



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 05:32 AM UTC

PDB ID : 5G4I / pdb_00005g4i
Title : PLP-dependent phospholyase A1RDF1 from *Arthrobacter aurescens* TC1
Authors : Cuetos, A.; Tuan, A.N.; Mangas Sanchez, J.; Grogan, G.
Deposited on : 2016-05-13
Resolution : 1.50 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

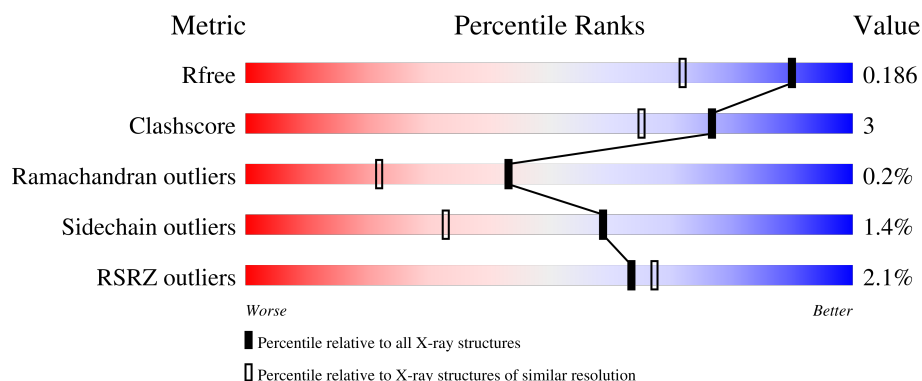
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

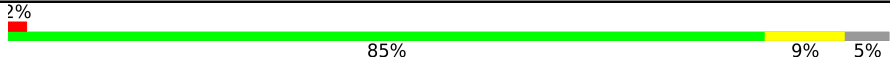
The reported resolution of this entry is 1.50 Å.

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	4037 (1.50-1.50)
Clashscore	190562	4235 (1.50-1.50)
Ramachandran outliers	187476	4153 (1.50-1.50)
Sidechain outliers	187428	4150 (1.50-1.50)
RSRZ outliers	180081	4039 (1.50-1.50)

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	446	
1	B	446	

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
3	PO4	A	1442	-	X	-	-
3	PO4	B	1442	-	X	-	-
4	ACT	A	1443	-	-	X	-
4	ACT	B	1445	-	-	X	-

2 Entry composition [i](#)

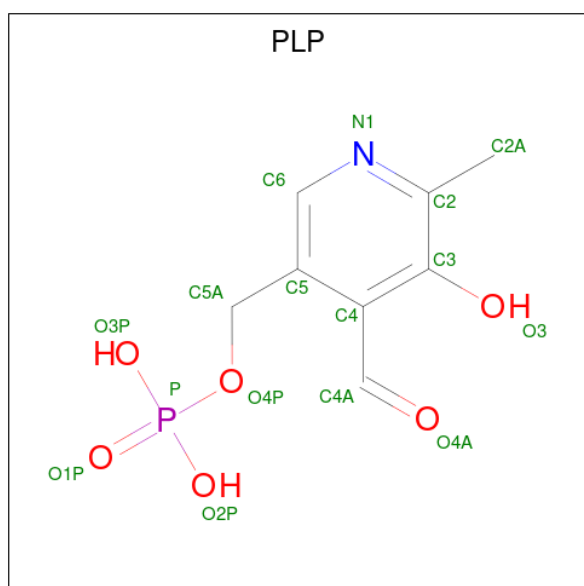
There are 6 unique types of molecules in this entry. The entry contains 7107 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 PHOSPHOLYASE.

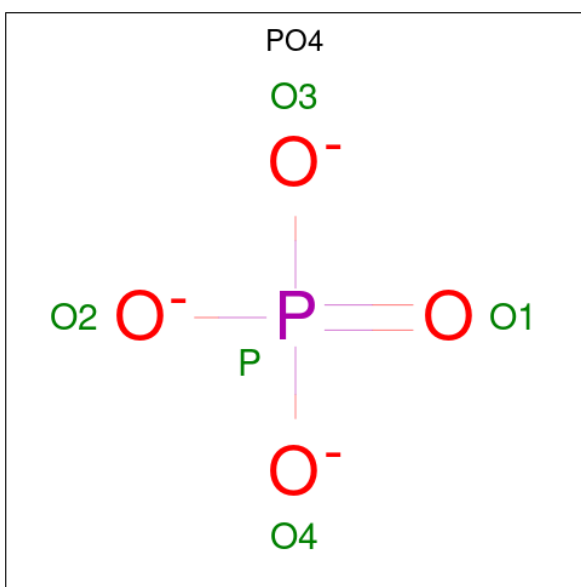
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	423	Total	C	N	O	S	0	5	0
			3198	2031	554	604	9			
1	B	422	Total	C	N	O	S	0	1	0
			3158	2006	550	593	9			

- Molecule 2 is PYRIDOXAL-5'-PHOSPHATE (CCD ID: PLP) (formula: $C_8H_{10}NO_6P$).



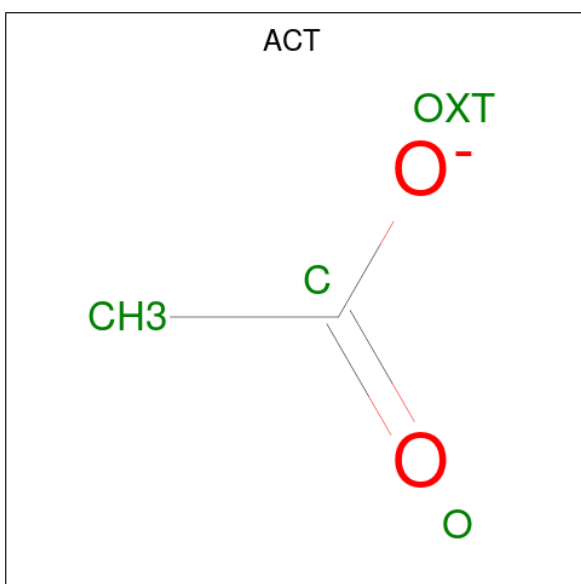
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			15	8	1	5	1		
2	B	1	Total	C	N	O	P	0	0
			15	8	1	5	1		

- Molecule 3 is PHOSPHATE ION (CCD ID: PO4) (formula: O_4P).



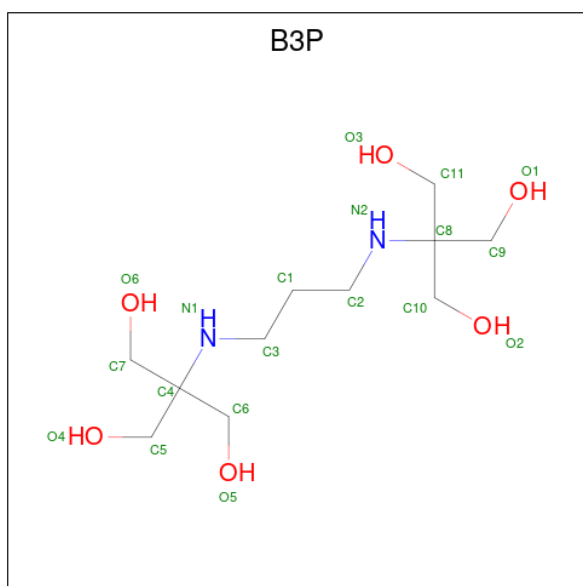
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	P	0	0
			5	4	1		
3	B	1	Total	O	P	0	0
			5	4	1		
3	B	1	Total	O	P	0	0
			5	4	1		
3	B	1	Total	O	P	0	0
			5	4	1		
3	B	1	Total	O	P	0	0
			5	4	1		

- Molecule 4 is ACETATE ION (CCD ID: ACT) (formula: $C_2H_3O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		
4	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 5 is 2-[3-(2-HYDROXY-1,1-DIHYDROXYMETHYL-ETHYLAMINO)-PROPYLAMINO]-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (CCD ID: B3P) (formula: $C_{11}H_{26}N_2O_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			19	11	2	6		

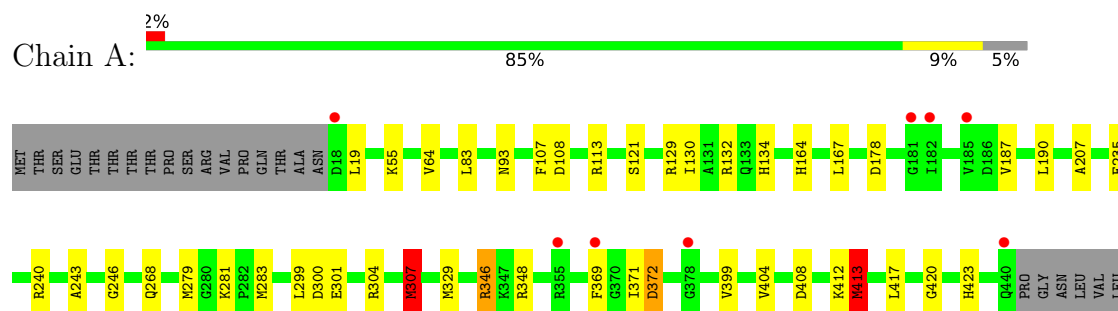
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	415	Total	O	0	0
			415	415		
6	B	250	Total	O	0	0
			250	250		

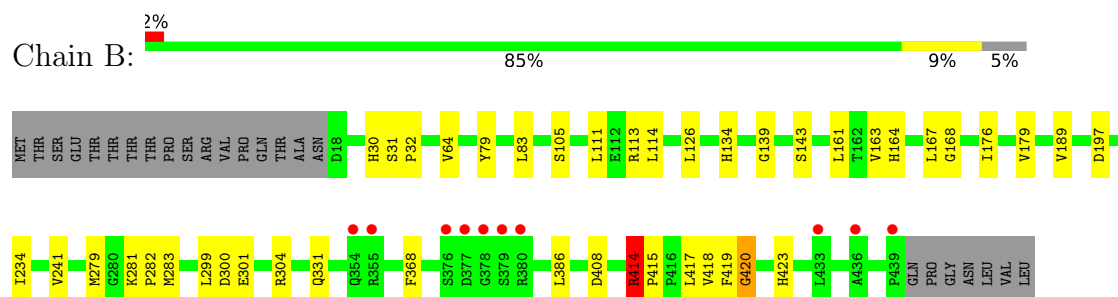
3 Residue-property plots

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: PHOSPHOLYASE



• Molecule 1: PHOSPHOLYASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	76.94Å 97.18Å 124.26Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	76.55 – 1.50 76.55 – 1.50	Depositor EDS
% Data completeness (in resolution range)	99.3 (76.55-1.50) 99.3 (76.55-1.50)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.00 (at 1.50Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.153 , 0.175 0.166 , 0.186	Depositor DCC
R_{free} test set	7219 reflections (4.87%)	wwPDB-VP
Wilson B-factor (Å ²)	14.0	Xtriage
Anisotropy	0.650	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 37.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	7107	wwPDB-VP
Average B, all atoms (Å ²)	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.03% 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, ACT, PLP, B3P

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	1.72	32/3278 (1.0%)	1.35	15/4458 (0.3%)
1	B	1.62	15/3225 (0.5%)	1.31	12/4388 (0.3%)
All	All	1.67	47/6503 (0.7%)	1.33	27/8846 (0.3%)

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	1

All (47) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	371	ILE	C-N	-17.67	1.08	1.33
1	A	279	MET	SD-CE	-12.80	1.47	1.79
1	B	279	MET	SD-CE	-9.45	1.55	1.79
1	B	420	GLY	N-CA	9.08	1.51	1.44
1	A	167	LEU	C-O	7.10	1.32	1.23
1	B	168	GLY	CA-C	7.06	1.58	1.52
1	B	64	VAL	CA-CB	6.83	1.58	1.54
1	A	132	ARG	N-CA	6.75	1.54	1.46
1	A	55	LYS	C-O	6.61	1.31	1.23
1	A	235	GLU	CD-OE1	-6.36	1.13	1.25
1	A	329	MET	CA-CB	6.36	1.63	1.53
1	A	83	LEU	N-CA	6.28	1.54	1.46
1	A	108	ASP	CB-CG	-6.28	1.36	1.52
1	B	415	PRO	CA-C	6.24	1.57	1.52
1	A	64	VAL	CA-CB	6.18	1.57	1.54
1	B	418	VAL	N-CA	6.06	1.53	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	197	ASP	CA-C	6.04	1.60	1.52
1	A	129	ARG	C-O	5.88	1.30	1.24
1	B	179	VAL	N-CA	5.86	1.54	1.46
1	A	246	GLY	N-CA	5.85	1.49	1.44
1	A	121	SER	N-CA	5.81	1.53	1.46
1	A	399	VAL	N-CA	5.77	1.53	1.46
1	A	413	MET	CA-CB	-5.76	1.44	1.53
1	A	283	MET	SD-CE	-5.67	1.65	1.79
1	A	132	ARG	C-O	5.67	1.30	1.24
1	A	107	PHE	C-O	-5.66	1.16	1.23
1	B	126	LEU	CA-C	5.66	1.60	1.52
1	A	207	ALA	N-CA	5.62	1.53	1.46
1	A	178	ASP	CA-C	5.55	1.58	1.52
1	A	190	LEU	C-O	5.55	1.30	1.24
1	A	187	VAL	CA-C	5.53	1.58	1.52
1	A	130	ILE	N-CA	5.50	1.52	1.46
1	B	163	VAL	N-CA	5.48	1.52	1.46
1	A	283	MET	CB-CG	-5.46	1.36	1.52
1	A	19	LEU	N-CA	5.27	1.52	1.46
1	B	167	LEU	C-O	5.24	1.30	1.23
1	B	241	VAL	N-CA	5.24	1.52	1.46
1	A	372	ASP	CB-CG	-5.22	1.39	1.52
1	A	93	ASN	CG-OD1	5.20	1.33	1.23
1	A	108	ASP	C-O	-5.19	1.16	1.23
1	A	240	ARG	N-CA	5.17	1.52	1.46
1	B	114	LEU	N-CA	5.13	1.52	1.46
1	A	268	GLN	CD-OE1	-5.11	1.13	1.23
1	B	105	SER	N-CA	5.11	1.52	1.46
1	B	111	LEU	N-CA	5.11	1.53	1.46
1	A	346	ARG	CZ-NH1	5.09	1.39	1.32
1	A	243	ALA	N-CA	5.08	1.52	1.46

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	414	ARG	NE-CZ-NH1	-12.99	108.51	121.50
1	B	414	ARG	NE-CZ-NH2	12.92	130.82	119.20
1	A	346	ARG	NE-CZ-NH2	-11.79	108.59	119.20
1	A	372	ASP	CB-CA-C	-10.58	92.94	110.29
1	A	413	MET	CG-SD-CE	10.56	124.14	100.90
1	A	346	ARG	NE-CZ-NH1	10.38	131.88	121.50
1	A	371	ILE	CA-C-N	10.12	136.39	121.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	371	ILE	C-N-CA	10.12	136.39	121.72
1	A	108	ASP	CB-CA-C	-8.73	93.58	111.22
1	B	414	ARG	CB-CG-CD	8.28	130.34	111.30
1	A	346	ARG	CD-NE-CZ	7.79	135.30	124.40
1	A	235	GLU	CG-CD-OE1	7.73	136.18	118.40
1	B	414	ARG	CD-NE-CZ	-7.62	113.74	124.40
1	A	235	GLU	OE1-CD-OE2	-7.42	105.08	122.90
1	A	235	GLU	CB-CG-CD	7.31	125.03	112.60
1	B	282	PRO	N-CA-C	6.84	122.74	113.84
1	A	348	ARG	CG-CD-NE	-6.77	97.10	112.00
1	A	372	ASP	CA-CB-CG	6.40	119.00	112.60
1	A	283	MET	N-CA-C	5.99	118.59	111.71
1	B	414	ARG	CG-CD-NE	5.93	125.05	112.00
1	B	283	MET	CG-SD-CE	-5.86	88.02	100.90
1	B	176	ILE	N-CA-C	-5.71	103.72	108.63
1	B	161	LEU	N-CA-CB	-5.67	102.19	111.66
1	A	307	MET	CG-SD-CE	-5.64	88.49	100.90
1	B	283	MET	N-CA-CB	-5.61	101.64	110.06
1	B	139	GLY	N-CA-C	5.52	117.82	111.36
1	B	419	PHE	CA-CB-CG	5.21	119.01	113.80

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	372	ASP	Sidechain

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	3198	0	3160	14	0
1	B	3158	0	3113	21	0
2	A	15	0	6	0	0
2	B	15	0	7	0	0
3	A	5	0	0	0	0
3	B	20	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	8	0	6	4	0
4	B	4	0	3	2	0
5	A	19	0	26	0	0
6	A	415	0	0	4	0
6	B	250	0	0	4	0
All	All	7107	0	6321	37	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 3.

All (37) 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:414:ARG:HG2	1:B:414:ARG:HH11	1.11	1.13
1:B:414:ARG:HH11	1:B:414:ARG:CG	1.53	1.10
1:B:414:ARG:CG	1:B:414:ARG:NH1	2.31	0.88
1:A:300[A]:ASP:OD2	1:A:304:ARG:NH1	2.10	0.85
1:B:300:ASP:OD2	1:B:304:ARG:NH1	2.13	0.81
1:B:368:PHE:CE2	1:B:414:ARG:HD3	2.15	0.81
1:B:414:ARG:HG2	1:B:414:ARG:NH1	1.92	0.80
1:A:420:GLY:H	1:A:423:HIS:HD2	1.30	0.80
1:B:420:GLY:H	1:B:423:HIS:HD2	1.30	0.76
1:A:369[B]:PHE:CE2	1:A:413:MET:HE3	2.28	0.68
1:A:134:HIS:HD2	1:A:301:GLU:OE1	1.76	0.68
1:A:164:HIS:HD2	4:A:1443:ACT:OXT	1.75	0.67
1:B:134:HIS:HD2	1:B:301:GLU:OE1	1.78	0.67
1:B:164:HIS:HD2	4:B:1445:ACT:O	1.81	0.64
1:B:414:ARG:HD2	6:B:2033:HOH:O	1.98	0.64
1:A:307:MET:HE1	1:B:32:PRO:HG3	1.80	0.63
1:B:420:GLY:H	1:B:423:HIS:CD2	2.15	0.62
1:A:420:GLY:H	1:A:423:HIS:CD2	2.16	0.60
1:A:369[B]:PHE:HE2	1:A:413:MET:HE3	1.65	0.59
1:A:404:VAL:HG13	1:A:412:LYS:HE3	1.87	0.56
1:B:164:HIS:CD2	4:B:1445:ACT:O	2.58	0.56
1:A:300[B]:ASP:CG	6:A:2154:HOH:O	2.49	0.54
4:A:1444:ACT:H1	6:A:2354:HOH:O	2.07	0.54
4:A:1443:ACT:H3	6:A:2231:HOH:O	2.07	0.54
1:B:79:TYR:CZ	1:B:83:LEU:HD11	2.46	0.52
1:A:346:ARG:HD2	6:A:2371:HOH:O	2.09	0.51
1:B:234:ILE:HB	3:B:1442:PO4:O1	2.11	0.49
1:B:368:PHE:CD2	1:B:414:ARG:HD3	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:412:LYS:C	1:A:413:MET:HG3	2.36	0.49
1:B:113:ARG:HB3	1:B:299:LEU:HD13	1.95	0.47
1:A:113:ARG:HB3	1:A:299:LEU:HD13	1.98	0.46
1:B:331:GLN:NE2	6:B:2222:HOH:O	2.50	0.44
1:A:369[B]:PHE:CD2	1:A:413:MET:HE3	2.55	0.42
1:B:189:VAL:HG23	6:B:2144:HOH:O	2.20	0.41
4:A:1443:ACT:H2	6:B:2034:HOH:O	2.21	0.41
1:B:30:HIS:O	1:B:31:SER:C	2.63	0.40
1:B:79:TYR:CE2	1:B:83:LEU:HD11	2.56	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	426/446 (96%)	412 (97%)	13 (3%)	1 (0%)	43	22
1	B	421/446 (94%)	410 (97%)	10 (2%)	1 (0%)	43	22
All	All	847/892 (95%)	822 (97%)	23 (3%)	2 (0%)	43	22

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	281	LYS
1	B	281	LYS

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	335/361 (93%)	331 (99%)	4 (1%)	63	38
1	B	326/361 (90%)	321 (98%)	5 (2%)	57	31
All	All	661/722 (92%)	652 (99%)	9 (1%)	59	33

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

Mol	Chain	Res	Type
1	A	307	MET
1	A	408	ASP
1	A	413	MET
1	A	417	LEU
1	B	143	SER
1	B	386	LEU
1	B	408	ASP
1	B	414	ARG
1	B	417	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	93	ASN
1	A	134	HIS
1	A	137	ASN
1	A	164	HIS
1	A	331	GLN
1	A	423	HIS
1	B	93	ASN
1	B	134	HIS
1	B	164	HIS
1	B	254	GLN
1	B	287	HIS
1	B	331	GLN
1	B	407	HIS
1	B	423	HIS

5.3.3 RNA ⓘ

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](#)

11 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	PO4	B	1442	-	4,4,4	1.76	1 (25%)	6,6,6	2.13	3 (50%)
4	ACT	A	1444	-	3,3,3	0.90	0	3,3,3	2.15	2 (66%)
4	ACT	B	1445	-	3,3,3	0.50	0	3,3,3	1.88	1 (33%)
3	PO4	A	1442	-	4,4,4	2.03	1 (25%)	6,6,6	2.71	3 (50%)
3	PO4	B	1443	-	4,4,4	1.09	0	6,6,6	1.05	1 (16%)
3	PO4	B	1441	-	4,4,4	1.64	1 (25%)	6,6,6	1.51	2 (33%)
2	PLP	B	1440	1	15,15,16	3.08	8 (53%)	21,22,23	2.04	7 (33%)
4	ACT	A	1443	-	3,3,3	1.55	0	3,3,3	2.19	1 (33%)
3	PO4	B	1444	-	4,4,4	0.72	0	6,6,6	0.78	0
5	B3P	A	1445	-	18,18,18	2.60	6 (33%)	23,23,23	1.70	5 (21%)
2	PLP	A	1441	1	15,15,16	3.03	3 (20%)	21,22,23	1.18	1 (4%)

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
2	PLP	B	1440	1	-	0/6/6/8	0/1/1/1
5	B3P	A	1445	-	-	0/28/28/28	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	PLP	A	1441	1	-	1/6/6/8	0/1/1/1

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1440	PLP	C3-C2	8.50	1.49	1.41
2	A	1441	PLP	C3-C2	7.64	1.48	1.41
2	A	1441	PLP	C5-C4	7.26	1.48	1.40
5	A	1445	B3P	C9-C8	7.09	1.61	1.53
2	A	1441	PLP	C3-C4	4.37	1.48	1.40
2	B	1440	PLP	C5-C4	4.32	1.45	1.40
5	A	1445	B3P	C3-N1	4.00	1.51	1.46
2	B	1440	PLP	C3-C4	3.99	1.47	1.40
3	A	1442	PO4	P-O1	-3.79	1.41	1.50
5	A	1445	B3P	C5-C4	-3.74	1.49	1.53
5	A	1445	B3P	C7-C4	3.36	1.57	1.53
3	B	1442	PO4	P-O3	-3.12	1.45	1.54
2	B	1440	PLP	C6-C5	3.12	1.43	1.37
5	A	1445	B3P	O5-C6	3.11	1.52	1.42
3	B	1441	PO4	P-O2	-3.11	1.45	1.54
5	A	1445	B3P	O2-C10	2.69	1.50	1.42
2	B	1440	PLP	C2-N1	-2.59	1.29	1.33
2	B	1440	PLP	C2A-C2	2.15	1.53	1.50
2	B	1440	PLP	C5A-C5	2.12	1.56	1.50
2	B	1440	PLP	P-O3P	-2.02	1.47	1.54

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1442	PO4	O3-P-O1	-4.43	95.27	110.95
2	B	1440	PLP	C3-C4-C5	-4.27	113.47	118.59
2	B	1440	PLP	O2P-P-O4P	-4.05	96.10	106.67
3	A	1442	PO4	O3-P-O2	-3.45	97.18	107.91
3	B	1442	PO4	O3-P-O1	-3.44	98.79	110.95
4	A	1443	ACT	OXT-C-O	-3.26	109.95	122.03
5	A	1445	B3P	O3-C11-C8	-3.20	105.18	111.68
5	A	1445	B3P	O2-C10-C8	-2.95	105.68	111.68
5	A	1445	B3P	C11-C8-C10	-2.88	103.71	110.02
4	A	1444	ACT	O-C-CH3	-2.77	111.16	122.53
5	A	1445	B3P	C10-C8-C9	-2.63	104.25	110.02
2	B	1440	PLP	C6-C5-C4	2.62	120.24	118.10
2	B	1440	PLP	O4P-P-O1P	2.60	113.46	106.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1442	PO4	O3-P-O2	2.58	115.94	107.91
4	B	1445	ACT	OXT-C-CH3	2.55	125.74	115.05
3	A	1442	PO4	O4-P-O1	2.49	119.75	110.95
2	B	1440	PLP	C5A-C5-C4	-2.46	117.77	122.64
4	A	1444	ACT	OXT-C-CH3	2.46	125.35	115.05
5	A	1445	B3P	C7-C4-N1	2.42	116.24	109.02
2	B	1440	PLP	C4A-C4-C5	2.28	123.29	120.94
3	B	1443	PO4	O4-P-O2	2.21	114.78	107.91
2	A	1441	PLP	C6-N1-C2	2.19	123.18	119.20
3	B	1441	PO4	O2-P-O1	2.16	118.59	110.95
2	B	1440	PLP	C6-N1-C2	2.13	123.06	119.20
3	B	1441	PO4	O4-P-O1	-2.03	103.77	110.95
3	B	1442	PO4	O4-P-O1	-2.02	103.80	110.95

There are no chirality outliers.

All (1) torsion outliers are listed below:

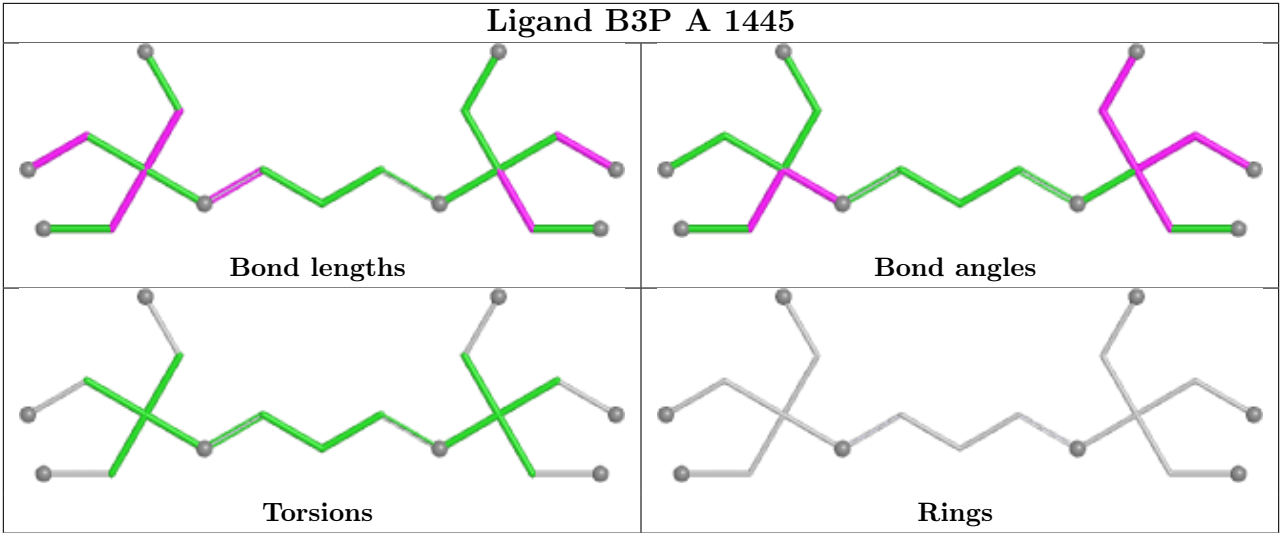
Mol	Chain	Res	Type	Atoms
2	A	1441	PLP	C5A-O4P-P-O1P

There are no ring outliers.

4 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1442	PO4	1	0
4	A	1444	ACT	1	0
4	B	1445	ACT	2	0
4	A	1443	ACT	3	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 ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	371:ILE	C	372:ASP	N	1.08

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	A	423/446 (94%)	-0.17	8 (1%) 66 70	7, 14, 27, 47	5 (1%)
1	B	422/446 (94%)	0.03	10 (2%) 59 63	10, 17, 32, 57	1 (0%)
All	All	845/892 (94%)	-0.07	18 (2%) 63 67	7, 15, 31, 57	6 (0%)

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	439	PRO	3.5
1	A	185	VAL	3.4
1	A	182	ILE	2.8
1	A	440	GLN	2.8
1	A	18	ASP	2.4
1	B	354	GLN	2.4
1	B	378	GLY	2.4
1	A	181	GLY	2.4
1	B	433	LEU	2.3
1	A	369[A]	PHE	2.2
1	B	376	SER	2.2
1	B	436	ALA	2.2
1	A	378	GLY	2.2
1	B	379	SER	2.2
1	B	355	ARG	2.1
1	B	380	ARG	2.1
1	A	355	ARG	2.1
1	B	377	ASP	2.1

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 ⓘ

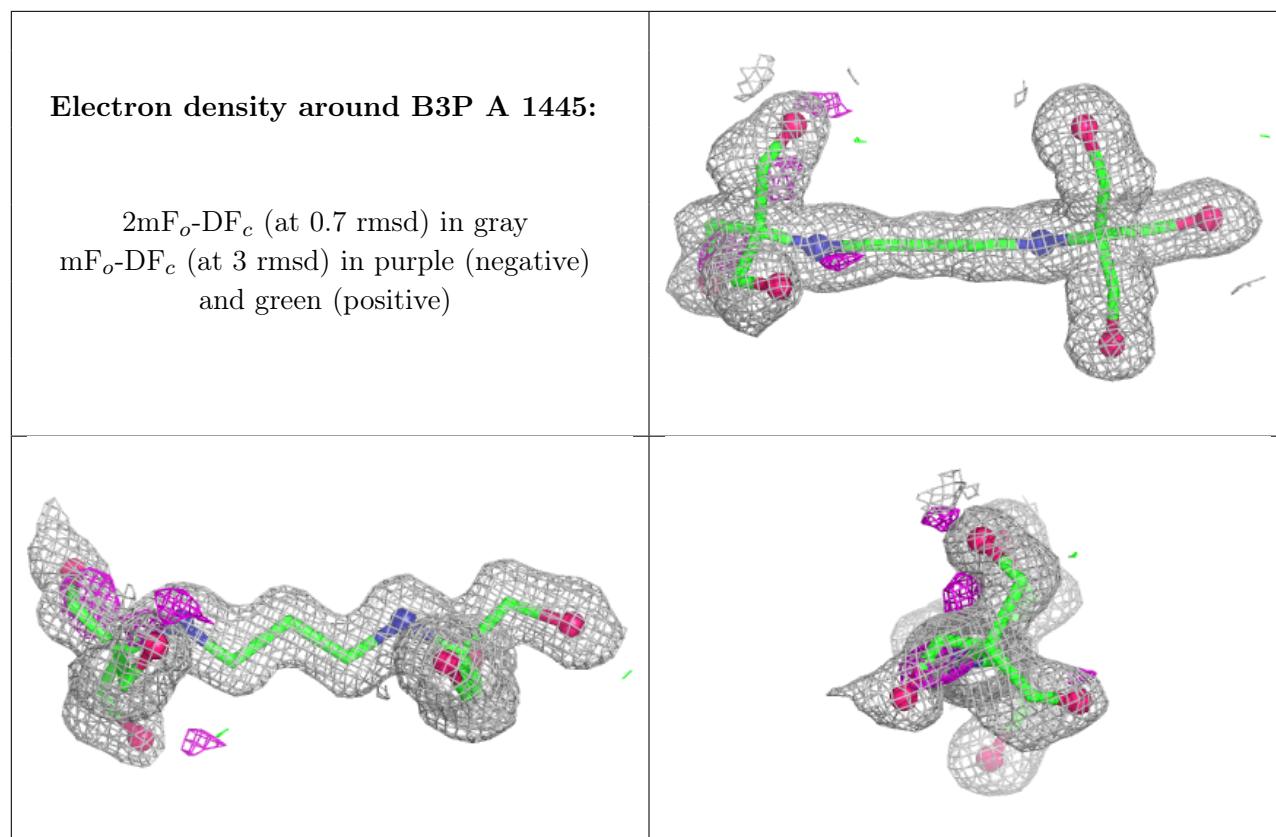
There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 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
4	ACT	A	1444	4/4	0.77	0.17	34,37,39,39	0
3	PO4	B	1444	5/5	0.79	0.10	47,57,62,64	0
4	ACT	B	1445	4/4	0.87	0.15	16,21,34,35	0
3	PO4	B	1443	5/5	0.88	0.12	50,53,60,65	0
4	ACT	A	1443	4/4	0.90	0.14	16,20,31,33	0
5	B3P	A	1445	19/19	0.93	0.09	14,19,32,34	0
3	PO4	B	1441	5/5	0.94	0.09	30,31,38,41	0
3	PO4	A	1442	5/5	0.95	0.09	28,29,36,43	0
3	PO4	B	1442	5/5	0.95	0.10	31,37,47,60	0
2	PLP	A	1441	15/16	0.96	0.09	12,16,22,23	0
2	PLP	B	1440	15/16	0.97	0.07	13,17,21,21	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.



6.5 Other polymers [i](#)

There are no such residues in this entry.