



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 10TH / pdb_000010th
Title : Crystal structure of the PXL domain of the human promyelocytic leukemia (PML) protein isoform 1
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Deposited on : 2026-02-07
Resolution : 3.01 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity	:	4-5-2 with Phenix2.2
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.2
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.51

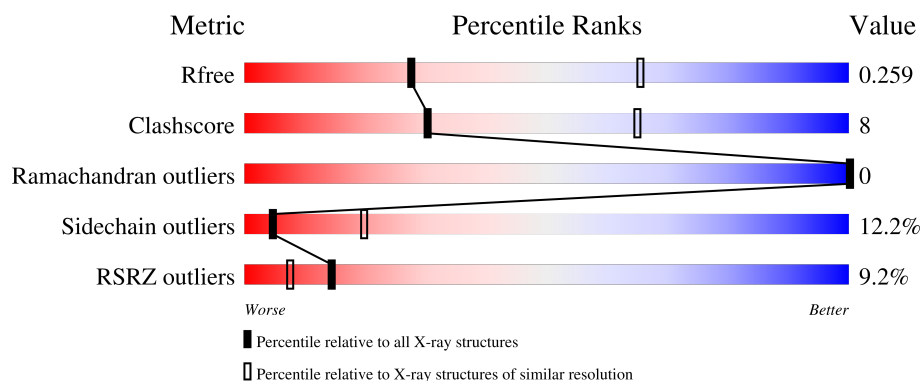
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.01 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3131 (3.04-3.00)
Clashscore	190562	3444 (3.04-3.00)
Ramachandran outliers	187476	3319 (3.04-3.00)
Sidechain outliers	187428	3322 (3.04-3.00)
RSRZ outliers	180081	3130 (3.04-3.00)

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

Mol	Chain	Length	Quality of chain
1	A	281	

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	PO4	A	901	-	X	-	-

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 2268 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.

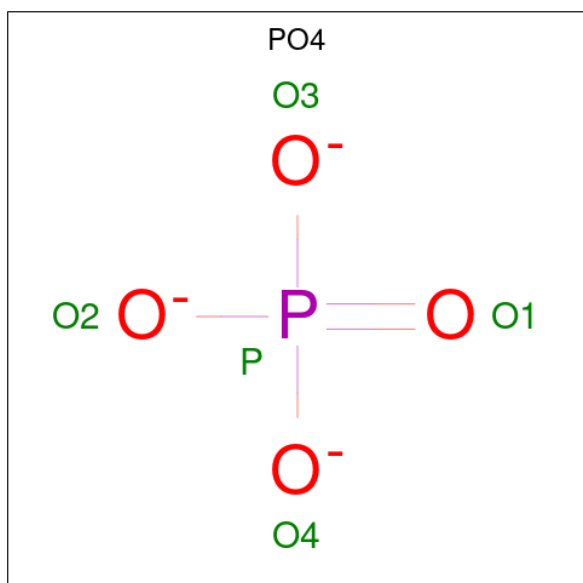
- Molecule 1 is a protein called Protein PML.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	262	2069	1325	368	368	8	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	576	SER	-	expression tag	UNP P29590
A	577	ASN	-	expression tag	UNP P29590
A	578	ALA	-	expression tag	UNP P29590
A	?	-	PHE	deletion	UNP P29590
A	?	-	PHE	deletion	UNP P29590
A	?	-	SER	deletion	UNP P29590
A	?	-	ILE	deletion	UNP P29590
A	?	-	TYR	deletion	UNP P29590

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



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

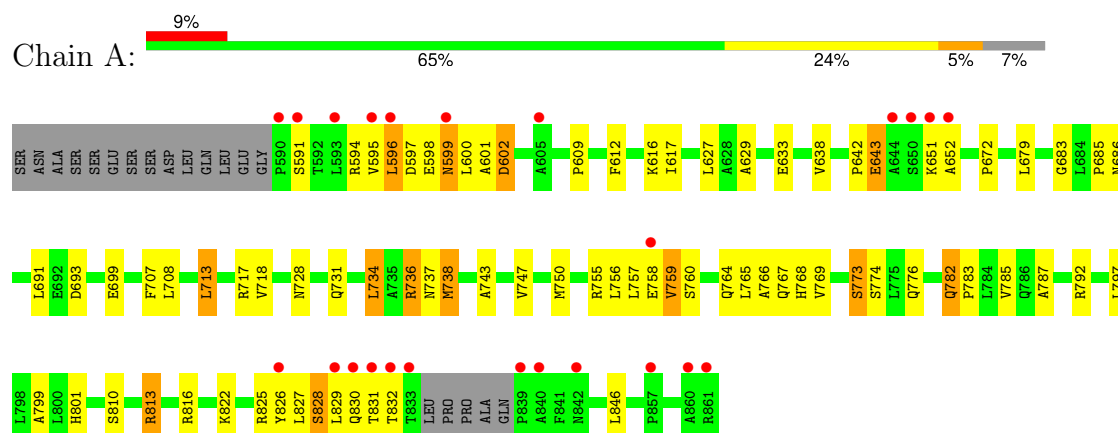
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	194	Total	O	0	0
			194	194		

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: Protein PML



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	114.34Å 114.34Å 102.68Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	21.61 – 3.01 21.61 – 3.01	Depositor EDS
% Data completeness (in resolution range)	61.9 (21.61-3.01) 61.7 (21.61-3.01)	Depositor EDS
R_{merge}	0.46	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.08 (at 3.00Å)	Xtriage
Refinement program	BUSTER 2.10.3	Depositor
R, R_{free}	0.167 , 0.241 0.195 , 0.259	Depositor DCC
R_{free} test set	538 reflections (3.41%)	wwPDB-VP
Wilson B-factor (Å ²)	72.1	Xtriage
Anisotropy	0.120	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 98.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.41$, $\langle L^2 \rangle = 0.23$	Xtriage
Estimated twinning fraction	0.088 for -h,-k,l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	2268	wwPDB-VP
Average B, all atoms (Å ²)	84.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.65% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	0.85	1/2112 (0.0%)	1.36	14/2857 (0.5%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	602	ASP	CA-C	5.29	1.59	1.52

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	828	SER	N-CA-C	-7.71	103.67	113.23
1	A	738	MET	N-CA-C	5.92	117.73	111.28
1	A	773	SER	CA-C-N	5.51	128.12	120.29
1	A	773	SER	C-N-CA	5.51	128.12	120.29
1	A	755	ARG	CA-C-N	5.27	127.29	120.44
1	A	755	ARG	C-N-CA	5.27	127.29	120.44
1	A	691	LEU	CA-C-N	5.26	127.33	120.28
1	A	691	LEU	C-N-CA	5.26	127.33	120.28
1	A	737	ASN	CA-C-N	5.17	127.21	120.28
1	A	737	ASN	C-N-CA	5.17	127.21	120.28
1	A	591	SER	CA-C-N	5.06	130.47	122.78
1	A	591	SER	C-N-CA	5.06	130.47	122.78

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	686	ASN	CA-C-N	5.03	127.02	120.28
1	A	686	ASN	C-N-CA	5.03	127.02	120.28

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	598	GLU	Peptide
1	A	759	VAL	Mainchain

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	2069	0	2114	35	3
2	A	5	0	0	0	0
3	A	194	0	0	0	0
All	All	2268	0	2114	35	3

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 8.

All (35) 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:643:GLU:HG3	1:A:652:ALA:HB1	1.61	0.80
1:A:713:LEU:HD11	1:A:765:LEU:HD22	1.73	0.69
1:A:596:LEU:HB2	1:A:600:LEU:HD22	1.83	0.61
1:A:595:VAL:HG21	1:A:601:ALA:O	2.05	0.57
1:A:731:GLN:HA	1:A:736:ARG:H	1.72	0.55
1:A:597:ASP:C	1:A:599:ASN:HA	2.32	0.54
1:A:717:ARG:HD2	1:A:759:VAL:HG22	1.90	0.54
1:A:717:ARG:NH1	1:A:759:VAL:HA	2.27	0.50
1:A:828:SER:HA	1:A:831:THR:HG23	1.94	0.49
1:A:600:LEU:HA	1:A:769:VAL:O	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:757:LEU:HB2	1:A:759:VAL:HG23	1.95	0.47
1:A:617:ILE:HD11	1:A:683:GLY:HA2	1.97	0.46
1:A:616:LYS:HG2	1:A:743:ALA:HB2	1.98	0.46
1:A:799:ALA:C	1:A:801:HIS:H	2.24	0.46
1:A:785:VAL:HG11	1:A:792:ARG:HG2	1.98	0.45
1:A:785:VAL:C	1:A:787:ALA:H	2.24	0.45
1:A:609:PRO:HG2	1:A:672:PRO:HB3	1.99	0.45
1:A:766:ALA:C	1:A:768:HIS:N	2.75	0.45
1:A:827:LEU:HD22	1:A:830:GLN:HE22	1.81	0.45
1:A:797:LEU:HD23	1:A:829:LEU:HD12	1.98	0.45
1:A:756:LEU:C	1:A:758:GLU:H	2.25	0.44
1:A:595:VAL:HG22	1:A:596:LEU:H	1.82	0.44
1:A:642:PRO:HA	1:A:643:GLU:HA	1.63	0.44
1:A:734:LEU:HB2	1:A:736:ARG:HG3	2.00	0.44
1:A:801:HIS:CG	1:A:826:TYR:CE1	3.05	0.44
1:A:747:VAL:HA	1:A:750:MET:HE2	2.01	0.42
1:A:627:LEU:HB2	1:A:638:VAL:HG12	2.00	0.42
1:A:810:SER:HA	1:A:813:ARG:HB2	2.01	0.42
1:A:685:PRO:HG2	1:A:782:GLN:HG3	2.01	0.41
1:A:612:PHE:O	1:A:629:ALA:HA	2.20	0.41
1:A:782:GLN:N	1:A:783:PRO:HD2	2.35	0.41
1:A:830:GLN:O	1:A:831:THR:C	2.63	0.41
1:A:828:SER:C	1:A:830:GLN:N	2.77	0.41
1:A:599:ASN:ND2	1:A:774:SER:OG	2.54	0.41
1:A:679:LEU:HD11	1:A:707:PHE:HB3	2.03	0.40

All (3) 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:758:GLU:OE2	1:A:822:LYS:CE[2_545]	1.72	0.48
1:A:758:GLU:OE2	1:A:822:LYS:CD[2_545]	2.06	0.14
1:A:758:GLU:OE2	1:A:822:LYS:NZ[2_545]	2.14	0.06

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

of similar resolution.

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	258/281 (92%)	218 (84%)	40 (16%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	221/237 (93%)	194 (88%)	27 (12%)	5	20

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	594	ARG
1	A	596	LEU
1	A	599	ASN
1	A	602	ASP
1	A	633	GLU
1	A	643	GLU
1	A	651	LYS
1	A	693	ASP
1	A	699	GLU
1	A	708	LEU
1	A	713	LEU
1	A	718	VAL
1	A	728	ASN
1	A	734	LEU
1	A	736	ARG
1	A	738	MET
1	A	760	SER
1	A	764	GLN
1	A	767	GLN

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Mol	Chain	Res	Type
1	A	773	SER
1	A	776	GLN
1	A	782	GLN
1	A	813	ARG
1	A	816	ARG
1	A	825	ARG
1	A	832	THR
1	A	846	LEU

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	830	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

1 ligand is 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
2	PO4	A	901	-	4,4,4	2.03	4 (100%)	6,6,6	0.59	0

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	901	PO4	P-O1	2.04	1.55	1.50
2	A	901	PO4	P-O2	2.03	1.60	1.54
2	A	901	PO4	P-O3	2.02	1.60	1.54
2	A	901	PO4	P-O4	2.01	1.60	1.54

There are no bond angle outliers.

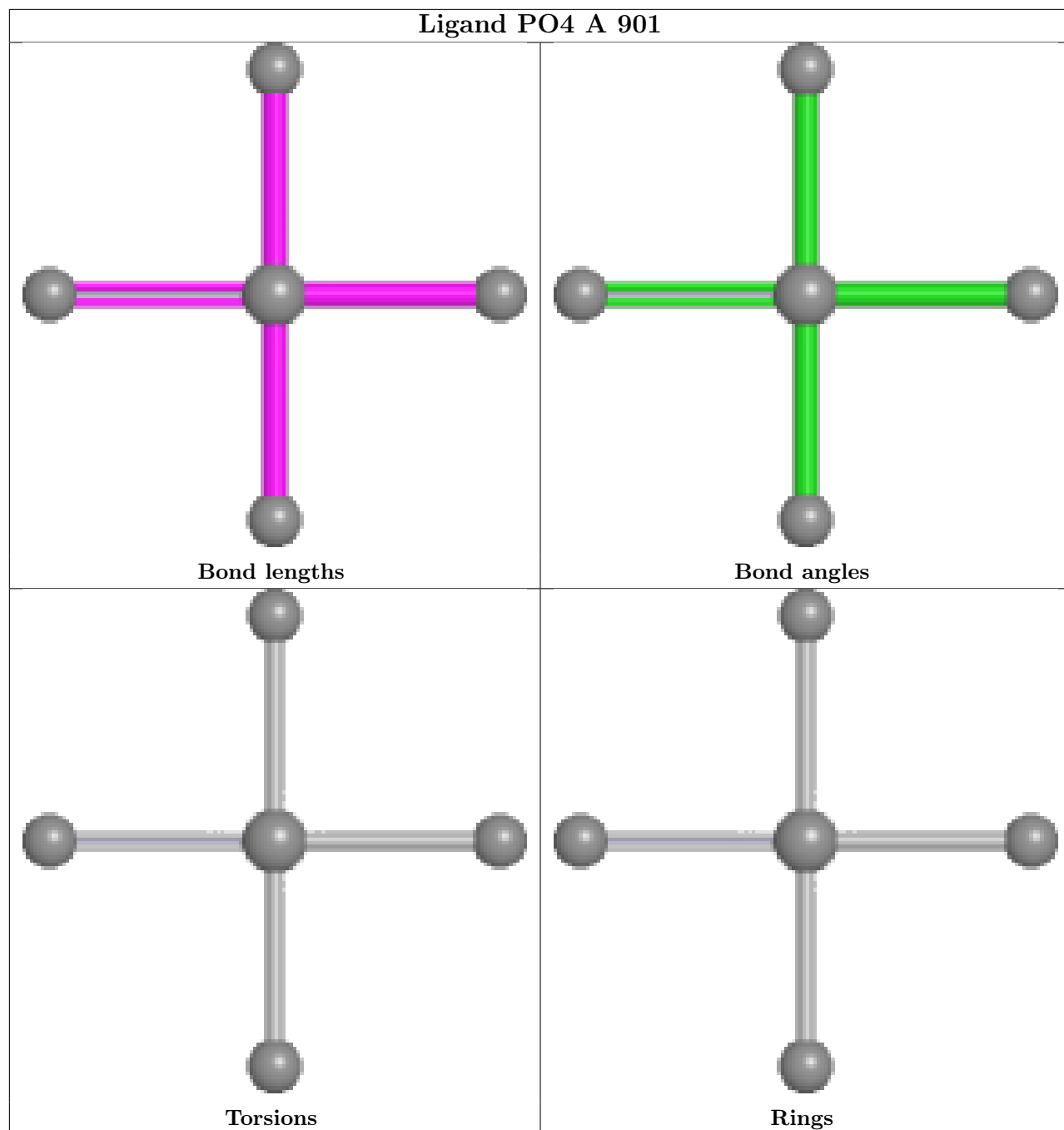
There are no chirality outliers.

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 than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	262/281 (93%)	0.16	24 (9%) 14 7	44, 78, 150, 206	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	595	VAL	5.3
1	A	590	PRO	5.1
1	A	599	ASN	4.8
1	A	758	GLU	3.9
1	A	842	ASN	3.8
1	A	831	THR	3.7
1	A	644	ALA	3.4
1	A	839	PRO	3.2
1	A	591	SER	3.2
1	A	833	THR	3.2
1	A	860	ALA	3.1
1	A	832	THR	2.7
1	A	830	GLN	2.6
1	A	652	ALA	2.6
1	A	829	LEU	2.6
1	A	651	LYS	2.5
1	A	861	ARG	2.5
1	A	650	SER	2.4
1	A	857	PRO	2.4
1	A	593	LEU	2.3
1	A	826	TYR	2.3
1	A	605	ALA	2.3
1	A	840	ALA	2.3
1	A	596	LEU	2.2

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

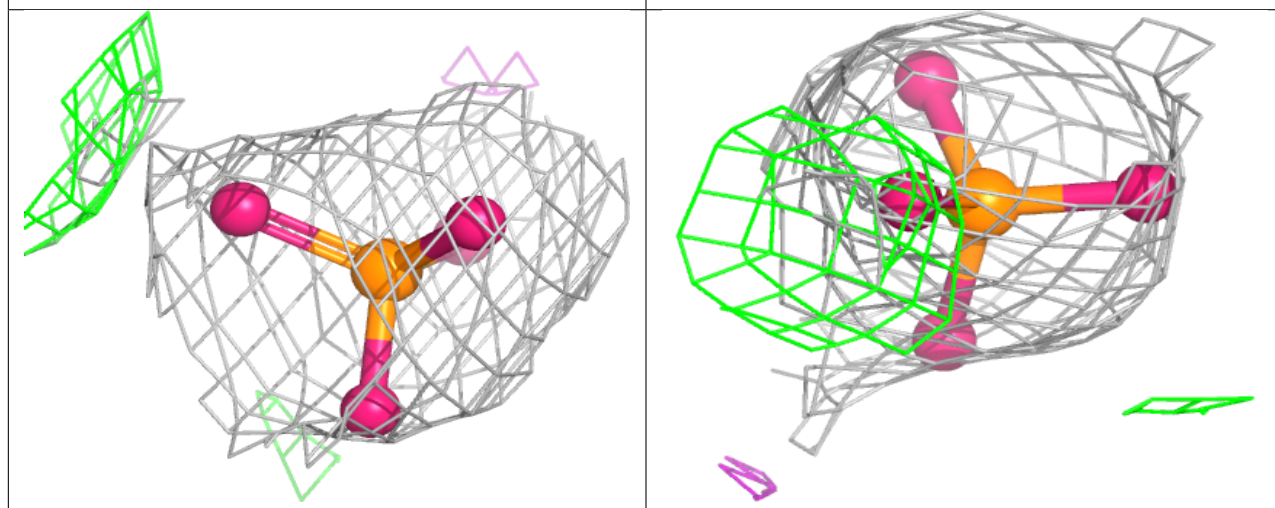
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	PO4	A	901	5/5	0.93	0.16	156,156,157,157	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 901:

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



6.5 Other polymers ⓘ

There are no such residues in this entry.