



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:13 PM UTC

PDB ID : 1W9L / pdb_00001w9l
Title : Myosin II Dictyostelium discoideum motor domain S456E bound with MgADP-AlF4
Authors : Morris, C.A.; Coureux, P.-D.; Wells, A.L.; Houdusse, A.; Sweeney, H.L.
Deposited on : 2004-10-13
Resolution : 1.95 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 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) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

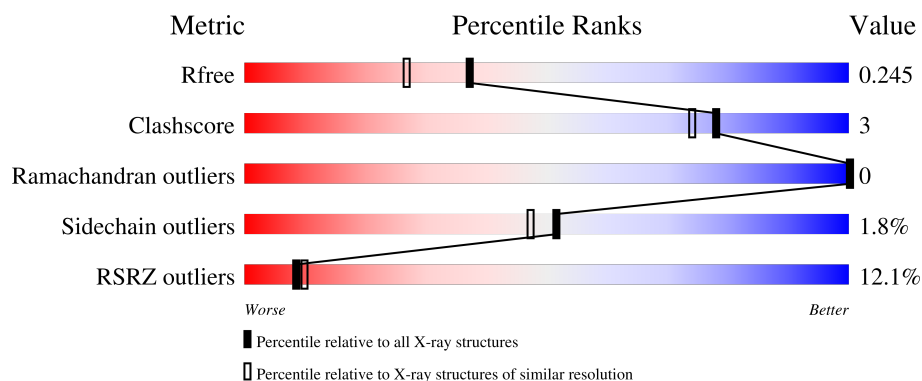
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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}	180053	3494 (1.96-1.96)
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)
RSRZ outliers	180081	3495 (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 $\leq 5\%$. 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	770	11% 88% 6% 6%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6228 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 II HEAVY CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	726	Total	C	N	O	S	0	1	0
			5736	3655	977	1088	16			

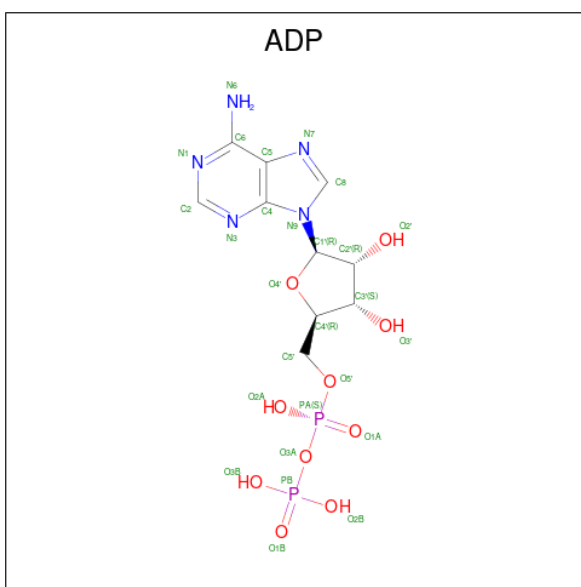
There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	249	SER	ASN	conflict	UNP P08799
A	312	CYS	TYR	conflict	UNP P08799
A	456	GLU	SER	engineered mutation	UNP P08799

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

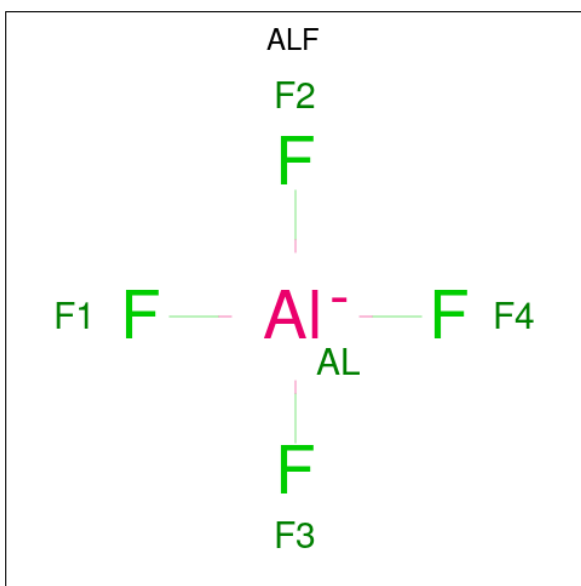
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mg	0	0
			1	1		

- Molecule 3 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: C₁₀H₁₅N₅O₁₀P₂).



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

- Molecule 4 is TETRAFLUOROALUMINATE ION (CCD ID: ALF) (formula: AlF_4).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total 5	Al 1	F 4	0	0

- Molecule 5 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $\text{C}_2\text{H}_6\text{O}_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		

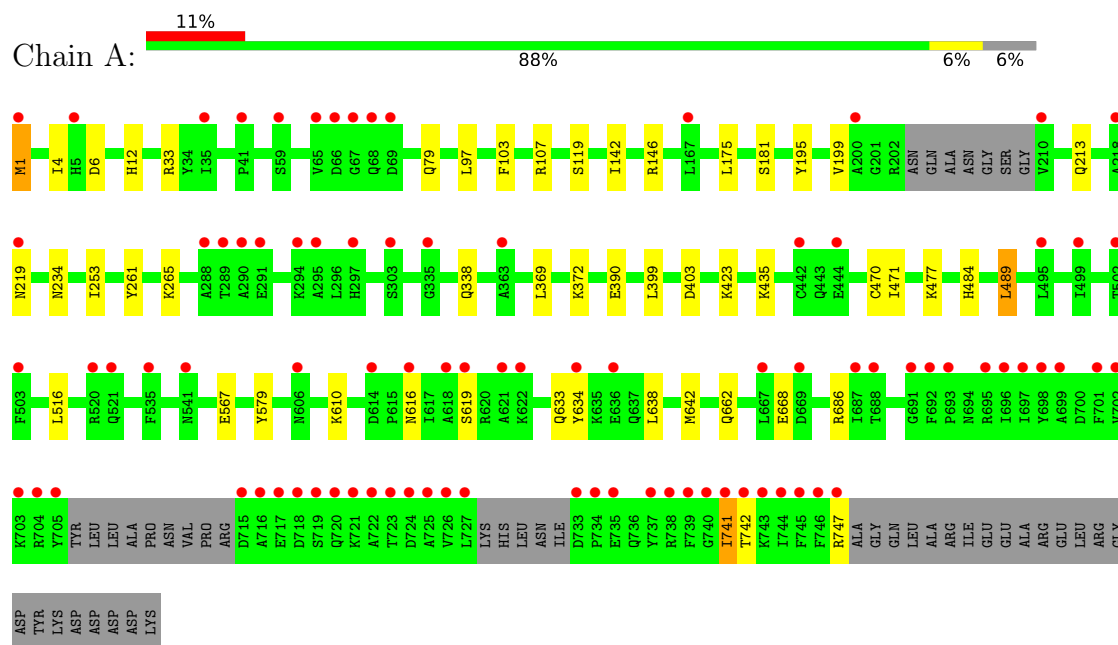
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	435	Total	O	0	0
			435	435		

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: MYOSIN II HEAVY CHAIN



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	88.04Å 145.98Å 152.37Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 1.95 40.00 – 1.95	Depositor EDS
% Data completeness (in resolution range)	97.8 (40.00-1.95) 97.8 (40.00-1.95)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.14 (at 1.95Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.193 , 0.225 (Not available) , 0.245	Depositor DCC
R_{free} test set	3525 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	28.9	Xtriage
Anisotropy	0.693	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 34.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.014 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.025 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6228	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.18% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.66	1/5850 (0.0%)	0.78	0/7910

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1	MET	SD-CE	-12.00	1.49	1.79

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	5736	0	5566	32	0
2	A	1	0	0	0	0
3	A	27	0	12	0	0
4	A	5	0	0	1	0
5	A	24	0	36	2	0
6	A	435	0	0	8	0
All	All	6228	0	5614	32	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 3.

All (32) 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:4:ILE:HD11	1:A:142:ILE:HG23	1.61	0.83
1:A:234:ASN:H	1:A:662:GLN:HE22	1.44	0.65
1:A:181:SER:HA	4:A:1750:ALF:F1	1.87	0.65
1:A:616:ASN:ND2	6:A:2371:HOH:O	2.30	0.62
1:A:1:MET:HE3	1:A:6:ASP:CB	2.38	0.53
1:A:103:PHE:CE2	1:A:107:ARG:NH2	2.78	0.52
1:A:103:PHE:HE2	1:A:107:ARG:NH2	2.08	0.51
1:A:265:LYS:HE3	1:A:423:LYS:HB3	1.94	0.49
1:A:97:LEU:HD23	1:A:686:ARG:HG3	1.95	0.48
1:A:261:TYR:C	6:A:2167:HOH:O	2.55	0.48
1:A:638:LEU:O	1:A:642:MET:HG2	2.14	0.48
1:A:33:ARG:HH21	1:A:79:GLN:HE22	1.61	0.47
1:A:741:ILE:HG13	1:A:742:THR:HG23	1.96	0.47
1:A:633:GLN:NE2	6:A:2381:HOH:O	2.45	0.47
1:A:1:MET:HE3	1:A:6:ASP:CG	2.39	0.47
1:A:477:LYS:HG3	1:A:638:LEU:HD21	1.97	0.47
1:A:668:GLU:OE1	6:A:2402:HOH:O	2.20	0.46
1:A:234:ASN:ND2	6:A:2151:HOH:O	2.50	0.44
1:A:747:ARG:C	6:A:2423:HOH:O	2.61	0.44
1:A:369:LEU:O	1:A:372:LYS:NZ	2.50	0.44
1:A:97:LEU:CD2	1:A:686:ARG:HG3	2.48	0.43
1:A:470:CYS:HB3	1:A:634[A]:TYR:CZ	2.54	0.43
1:A:399:LEU:HA	1:A:403:ASP:O	2.19	0.42
1:A:372:LYS:NZ	1:A:390:GLU:OE2	2.51	0.42
1:A:1:MET:HE3	1:A:6:ASP:HB2	2.01	0.42
1:A:516:LEU:HA	5:A:1755:EDO:H11	2.00	0.42
1:A:213:GLN:HG2	6:A:2135:HOH:O	2.19	0.42
1:A:195:TYR:CZ	1:A:199:VAL:HG11	2.55	0.42
1:A:567:GLU:HA	1:A:579:TYR:O	2.20	0.41
1:A:471:ILE:HG13	5:A:1752:EDO:H12	2.03	0.41
1:A:610:LYS:HD2	6:A:2369:HOH:O	2.21	0.40
1:A:484:HIS:CE1	1:A:489:LEU:HD13	2.56	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	719/770 (93%)	707 (98%)	12 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	608/673 (90%)	597 (98%)	11 (2%)	51	47

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	HIS
1	A	119	SER
1	A	146	ARG
1	A	175	LEU
1	A	219	ASN
1	A	253	ILE
1	A	338	GLN
1	A	435	LYS
1	A	489	LEU
1	A	619	SER
1	A	741	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	79	GLN
1	A	104	HIS
1	A	188	ASN
1	A	234	ASN
1	A	283	GLN
1	A	338	GLN
1	A	407	GLN
1	A	500	ASN
1	A	582	GLN
1	A	606	ASN
1	A	662	GLN
1	A	679	ASN

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 9 ligands modelled in this entry, 1 is monoatomic - leaving 8 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	ALF	A	1750	6,2,3	4,4,4	1.29	0	-		
3	ADP	A	1749	2,4	28,29,29	2.07	4 (14%)	43,45,45	1.97	11 (25%)
5	EDO	A	1756	-	3,3,3	0.28	0	2,2,2	0.35	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
5	EDO	A	1753	-	3,3,3	0.26	0	2,2,2	0.52	0
5	EDO	A	1752	-	3,3,3	0.25	0	2,2,2	0.51	0
5	EDO	A	1754	-	3,3,3	0.25	0	2,2,2	0.35	0
5	EDO	A	1755	-	3,3,3	0.26	0	2,2,2	0.46	0
5	EDO	A	1751	-	3,3,3	0.24	0	2,2,2	0.25	0

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	ADP	A	1749	2,4	-	1/16/32/32	0/3/3/3
5	EDO	A	1756	-	-	1/1/1/1	-
5	EDO	A	1753	-	-	0/1/1/1	-
5	EDO	A	1752	-	-	0/1/1/1	-
5	EDO	A	1754	-	-	0/1/1/1	-
5	EDO	A	1755	-	-	0/1/1/1	-
5	EDO	A	1751	-	-	0/1/1/1	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1749	ADP	C8-N7	6.84	1.44	1.31
3	A	1749	ADP	C8-N9	6.65	1.48	1.37
3	A	1749	ADP	PA-O1A	2.76	1.60	1.50
3	A	1749	ADP	C6-N6	2.28	1.40	1.34

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1749	ADP	N9-C8-N7	-6.13	105.24	113.94
3	A	1749	ADP	N3-C2-N1	-4.75	121.39	128.58
3	A	1749	ADP	C5-C4-N3	-3.67	121.66	126.72
3	A	1749	ADP	C5-C4-N9	3.37	109.48	105.81
3	A	1749	ADP	C2'-C1'-N9	-3.21	105.32	113.30
3	A	1749	ADP	C4-N9-C1'	-2.80	120.09	126.63
3	A	1749	ADP	C6-C5-C4	2.66	120.81	117.18
3	A	1749	ADP	O2B-PB-O1B	2.35	120.01	110.83
3	A	1749	ADP	C2-N1-C6	2.33	122.56	118.73
3	A	1749	ADP	C6-C5-N7	-2.23	127.78	132.09
3	A	1749	ADP	C2-N3-C4	2.20	117.21	111.83

There are no chirality outliers.

All (2) torsion outliers are listed below:

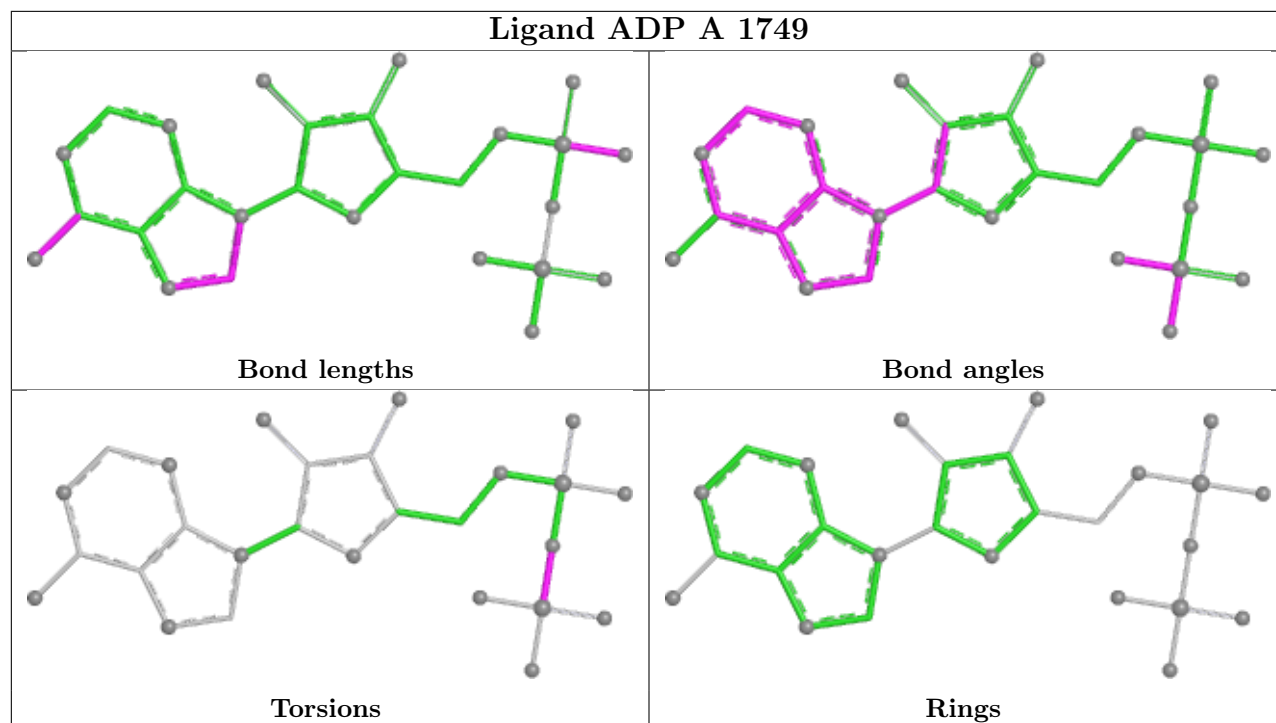
Mol	Chain	Res	Type	Atoms
3	A	1749	ADP	PA-O3A-PB-O3B
5	A	1756	EDO	O1-C1-C2-O2

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1750	ALF	1	0
5	A	1752	EDO	1	0
5	A	1755	EDO	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 ⓘ

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	A	726/770 (94%)	0.77	88 (12%) 8 10	16, 36, 65, 94	1 (0%)

All (88) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	65	VAL	6.9
1	A	704	ARG	6.8
1	A	67	GLY	6.4
1	A	727	LEU	6.1
1	A	722	ALA	5.5
1	A	702	VAL	5.4
1	A	718	ASP	5.4
1	A	725	ALA	5.4
1	A	726	VAL	5.2
1	A	715	ASP	5.0
1	A	741	ILE	5.0
1	A	703	LYS	5.0
1	A	696	ILE	4.8
1	A	716	ALA	4.7
1	A	719	SER	4.6
1	A	495	LEU	4.4
1	A	68	GLN	4.3
1	A	502	THR	4.3
1	A	442	CYS	4.2
1	A	705	TYR	4.0
1	A	746	PHE	3.9
1	A	744	ILE	3.8
1	A	210	VAL	3.8
1	A	616	ASN	3.8
1	A	739	PHE	3.8
1	A	737	TYR	3.7
1	A	717	GLU	3.6

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Mol	Chain	Res	Type	RSRZ
1	A	219	ASN	3.6
1	A	698	TYR	3.5
1	A	734	PRO	3.4
1	A	721	LYS	3.3
1	A	723	THR	3.2
1	A	692	PHE	3.2
1	A	521	GLN	3.1
1	A	614	ASP	3.1
1	A	499	ILE	3.1
1	A	363	ALA	3.1
1	A	687	ILE	3.0
1	A	747	ARG	3.0
1	A	745	PHE	3.0
1	A	200	ALA	3.0
1	A	291	GLU	3.0
1	A	697	ILE	2.9
1	A	699	ALA	2.9
1	A	66	ASP	2.8
1	A	303	SER	2.8
1	A	733	ASP	2.8
1	A	740	GLY	2.8
1	A	297	HIS	2.8
1	A	622	LYS	2.8
1	A	619	SER	2.7
1	A	720	GLN	2.7
1	A	742	THR	2.7
1	A	295	ALA	2.6
1	A	691	GLY	2.6
1	A	738	ARG	2.6
1	A	695	ARG	2.6
1	A	41	PRO	2.6
1	A	669	ASP	2.5
1	A	5	HIS	2.5
1	A	69	ASP	2.5
1	A	618	ALA	2.5
1	A	621	ALA	2.5
1	A	289	THR	2.4
1	A	636	GLU	2.4
1	A	167	LEU	2.4
1	A	693	PRO	2.4
1	A	735	GLU	2.4
1	A	541	ASN	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	667	LEU	2.4
1	A	724	ASP	2.4
1	A	1	MET	2.4
1	A	290	ALA	2.3
1	A	520	ARG	2.3
1	A	634[A]	TYR	2.3
1	A	503	PHE	2.3
1	A	294	LYS	2.3
1	A	701	PHE	2.2
1	A	688	THR	2.2
1	A	59	SER	2.2
1	A	218	ALA	2.2
1	A	535	PHE	2.2
1	A	743	LYS	2.2
1	A	288	ALA	2.2
1	A	335	GLY	2.1
1	A	606	ASN	2.1
1	A	444	GLU	2.0
1	A	35	ILE	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 oligosaccharides 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.

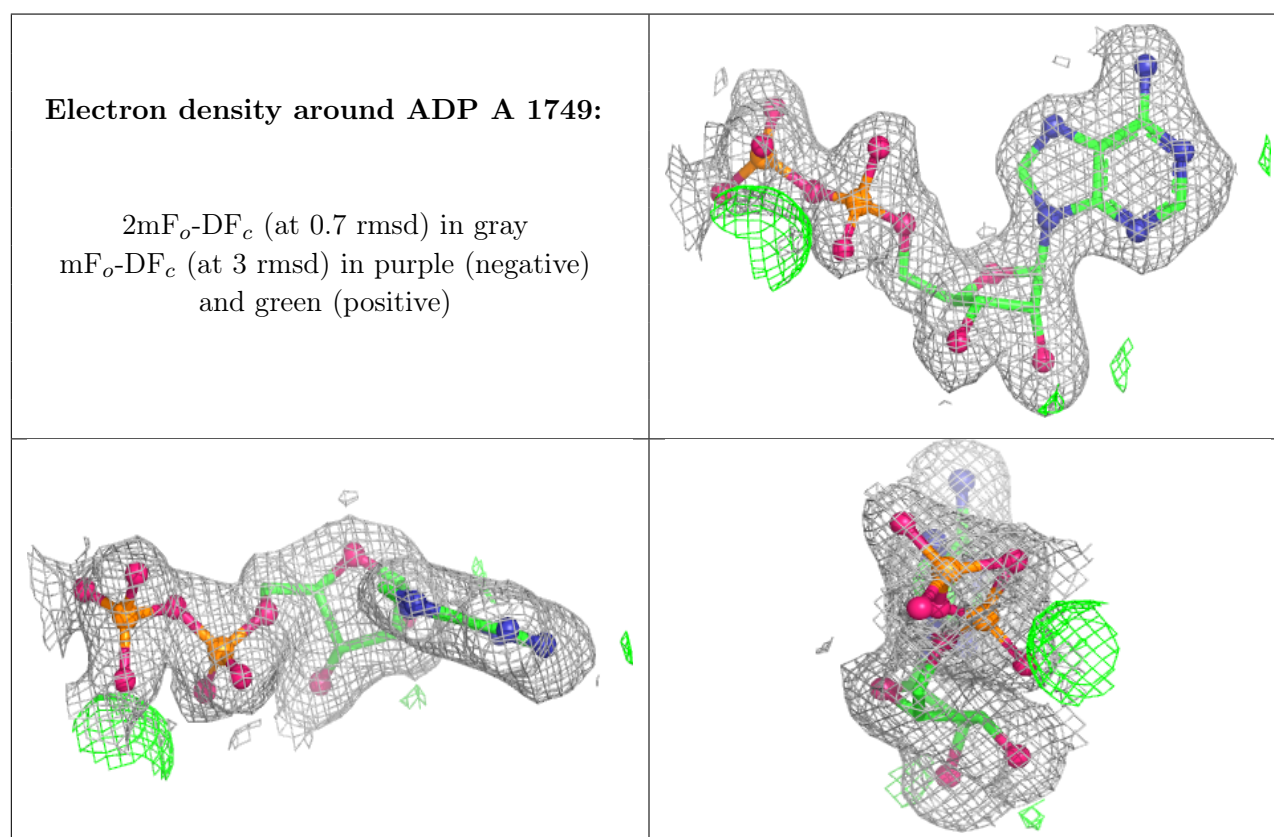
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	EDO	A	1756	4/4	0.76	0.19	69,70,70,71	0
5	EDO	A	1754	4/4	0.82	0.17	58,60,61,62	0
2	MG	A	1748	1/1	0.90	0.07	28,28,28,28	0
5	EDO	A	1755	4/4	0.90	0.13	58,58,58,60	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	EDO	A	1752	4/4	0.90	0.18	59,61,61,62	0
5	EDO	A	1751	4/4	0.96	0.09	47,48,48,48	0
5	EDO	A	1753	4/4	0.97	0.06	40,40,41,43	0
3	ADP	A	1749	27/27	0.98	0.06	22,29,31,31	0
4	ALF	A	1750	5/5	0.99	0.04	24,24,25,26	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.



6.5 Other polymers [i](#)

There are no such residues in this entry.