Full wwPDB X-ray Structure Validation Report (i)

May 1, 2024 – 12:43 am BST

PDB ID	:	4CBU
Title	:	Crystal structure of Plasmodium falciparum actin I
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Deposited on	:	2013-10-16
Resolution	:	1.30 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 (i)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36.2
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)	:	2.36.2

PERCENTILES INFOmissingINFO

1 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 9588 atoms, of which 4348 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 Actin-1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	357	6463	2038	3219	550	639	17	0	64	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP Q8I4X0
A	0	ALA	-	expression tag	UNP Q8I4X0

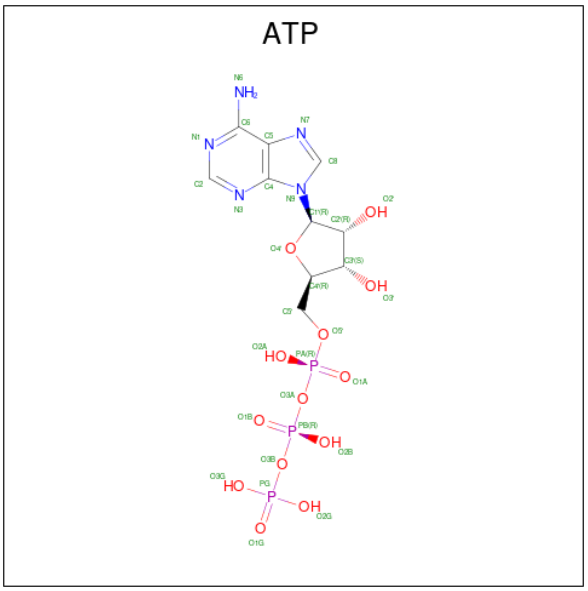
- Molecule 2 is a protein called Gelsolin.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
2	G	125	2260	741	1117	190	210	2	0	18	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
G	23	GLY	-	expression tag	UNP P13020
G	24	PRO	-	expression tag	UNP P13020

- Molecule 3 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
			Total	C	H	N	O	P		
3	A	1	43	10	12	5	13	3	0	0

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	A	549	584	584	0	44
5	G	225	235	235	0	15

SEQUENCE-PLOTS INFOmissingINFO

2 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α, β, γ	68.98Å 110.30Å 71.48Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	55.15 – 1.30 71.48 – 1.30	Depositor EDS
% Data completeness (in resolution range)	99.9 (55.15-1.30) 99.9 (71.48-1.30)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.75 (at 1.30Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.120 , 0.153 0.117 , 0.151	Depositor DCC
R_{free} test set	4041 reflections (3.01%)	wwPDB-VP
Wilson B-factor (Å^2)	10.7	Xtriage
Anisotropy	0.263	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.39 , 55.9	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	0.016 for l,-k,h	Xtriage
F_o,F_c correlation	0.98	EDS
Total number of atoms	9588	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 4.88% 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.

3 Model quality (i)

3.1 Standard geometry (i)

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

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	0.58	0/3385	0.74	0/4574
2	G	0.58	0/1178	0.70	0/1586
All	All	0.58	0/4563	0.73	0/6160

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

3.2 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	3244	3219	3098	31	2
2	G	1143	1117	1098	12	0
3	A	31	12	12	0	0
4	A	1	0	0	0	0
4	G	2	0	0	0	0
5	A	584	0	0	19	1
5	G	235	0	0	4	0
All	All	5240	4348	4208	42	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 5.

All (42) 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:285[B]:LYS:NZ	5:A:2773[B]:HOH:O	1.68	0.99
1:A:338[B]:TYR:OH	5:A:2544:HOH:O	1.93	0.87
1:A:101:GLU:OE1	5:A:2683:HOH:O	2.02	0.77
1:A:289:ASP:OD1	5:A:2637:HOH:O	2.05	0.74
1:A:254:GLU:OE2	5:A:2279:HOH:O	2.07	0.73
1:A:84[B]:GLU:OE2	5:A:2276:HOH:O	2.08	0.71
2:G:75:GLN:OE1	5:G:2496:HOH:O	2.12	0.67
1:A:122[B]:GLN:OE1	5:A:2559:HOH:O	2.14	0.65
1:A:329[A]:LYS:NZ	5:A:2322[A]:HOH:O	2.29	0.64
1:A:335[A]:GLU:O	5:A:2292:HOH:O	2.07	0.63
2:G:75:GLN:NE2	2:G:79:GLY:O	2.31	0.63
1:A:122[B]:GLN:NE2	5:A:2612:HOH:O	2.23	0.62
2:G:139[B]:LYS:NZ	5:G:2402:HOH:O	2.33	0.62
2:G:135[A]:LYS:H	2:G:135[A]:LYS:CD	2.15	0.59
1:A:363:TYR:O	1:A:368:PRO:HB3	2.03	0.59
1:A:355:GLN:NE2	2:G:28:GLU:O	2.30	0.58
2:G:29:HIS:HB3	2:G:32[B]:PHE:CD1	2.40	0.57
1:A:292[B]:LYS:HG2	1:A:326[B]:MET:SD	2.46	0.56
1:A:336:ARG:HA	1:A:339[B]:SER:OG	2.08	0.54
1:A:22[B]:VAL:HG11	5:A:2735:HOH:O	2.07	0.54
1:A:120[B]:MET:HE1	5:A:2334[B]:HOH:O	2.08	0.53
1:A:35[B]:ILE:CD1	1:A:60:GLN:HG3	2.40	0.52
1:A:106:LEU:HD12	1:A:133[B]:MET:SD	2.50	0.52
1:A:366:SER:N	1:A:367:GLY:CA	2.74	0.50
2:G:78:ASN:ND2	2:G:80:ASN:OD1	2.45	0.50
1:A:118[A]:GLU:OE1	5:A:2317:HOH:O	2.20	0.47
2:G:72:LYS:NZ	5:G:2468:HOH:O	2.45	0.47
1:A:338[B]:TYR:HB2	5:A:2292:HOH:O	2.14	0.47
1:A:13[B]:ASN:HD22	1:A:18:VAL:HG22	1.81	0.46
1:A:22[B]:VAL:CG1	5:A:2735:HOH:O	2.61	0.46
1:A:125:PHE:CZ	1:A:133[A]:MET:HG2	2.51	0.45
2:G:135[A]:LYS:H	2:G:135[A]:LYS:HD2	1.80	0.45
1:A:29:ARG:HD3	1:A:29:ARG:HA	1.84	0.45
1:A:71[A]:PRO:HG3	1:A:82:ASP:HB3	1.99	0.45
1:A:57:ASP:OD2	5:A:2632:HOH:O	2.21	0.44
1:A:180[A]:ASP:OD2	5:A:2635:HOH:O	2.21	0.44
2:G:85:LEU:HD21	2:G:108[A]:LEU:HD23	2.01	0.43
1:A:208:GLU:OE2	5:A:2510:HOH:O	2.21	0.42
2:G:135[B]:LYS:HE2	5:G:2772:HOH:O	2.18	0.42
1:A:18:VAL:HG22	5:A:2545:HOH:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:347[A]:LEU:HD23	1:A:347[A]:LEU:C	2.41	0.40
2:G:25:MET:O	2:G:25:MET:HG3	2.22	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:101:GLU:OE2	1:A:324[B]:SER:OG[3_556]	2.11	0.09
1:A:5[B]:ASP:OD1	5:A:2724:HOH:O[4_456]	2.19	0.01

3.3 Torsion angles (i)

3.3.1 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	418/378 (111%)	407 (97%)	11 (3%)	0	100	100
2	G	141/127 (111%)	138 (98%)	3 (2%)	0	100	100
All	All	559/505 (111%)	545 (98%)	14 (2%)	0	100	100

There are no Ramachandran outliers to report.

3.3.2 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	369/322 (115%)	360 (98%)	9 (2%)	49	12

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
2	G	121/105 (115%)	118 (98%)	3 (2%)	47 10
All	All	490/427 (115%)	478 (98%)	12 (2%)	52 12

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	51	LYS
1	A	52	ASP
1	A	63	ARG
1	A	239	LYS
1	A	292[A]	LYS
1	A	292[B]	LYS
1	A	335[A]	GLU
1	A	335[B]	GLU
1	A	357	TRP
2	G	78	ASN
2	G	135[A]	LYS
2	G	135[B]	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
2	G	78	ASN
2	G	80	ASN

3.3.3 RNA (i)

There are no RNA molecules in this entry.

3.4 Non-standard residues in protein, DNA, RNA chains (i)

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

3.5 Carbohydrates (i)

There are no monosaccharides in this entry.

3.6 Ligand geometry (i)

Of 4 ligands modelled in this entry, 3 are 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	ATP	A	1369	4	26,33,33	0.68	0	31,52,52	1.25	3 (9%)

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	ATP	A	1369	4	-	3/18/38/38	0/3/3/3

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	1369	ATP	N3-C2-N1	-3.28	123.55	128.68
3	A	1369	ATP	C2-N1-C6	2.77	123.50	118.75
3	A	1369	ATP	N6-C6-N1	2.24	123.23	118.57

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1369	ATP	PG-O3B-PB-O1B
3	A	1369	ATP	PB-O3B-PG-O1G
3	A	1369	ATP	PG-O3B-PB-O2B

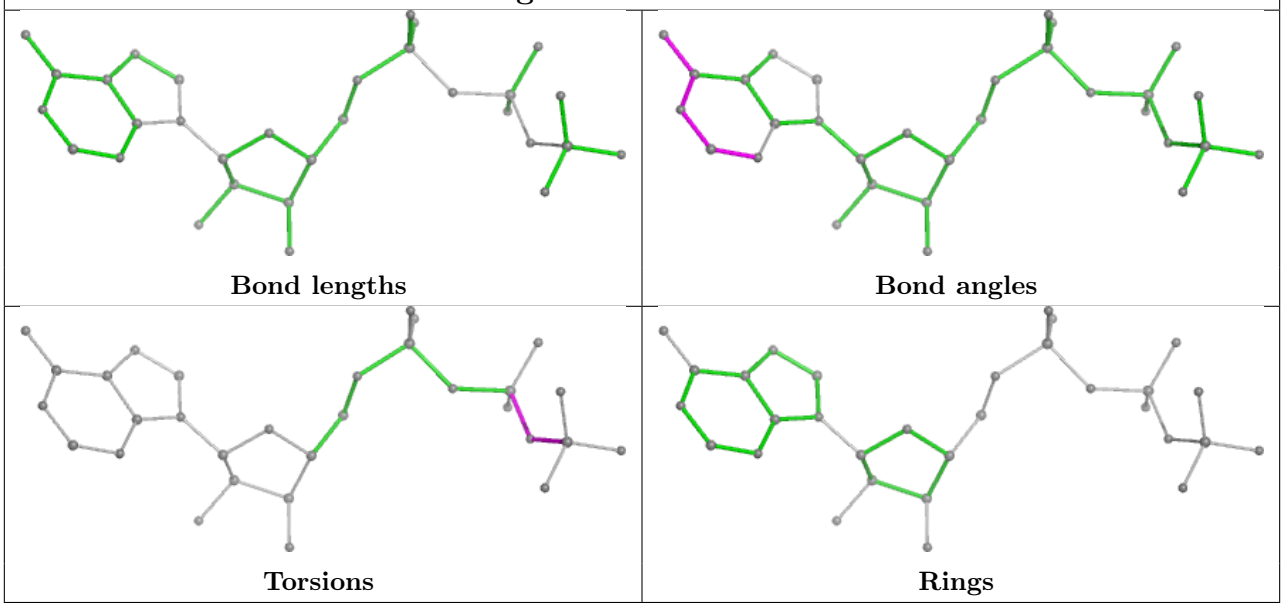
There are no ring outliers.

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 then 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.

Ligand ATP A 1369



3.7 Other polymers (i)

There are no such residues in this entry.

3.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

4 Fit of model and data (i)

4.1 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	357/378 (94%)	-0.49	13 (3%)	42	39	6, 13, 37, 61	1 (0%)
2	G	125/127 (98%)	-0.42	5 (4%)	38	35	6, 12, 34, 56	0
All	All	482/505 (95%)	-0.47	18 (3%)	41	38	6, 12, 37, 61	1 (0%)

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	367	GLY	10.6
1	A	51	LYS	8.3
2	G	26	VAL	8.2
1	A	368	PRO	8.2
2	G	25	MET	8.0
1	A	42	PRO	5.9
2	G	77	ARG	5.6
1	A	5[A]	ASP	5.0
1	A	41	ASN	4.7
2	G	78	ASN	4.6
1	A	52	ASP	3.5
1	A	65	ILE	3.3
1	A	4	GLU	2.9
1	A	61	THR	2.8
1	A	53	ALA	2.8
1	A	54	PHE	2.5
2	G	79	GLY	2.2
1	A	39	PRO	2.0

4.2 Non-standard residues in protein, DNA, RNA chains (i)

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

4.3 Carbohydrates (i)

There are no monosaccharides in this entry.

4.4 Ligands (i)

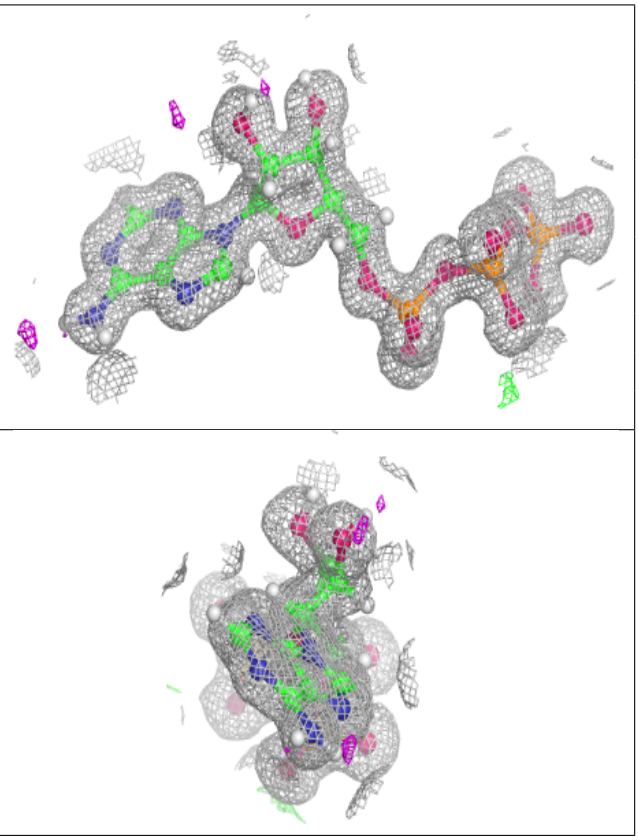
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	ATP	A	1369	31/31	1.00	0.04	8,10,13,15	0
4	CA	A	1370	1/1	1.00	0.05	9,9,9,9	0
4	CA	G	1150	1/1	1.00	0.07	7,7,7,7	0
4	CA	G	1151	1/1	1.00	0.06	10,10,10,10	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 ATP A 1369:

2mF$_o$-DF$_c$ (at 0.7 rmsd) in gray
mF$_o$-DF$_c$ (at 3 rmsd) in purple (negative) and green (positive)



4.5 Other polymers (i)

There are no such residues in this entry.