Full wwPDB X-ray Structure Validation Report (i)

Jun 23, 2024 – 12:39 AM EDT

PDB ID	:	6KZ9
Title	:	Crystal structure of plant Phospholipase D alpha
Authors	:	Li, J.X.; Yu, F.; Zhang, P.
Deposited on	:	2019-09-23
Resolution	:	1.80 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 (i) 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 (1)) were used in the production of this report:

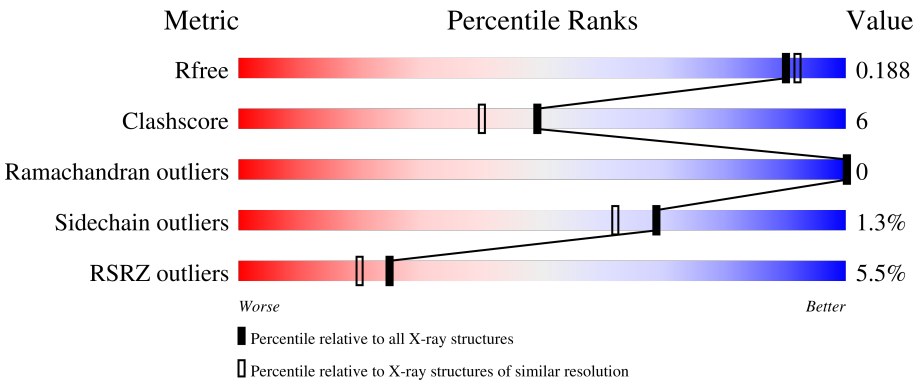
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

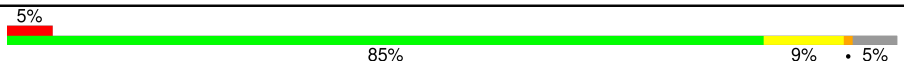
The reported resolution of this entry is 1.80 Å.

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	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

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 <=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	810	

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 6828 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 Phospholipase D alpha 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	770	6199	3941	1077	1154	27	0	0	0

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

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

- Molecule 3 is water.

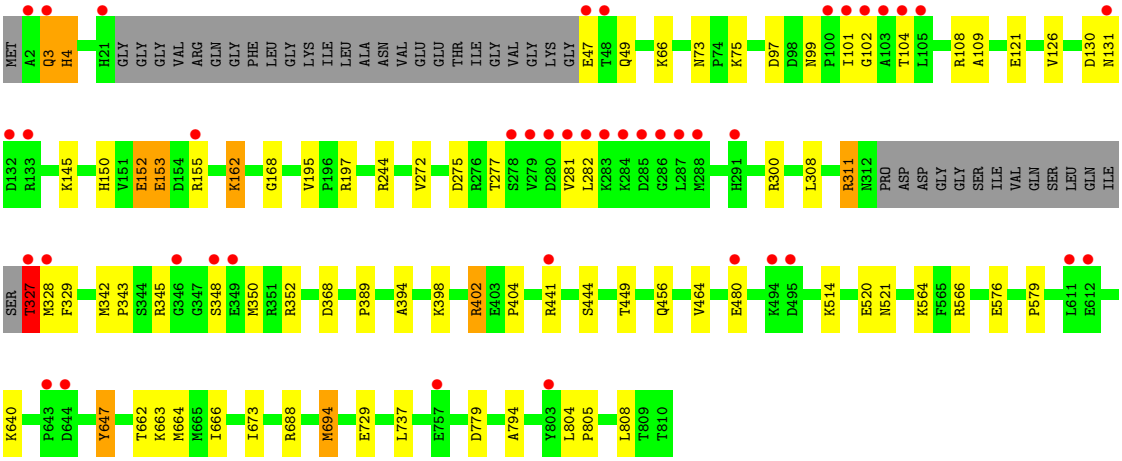
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	628	628	628	0	0

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: Phospholipase D alpha 1

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	118.89Å 127.44Å 55.36Å 90.00° 95.47° 90.00°	Depositor
Resolution (Å)	42.40 – 1.80 42.39 – 1.80	Depositor EDS
% Data completeness (in resolution range)	96.3 (42.40-1.80) 96.1 (42.39-1.80)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.54 (at 1.81Å)	Xtriage
Refinement program	PHENIX 1.10_2155	Depositor
R, R_{free}	0.167 , 0.187 0.168 , 0.188	Depositor DCC
R_{free} test set	2000 reflections (2.66%)	wwPDB-VP
Wilson B-factor (Å^2)	18.0	Xtriage
Anisotropy	0.055	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 52.3	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	6828	wwPDB-VP
Average B, all atoms (Å^2)	14.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.83% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

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

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	0.50	2/6361 (0.0%)	0.69	10/8620 (0.1%)

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	3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	152	GLU	CG-CD	6.13	1.61	1.51
1	A	153	GLU	CG-CD	5.91	1.60	1.51

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	4	HIS	N-CA-CB	14.29	136.31	110.60
1	A	3	GLN	CB-CA-C	14.04	138.47	110.40
1	A	3	GLN	C-N-CA	8.95	144.06	121.70
1	A	162	LYS	CB-CA-C	-8.74	92.92	110.40
1	A	402	ARG	NE-CZ-NH2	-8.21	116.20	120.30
1	A	402	ARG	NE-CZ-NH1	7.08	123.84	120.30
1	A	152	GLU	OE1-CD-OE2	-6.86	115.06	123.30
1	A	3	GLN	O-C-N	-5.42	114.04	122.70
1	A	275	ASP	CB-CG-OD2	-5.35	113.49	118.30
1	A	694	MET	CG-SD-CE	5.34	108.74	100.20

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	3	GLN	Mainchain,Peptide
1	A	327	THR	Peptide

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	6199	0	6017	72	0
2	A	1	0	0	0	0
3	A	628	0	0	24	2
All	All	6828	0	6017	72	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 6.

All (72) 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:47:GLU:HG3	3:A:1002:HOH:O	1.54	1.07
1:A:47:GLU:CG	3:A:1002:HOH:O	2.05	1.04
1:A:130:ASP:OD2	3:A:1001:HOH:O	1.81	0.96
1:A:345:ARG:NH2	3:A:1004:HOH:O	2.00	0.93
1:A:345:ARG:NH1	3:A:1006:HOH:O	2.03	0.90
1:A:47:GLU:OE2	3:A:1002:HOH:O	1.90	0.90
1:A:277:THR:O	1:A:327:THR:HG21	1.72	0.90
1:A:398:LYS:NZ	3:A:1005:HOH:O	2.02	0.88
1:A:345:ARG:CZ	3:A:1004:HOH:O	2.29	0.76
1:A:99:ASN:HB3	1:A:102:GLY:O	1.86	0.75
1:A:664:MET:SD	1:A:666:ILE:HG13	2.28	0.73
1:A:564:LYS:HZ3	1:A:737:LEU:H	1.39	0.70
1:A:464:VAL:HG13	1:A:694:MET:HE2	1.77	0.67
1:A:47:GLU:CD	3:A:1002:HOH:O	2.22	0.65
1:A:640:LYS:NZ	3:A:1007:HOH:O	2.10	0.64
1:A:155:ARG:NE	3:A:1018:HOH:O	2.32	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:464:VAL:HG13	1:A:694:MET:CE	2.30	0.61
1:A:244:ARG:NH1	3:A:1020:HOH:O	2.34	0.61
1:A:155:ARG:NH1	1:A:162:LYS:NZ	2.50	0.60
1:A:300:ARG:O	3:A:1009:HOH:O	2.17	0.58
1:A:150:HIS:HD2	1:A:152:GLU:HG3	1.68	0.58
1:A:328:MET:HE2	1:A:808:LEU:HG	1.85	0.58
1:A:345:ARG:NE	3:A:1004:HOH:O	2.33	0.57
1:A:779:ASP:OD2	3:A:1008:HOH:O	2.17	0.57
1:A:49:GLN:HB3	1:A:66:LYS:HD2	1.87	0.55
1:A:155:ARG:NH2	3:A:1026:HOH:O	2.39	0.55
1:A:155:ARG:NH1	1:A:162:LYS:HZ1	2.08	0.52
1:A:328:MET:CE	1:A:808:LEU:HG	2.40	0.52
1:A:162:LYS:O	1:A:162:LYS:HG2	2.10	0.50
1:A:329:PHE:CB	1:A:688:ARG:O	2.59	0.50
1:A:564:LYS:HE2	1:A:566:ARG:HH21	1.75	0.50
1:A:121:GLU:OE1	3:A:1010:HOH:O	2.20	0.50
1:A:329:PHE:HB3	1:A:688:ARG:O	2.11	0.50
1:A:343:PRO:HD3	1:A:352:ARG:NE	2.26	0.50
1:A:121:GLU:OE2	1:A:145:LYS:HD2	2.13	0.48
1:A:109:ALA:HB1	1:A:126:VAL:HG23	1.95	0.48
1:A:150:HIS:CD2	1:A:152:GLU:HG3	2.49	0.48
1:A:729:GLU:H	1:A:729:GLU:CD	2.17	0.48
1:A:97:ASP:HB3	1:A:104:THR:HG22	1.96	0.47
1:A:272:VAL:HG23	1:A:308:LEU:HD23	1.97	0.47
1:A:368:ASP:CG	1:A:402:ARG:HH22	2.17	0.47
1:A:327:THR:HG23	1:A:328:MET:HA	1.96	0.47
1:A:73:ASN:OD1	1:A:75:LYS:HE3	2.16	0.46
1:A:282:LEU:HD11	1:A:808:LEU:HD22	1.96	0.46
1:A:277:THR:HB	1:A:327:THR:HB	1.97	0.46
1:A:394:ALA:HA	1:A:647:TYR:CE2	2.51	0.46
1:A:108:ARG:NH2	1:A:131:ASN:OD1	2.50	0.45
1:A:343:PRO:HD3	1:A:352:ARG:CD	2.46	0.45
1:A:4:HIS:HD2	3:A:1572:HOH:O	1.99	0.45
1:A:150:HIS:CD2	1:A:152:GLU:CG	3.01	0.44
1:A:514:LYS:HE2	3:A:1258:HOH:O	2.17	0.44
1:A:343:PRO:HD3	1:A:352:ARG:HD3	2.00	0.44
1:A:456:GLN:HG3	3:A:1480:HOH:O	2.16	0.44
1:A:311:ARG:CZ	3:A:1111:HOH:O	2.65	0.44
1:A:521:ASN:HA	1:A:662:THR:O	2.17	0.44
1:A:389:PRO:HG2	1:A:576:GLU:HG3	2.00	0.43
1:A:342:MET:HG2	1:A:343:PRO:HD2	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:402:ARG:O	1:A:404:PRO:HD3	2.18	0.43
1:A:449:THR:HG22	3:A:1347:HOH:O	2.18	0.43
1:A:520:GLU:HB2	1:A:662:THR:HG21	1.99	0.43
1:A:168:GLY:HA2	3:A:1138:HOH:O	2.19	0.43
1:A:805:PRO:HD2	1:A:808:LEU:HD12	2.00	0.43
1:A:342:MET:HA	1:A:352:ARG:HD3	2.00	0.42
1:A:155:ARG:NH1	1:A:162:LYS:HZ2	2.18	0.42
1:A:348:SER:C	1:A:350:MET:H	2.24	0.42
1:A:97:ASP:HB3	1:A:104:THR:CG2	2.50	0.41
1:A:99:ASN:OD1	1:A:101:ILE:HG22	2.20	0.41
1:A:579:PRO:HB2	1:A:794:ALA:CB	2.51	0.41
1:A:480:GLU:HG3	3:A:1189:HOH:O	2.21	0.41
1:A:348:SER:O	1:A:350:MET:HG2	2.21	0.40
1:A:664:MET:HG2	1:A:673:ILE:O	2.21	0.40
1:A:281:VAL:HG21	1:A:804:LEU:HD21	2.02	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)
3:A:1428:HOH:O	3:A:1520:HOH:O[1_554]	1.94	0.26
3:A:1169:HOH:O	3:A:1169:HOH:O[2_8510]	2.13	0.07

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	764/810 (94%)	739 (97%)	25 (3%)	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	668/697 (96%)	659 (99%)	9 (1%)	69 62

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

Mol	Chain	Res	Type
1	A	153	GLU
1	A	195	VAL
1	A	197	ARG
1	A	311	ARG
1	A	327	THR
1	A	441	ARG
1	A	444	SER
1	A	647	TYR
1	A	663	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	4	HIS
1	A	81	HIS
1	A	150	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

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.

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, 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	770/810 (95%)	0.04	42 (5%) 25 20	3, 12, 28, 40	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	101	ILE	7.2
1	A	327	THR	6.7
1	A	803	TYR	6.4
1	A	100	PRO	5.3
1	A	102	GLY	4.8
1	A	103	ALA	4.7
1	A	104	THR	4.5
1	A	441	ARG	4.1
1	A	280	ASP	4.0
1	A	287	LEU	3.8
1	A	281	VAL	3.8
1	A	2	ALA	3.7
1	A	282	LEU	3.7
1	A	643	PRO	3.7
1	A	284	LYS	3.4
1	A	21	HIS	3.4
1	A	346	GLY	3.3
1	A	131	ASN	3.2
1	A	47	GLU	3.2
1	A	279	VAL	3.1
1	A	288	MET	2.9
1	A	133	ARG	2.9
1	A	644	ASP	2.8
1	A	132	ASP	2.8
1	A	612	GLU	2.8
1	A	283	LYS	2.7
1	A	291	HIS	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	611	LEU	2.6
1	A	285	ASP	2.5
1	A	757	GLU	2.5
1	A	349	GLU	2.5
1	A	3	GLN	2.4
1	A	480	GLU	2.4
1	A	348	SER	2.3
1	A	495	ASP	2.3
1	A	328	MET	2.3
1	A	105	LEU	2.3
1	A	494	LYS	2.2
1	A	278	SER	2.2
1	A	286	GLY	2.1
1	A	48	THR	2.0
1	A	155	ARG	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 monosaccharides in this entry.

6.4 Ligands (i)

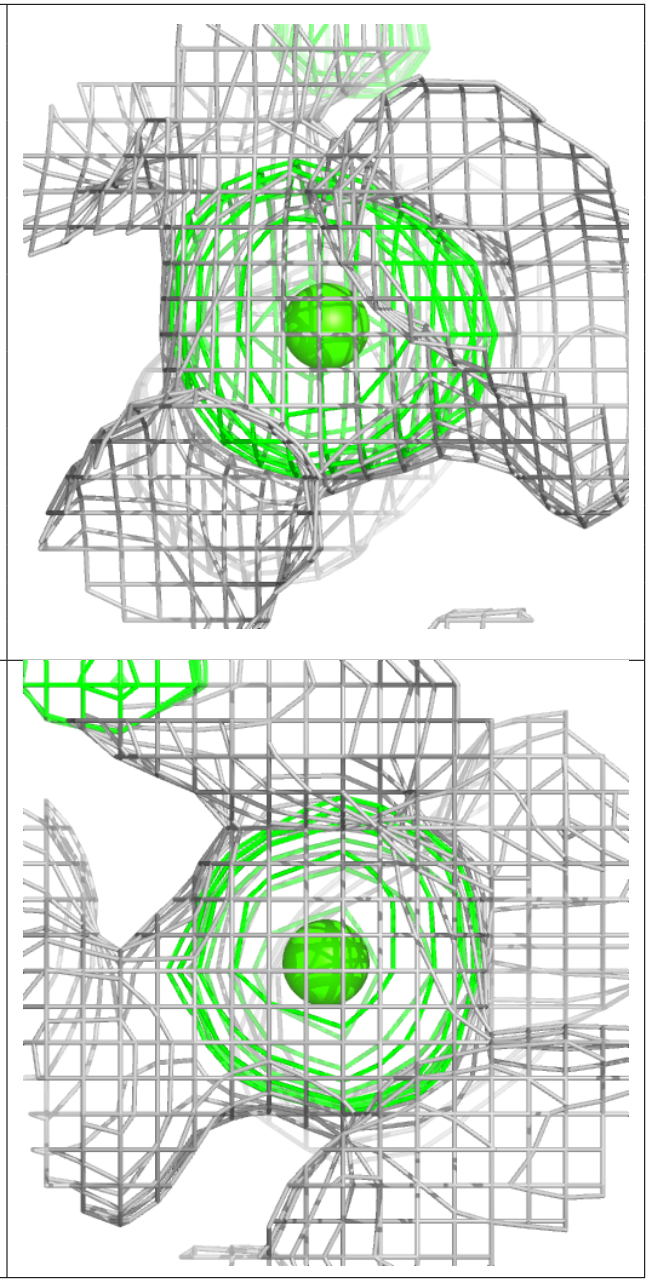
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	CA	A	901	1/1	1.00	0.28	5,5,5,5	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 CA 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 (i)

There are no such residues in this entry.