



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 9, 2023 – 07:06 PM EDT

PDB ID : 6WBO  
Title : DNA-Ligase from *Thermococcus gammatolerans*  
Authors : Flores-Hernandez, E.; Cardona-Felix, C.; Miranda-Blancas, R.; Rudino-Pinera, E.  
Deposited on : 2020-03-26  
Resolution : 2.65 Å(reported)

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

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.35.1  
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) : 2.35.1

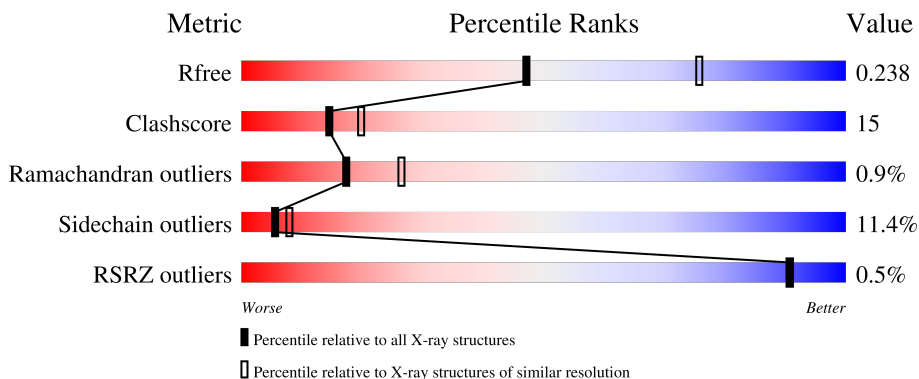
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*


The reported resolution of this entry is 2.65 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	1332 (2.68-2.64)
Clashscore	141614	1374 (2.68-2.64)
Ramachandran outliers	138981	1349 (2.68-2.64)
Sidechain outliers	138945	1349 (2.68-2.64)
RSRZ outliers	127900	1318 (2.68-2.64)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	555	

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 4644 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 DNA ligase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	555	4438	2839	759	828	12	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	11	HIS	-	expression tag	UNP C5A7K8
A	544	ASN	ASP	conflict	UNP C5A7K8

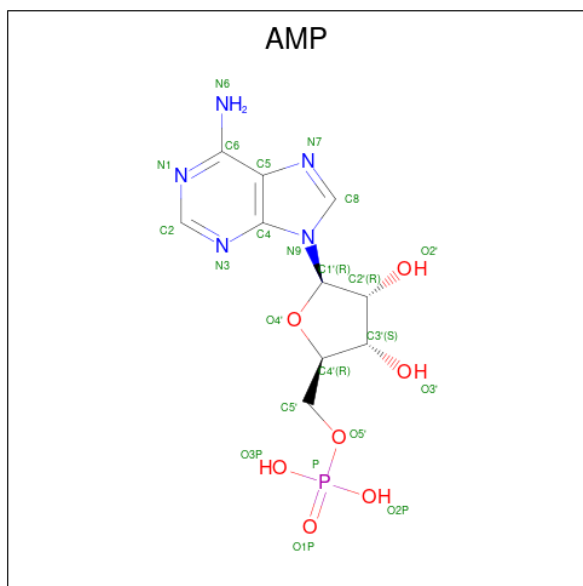
- Molecule 2 is PHOSPHATE ION (three-letter code: PO4) (formula: O<sub>4</sub>P) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 3 is ADENOSINE MONOPHOSPHATE (three-letter code: AMP) (formula: C<sub>10</sub>H<sub>14</sub>N<sub>5</sub>O<sub>7</sub>P)

(labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
3	A	1	22	10	5	6	1	0	0

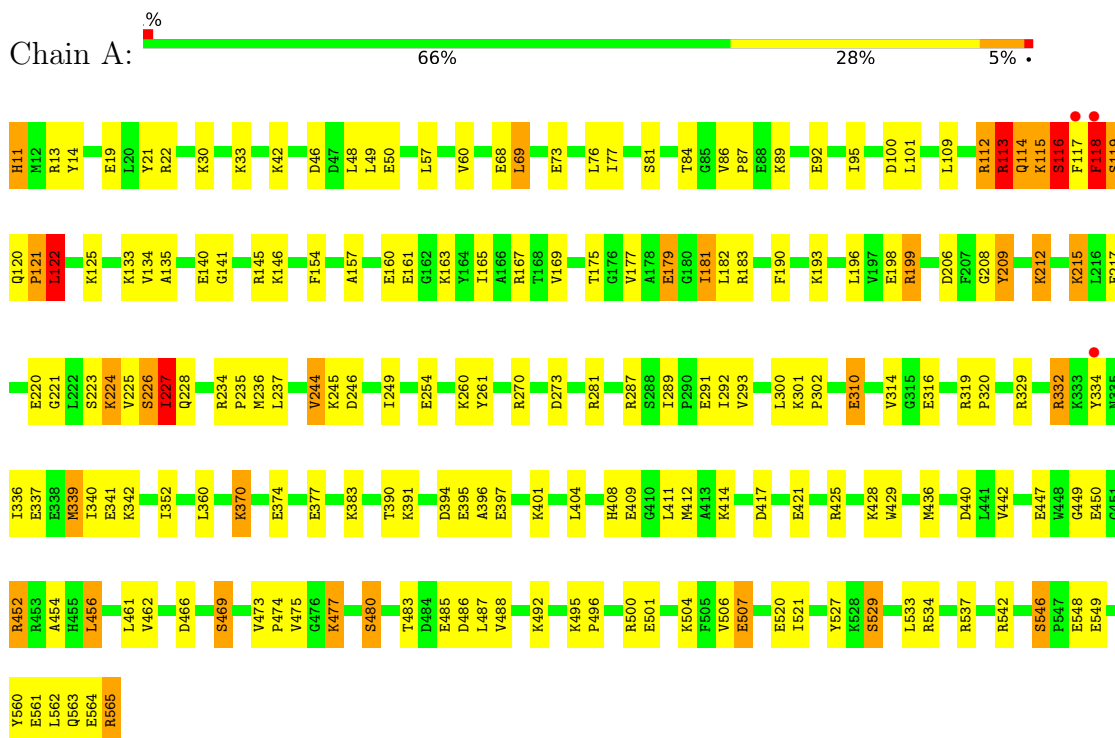
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	179	179	179	0	0

### 3 Residue-property plots

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

- Molecule 1: DNA ligase



## 4 Data and refinement statistics i

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	91.67Å 92.75Å 98.61Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.33 – 2.65 39.33 – 2.65	Depositor EDS
% Data completeness (in resolution range)	94.5 (39.33-2.65) 94.9 (39.33-2.65)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.54 (at 2.65Å)	Xtrriage
Refinement program	PHENIX 1.10.1_2155	Depositor
R, $R_{free}$	0.233 , 0.238 0.233 , 0.238	Depositor DCC
$R_{free}$ test set	1151 reflections (4.85%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	32.5	Xtrriage
Anisotropy	0.222	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 39.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.006 for k,h,-l	Xtrriage
$F_o, F_c$ correlation	0.87	EDS
Total number of atoms	4644	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	36.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.47% of the height of the origin peak. No significant pseudotranslation is detected.*

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

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

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

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	1.45	3/4511 (0.1%)	0.95	6/6062 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	209	TYR	CE1-CZ	-5.94	1.30	1.38
1	A	310	GLU	CD-OE2	-5.69	1.19	1.25
1	A	21	TYR	CE1-CZ	-5.51	1.31	1.38

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	113	ARG	N-CA-C	7.23	130.51	111.00
1	A	115	LYS	N-CA-C	6.14	127.59	111.00
1	A	114	GLN	N-CA-C	5.61	126.15	111.00
1	A	116	SER	N-CA-C	5.39	125.56	111.00
1	A	69	LEU	N-CA-C	-5.24	96.84	111.00
1	A	227	ILE	N-CA-C	-5.12	97.17	111.00

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4438	0	4576	139	2
2	A	5	0	0	0	0
3	A	22	0	12	1	0
4	A	179	0	0	12	0
All	All	4644	0	4588	139	2

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 15.

All (139) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:421:GLU:OE1	1:A:428:LYS:CE	1.78	1.29
1:A:118:PHE:O	1:A:120:GLN:HG2	1.22	1.29
1:A:421:GLU:OE1	1:A:428:LYS:HE3	1.17	1.26
1:A:377:GLU:OE2	4:A:701:HOH:O	1.56	1.21
1:A:118:PHE:O	1:A:120:GLN:N	1.73	1.21
1:A:112:ARG:NH2	1:A:157:ALA:O	1.73	1.20
1:A:121:PRO:O	1:A:122:LEU:HB2	1.32	1.11
1:A:118:PHE:O	1:A:120:GLN:CG	1.99	1.10
1:A:440:ASP:OD2	1:A:542:ARG:NH1	1.91	1.03
1:A:118:PHE:C	1:A:120:GLN:HG2	1.80	1.01
1:A:46:ASP:OD1	1:A:215:LYS:NZ	1.93	0.99
1:A:370:LYS:O	1:A:374:GLU:HG3	1.66	0.94
1:A:401:LYS:NZ	4:A:705:HOH:O	2.00	0.93
1:A:121:PRO:O	1:A:122:LEU:CB	2.16	0.87
1:A:118:PHE:O	1:A:119:SER:C	2.07	0.86
1:A:421:GLU:OE1	1:A:428:LYS:HE2	1.74	0.86
1:A:118:PHE:C	1:A:120:GLN:N	2.28	0.85
1:A:69:LEU:HD23	1:A:135:ALA:HB2	1.59	0.84
1:A:228:GLN:O	4:A:704:HOH:O	1.97	0.81
1:A:440:ASP:CG	1:A:542:ARG:NH1	2.33	0.81
1:A:175:THR:OG1	4:A:702:HOH:O	1.68	0.78
1:A:564:GLU:C	1:A:565:ARG:HG3	2.04	0.77
1:A:118:PHE:O	1:A:120:GLN:CB	2.32	0.77
1:A:565:ARG:HE	1:A:565:ARG:C	1.89	0.76
1:A:440:ASP:CG	1:A:542:ARG:HH12	1.88	0.74
1:A:440:ASP:OD1	1:A:542:ARG:NH1	2.19	0.74
1:A:474:PRO:O	1:A:529:SER:OG	2.04	0.74
1:A:332:ARG:NE	1:A:334:TYR:O	2.20	0.73
1:A:501:GLU:HG2	1:A:506:VAL:HG12	1.71	0.72
1:A:92:GLU:OE2	4:A:706:HOH:O	2.06	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:165:ILE:O	1:A:169:VAL:HG23	1.88	0.71
1:A:118:PHE:CA	1:A:120:GLN:HG2	2.19	0.71
1:A:390:THR:OG1	1:A:395:GLU:OE2	2.06	0.71
1:A:246:ASP:OD2	1:A:537:ARG:NH2	2.24	0.71
1:A:226:SER:O	1:A:227:ILE:HG13	1.90	0.70
1:A:261:TYR:HB2	1:A:310:GLU:HG2	1.72	0.70
1:A:134:VAL:HG13	1:A:146:LYS:HG2	1.74	0.70
1:A:118:PHE:C	1:A:120:GLN:H	1.97	0.68
1:A:19:GLU:OE2	1:A:22:ARG:NE	2.27	0.67
1:A:183:ARG:NH2	1:A:198:GLU:OE2	2.27	0.67
1:A:68:GLU:O	4:A:707:HOH:O	2.12	0.66
1:A:177:VAL:HG13	1:A:181:ILE:HG22	1.79	0.65
1:A:452:ARG:HG2	1:A:452:ARG:NH1	2.12	0.64
1:A:564:GLU:O	1:A:565:ARG:HG3	1.96	0.64
1:A:560:TYR:O	1:A:563:GLN:HB3	1.97	0.64
1:A:329:ARG:HD2	1:A:339:MET:HG2	1.78	0.64
1:A:542:ARG:HH11	1:A:542:ARG:HG3	1.62	0.64
1:A:115:LYS:HB2	4:A:788:HOH:O	1.98	0.63
1:A:73:GLU:O	1:A:77:ILE:HG12	1.97	0.63
1:A:452:ARG:CG	1:A:452:ARG:HH11	2.12	0.63
1:A:452:ARG:HG2	1:A:452:ARG:HH11	1.65	0.62
1:A:190:PHE:CE1	1:A:215:LYS:HG3	2.34	0.61
1:A:118:PHE:O	1:A:120:GLN:CA	2.49	0.61
1:A:177:VAL:HG13	1:A:181:ILE:CG2	2.31	0.61
1:A:520:GLU:HG2	1:A:521:ILE:N	2.15	0.60
1:A:225:VAL:O	4:A:709:HOH:O	2.17	0.59
1:A:48:LEU:HD21	1:A:125:LYS:HG3	1.84	0.59
1:A:14:TYR:HD2	1:A:161:GLU:HG2	1.67	0.58
1:A:254:GLU:HG3	1:A:391:LYS:HG2	1.85	0.58
1:A:260:LYS:HD3	1:A:409:GLU:HB2	1.85	0.58
1:A:46:ASP:HA	1:A:49:LEU:HD12	1.84	0.58
1:A:84:THR:CG2	1:A:109:LEU:HD23	2.35	0.57
1:A:456:LEU:HD11	1:A:488:VAL:HG22	1.86	0.56
1:A:495:LYS:N	1:A:496:PRO:HD2	2.20	0.56
1:A:483:THR:H	1:A:486:ASP:HB2	1.71	0.55
1:A:475:VAL:HG12	1:A:533:LEU:HG	1.90	0.54
1:A:449:GLY:O	1:A:454:ALA:HA	2.07	0.54
1:A:520:GLU:OE1	1:A:534:ARG:HD3	2.09	0.53
1:A:220:GLU:O	1:A:224:LYS:HG3	2.08	0.53
1:A:337:GLU:HA	1:A:340:ILE:HD12	1.91	0.53
1:A:483:THR:O	1:A:486:ASP:HB2	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:289:ILE:O	1:A:293:VAL:HG23	2.09	0.52
1:A:316:GLU:HA	1:A:316:GLU:OE2	2.10	0.52
1:A:245:LYS:O	1:A:249:ILE:HG12	2.11	0.51
1:A:244:VAL:HG21	1:A:396:ALA:HB3	1.92	0.50
1:A:33:LYS:NZ	4:A:715:HOH:O	2.38	0.50
1:A:352:ILE:HG23	1:A:360:LEU:HD12	1.94	0.50
1:A:474:PRO:HB2	1:A:527:TYR:HB3	1.94	0.50
1:A:237:LEU:HB2	1:A:425:ARG:HB2	1.93	0.50
1:A:134:VAL:O	1:A:146:LYS:NZ	2.43	0.50
1:A:461:LEU:HD13	1:A:477:LYS:HE3	1.93	0.49
1:A:112:ARG:NH1	1:A:154:PHE:O	2.42	0.49
1:A:473:VAL:HG13	1:A:473:VAL:O	2.12	0.48
1:A:160:GLU:OE2	1:A:163:LYS:NZ	2.43	0.48
1:A:226:SER:OG	4:A:708:HOH:O	2.13	0.48
1:A:92:GLU:O	1:A:95:ILE:HG12	2.13	0.48
1:A:565:ARG:HE	1:A:565:ARG:CA	2.26	0.47
1:A:30:LYS:HG2	1:A:181:ILE:HD11	1.95	0.47
1:A:14:TYR:CD2	1:A:161:GLU:HG2	2.48	0.47
1:A:301:LYS:N	1:A:302:PRO:HD2	2.30	0.47
1:A:49:LEU:HD11	1:A:215:LYS:HD2	1.97	0.47
1:A:270:ARG:NH1	1:A:300:LEU:O	2.48	0.47
1:A:316:GLU:HB3	1:A:319:ARG:HH21	1.81	0.46
1:A:417:ASP:OD1	1:A:417:ASP:N	2.41	0.46
1:A:114:GLN:H	1:A:114:GLN:HG2	1.52	0.46
1:A:291:GLU:HG2	1:A:292:ILE:N	2.30	0.46
1:A:480:SER:HB3	1:A:537:ARG:HG3	1.98	0.46
1:A:466:ASP:HB3	1:A:469:SER:HB2	1.98	0.46
1:A:483:THR:OG1	1:A:486:ASP:HB2	2.16	0.45
1:A:57:LEU:HD21	1:A:182:LEU:HD22	1.99	0.45
1:A:11:HIS:N	1:A:125:LYS:HE3	2.32	0.44
1:A:500:ARG:HB3	1:A:507:GLU:HB2	1.99	0.44
1:A:236:MET:HE3	1:A:429:TRP:CE2	2.53	0.44
1:A:370:LYS:HA	1:A:370:LYS:HD2	1.69	0.44
1:A:404:LEU:HD23	1:A:404:LEU:HA	1.82	0.44
1:A:542:ARG:NH1	1:A:542:ARG:HG3	2.30	0.44
1:A:500:ARG:HE	1:A:500:ARG:HB2	1.57	0.44
1:A:495:LYS:N	1:A:496:PRO:CD	2.80	0.43
1:A:320:PRO:HG3	1:A:408:HIS:CE1	2.53	0.43
1:A:212:LYS:HE3	1:A:212:LYS:HB3	1.57	0.43
1:A:414:LYS:HD3	1:A:429:TRP:CD1	2.53	0.42
1:A:483:THR:N	1:A:486:ASP:HB2	2.32	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:301:LYS:HB2	1:A:302:PRO:HD3	2.00	0.42
1:A:100:ASP:OD2	1:A:167:ARG:NH2	2.50	0.42
1:A:442:VAL:O	1:A:462:VAL:HA	2.19	0.42
1:A:196:LEU:HD23	1:A:196:LEU:HA	1.87	0.42
1:A:199:ARG:NH2	4:A:709:HOH:O	2.52	0.42
1:A:301:LYS:HB2	1:A:302:PRO:CD	2.49	0.42
1:A:236:MET:HE3	1:A:429:TRP:NE1	2.35	0.42
1:A:336:ILE:O	1:A:339:MET:N	2.52	0.41
1:A:520:GLU:CG	1:A:521:ILE:N	2.82	0.41
1:A:487:LEU:HD23	1:A:487:LEU:HA	1.88	0.41
1:A:436:MET:SD	1:A:436:MET:O	2.78	0.41
1:A:113:ARG:HG2	4:A:782:HOH:O	2.19	0.41
1:A:141:GLY:O	1:A:145:ARG:HB2	2.20	0.41
1:A:546:SER:HB3	1:A:549:GLU:HG3	2.02	0.41
1:A:86:VAL:HA	1:A:87:PRO:HD3	1.88	0.41
1:A:13:ARG:HH21	1:A:13:ARG:HD2	1.73	0.41
1:A:50:GLU:HA	1:A:208:GLY:HA3	2.03	0.41
1:A:227:ILE:HG21	1:A:227:ILE:HD13	1.83	0.41
1:A:183:ARG:HH21	1:A:183:ARG:HD3	1.75	0.40
1:A:179:GLU:H	1:A:179:GLU:HG2	1.54	0.40
1:A:292:ILE:HA	1:A:292:ILE:HD13	1.79	0.40
1:A:412:MET:HG3	3:A:602:AMP:C6	2.55	0.40
1:A:206:ASP:HB3	1:A:209:TYR:HB3	2.03	0.40
1:A:217:GLU:HB3	1:A:221:GLY:HA3	2.04	0.40
1:A:428:LYS:H	1:A:428:LYS:HG2	1.69	0.40
1:A:115:LYS:HB2	1:A:116:SER:H	1.45	0.40
1:A:234:ARG:HA	1:A:235:PRO:HD3	1.78	0.40

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:337:GLU:OE2	1:A:450:GLU:OE2[2_575]	1.17	1.03
1:A:337:GLU:CD	1:A:450:GLU:OE2[2_575]	1.89	0.31

## 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	553/555 (100%)	530 (96%)	18 (3%)	5 (1%)	17 26

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	116	SER
1	A	118	PHE
1	A	119	SER
1	A	227	ILE
1	A	122	LEU

### 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	475/475 (100%)	421 (89%)	54 (11%)	5 8

All (54) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	HIS
1	A	42	LYS
1	A	60	VAL
1	A	76	LEU
1	A	81	SER
1	A	89	LYS

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	A	101	LEU
1	A	112	ARG
1	A	113	ARG
1	A	117	PHE
1	A	118	PHE
1	A	121	PRO
1	A	122	LEU
1	A	133	LYS
1	A	140	GLU
1	A	179	GLU
1	A	181	ILE
1	A	193	LYS
1	A	199	ARG
1	A	212	LYS
1	A	215	LYS
1	A	223	SER
1	A	224	LYS
1	A	226	SER
1	A	244	VAL
1	A	273	ASP
1	A	281	ARG
1	A	287	ARG
1	A	314	VAL
1	A	332	ARG
1	A	339	MET
1	A	341	GLU
1	A	342	LYS
1	A	370	LYS
1	A	383	LYS
1	A	394	ASP
1	A	397	GLU
1	A	411	LEU
1	A	447	GLU
1	A	452	ARG
1	A	456	LEU
1	A	469	SER
1	A	477	LYS
1	A	480	SER
1	A	485	GLU
1	A	492	LYS
1	A	504	LYS
1	A	507	GLU

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Mol	Chain	Res	Type
1	A	529	SER
1	A	546	SER
1	A	548	GLU
1	A	561	GLU
1	A	562	LEU
1	A	565	ARG

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

Mol	Chain	Res	Type
1	A	228	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 monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

2 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	$\# Z  > 2$
3	AMP	A	602	1	18,24,25	1.13	3 (16%)	18,35,38	1.30	3 (16%)
2	PO4	A	601	-	4,4,4	0.72	0	6,6,6	0.62	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	AMP	A	602	1	-	0/3/25/26	0/3/3/3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	602	AMP	C2-N3	2.25	1.35	1.32
3	A	602	AMP	C5-C4	2.20	1.46	1.40
3	A	602	AMP	O4'-C1'	2.02	1.43	1.41

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	602	AMP	C4-C5-N7	-2.77	106.51	109.40
3	A	602	AMP	N3-C2-N1	-2.46	124.83	128.68
3	A	602	AMP	C2'-C3'-C4'	2.18	106.89	102.64

There are no chirality outliers.

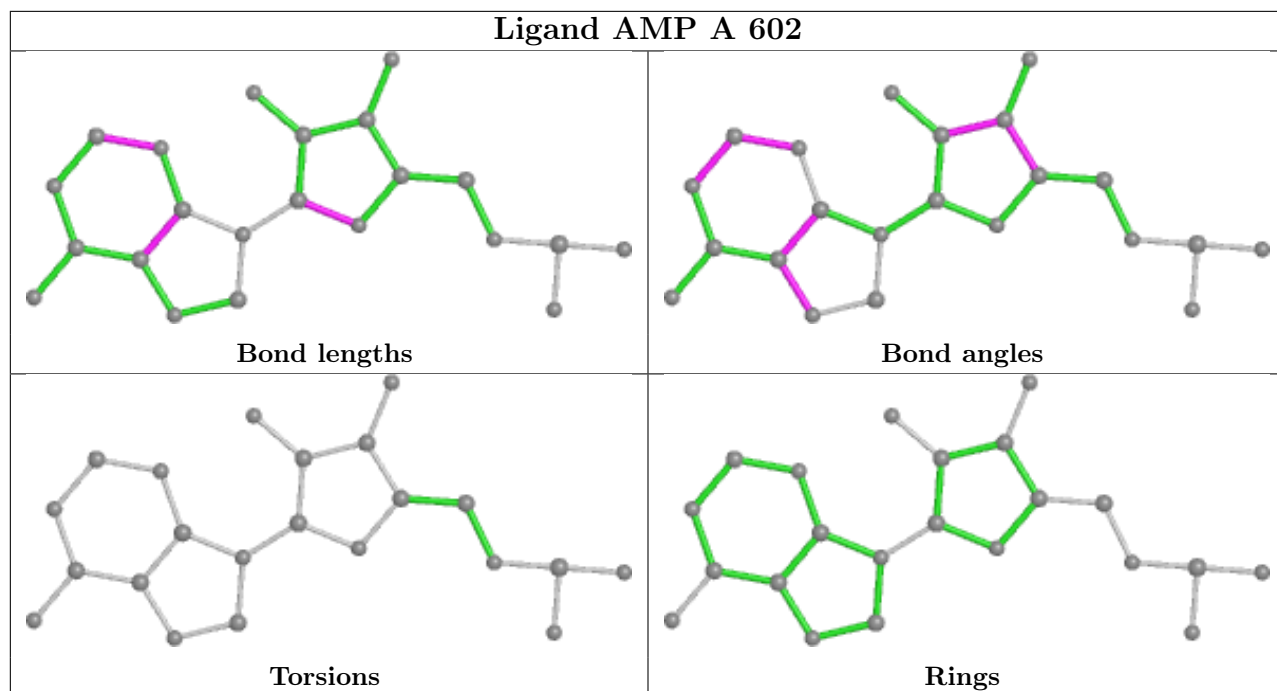
There are no torsion outliers.

There are no ring outliers.

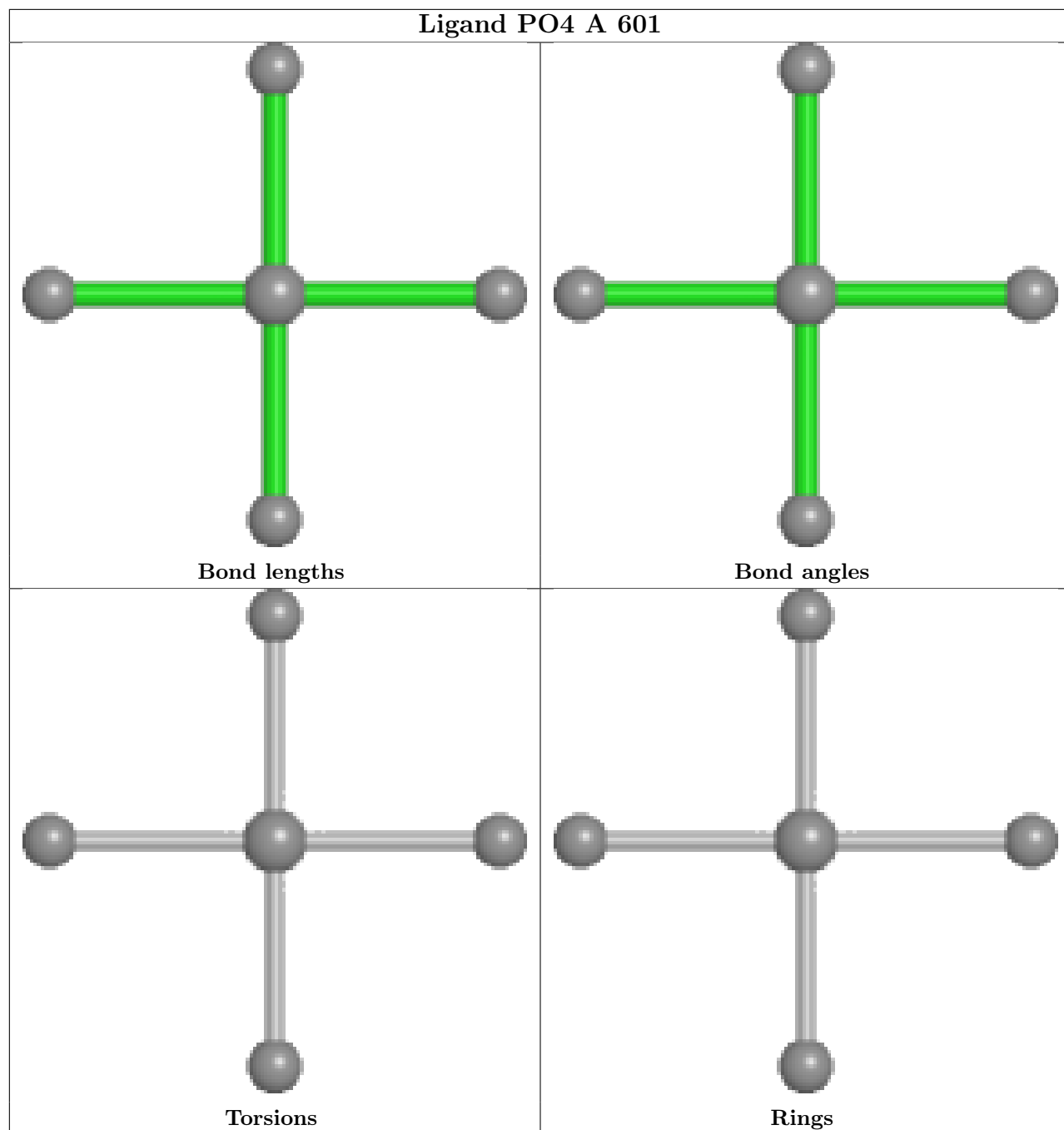
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	602	AMP	1	0

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







### 5.7 Other polymers [i](#)

There are no such residues in this entry.

### 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data [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, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	555/555 (100%)	-0.14	3 (0%) 91   91	21, 34, 56, 78	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	118	PHE	5.7
1	A	117	PHE	3.0
1	A	334	TYR	2.4

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

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

### 6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

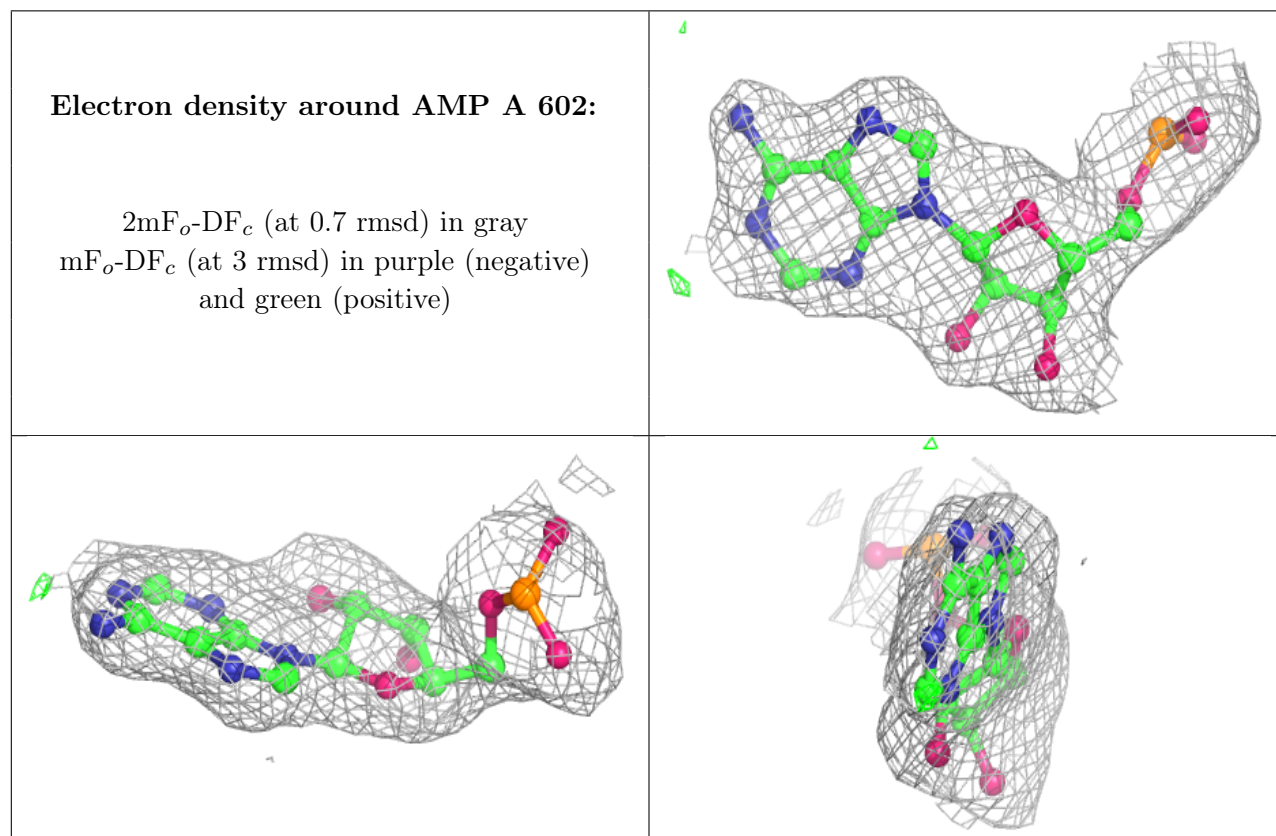
### 6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
3	AMP	A	602	22/23	0.95	0.16	25,28,32,35	0
2	PO4	A	601	5/5	0.98	0.14	22,24,27,31	0

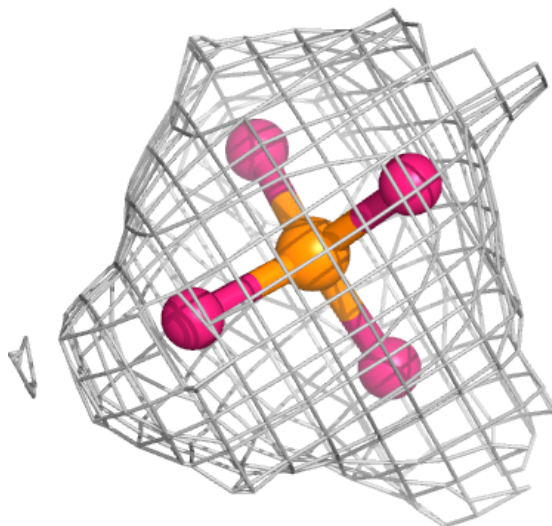
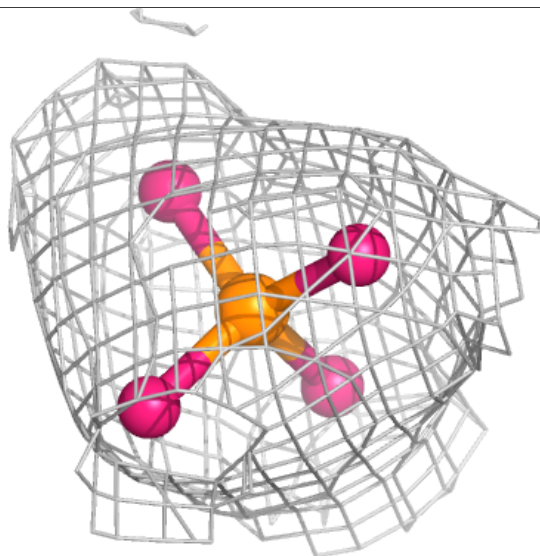
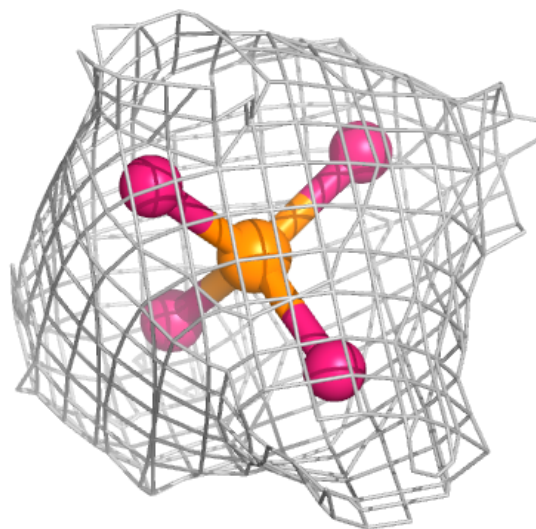
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers

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 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)



## 6.5 Other polymers [i](#)

There are no such residues in this entry.