



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 5, 2026 – 06:38 PM UTC

PDB ID : 1DFL / pdb_00001dfL
Title : SCALLOP MYOSIN S1 COMPLEXED WITH MGADP:VANADATE-TRANSITION STATE
Authors : Houdusse, A.; Szent-Gyorgyi, A.G.; Cohen, C.
Deposited on : 1999-11-19
Resolution : 4.20 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

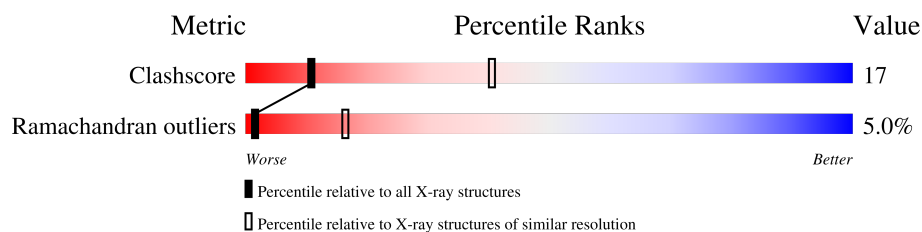
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 4.20 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))
Clashscore	190562	1019 (4.50-3.90)
Ramachandran outliers	187476	1020 (4.56-3.84)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	831	71% 18% • 7%
1	B	831	71% 18% • 8%
2	W	139	70% 24% • •
2	Y	139	71% 24% • •
3	X	152	70% 27% • •
3	Z	152	72% 25% • •

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 10500 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 MYOSIN HEAD.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	A	772	Total	C	N	O	0	0	0
			3816	2271	772	773			
1	B	766	Total	C	N	O	0	0	0
			3786	2253	766	767			

- Molecule 2 is a protein called MYOSIN HEAD.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	Y	136	Total	C	N	O	0	0	0
			673	400	136	137			
2	W	136	Total	C	N	O	0	0	0
			673	400	136	137			

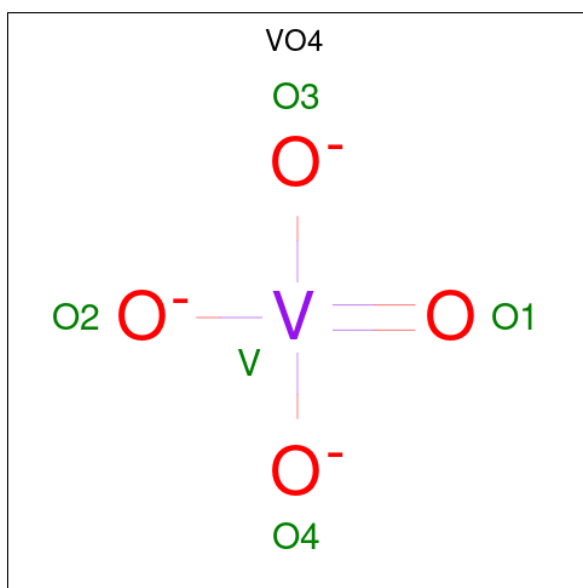
- Molecule 3 is a protein called MYOSIN HEAD.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	Z	151	Total	C	N	O	0	0	0
			741	438	151	152			
3	X	151	Total	C	N	O	0	0	0
			741	438	151	152			

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

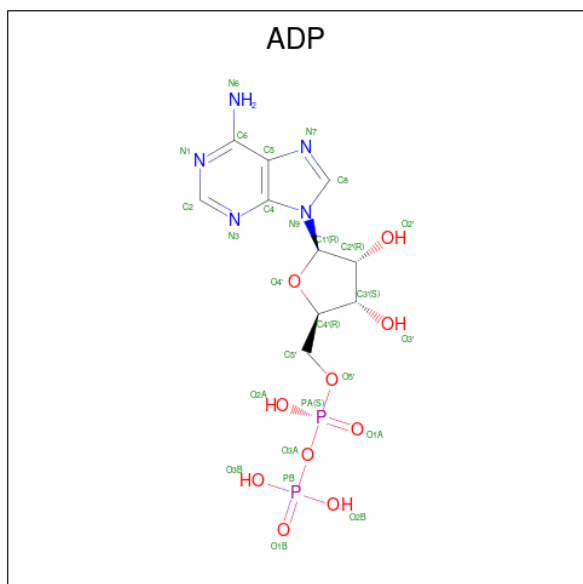
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Mg	0	0
			1	1		
4	Y	1	Total	Mg	0	0
			1	1		
4	B	1	Total	Mg	0	0
			1	1		
4	W	1	Total	Mg	0	0
			1	1		

- Molecule 5 is VANADATE ION (CCD ID: VO4) (formula: O_4V).



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

- Molecule 6 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



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

Continued on next page...

Continued from previous page...

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

- Molecule 7 is CALCIUM ION (CCD ID: CA) (formula: Ca).

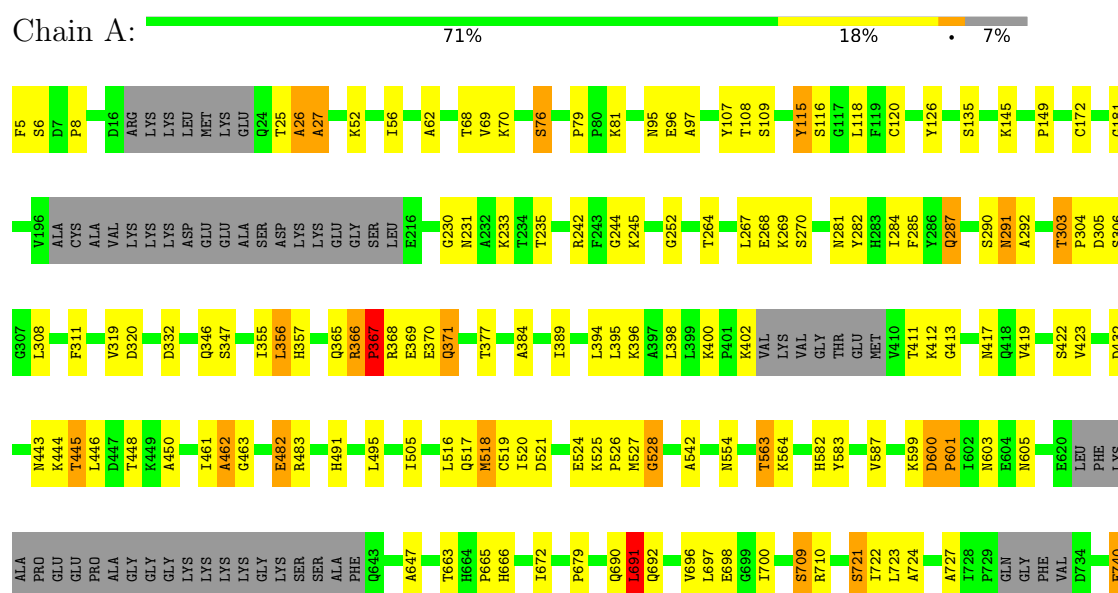
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	Z	1	Total	Ca	0	0
			1	1		
7	X	1	Total	Ca	0	0
			1	1		

3 Residue-property plots

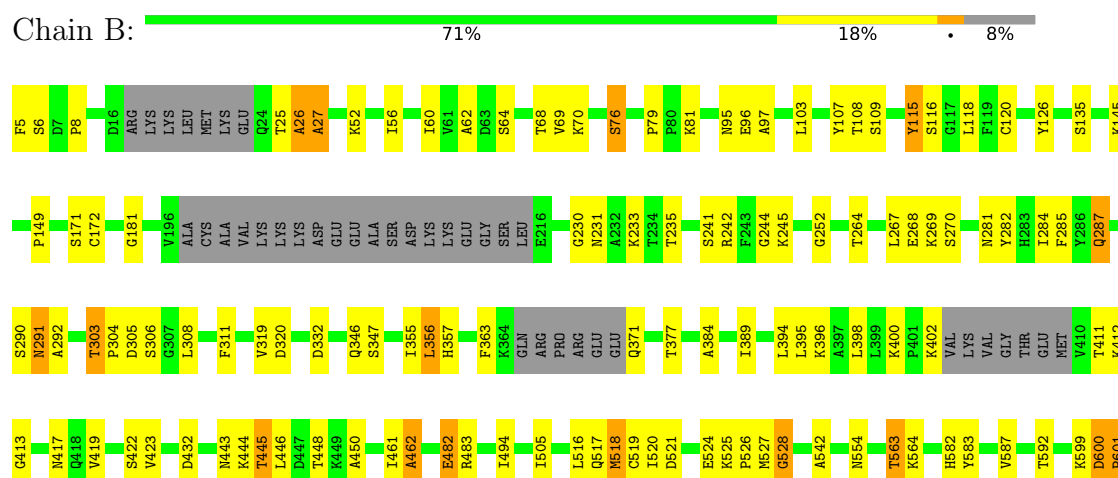
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. 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.

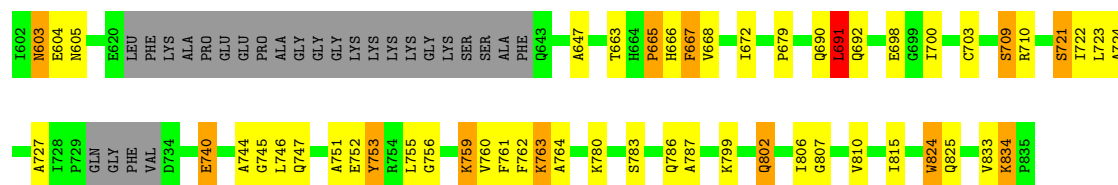
Note EDS was not executed.

• Molecule 1: MYOSIN HEAD



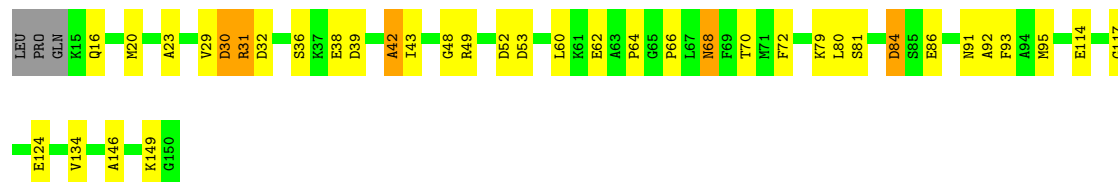
• Molecule 1: MYOSIN HEAD





• Molecule 2: MYOSIN HEAD

Chain Y: 71% 24% . .



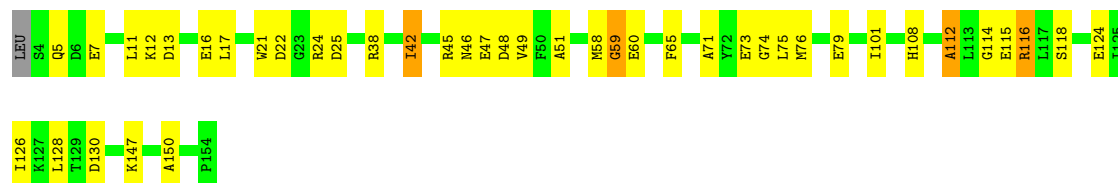
• Molecule 2: MYOSIN HEAD

Chain W: 70% 24% . .



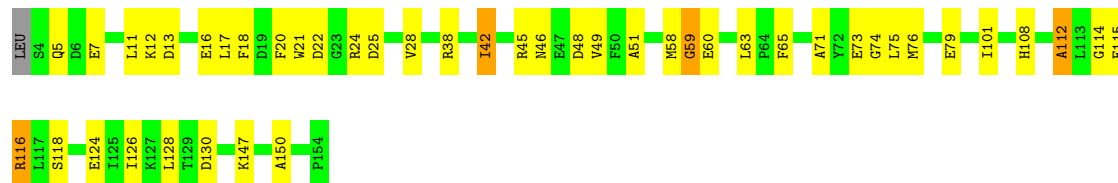
• Molecule 3: MYOSIN HEAD

Chain Z: 72% 25% . .



• Molecule 3: MYOSIN HEAD

Chain X: 70% 27% . .



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	53.03Å 243.47Å 124.65Å 90.00° 100.20° 90.00°	Depositor
Resolution (Å)	20.00 – 4.20	Depositor
% Data completeness (in resolution range)	95.0 (20.00-4.20)	Depositor
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.394 , 0.400	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	10500	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ADP, MG, CA, VO4

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	1.64	15/3809 (0.4%)	1.69	72/5293 (1.4%)
1	B	1.66	18/3778 (0.5%)	1.86	88/5248 (1.7%)
2	W	0.73	2/672 (0.3%)	1.69	12/933 (1.3%)
2	Y	0.89	2/672 (0.3%)	1.36	10/933 (1.1%)
3	X	0.58	0/740	1.32	10/1024 (1.0%)
3	Z	0.58	0/740	1.32	10/1024 (1.0%)
All	All	1.45	37/10411 (0.4%)	1.69	202/14455 (1.4%)

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	4
1	B	0	5
2	W	0	1
2	Y	0	1
All	All	0	11

The worst 5 of 37 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	462	ALA	C-N	47.04	1.92	1.33
1	A	709	SER	C-N	45.57	1.94	1.33
1	B	667	PHE	C-N	-43.92	0.75	1.33
1	B	709	SER	C-N	30.99	1.75	1.33
1	B	482	GLU	C-N	-27.83	0.94	1.33

The worst 5 of 202 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	824	TRP	O-C-N	-36.73	73.73	122.59
1	A	709	SER	O-C-N	34.79	166.03	123.24
1	B	709	SER	O-C-N	29.72	158.04	123.27
1	A	482	GLU	O-C-N	22.07	151.94	122.59
1	B	482	GLU	O-C-N	22.06	151.93	122.59

There are no chirality outliers.

5 of 11 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	115	TYR	Mainchain
1	A	691	LEU	Mainchain
1	A	76	SER	Mainchain
1	A	824	TRP	Mainchain
2	Y	84	ASP	Mainchain

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	3816	0	1691	88	0
1	B	3786	0	1675	89	0
2	W	673	0	302	22	0
2	Y	673	0	302	21	0
3	X	741	0	338	18	0
3	Z	741	0	338	18	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	W	1	0	0	0	0
4	Y	1	0	0	0	0
5	A	5	0	0	0	0
5	B	5	0	0	0	0
6	A	27	0	12	0	0
6	B	27	0	12	0	0
7	X	1	0	0	0	0
7	Z	1	0	0	0	0
All	All	10500	0	4670	252	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 17.

The worst 5 of 252 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:445:THR:C	1:B:446:LEU:N	1.68	1.52
1:B:667:PHE:C	1:B:668:VAL:CA	1.77	1.51
1:A:445:THR:C	1:A:446:LEU:N	1.68	1.49
1:B:667:PHE:CA	1:B:668:VAL:N	1.73	1.49
1:B:709:SER:C	1:B:710:ARG:N	1.75	1.45

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	758/831 (91%)	606 (80%)	114 (15%)	38 (5%)	1	16
1	B	750/831 (90%)	606 (81%)	108 (14%)	36 (5%)	2	17
2	W	134/139 (96%)	94 (70%)	34 (25%)	6 (4%)	2	18
2	Y	134/139 (96%)	95 (71%)	33 (25%)	6 (4%)	2	18
3	X	149/152 (98%)	104 (70%)	36 (24%)	9 (6%)	1	14
3	Z	149/152 (98%)	104 (70%)	36 (24%)	9 (6%)	1	14
All	All	2074/2244 (92%)	1609 (78%)	361 (17%)	104 (5%)	1	16

5 of 104 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	27	ALA
1	A	366	ARG
1	A	368	ARG

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	371	GLN
1	A	542	ALA

5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 10 ligands modelled in this entry, 6 are monoatomic - leaving 4 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	$\# Z > 2$
5	VO4	B	2998	6,4	0,4,4	-	-	-		
5	VO4	A	1998	6,4	0,4,4	-	-	-		
6	ADP	A	1999	4,5	28,29,29	0.95	0	43,45,45	0.84	1 (2%)
6	ADP	B	2999	4,5	28,29,29	0.95	1 (3%)	43,45,45	0.84	1 (2%)

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	ADP	A	1999	4,5	-	3/16/32/32	0/3/3/3
6	ADP	B	2999	4,5	-	3/16/32/32	0/3/3/3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	2999	ADP	C2-N1	2.01	1.37	1.33

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	1999	ADP	C2'-C3'-C4'	-2.13	98.49	102.61
6	B	2999	ADP	C2'-C3'-C4'	-2.13	98.50	102.61

There are no chirality outliers.

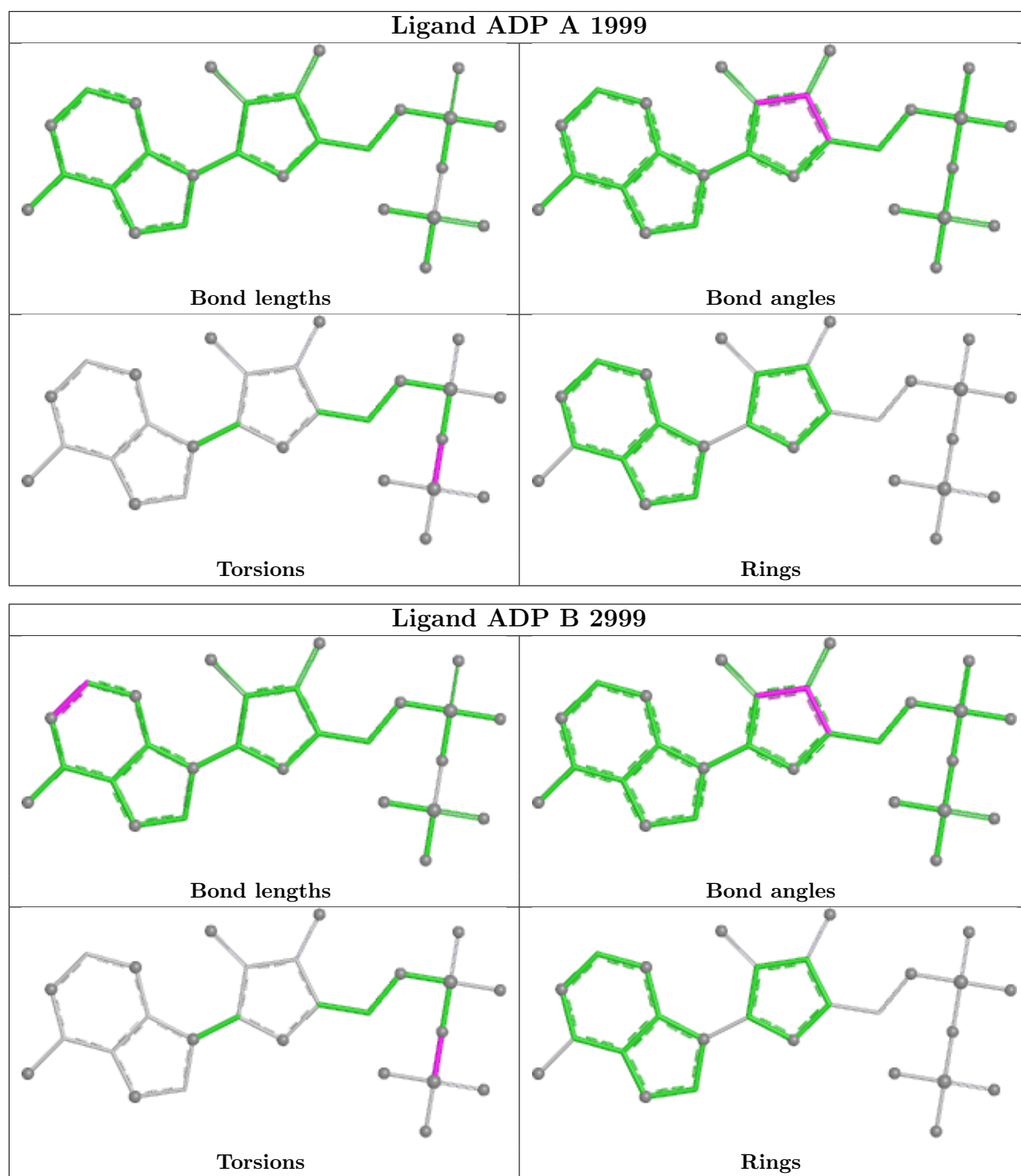
5 of 6 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	1999	ADP	PA-O3A-PB-O1B
6	B	2999	ADP	PA-O3A-PB-O1B
6	A	1999	ADP	PA-O3A-PB-O2B
6	A	1999	ADP	PA-O3A-PB-O3B
6	B	2999	ADP	PA-O3A-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 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	B	12
1	A	10

The worst 5 of 22 chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	461:ILE	C	462:ALA	N	2.12
1	B	461:ILE	C	462:ALA	N	2.12
1	A	709:SER	C	710:ARG	N	1.94
1	A	462:ALA	C	463:GLY	N	1.92
1	B	709:SER	C	710:ARG	N	1.75

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.