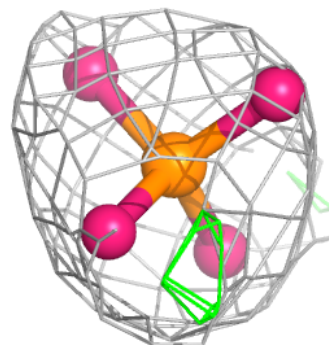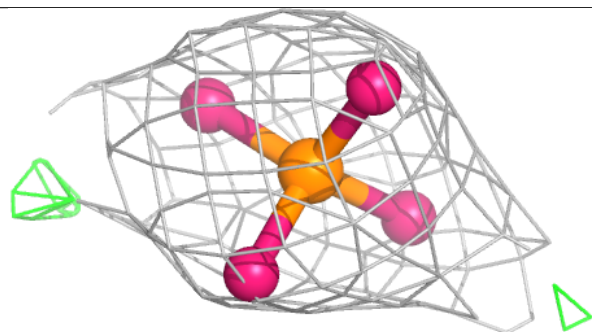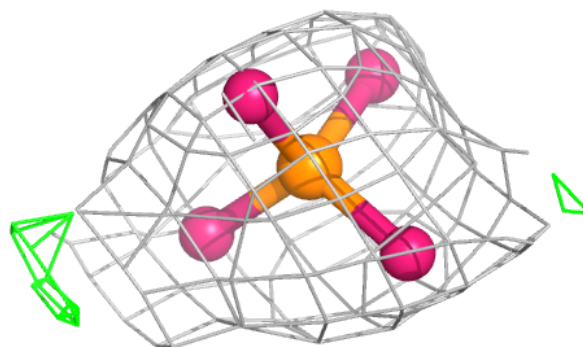
Electron density around PO4 I 301:

$2mF_o$-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 H 301:

$2mF_o$-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.