Full wwPDB X-ray Structure Validation Report ⓘ

Jun 13, 2022 – 04:37 pm BST

PDB ID	:	7B57
Title	:	Crystal structure of CaMKII-actinin complex bound to ADP
Authors	:	Zhu, J.; Gold, M.
Deposited on	:	2020-12-03
Resolution	:	1.95 Å(reported)

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

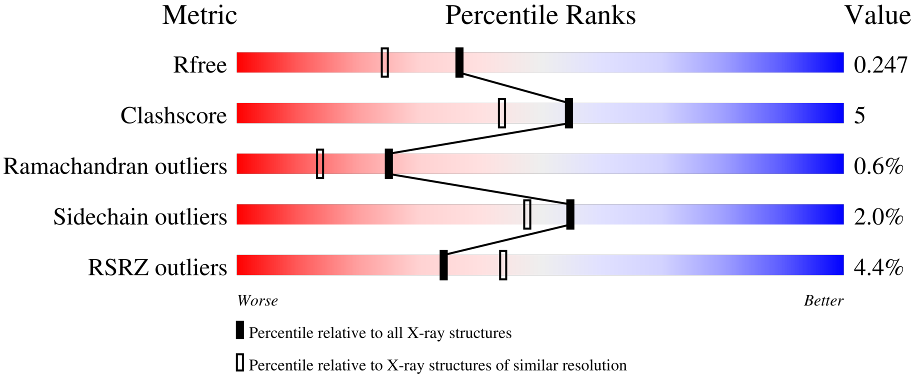
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.28.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0267
CCP4	:	7.1.010 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.28.1

1 Overall quality at a glance ⓘ

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

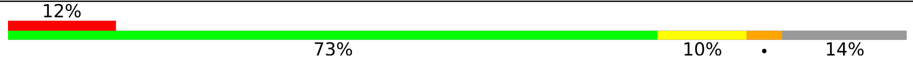
The reported resolution of this entry is 1.95 Å.

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	2580 (1.96-1.96)
Clashscore	141614	2705 (1.96-1.96)
Ramachandran outliers	138981	2678 (1.96-1.96)
Sidechain outliers	138945	2678 (1.96-1.96)
RSRZ outliers	127900	2539 (1.96-1.96)

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	B	317	
2	A	73	

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 3227 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 Calcium/calmodulin-dependent protein kinase type II subunit alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	B	303	2402	1537	420	431	14	0	3	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-1	SER	-	expression tag	UNP P11798
B	0	MET	-	expression tag	UNP P11798
B	305	ALA	THR	engineered mutation	UNP P11798
B	306	ALA	THR	engineered mutation	UNP P11798

- Molecule 2 is a protein called Alpha-actinin-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	A	63	483	311	80	90	2	0	1	0

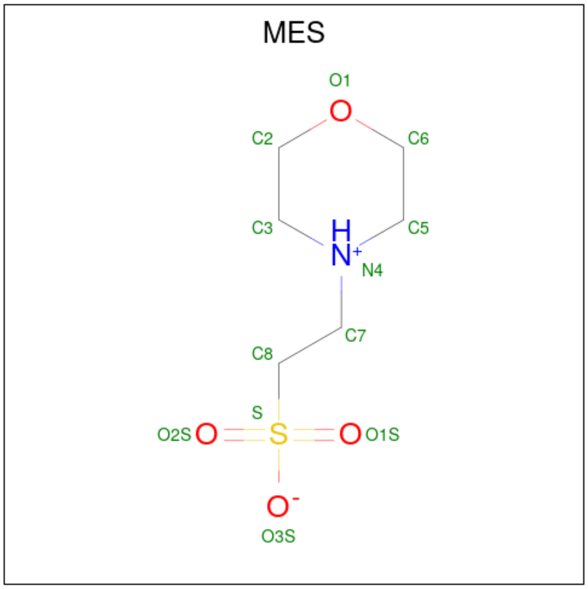
There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	822	SER	-	expression tag	UNP P35609
A	823	ASN	-	expression tag	UNP P35609
A	824	ALA	-	expression tag	UNP P35609

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

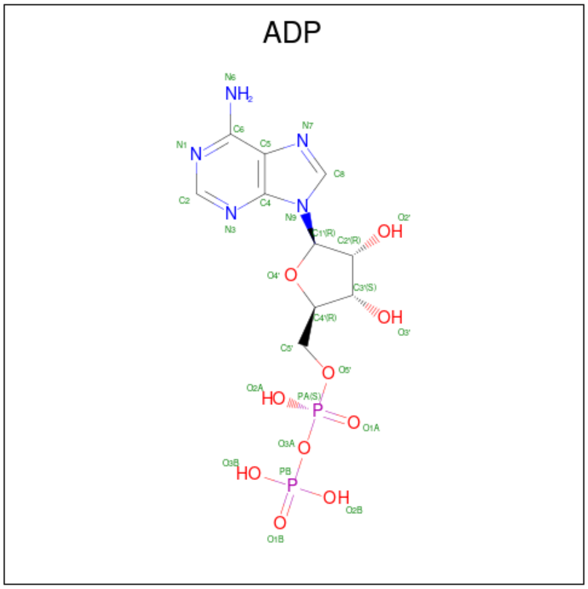
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Mg		
3	B	1	1	1	0	0

- Molecule 4 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES) (formula: $C_6H_{13}NO_4S$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	S		
4	B	1	12	6	1	4	1	0	0

- Molecule 5 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
5	B	1	27	10	5	10	2	0	0

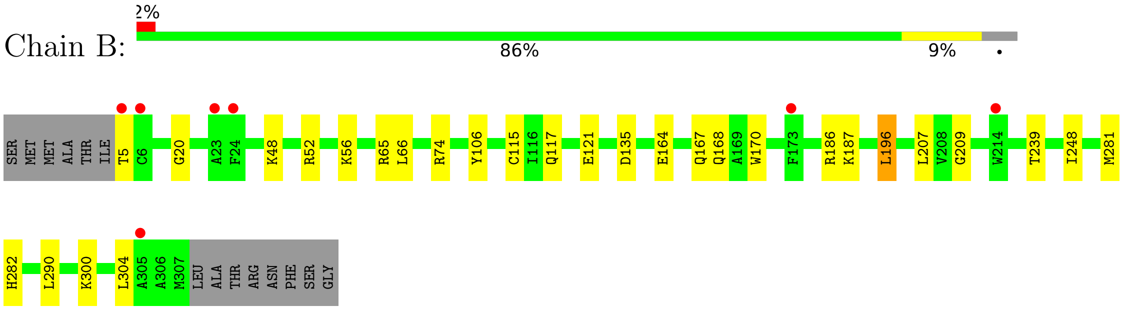
- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	B	275	Total O 275 275	0	0
6	A	27	Total O 27 27	0	0

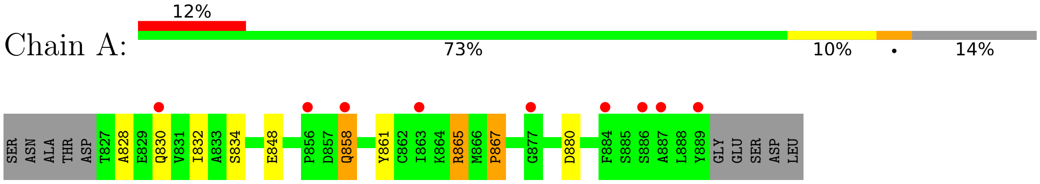
3 Residue-property plots ⓘ

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: Calcium/calmodulin-dependent protein kinase type II subunit alpha



- Molecule 2: Alpha-actinin-2



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	65.92Å 71.85Å 91.60Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	56.53 – 1.95 56.53 – 1.95	Depositor EDS
% Data completeness (in resolution range)	99.8 (56.53-1.95) 91.8 (56.53-1.95)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.14 (at 1.95Å)	Xtriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, R_{free}	0.203 , 0.247 0.203 , 0.247	Depositor DCC
R_{free} test set	1996 reflections (6.17%)	wwPDB-VP
Wilson B-factor (Å^2)	31.4	Xtriage
Anisotropy	0.562	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	(Not available) , (Not available)	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	0.95	EDS
Total number of atoms	3227	wwPDB-VP
Average B, all atoms (Å^2)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.40% of the height of the origin peak. No significant pseudotranslation is detected.*

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[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 ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: MES, MG, ADP

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	B	0.36	0/2463	0.52	1/3338 (0.0%)
2	A	0.29	0/495	0.42	0/673
All	All	0.35	0/2958	0.51	1/4011 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	196	LEU	CA-CB-CG	-7.15	98.85	115.30

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	B	2402	0	2368	19	0
2	A	483	0	471	11	0
3	B	1	0	0	0	0
4	B	12	0	12	0	0
5	B	27	0	12	1	0
6	A	27	0	0	9	2
6	B	275	0	0	10	2
All	All	3227	0	2863	30	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 (30) 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)
2:A:865:ARG:NH1	6:A:901:HOH:O	2.03	0.90
1:B:66:LEU:O	6:B:701:HOH:O	1.91	0.89
1:B:300:LYS:O	6:B:702:HOH:O	2.08	0.70
2:A:832:ILE:N	6:A:902:HOH:O	2.20	0.69
2:A:858:GLN:N	6:A:906:HOH:O	2.20	0.68
2:A:828:ALA:O	6:A:902:HOH:O	2.10	0.68
1:B:135:ASP:OD1	6:B:703:HOH:O	2.14	0.66
1:B:168:GLN:NE2	6:B:705:HOH:O	2.29	0.65
1:B:115[A]:CYS:SG	6:B:917:HOH:O	2.55	0.65
1:B:164:GLU:H	1:B:167:GLN:HE21	1.45	0.64
1:B:74:ARG:NH2	6:B:708:HOH:O	2.32	0.62
2:A:828:ALA:O	6:A:903:HOH:O	2.15	0.61
1:B:209:GLY:HA2	1:B:290:LEU:HD13	1.86	0.57
1:B:164:GLU:O	1:B:167:GLN:HG3	2.05	0.57
1:B:304:LEU:N	6:B:702:HOH:O	2.12	0.57
2:A:834:SER:OG	6:A:904:HOH:O	2.18	0.57
1:B:239:THR:HB	1:B:281:MET:HE1	1.87	0.56
2:A:848:GLU:OE1	6:A:905:HOH:O	2.18	0.55
1:B:52:ARG:NE	1:B:56:LYS:HE3	2.24	0.53
1:B:106:TYR:HA	1:B:282:HIS:CE1	2.47	0.50
1:B:65:ARG:NH2	6:B:714:HOH:O	2.39	0.47
1:B:117:GLN:NE2	6:B:726:HOH:O	2.47	0.47
1:B:170:TRP:CZ3	1:B:187:LYS:HA	2.49	0.47
2:A:858:GLN:HA	2:A:861:TYR:HB3	1.98	0.46
1:B:207:LEU:HD12	1:B:248:ILE:HD11	1.97	0.45
2:A:867:PRO:HG2	2:A:880:ASP:HB3	1.98	0.45
1:B:186:ARG:NH2	6:B:732:HOH:O	2.50	0.44
2:A:865:ARG:HD2	6:A:911:HOH:O	2.20	0.42
1:B:20:GLY:HA3	5:B:603:ADP:H4’	2.02	0.41
2:A:865:ARG:HB2	6:A:911:HOH:O	2.20	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)
6:B:767:HOH:O	6:A:915:HOH:O[2_555]	1.98	0.22

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:B:892:HOH:O	6:A:908:HOH:O[4_445]	2.11	0.09

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	B	304/317 (96%)	296 (97%)	8 (3%)	0	100	100
2	A	62/73 (85%)	55 (89%)	5 (8%)	2 (3%)	4	0
All	All	366/390 (94%)	351 (96%)	13 (4%)	2 (0%)	25	17

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	A	867	PRO
2	A	865	ARG

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	B	249/267 (93%)	245 (98%)	4 (2%)	62	58
2	A	48/58 (83%)	46 (96%)	2 (4%)	30	17
All	All	297/325 (91%)	291 (98%)	6 (2%)	55	48

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	5	THR
1	B	48	LYS
1	B	121	GLU
1	B	196	LEU
2	A	830	GLN
2	A	858	GLN

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

Mol	Chain	Res	Type
1	B	14	GLN
1	B	55	GLN
1	B	84	HIS
1	B	85	HIS
1	B	167	GLN
1	B	168	GLN
2	A	830	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 3 ligands modelled in this entry, 1 is monoatomic - leaving 2 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
4	MES	B	602	-	12,12,12	2.35	1 (8%)	14,16,16	2.00	3 (21%)
5	ADP	B	603	3	24,29,29	0.95	1 (4%)	29,45,45	1.35	3 (10%)

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
4	MES	B	602	-	-	1/6/14/14	0/1/1/1
5	ADP	B	603	3	-	2/12/32/32	0/3/3/3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	602	MES	C8-S	-7.87	1.66	1.77
5	B	603	ADP	C5-C4	2.33	1.47	1.40

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	B	602	MES	C5-N4-C3	4.60	119.18	108.83
5	B	603	ADP	N3-C2-N1	-3.54	123.14	128.68
4	B	602	MES	O3S-S-C8	3.37	111.22	105.77
5	B	603	ADP	PA-O3A-PB	-3.05	122.36	132.83
4	B	602	MES	C6-C5-N4	-2.45	106.38	110.10
5	B	603	ADP	N6-C6-N1	2.16	123.07	118.57

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	603	ADP	C5'-O5'-PA-O3A
4	B	602	MES	C8-C7-N4-C3
5	B	603	ADP	C5'-O5'-PA-O1A

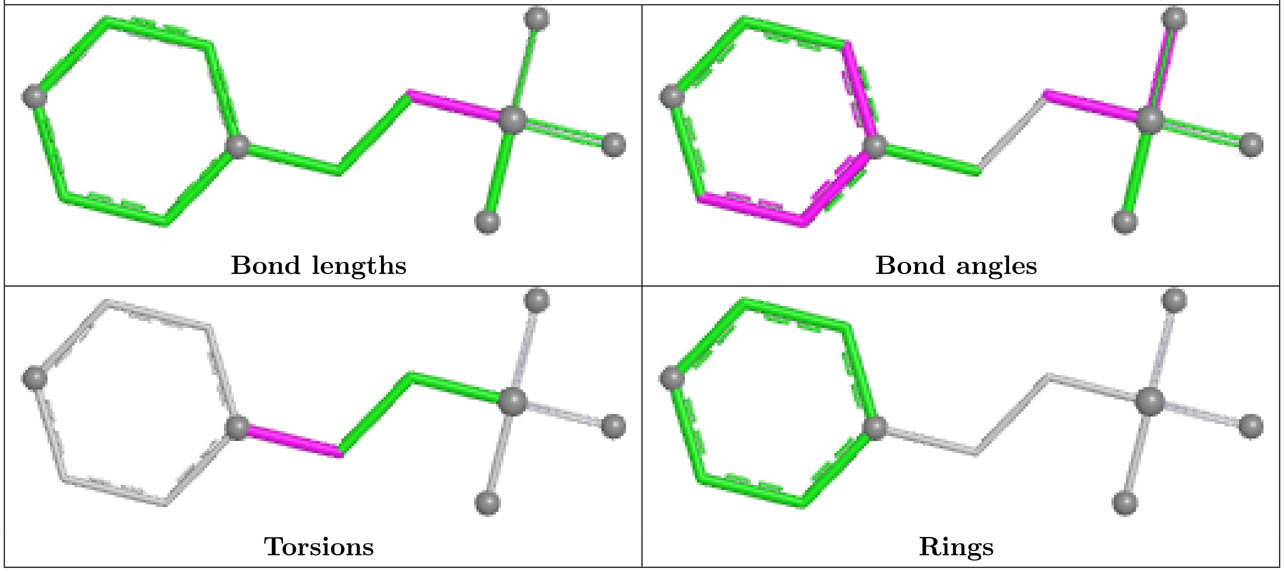
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	603	ADP	1	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 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 MES B 602

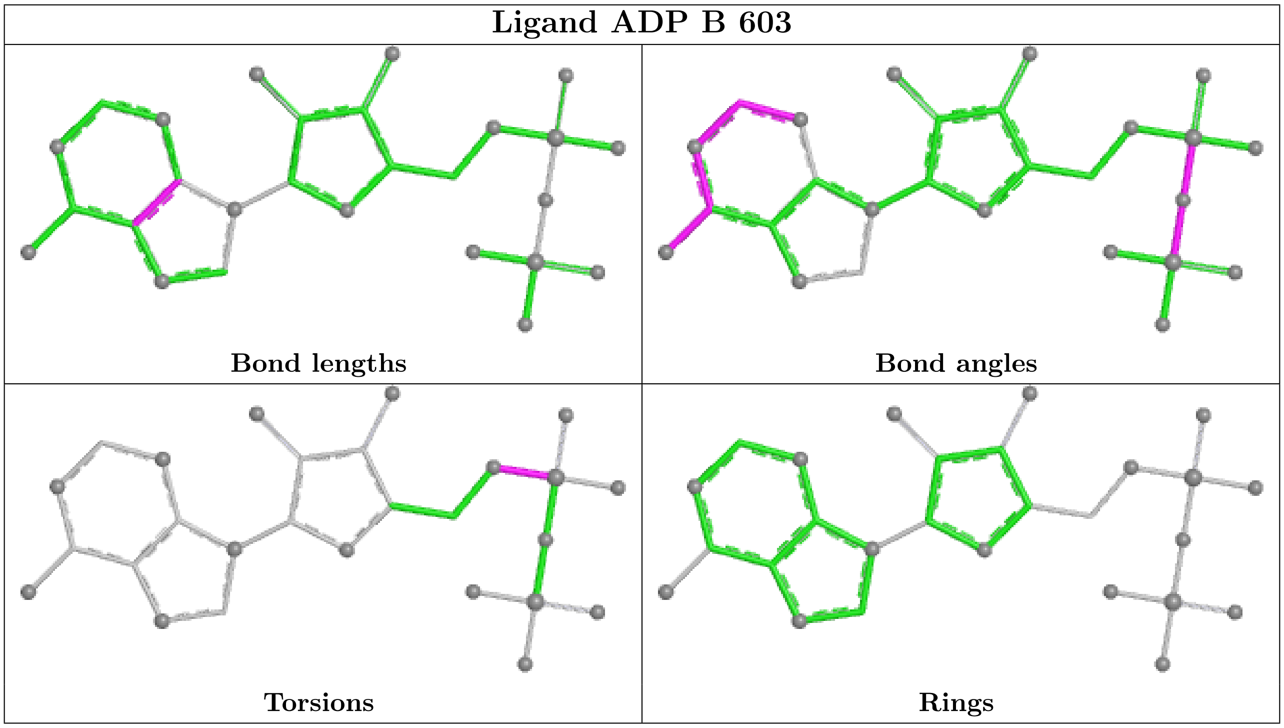


Bond lengths

Bond angles

Torsions

Rings


Ligand ADP B 603

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	B	303/317 (95%)	0.00	7 (2%)	60	69	27, 37, 57, 88	0
2	A	63/73 (86%)	0.81	9 (14%)	2	4	40, 64, 80, 90	0
All	All	366/390 (93%)	0.14	16 (4%)	34	44	27, 39, 74, 90	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	23	ALA	4.7
1	B	24	PHE	4.4
2	A	889	TYR	4.4
1	B	173	PHE	4.2
2	A	887	ALA	3.4
2	A	863	ILE	3.2
2	A	830	GLN	2.9
1	B	5	THR	2.8
2	A	886	SER	2.7
2	A	856	PRO	2.6
2	A	884	PHE	2.5
1	B	214	TRP	2.4
2	A	877	GLY	2.1
2	A	858	GLN	2.1
1	B	305	ALA	2.0
1	B	6	CYS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates ⓘ

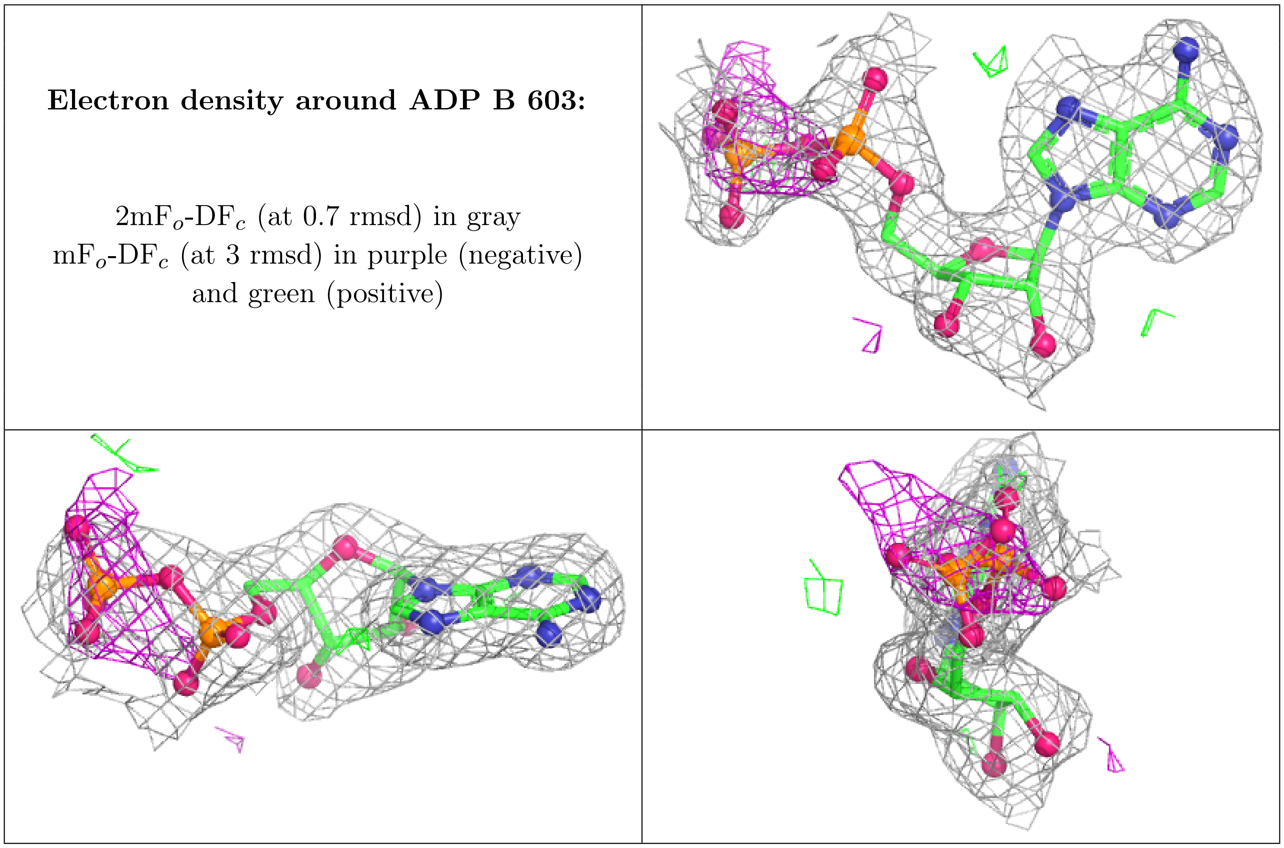
There are no monosaccharides in this entry.

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, 95^{th} 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
5	ADP	B	603	27/27	0.88	0.14	32,43,62,72	0
4	MES	B	602	12/12	0.94	0.20	44,56,65,69	0
3	MG	B	601	1/1	0.94	0.31	58,58,58,58	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.

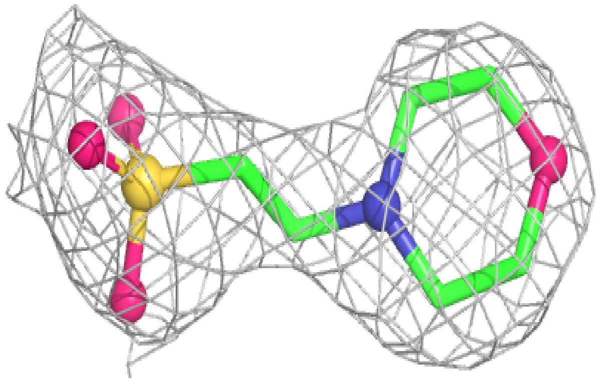


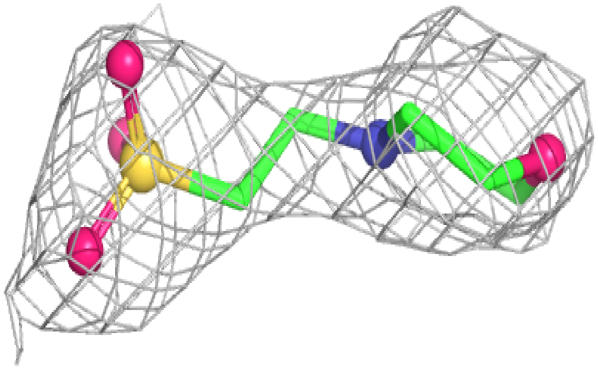
Electron density around ADP B 603:

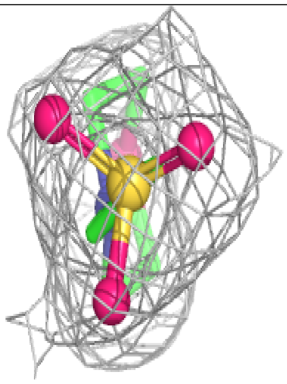
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

Electron density around MES B 602:

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

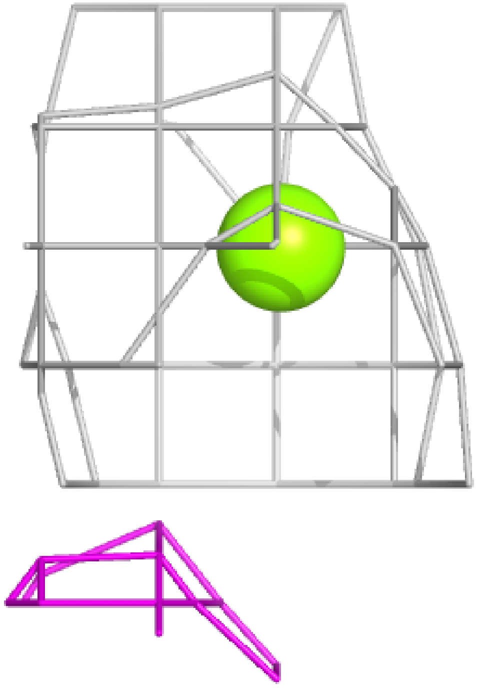


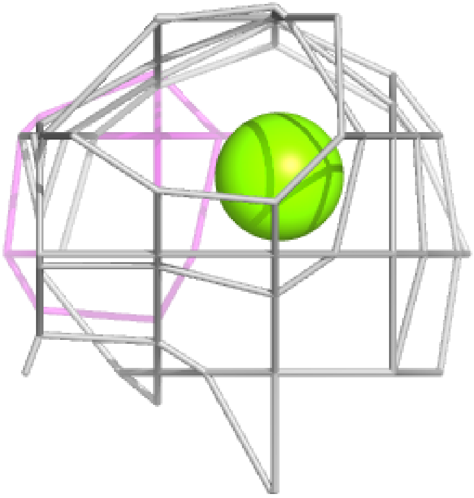


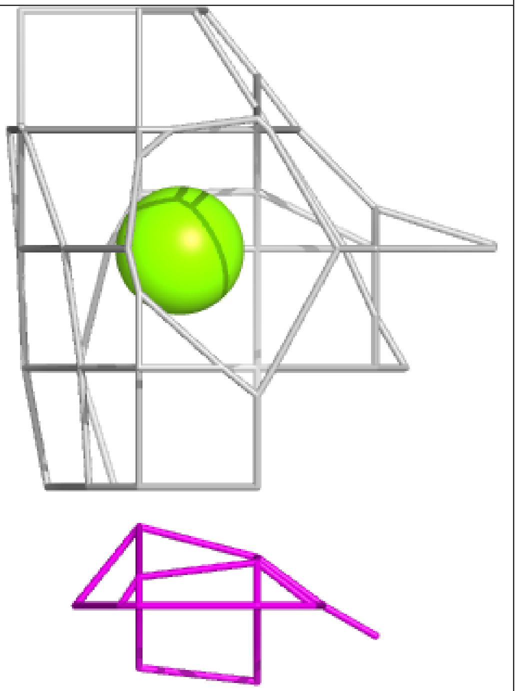


Electron density around MG B 601:

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.