



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 01:30 AM UTC

PDB ID : 9JZ1 / pdb_00009jz1
Title : Crystal structure of Nir2 C-terminal domain in complex with phosphate
Authors : Kim, D.; Lee, C.
Deposited on : 2024-10-13
Resolution : 2.40 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

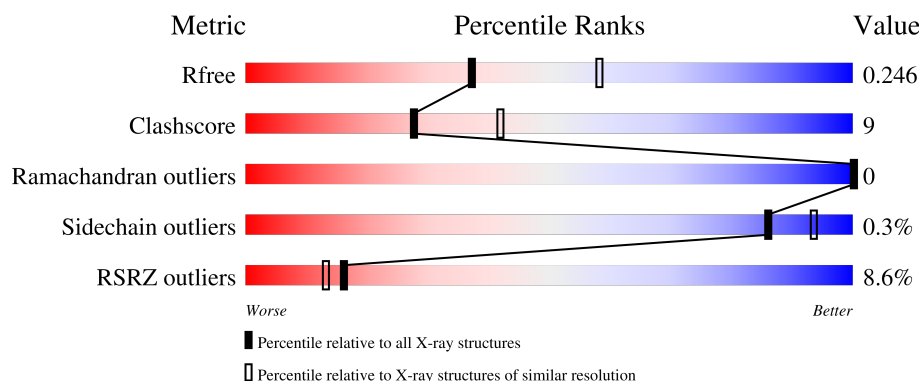
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	284	7% 71% 16% 13%
1	B	284	8% 70% 18% 11%
1	C	284	8% 74% 12% 14%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5858 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 Membrane-associated phosphatidylinositol transfer protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	248	Total	C	N	O	S	0	0	0
			1912	1209	341	352	10			
1	B	253	Total	C	N	O	S	0	0	0
			1948	1233	347	358	10			
1	C	243	Total	C	N	O	S	0	0	0
			1877	1191	332	344	10			

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	907	SER	-	expression tag	UNP O35954
A	908	GLY	-	expression tag	UNP O35954
A	909	SER	-	expression tag	UNP O35954
B	907	SER	-	expression tag	UNP O35954
B	908	GLY	-	expression tag	UNP O35954
B	909	SER	-	expression tag	UNP O35954
C	907	SER	-	expression tag	UNP O35954
C	908	GLY	-	expression tag	UNP O35954
C	909	SER	-	expression tag	UNP O35954

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Ca	0	0
			1	1		

- Molecule 3 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P) (labeled as "Ligand of Interest" by depositor).



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

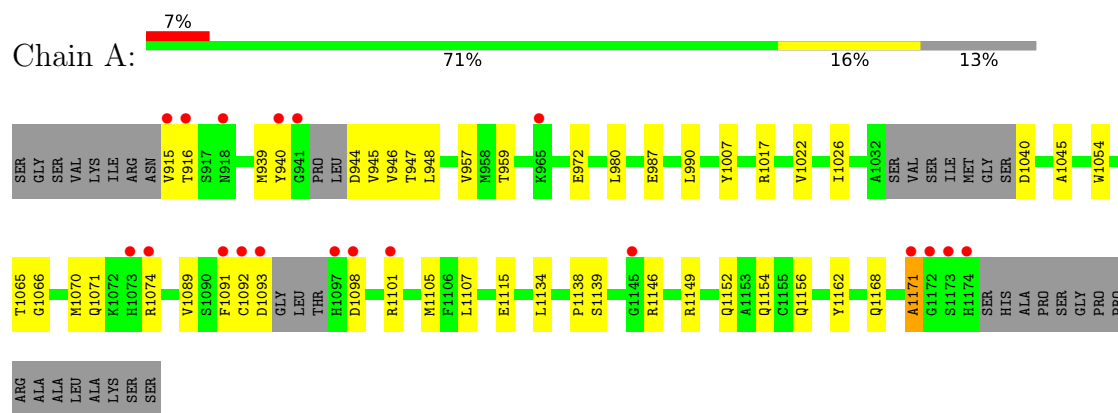
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	46	Total	O	0	0
			46	46		
4	B	39	Total	O	0	0
			39	39		
4	C	30	Total	O	0	0
			30	30		

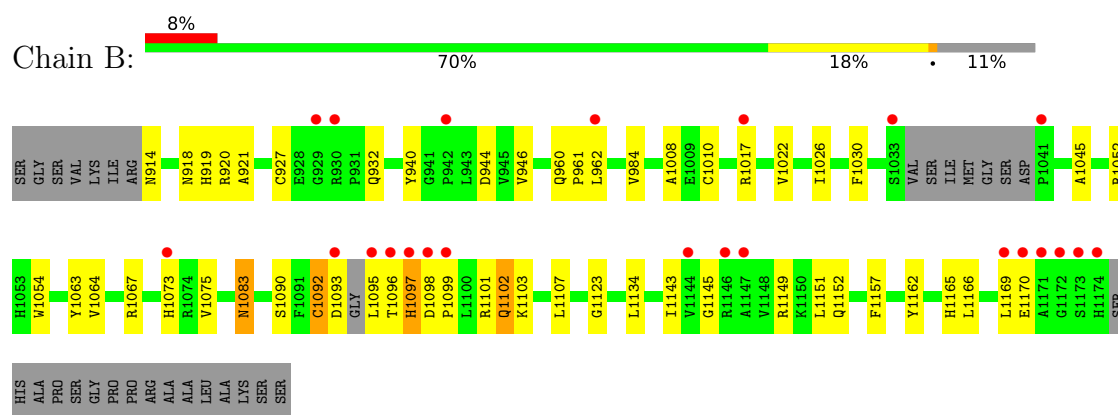
3 Residue-property plots [i](#)

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.

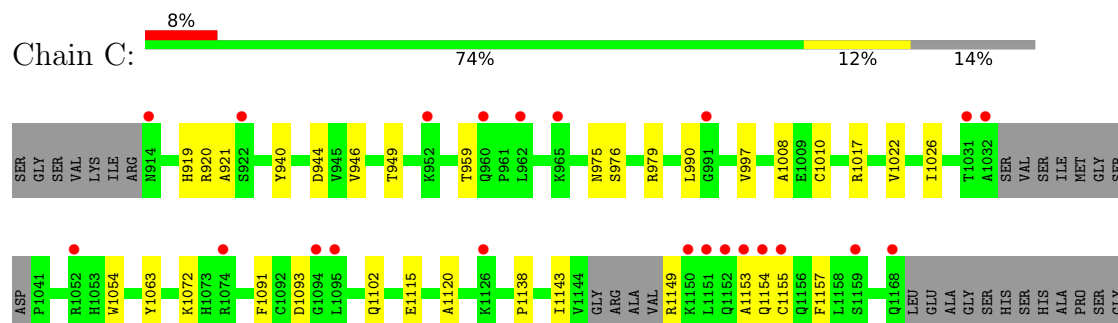
- Molecule 1: Membrane-associated phosphatidylinositol transfer protein 1



- Molecule 1: Membrane-associated phosphatidylinositol transfer protein 1



- Molecule 1: Membrane-associated phosphatidylinositol transfer protein 1



PRO
PRO
ARG
ALA
ALA
LEU
ALA
LYS
SER
SER

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	69.98Å 108.84Å 120.58Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	33.60 – 2.40 33.60 – 2.40	Depositor EDS
% Data completeness (in resolution range)	98.7 (33.60-2.40) 98.6 (33.60-2.40)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.44 (at 2.39Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.211 , 0.260 0.240 , 0.246	Depositor DCC
R_{free} test set	2000 reflections (5.53%)	wwPDB-VP
Wilson B-factor (Å ²)	36.4	Xtriage
Anisotropy	0.150	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 45.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	5858	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.42% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: CA, 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	0.39	0/1954	0.85	3/2654 (0.1%)
1	B	0.47	1/1992 (0.1%)	0.86	7/2708 (0.3%)
1	C	0.45	1/1920 (0.1%)	0.85	2/2611 (0.1%)
All	All	0.44	2/5866 (0.0%)	0.86	12/7973 (0.2%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	1153	ALA	CA-CB	-7.02	1.44	1.53
1	B	1093	ASP	CA-C	-5.38	1.41	1.52

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1171	ALA	N-CA-C	-7.40	97.92	109.25
1	C	1153	ALA	N-CA-C	-7.00	105.55	112.97
1	A	1026	ILE	N-CA-C	5.88	116.35	110.23
1	B	1145	GLY	N-CA-C	5.56	121.09	110.86
1	B	1151	LEU	N-CA-C	5.47	119.58	113.02
1	B	1083	ASN	CB-CA-C	-5.38	104.10	111.73
1	B	1092	CYS	CA-C-N	-5.33	112.11	121.70
1	B	1092	CYS	C-N-CA	-5.33	112.11	121.70
1	A	1007	TYR	N-CA-C	5.27	116.29	108.60
1	B	1030	PHE	N-CA-C	5.25	119.82	113.41
1	C	1102	GLN	N-CA-C	5.08	116.81	111.28
1	B	1102	GLN	N-CA-C	-5.00	105.72	111.07

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	1912	0	1879	38	0
1	B	1948	0	1924	44	0
1	C	1877	0	1855	22	0
2	A	1	0	0	0	0
3	A	5	0	0	0	0
4	A	46	0	0	1	0
4	B	39	0	0	4	0
4	C	30	0	0	0	0
All	All	5858	0	5658	102	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 9.

All (102) 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:B:1095:LEU:HA	1:B:1102:GLN:HE22	1.19	1.04
1:B:1165:HIS:NE2	1:B:1169:LEU:HD11	1.71	1.03
1:B:1165:HIS:CE1	1:B:1169:LEU:HD21	1.95	1.00
1:B:1165:HIS:CD2	1:B:1169:LEU:HD11	1.97	1.00
1:A:1101:ARG:HE	1:A:1105:MET:HE1	1.37	0.89
1:A:1101:ARG:HE	1:A:1105:MET:CE	1.85	0.88
1:A:939:MET:SD	1:A:947:THR:HA	2.14	0.88
1:B:1095:LEU:HA	1:B:1102:GLN:NE2	1.92	0.85
1:A:940:TYR:HE2	1:A:948:LEU:HD13	1.42	0.85
1:A:939:MET:HE1	1:A:947:THR:HG23	1.57	0.84
1:B:918:ASN:OD1	1:B:1073:HIS:CD2	2.33	0.81
1:B:1098:ASP:OD1	1:B:1099:PRO:HD2	1.80	0.81
1:A:940:TYR:CE2	1:A:948:LEU:HD13	2.16	0.80
1:A:1065:THR:HG21	1:A:1091:PHE:CE1	2.19	0.78
1:A:1065:THR:HG21	1:A:1091:PHE:CD1	2.23	0.74
1:A:1065:THR:CG2	1:A:1091:PHE:CD1	2.70	0.73
1:B:1092:CYS:SG	4:B:1235:HOH:O	2.45	0.73
1:B:1165:HIS:O	1:B:1169:LEU:HG	1.89	0.72
1:A:915:VAL:HG23	1:A:916:THR:H	1.54	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:927:CYS:SG	1:B:1017:ARG:HG2	2.31	0.70
1:A:1092:CYS:O	1:A:1093:ASP:C	2.34	0.70
1:A:1101:ARG:NE	1:A:1105:MET:HE1	2.06	0.69
1:B:918:ASN:OD1	1:B:1073:HIS:HD2	1.75	0.68
1:C:1026:ILE:HD11	1:C:1063:TYR:HB3	1.78	0.66
1:C:1138:PRO:HB3	1:C:1154:GLN:O	1.95	0.65
1:A:945:VAL:HG23	1:A:946:VAL:HG23	1.79	0.64
1:B:1166:LEU:O	1:B:1170:GLU:HG3	1.97	0.64
1:C:1138:PRO:CB	1:C:1154:GLN:O	2.45	0.64
1:B:1067:ARG:HH11	1:B:1075:VAL:HG22	1.63	0.63
1:B:1052:ARG:HH12	1:B:1083:ASN:HD22	1.47	0.63
1:A:939:MET:CE	1:A:947:THR:HG23	2.28	0.63
1:A:1089:VAL:HG12	1:A:1091:PHE:CE2	2.34	0.62
1:B:914:ASN:N	4:B:1201:HOH:O	2.33	0.62
1:A:1022:VAL:HG21	1:A:1054:TRP:CE3	2.36	0.61
1:B:1067:ARG:NH2	4:B:1202:HOH:O	2.33	0.60
1:C:1143:ILE:HG12	1:C:1155:CYS:HB2	1.84	0.60
1:A:1071:GLN:HG2	1:A:1074:ARG:HD2	1.84	0.60
1:C:1149:ARG:HA	1:C:1157:PHE:CZ	2.36	0.59
1:C:979:ARG:NH2	1:C:1093:ASP:O	2.35	0.59
1:B:1165:HIS:CD2	1:B:1169:LEU:CD1	2.81	0.59
1:B:1098:ASP:OD1	1:B:1099:PRO:CD	2.49	0.58
1:A:1040:ASP:N	4:A:1302:HOH:O	2.35	0.58
1:B:1097:HIS:O	1:B:1098:ASP:C	2.46	0.58
1:C:1017:ARG:HG3	1:C:1115:GLU:OE1	2.04	0.57
1:B:920:ARG:NH2	1:B:944:ASP:OD2	2.35	0.57
1:C:959:THR:HG22	1:C:990:LEU:HD21	1.88	0.56
1:A:1107:LEU:HB3	1:A:1134:LEU:HD11	1.89	0.55
1:A:939:MET:HE1	1:A:947:THR:CG2	2.34	0.54
1:B:927:CYS:SG	1:B:1017:ARG:CG	2.95	0.54
1:B:1152:GLN:HA	1:B:1157:PHE:HE2	1.72	0.54
1:B:1165:HIS:CE1	1:B:1169:LEU:CD2	2.83	0.53
1:A:1101:ARG:NE	1:A:1105:MET:CE	2.62	0.53
1:B:919:HIS:O	1:B:1073:HIS:CE1	2.62	0.52
1:A:1098:ASP:CG	1:B:962:LEU:HD13	2.34	0.52
1:B:1098:ASP:HB3	1:B:1101:ARG:HD2	1.90	0.52
1:C:921:ALA:HB3	1:C:1010:CYS:HB3	1.93	0.51
1:B:1096:THR:H	1:B:1102:GLN:HE21	1.60	0.50
1:A:944:ASP:OD1	1:A:1070:MET:HG3	2.11	0.50
1:C:920:ARG:HH22	1:C:944:ASP:CG	2.20	0.50
1:A:959:THR:HG22	1:A:990:LEU:HD21	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1022:VAL:HG21	1:C:1054:TRP:CE3	2.46	0.49
1:B:1064:VAL:HG11	1:B:1107:LEU:HD21	1.95	0.49
1:B:940:TYR:HB2	1:B:946:VAL:HB	1.95	0.49
1:A:1098:ASP:OD2	1:B:962:LEU:HD13	2.12	0.49
1:A:1066:GLY:HA2	1:A:1092:CYS:SG	2.52	0.49
1:C:997:VAL:HB	1:C:1010:CYS:O	2.13	0.48
1:C:975:ASN:O	1:C:976:SER:C	2.54	0.48
1:B:1022:VAL:HG21	1:B:1054:TRP:CE3	2.48	0.48
1:A:1045:ALA:O	1:A:1162:TYR:HB2	2.14	0.48
1:B:1103:LYS:NZ	4:B:1205:HOH:O	2.48	0.47
1:B:1123:GLY:O	1:B:1143:ILE:HG23	2.14	0.47
1:A:1089:VAL:CG1	1:A:1091:PHE:CE2	2.99	0.46
1:A:1138:PRO:HB3	1:A:1154:GLN:C	2.41	0.46
1:A:1017:ARG:HG3	1:A:1115:GLU:OE1	2.17	0.45
1:B:1165:HIS:HE1	1:B:1169:LEU:HD21	1.73	0.45
1:A:957:VAL:HG11	1:A:990:LEU:HD11	1.99	0.45
1:C:1138:PRO:CA	1:C:1154:GLN:O	2.64	0.45
1:C:940:TYR:HB2	1:C:946:VAL:HB	1.97	0.45
1:B:1064:VAL:HG12	1:B:1090:SER:OG	2.17	0.44
1:A:1168:GLN:O	1:A:1171:ALA:O	2.34	0.44
1:B:1045:ALA:O	1:B:1162:TYR:HB2	2.18	0.44
1:C:949:THR:HG22	1:C:976:SER:O	2.18	0.44
1:B:921:ALA:HB3	1:B:1010:CYS:HB3	1.99	0.43
1:B:1107:LEU:HB3	1:B:1134:LEU:HD11	1.99	0.43
1:A:972:GLU:HB2	1:A:980:LEU:HD11	2.00	0.43
1:A:1101:ARG:HG3	1:A:1105:MET:HE3	2.00	0.43
1:B:932:GLN:HB2	1:B:984:VAL:HB	2.01	0.43
1:C:1022:VAL:HG13	1:C:1120:ALA:HB3	1.99	0.43
1:A:1149:ARG:O	1:A:1152:GLN:HB2	2.19	0.42
1:C:920:ARG:NH2	1:C:944:ASP:OD1	2.52	0.42
1:B:1165:HIS:NE2	1:B:1169:LEU:CD1	2.62	0.42
1:C:1143:ILE:HG13	1:C:1155:CYS:SG	2.60	0.42
1:A:1139:SER:O	1:A:1156:GLN:NE2	2.52	0.42
1:B:919:HIS:HB3	1:B:1008:ALA:HB2	2.02	0.42
1:C:1026:ILE:HD11	1:C:1063:TYR:CD1	2.55	0.42
1:B:1026:ILE:HD11	1:B:1063:TYR:HB3	2.02	0.41
1:B:1149:ARG:O	1:B:1152:GLN:HG2	2.20	0.41
1:C:1072:LYS:HA	1:C:1091:PHE:CZ	2.55	0.41
1:A:945:VAL:HG23	1:A:946:VAL:N	2.34	0.41
1:A:987:GLU:CD	1:A:987:GLU:H	2.29	0.41
1:B:960:GLN:O	1:B:961:PRO:C	2.64	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:919:HIS:HB3	1:C:1008:ALA:HB2	2.02	0.40

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	A	240/284 (84%)	239 (100%)	1 (0%)	0	100	100
1	B	247/284 (87%)	245 (99%)	2 (1%)	0	100	100
1	C	237/284 (84%)	235 (99%)	2 (1%)	0	100	100
All	All	724/852 (85%)	719 (99%)	5 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	208/236 (88%)	207 (100%)	1 (0%)	81	91
1	B	213/236 (90%)	212 (100%)	1 (0%)	81	91
1	C	206/236 (87%)	206 (100%)	0	100	100
All	All	627/708 (89%)	625 (100%)	2 (0%)	86	93

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1146	ARG
1	B	1097	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1156	GLN
1	B	1073	HIS
1	B	1083	ASN
1	B	1102	GLN
1	C	1071	GLN
1	C	1083	ASN
1	C	1086	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	PO4	A	1202	-	4,4,4	0.97	0	6,6,6	0.36	0

There are no bond length outliers.

There are no bond angle outliers.

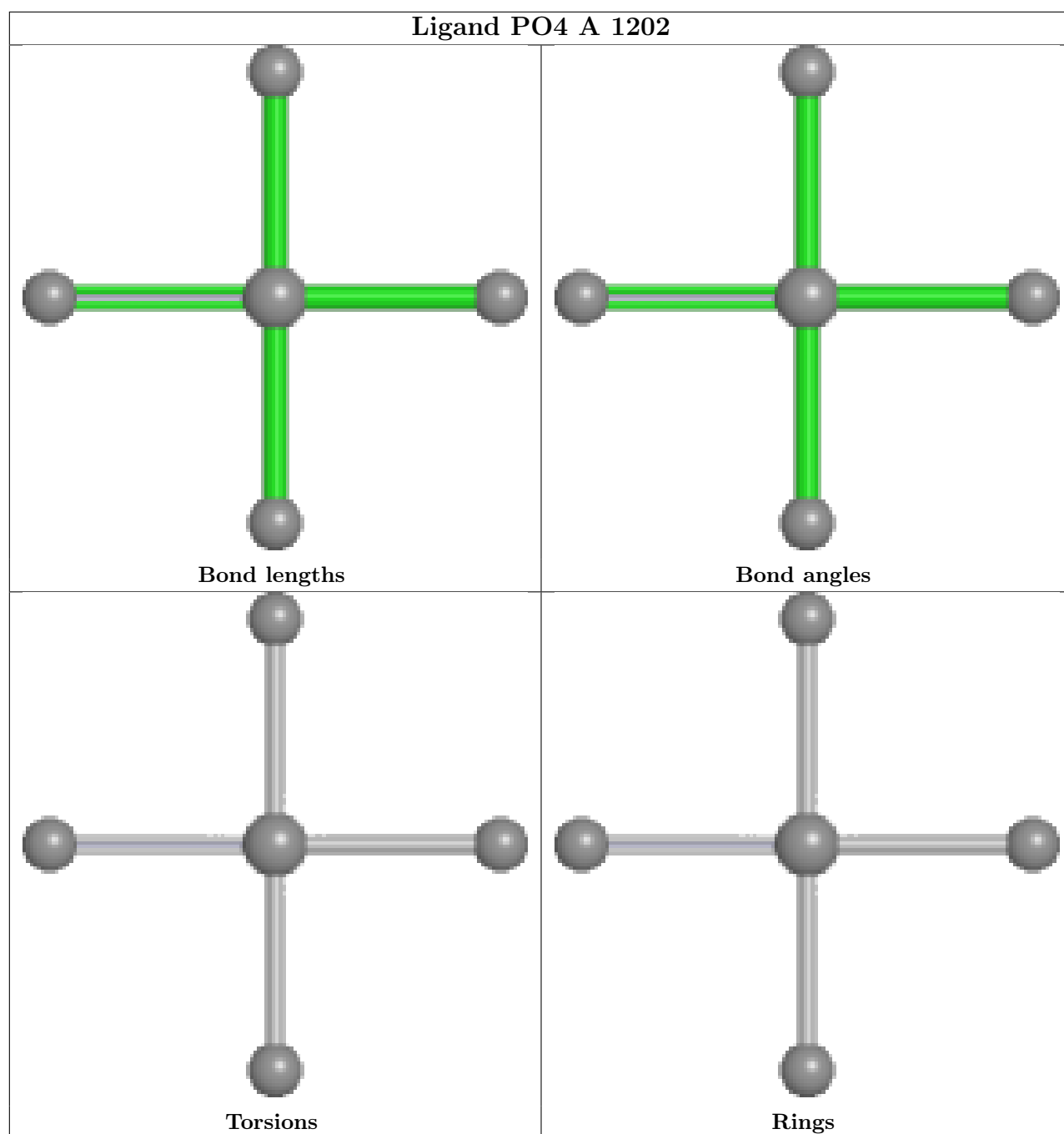
There are no chirality outliers.

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 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	248/284 (87%)	0.25	19 (7%)	19 16	26, 44, 77, 100	0
1	B	253/284 (89%)	0.33	23 (9%)	15 12	25, 44, 76, 92	0
1	C	243/284 (85%)	0.56	22 (9%)	15 12	30, 55, 80, 101	0
All	All	744/852 (87%)	0.38	64 (8%)	16 13	25, 48, 78, 101	0

All (64) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	941	GLY	4.8
1	A	1093	ASP	4.6
1	B	1095	LEU	4.6
1	C	1074	ARG	4.6
1	A	1073	HIS	4.4
1	C	1153	ALA	4.4
1	B	1073	HIS	4.2
1	B	1098	ASP	4.2
1	B	1096	THR	4.0
1	A	1101	ARG	3.8
1	A	1173	SER	3.7
1	A	1172	GLY	3.7
1	A	915	VAL	3.6
1	A	1171	ALA	3.6
1	C	1151	LEU	3.5
1	B	942	PRO	3.5
1	C	1155	CYS	3.5
1	B	1033	SER	3.5
1	B	962	LEU	3.4
1	C	914	ASN	3.3
1	C	1095	LEU	3.3
1	B	1097	HIS	3.3
1	C	1052	ARG	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	1074	ARG	3.1
1	B	1174	HIS	3.0
1	B	1169	LEU	3.0
1	B	1093	ASP	3.0
1	C	1150	LYS	3.0
1	B	929	GLY	3.0
1	C	960	GLN	3.0
1	B	930	ARG	2.9
1	C	1152	GLN	2.9
1	B	1041	PRO	2.9
1	C	1094	GLY	2.7
1	B	1144	VAL	2.6
1	B	1170	GLU	2.6
1	B	1146	ARG	2.6
1	C	952	LYS	2.6
1	C	1168	GLN	2.5
1	A	1097	HIS	2.5
1	C	1126	LYS	2.5
1	C	962	LEU	2.5
1	A	965	LYS	2.4
1	B	1017	ARG	2.4
1	B	1173	SER	2.4
1	C	965	LYS	2.4
1	C	1154	GLN	2.3
1	A	1091	PHE	2.3
1	A	1145	GLY	2.3
1	C	1032	ALA	2.2
1	B	1172	GLY	2.2
1	B	1171	ALA	2.1
1	C	1031	THR	2.1
1	A	940	TYR	2.1
1	A	1092	CYS	2.1
1	A	918	ASN	2.1
1	B	1099	PRO	2.1
1	A	916	THR	2.1
1	C	1159	SER	2.1
1	B	1147	ALA	2.0
1	C	991	GLY	2.0
1	A	1098	ASP	2.0
1	A	1174	HIS	2.0
1	C	922	SER	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

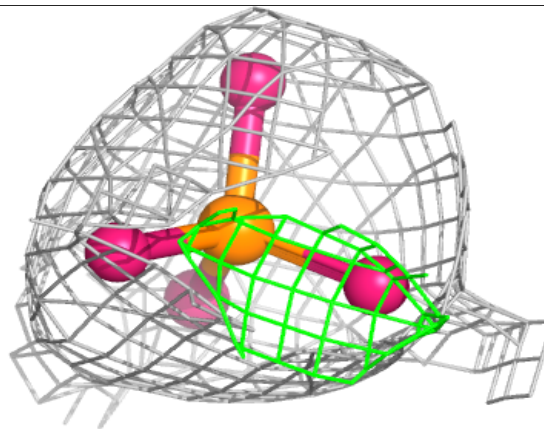
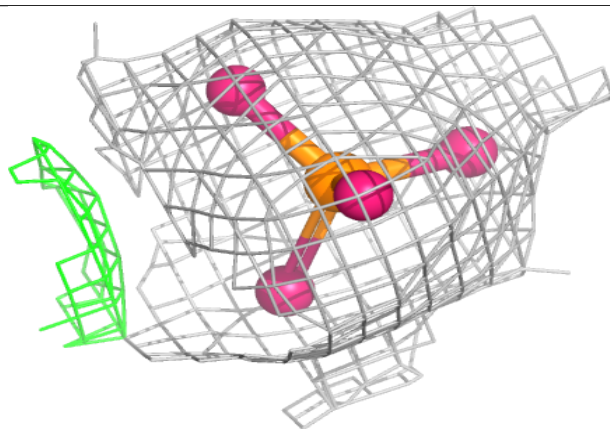
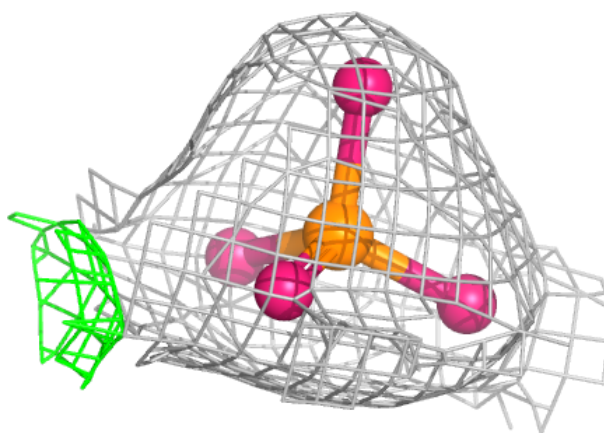
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

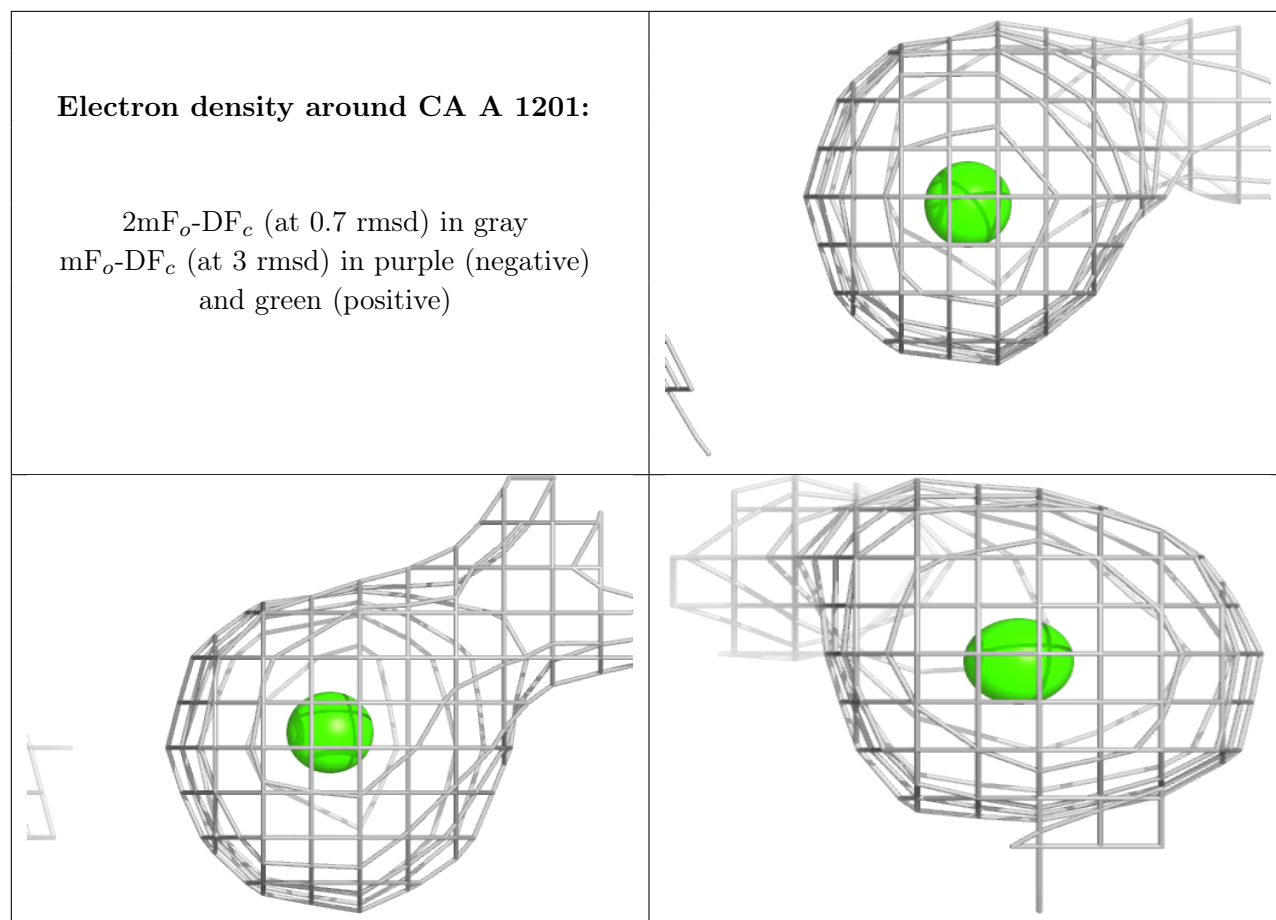
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	PO4	A	1202	5/5	0.95	0.07	61,61,63,64	0
2	CA	A	1201	1/1	0.96	0.09	80,80,80,80	0

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 1202:

$2mF_o-DF_c$ (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.