



Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 04:51 am BST

PDB ID : 2A5X
Title : Crystal Structure of a Cross-linked Actin Dimer
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Deposited on : 2005-07-01
Resolution : 2.49 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
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) : 2.11

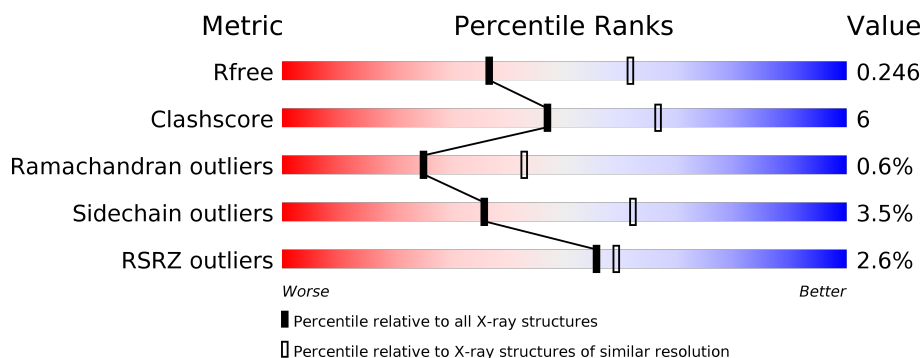
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.49 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	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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 on 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	375	<div> <div>2%</div> <div> <div></div> <div>78%</div> <div>13%</div> <div>9%</div> </div> </div>

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 2812 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.

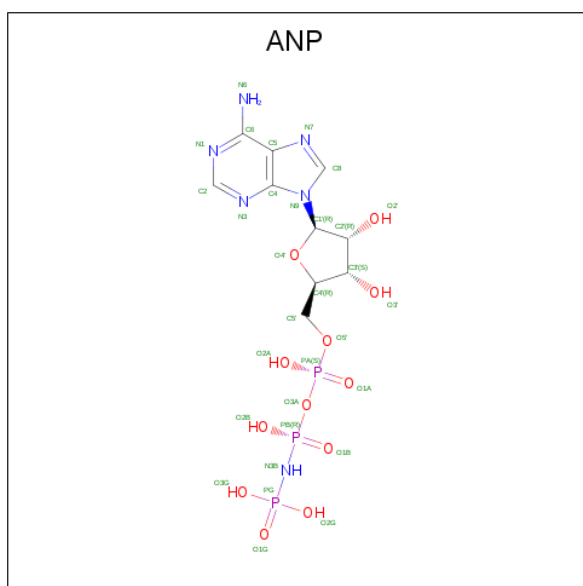
- Molecule 1 is a protein called Actin, alpha skeletal muscle.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	342	Total	C	N	O	S	0	2	0
			2671	1698	445	509	19			

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

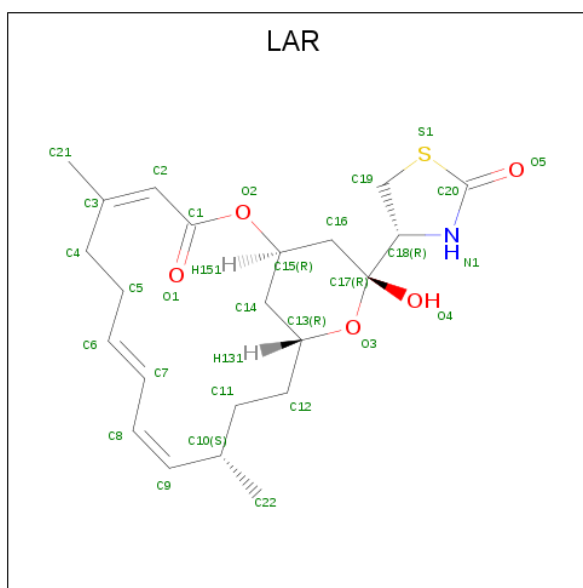
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Ca	0	0
			2	2		

- Molecule 3 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (three-letter code: ANP) (formula: C₁₀H₁₇N₆O₁₂P₃).



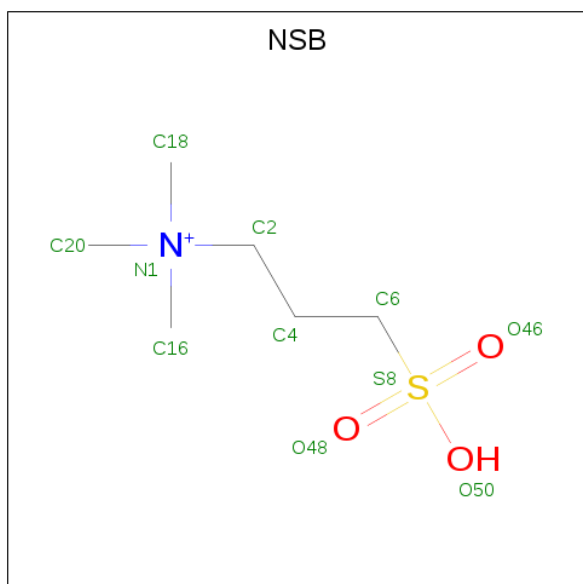
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			31	10	6	12	3		

- Molecule 4 is LATRUNCULIN A (three-letter code: LAR) (formula: $C_{22}H_{31}NO_5S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	S	0	0
			29	22	1	5	1		

- Molecule 5 is N,N,N-TRIMETHYL-3-SULFOPROPAN-1-AMINIUM (three-letter code: NSB) (formula: $C_6H_{16}NO_3S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	S	0	0
			11	6	1	3	1		

- Molecule 6 is (4S)-2-METHYL-2,4-PENTANEDIOL (three-letter code: MPD) (formula: $C_6H_{14}O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			8	6	2		
6	A	1	Total	C	O	0	0
			8	6	2		

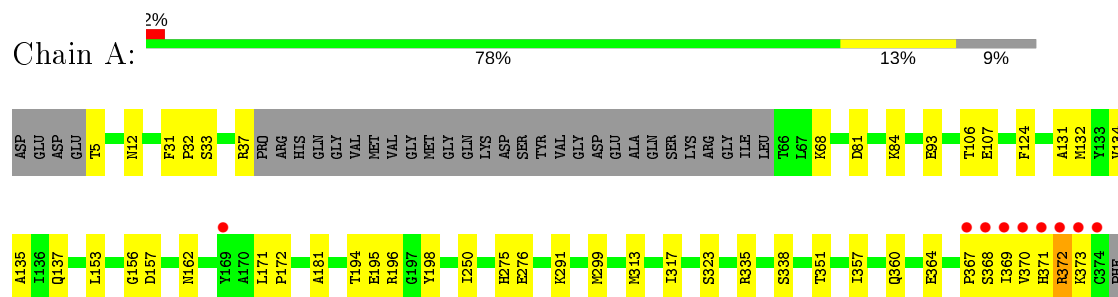
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	52	Total	O	0	0
			52	52		

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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: Actin, alpha skeletal muscle



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	207.38Å 54.37Å 36.20Å 90.00° 98.62° 90.00°	Depositor
Resolution (Å)	90.00 – 2.49 51.26 – 2.49	Depositor EDS
% Data completeness (in resolution range)	98.8 (90.00-2.49) 98.8 (51.26-2.49)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.15	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.08 (at 2.48Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.194 , 0.250 0.195 , 0.246	Depositor DCC
R_{free} test set	686 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	32.7	Xtriage
Anisotropy	0.748	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 34.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.038 for -h-2*k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	2812	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.68% 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: MPD, ANP, CA, LAR, NSB

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	0.53	0/2733	0.60	0/3708

There are no bond length outliers.

There are no bond angle outliers.

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	2671	0	2629	33	0
2	A	2	0	0	0	0
3	A	31	0	13	1	0
4	A	29	0	31	1	0
5	A	11	0	16	0	0
6	A	16	0	28	1	0
7	A	52	0	0	0	0
All	All	2812	0	2717	34	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 6.

All (34) 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:368:SER:HA	1:A:371:HIS:CE1	2.19	0.78
1:A:371:HIS:C	1:A:373:LYS:H	1.89	0.75
1:A:194:THR:HA	1:A:198:TYR:O	1.88	0.74
1:A:37:ARG:NH1	1:A:68:LYS:HG3	2.03	0.73
1:A:156:GLY:O	1:A:181:ALA:HB1	1.87	0.73
1:A:368:SER:O	1:A:370:VAL:N	2.30	0.63
1:A:299:MET:CE	1:A:313:MET:HG3	2.28	0.63
1:A:368:SER:C	1:A:370:VAL:H	2.03	0.62
1:A:31:PHE:HB2	1:A:32:PRO:HD2	1.84	0.58
1:A:368:SER:C	1:A:370:VAL:N	2.58	0.56
1:A:196:ARG:O	1:A:196:ARG:HG2	2.07	0.55
1:A:371:HIS:C	1:A:373:LYS:N	2.57	0.53
1:A:299:MET:HE1	1:A:313:MET:HG3	1.90	0.53
1:A:367:PRO:O	1:A:370:VAL:HG12	2.09	0.51
1:A:31:PHE:CE1	1:A:93:GLU:HG3	2.45	0.51
1:A:37:ARG:HH11	1:A:68:LYS:HG3	1.75	0.51
1:A:157:ASP:HB2	3:A:378:ANP:H5'1	1.93	0.51
1:A:368:SER:O	1:A:371:HIS:ND1	2.44	0.50
1:A:335:ARG:HA	1:A:338:SER:OG	2.12	0.49
1:A:81:ASP:HA	1:A:84:LYS:HD2	1.94	0.49
1:A:153:LEU:HD13	1:A:162:ASN:ND2	2.27	0.49
1:A:313:MET:O	1:A:317:ILE:HG12	2.16	0.46
4:A:379:LAR:O2	4:A:379:LAR:H42	2.16	0.46
1:A:131:ALA:HA	1:A:357:ILE:O	2.15	0.45
1:A:368:SER:CA	1:A:371:HIS:CE1	2.94	0.45
1:A:135:ALA:HB1	6:A:382:MPD:H51	1.99	0.45
1:A:153:LEU:HD13	1:A:162:ASN:HD21	1.84	0.43
1:A:107:GLU:HG2	1:A:134:VAL:HG12	2.01	0.43
1:A:106:THR:HB	1:A:137:GLN:HG3	2.01	0.42
1:A:275:HIS:CE1	1:A:276:GLU:HG3	2.55	0.42
1:A:171:LEU:HA	1:A:172:PRO:HD3	1.95	0.41
1:A:372:ARG:O	1:A:373:LYS:HG2	2.20	0.41
1:A:368:SER:HA	1:A:371:HIS:HE1	1.79	0.41
1:A:124:PHE:CZ	1:A:132:MET:HG3	2.57	0.40

There are no symmetry-related clashes.

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	A	339/375 (90%)	325 (96%)	12 (4%)	2 (1%)	25	43

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	372	ARG
1	A	369	ILE

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	A	289/318 (91%)	279 (96%)	10 (4%)	36	62

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	THR
1	A	12	ASN
1	A	33	SER
1	A	195	GLU
1	A	250	ILE
1	A	291	LYS
1	A	323	SER
1	A	351	THR
1	A	360	GLN

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Mol	Chain	Res	Type
1	A	364	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	275	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 2 are monoatomic - leaving 5 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
6	MPD	A	382	-	7,7,7	0.40	0	9,10,10	0.71	0
5	NSB	A	380	-	10,10,10	2.41	1 (10%)	14,15,15	1.43	2 (14%)
6	MPD	A	381	-	7,7,7	0.34	0	9,10,10	0.49	0
4	LAR	A	379	-	30,31,31	1.65	4 (13%)	32,43,43	2.08	13 (40%)
3	ANP	A	378	2	29,33,33	1.53	7 (24%)	31,52,52	2.31	8 (25%)

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
6	MPD	A	382	-	-	1/5/5/5	-
5	NSB	A	380	-	-	2/8/8/8	-
6	MPD	A	381	-	-	4/5/5/5	-
4	LAR	A	379	-	-	3/23/51/51	0/2/3/3
3	ANP	A	378	2	-	3/14/38/38	0/3/3/3

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	380	NSB	C6-S8	-7.20	1.67	1.77
4	A	379	LAR	O2-C1	5.72	1.46	1.34
4	A	379	LAR	C20-S1	-4.44	1.67	1.77
4	A	379	LAR	C8-C7	-3.91	1.32	1.44
3	A	378	ANP	PB-N3B	3.24	1.71	1.63
3	A	378	ANP	PG-O1G	3.04	1.51	1.46
3	A	378	ANP	PB-O1B	2.92	1.50	1.46
3	A	378	ANP	PG-N3B	2.74	1.70	1.63
3	A	378	ANP	C5-C4	2.52	1.47	1.40
3	A	378	ANP	PB-O3A	2.20	1.61	1.59
4	A	379	LAR	O4-C17	2.05	1.44	1.40
3	A	378	ANP	PG-O2G	-2.03	1.51	1.56

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	378	ANP	O1G-PG-N3B	-8.93	98.61	111.77
4	A	379	LAR	O2-C1-C2	5.17	123.41	111.27
4	A	379	LAR	C8-C7-C6	4.28	150.15	125.51
3	A	378	ANP	N3-C2-N1	-3.86	122.65	128.68
4	A	379	LAR	O1-C1-C2	-3.58	117.22	126.23
5	A	380	NSB	O46-S8-C6	3.43	111.04	106.92
3	A	378	ANP	O2B-PB-O1B	3.30	116.83	109.92
4	A	379	LAR	O3-C17-C18	3.17	108.32	104.25
3	A	378	ANP	O4'-C1'-C2'	-3.17	102.30	106.93
3	A	378	ANP	O3A-PB-N3B	-2.95	98.40	106.59
3	A	378	ANP	O2B-PB-O3A	2.92	114.39	104.64
4	A	379	LAR	C7-C8-C9	2.82	144.58	124.42
4	A	379	LAR	C4-C5-C6	-2.68	106.80	112.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	379	LAR	C16-C17-C18	-2.61	108.67	113.75
4	A	379	LAR	O2-C15-C16	2.49	113.80	107.59
4	A	379	LAR	O2-C1-O1	-2.44	119.38	123.35
5	A	380	NSB	O48-S8-C6	2.39	109.80	106.92
4	A	379	LAR	C14-C15-C16	-2.33	105.88	111.00
3	A	378	ANP	C4-C5-N7	-2.30	107.00	109.40
3	A	378	ANP	C2-N1-C6	2.29	122.66	118.75
4	A	379	LAR	O5-C20-N1	-2.12	124.49	126.81
4	A	379	LAR	C14-C13-C12	-2.12	109.23	113.24
4	A	379	LAR	C21-C3-C4	2.11	118.82	115.27

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	378	ANP	PB-N3B-PG-O1G
4	A	379	LAR	C7-C8-C9-C10
5	A	380	NSB	C2-C4-C6-S8
3	A	378	ANP	C3'-C4'-C5'-O5'
3	A	378	ANP	O4'-C4'-C5'-O5'
6	A	381	MPD	O2-C2-C3-C4
5	A	380	NSB	N1-C2-C4-C6
6	A	381	MPD	C1-C2-C3-C4
6	A	381	MPD	CM-C2-C3-C4
4	A	379	LAR	O2-C1-C2-C3
4	A	379	LAR	C3-C4-C5-C6
6	A	381	MPD	C2-C3-C4-C5
6	A	382	MPD	C2-C3-C4-C5

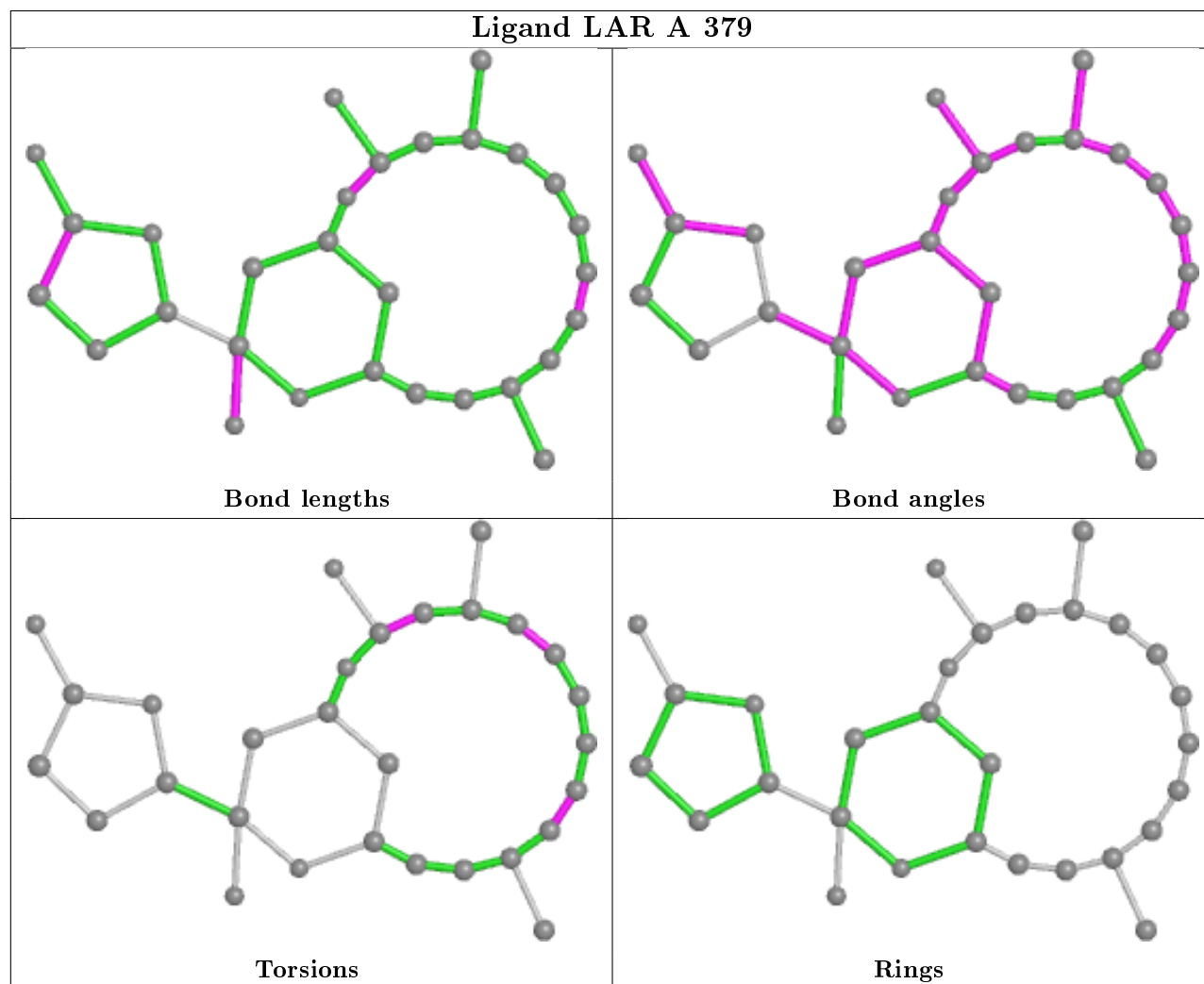
There are no ring outliers.

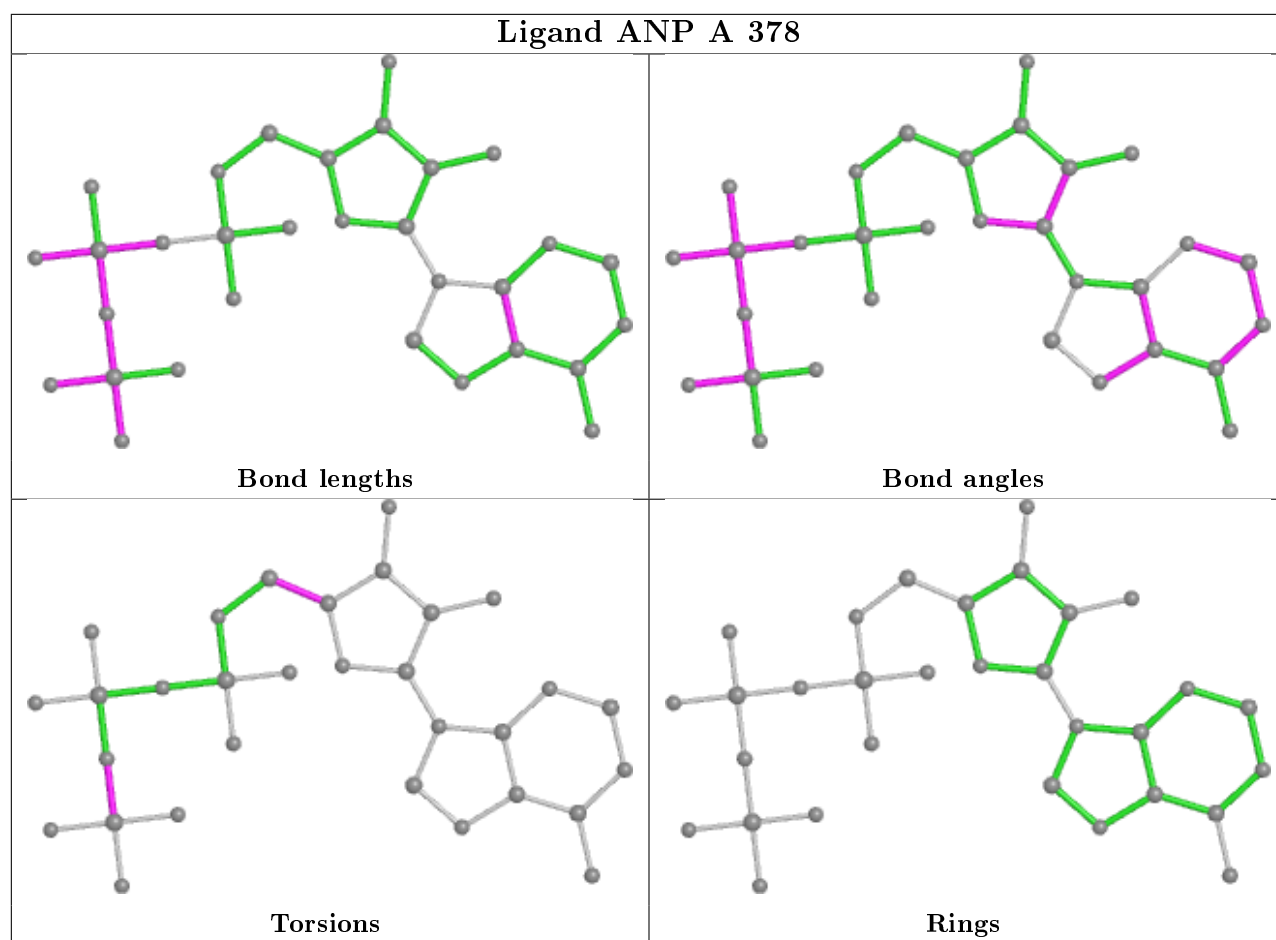
3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	382	MPD	1	0
4	A	379	LAR	1	0
3	A	378	ANP	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 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 [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	342/375 (91%)	-0.14	9 (2%) 56 59	22, 30, 40, 63	8 (2%)

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	374	CYS	12.5
1	A	372	ARG	8.3
1	A	373	LYS	6.0
1	A	370	VAL	5.5
1	A	371	HIS	4.7
1	A	169	TYR	3.9
1	A	369	ILE	3.5
1	A	367	PRO	2.9
1	A	368	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 carbohydrates in this entry.

6.4 Ligands [i](#)

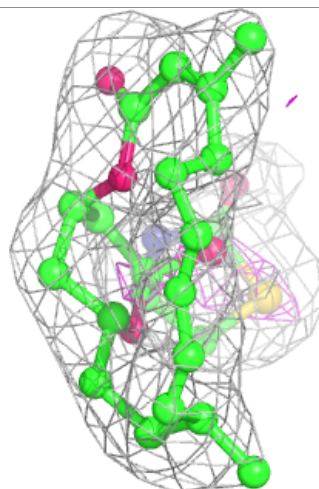
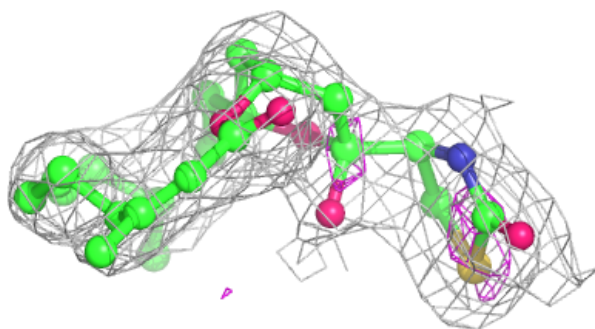
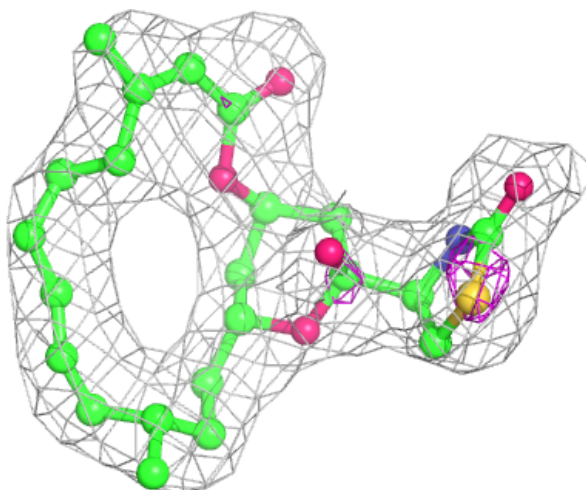
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
5	NSB	A	380	11/11	0.87	0.20	73,75,79,80	0
6	MPD	A	381	8/8	0.87	0.24	70,70,71,71	0
6	MPD	A	382	8/8	0.92	0.19	49,52,54,54	0
4	LAR	A	379	29/29	0.95	0.17	31,34,37,37	0
2	CA	A	377	1/1	0.96	0.06	44,44,44,44	0
2	CA	A	376	1/1	0.98	0.13	27,27,27,27	0
3	ANP	A	378	31/31	0.98	0.11	17,22,25,25	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.

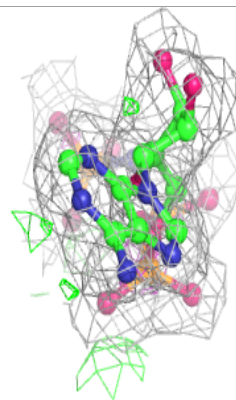
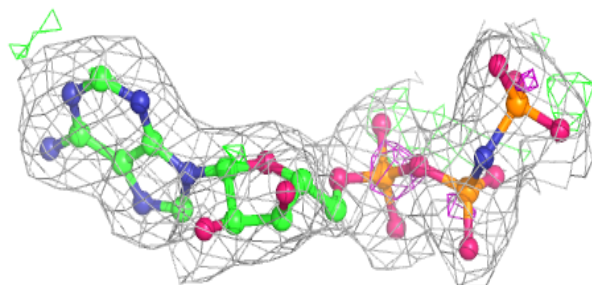
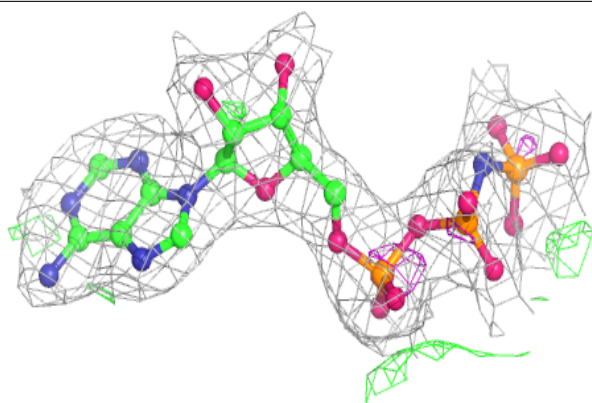
Electron density around LAR A 379:

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 ANP A 378:

$2mF_o-DF_c$ (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.