



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 02:17 AM UTC

PDB ID : 3DAW / pdb_00003daw
Title : Structure of the actin-depolymerizing factor homology domain in complex with actin
Authors : Paavilainen, V.O.; Oksanen, E.; Goldman, A.; Lappalainen, P.
Deposited on : 2008-05-30
Resolution : 2.55 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

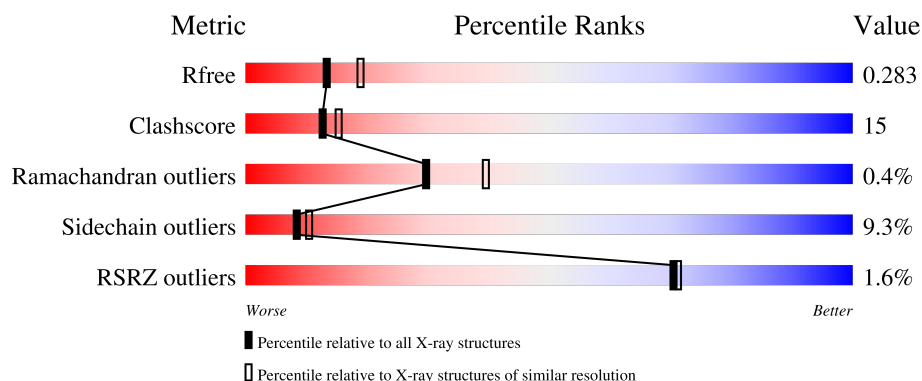
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

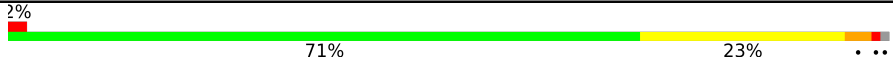
The reported resolution of this entry is 2.55 Å.

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	1091 (2.54-2.54)
Clashscore	190562	1120 (2.54-2.54)
Ramachandran outliers	187476	1106 (2.54-2.54)
Sidechain outliers	187428	1106 (2.54-2.54)
RSRZ outliers	180081	1091 (2.54-2.54)

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	377	 2% 71% 23% . .
2	B	164	 59% 24% . . 15%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4211 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	372	Total	C	N	O	S	0	0	0
			2901	1837	489	554	21			

- Molecule 2 is a protein called Twinfilin-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	140	Total	C	N	O	S	0	0	0
			1155	737	193	220	5			

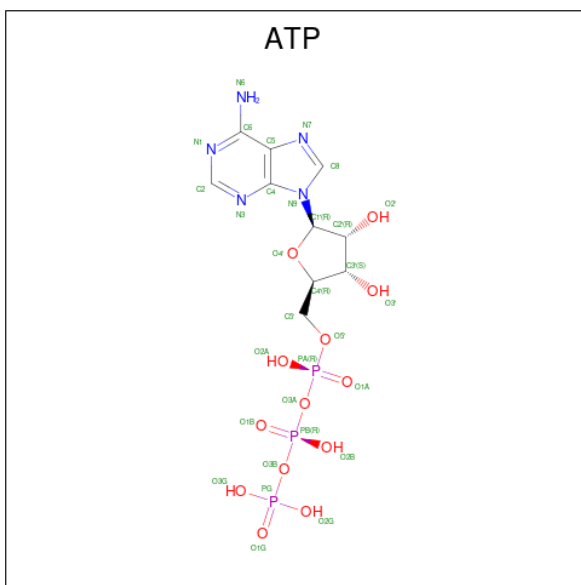
There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	159	MET	-	expression tag	UNP Q91YR1
B	160	SER	-	expression tag	UNP Q91YR1
B	161	HIS	-	expression tag	UNP Q91YR1
B	162	HIS	-	expression tag	UNP Q91YR1
B	163	HIS	-	expression tag	UNP Q91YR1
B	164	HIS	-	expression tag	UNP Q91YR1
B	165	HIS	-	expression tag	UNP Q91YR1
B	166	HIS	-	expression tag	UNP Q91YR1
B	168	MET	VAL	engineered mutation	UNP Q91YR1

- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Ca	0	0
			1	1		

- Molecule 4 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: C₁₀H₁₆N₅O₁₃P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			31	10	5	13	3		

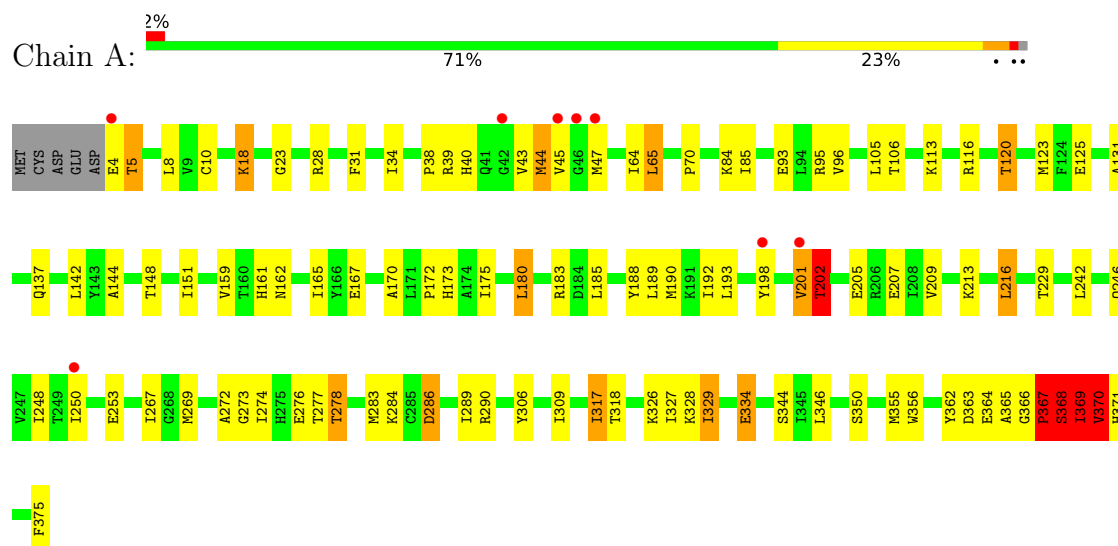
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	83	Total O 83 83	0	0
5	B	40	Total O 40 40	0	0

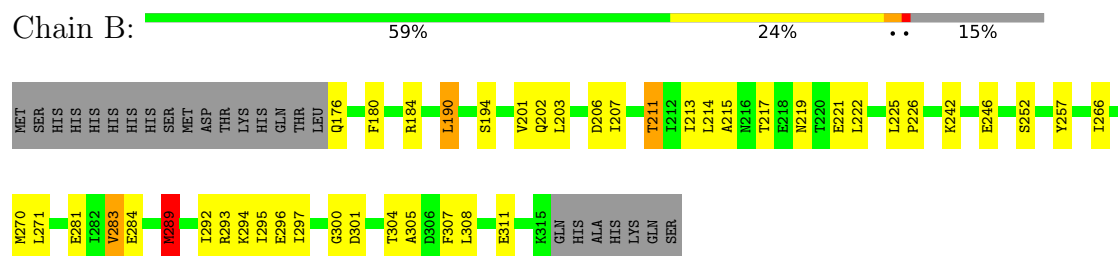
3 Residue-property plots

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: Actin, alpha skeletal muscle



- Molecule 2: Twinfilin-1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	52.78Å 72.99Å 168.88Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.76 – 2.55 42.76 – 2.55	Depositor EDS
% Data completeness (in resolution range)	99.3 (42.76-2.55) 99.3 (42.76-2.55)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	0.13	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.91 (at 2.54Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.208 , 0.279 0.212 , 0.283	Depositor DCC
R_{free} test set	1125 reflections (5.11%)	wwPDB-VP
Wilson B-factor (Å ²)	45.2	Xtriage
Anisotropy	0.127	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 40.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4211	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.41% 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: ATP, CA

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	0.74	0/2964	0.97	10/4016 (0.2%)
2	B	0.77	0/1178	0.94	2/1589 (0.1%)
All	All	0.75	0/4142	0.97	12/5605 (0.2%)

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	6
2	B	0	1
All	All	0	7

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	364	GLU	N-CA-C	8.74	122.69	112.72
1	A	202	THR	N-CA-C	-7.79	95.54	108.32
1	A	5	THR	N-CA-C	-7.25	98.49	110.17
1	A	368	SER	N-CA-CB	-6.63	99.28	110.49
1	A	368	SER	CB-CA-C	-5.96	98.57	110.42
1	A	286	ASP	CB-CA-C	-5.51	100.91	109.61
1	A	367	PRO	N-CA-C	-5.48	101.18	112.47
2	B	301	ASP	N-CA-C	-5.18	105.76	111.71
1	A	202	THR	N-CA-CB	5.15	119.27	110.57
1	A	369	ILE	N-CA-C	5.12	120.00	109.34
1	A	370	VAL	N-CA-C	5.10	115.82	110.62
2	B	289	MET	N-CA-C	-5.09	103.43	110.35

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	201	VAL	Peptide
1	A	367	PRO	Peptide
1	A	368	SER	Peptide
1	A	4	GLU	Peptide
1	A	40	HIS	Peptide
1	A	45	VAL	Peptide
2	B	300	GLY	Peptide

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	2901	0	2864	96	0
2	B	1155	0	1149	39	0
3	A	1	0	0	0	0
4	A	31	0	12	1	0
5	A	83	0	0	2	0
5	B	40	0	0	1	0
All	All	4211	0	4025	123	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 15.

All (123) 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:250:ILE:HD11	1:A:253:GLU:HB2	1.34	1.09
1:A:274:ILE:O	1:A:278:THR:HG23	1.54	1.07
2:B:206:ASP:HB3	2:B:211:THR:HG22	1.41	0.99
1:A:286:ASP:HB2	1:A:289:ILE:HG22	1.54	0.90
1:A:355:MET:HE2	5:A:410:HOH:O	1.75	0.85
1:A:193:LEU:HD21	1:A:250:ILE:HD13	1.59	0.83
1:A:229:THR:HG22	5:A:447:HOH:O	1.79	0.82
1:A:144:ALA:HA	2:B:270:MET:CE	2.11	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:120:THR:HG21	1:A:370:VAL:HG21	1.63	0.79
1:A:120:THR:HG21	1:A:370:VAL:CG2	2.14	0.77
1:A:274:ILE:O	1:A:278:THR:CG2	2.35	0.74
1:A:144:ALA:HA	2:B:270:MET:HE2	1.71	0.73
1:A:47:MET:HE2	1:A:65:LEU:HD12	1.72	0.71
1:A:144:ALA:CA	2:B:270:MET:CE	2.70	0.70
1:A:47:MET:CE	1:A:65:LEU:HD12	2.21	0.70
1:A:38:PRO:CG	1:A:47:MET:HE3	2.22	0.69
1:A:131:ALA:HB1	1:A:356:TRP:HB3	1.73	0.69
1:A:70:PRO:HG2	1:A:85:ILE:HD12	1.74	0.69
1:A:116:ARG:O	1:A:120:THR:HG23	1.91	0.69
1:A:242:LEU:HD12	1:A:246:GLN:HB3	1.75	0.69
1:A:202:THR:HG22	1:A:205:GLU:HB2	1.73	0.68
1:A:38:PRO:CA	1:A:47:MET:HE3	2.24	0.68
1:A:216:LEU:HD23	1:A:250:ILE:HG21	1.77	0.66
2:B:293:ARG:HD3	2:B:311:GLU:HB3	1.78	0.65
1:A:38:PRO:HG3	1:A:47:MET:HE3	1.79	0.65
2:B:202:GLN:HG2	2:B:215:ALA:HB3	1.79	0.64
1:A:250:ILE:HD11	1:A:253:GLU:CB	2.19	0.64
2:B:266:ILE:HD12	5:B:10:HOH:O	1.98	0.63
2:B:206:ASP:CB	2:B:211:THR:HG22	2.24	0.63
1:A:64:ILE:HG22	1:A:65:LEU:HD12	1.80	0.62
2:B:211:THR:HG23	2:B:213:ILE:HG23	1.82	0.61
1:A:368:SER:HB2	1:A:371:HIS:ND1	2.15	0.61
1:A:162:ASN:HD21	1:A:278:THR:HG22	1.65	0.61
1:A:31:PHE:CE1	1:A:93:GLU:HG3	2.36	0.60
1:A:173:HIS:NE2	1:A:284:LYS:O	2.27	0.60
2:B:304:THR:HG23	2:B:307:PHE:H	1.66	0.60
1:A:144:ALA:CA	2:B:270:MET:HE1	2.32	0.59
2:B:194:SER:HA	2:B:289:MET:HE2	1.83	0.59
1:A:64:ILE:HG22	1:A:65:LEU:CD1	2.32	0.59
1:A:198:TYR:CE1	1:A:248:ILE:HG22	2.37	0.59
1:A:144:ALA:HB2	2:B:270:MET:HE1	1.84	0.59
1:A:162:ASN:ND2	1:A:277:THR:OG1	2.36	0.59
2:B:214:LEU:HD11	2:B:217:THR:HG23	1.83	0.59
1:A:47:MET:HE2	1:A:64:ILE:HG22	1.85	0.57
1:A:202:THR:HG22	1:A:205:GLU:CB	2.34	0.57
1:A:120:THR:CG2	1:A:370:VAL:HG21	2.34	0.57
1:A:38:PRO:HA	1:A:47:MET:HE3	1.85	0.57
1:A:365:ALA:HB1	1:A:366:GLY:CA	2.35	0.56
2:B:222:LEU:HD21	2:B:305:ALA:HA	1.86	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:368:SER:HA	1:A:370:VAL:H	1.71	0.55
2:B:190:LEU:HD13	2:B:201:VAL:HG11	1.87	0.55
2:B:295:ILE:HG22	2:B:297:ILE:HG23	1.88	0.55
1:A:317:ILE:HG22	1:A:327:ILE:HD13	1.89	0.54
1:A:334:GLU:HG2	2:B:266:ILE:HD11	1.88	0.54
1:A:180:LEU:HB2	1:A:269:MET:HE1	1.90	0.54
1:A:185:LEU:HD23	1:A:306:TYR:OH	2.09	0.53
1:A:172:PRO:HA	1:A:175:ILE:HD12	1.90	0.53
1:A:44:MET:HB2	1:A:47:MET:CG	2.40	0.52
1:A:105:LEU:HD11	1:A:123:MET:HE3	1.91	0.51
1:A:273:GLY:O	1:A:277:THR:HG23	2.11	0.51
2:B:283:VAL:HG13	2:B:289:MET:CB	2.40	0.51
1:A:193:LEU:CD2	1:A:250:ILE:HD13	2.37	0.51
2:B:194:SER:CB	2:B:289:MET:CE	2.89	0.50
1:A:362:TYR:HE1	1:A:367:PRO:HB3	1.77	0.50
1:A:162:ASN:ND2	1:A:278:THR:HG22	2.26	0.49
2:B:225:LEU:N	2:B:226:PRO:CD	2.75	0.49
1:A:318:THR:HA	1:A:327:ILE:CD1	2.43	0.49
1:A:362:TYR:O	1:A:366:GLY:HA2	2.13	0.48
1:A:144:ALA:CB	2:B:270:MET:HE1	2.43	0.48
1:A:38:PRO:HA	1:A:47:MET:CE	2.43	0.48
1:A:272:ALA:HB1	1:A:276:GLU:HB2	1.97	0.47
1:A:38:PRO:HB3	1:A:44:MET:HG3	1.95	0.47
1:A:198:TYR:CE1	1:A:248:ILE:CG2	2.97	0.47
2:B:221:GLU:O	2:B:222:LEU:C	2.56	0.47
1:A:306:TYR:O	1:A:309:ILE:HG22	2.14	0.47
1:A:180:LEU:HD22	1:A:269:MET:CE	2.45	0.46
1:A:188:TYR:CD1	1:A:267:ILE:HG22	2.50	0.46
1:A:39:ARG:HD2	1:A:65:LEU:O	2.16	0.46
1:A:328:LYS:C	1:A:329:ILE:HD12	2.41	0.46
1:A:329:ILE:HD12	1:A:329:ILE:N	2.31	0.46
2:B:270:MET:HE2	2:B:270:MET:HA	1.97	0.46
1:A:142:LEU:HD22	1:A:165:ILE:HB	1.97	0.45
2:B:194:SER:HB3	2:B:289:MET:CE	2.46	0.45
1:A:47:MET:CE	1:A:65:LEU:CD1	2.93	0.45
1:A:47:MET:HE1	1:A:65:LEU:HD12	1.97	0.45
1:A:190:MET:HG2	1:A:209:VAL:HG21	1.98	0.45
1:A:38:PRO:CB	1:A:47:MET:HE3	2.46	0.45
1:A:144:ALA:N	2:B:270:MET:HE1	2.32	0.45
1:A:106:THR:HB	1:A:137:GLN:HG3	1.99	0.44
1:A:144:ALA:HB2	2:B:270:MET:CE	2.48	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:216:LEU:HD23	1:A:250:ILE:CG2	2.45	0.44
2:B:194:SER:CA	2:B:289:MET:CE	2.95	0.44
2:B:180:PHE:CZ	2:B:271:LEU:HD22	2.52	0.44
1:A:23:GLY:N	1:A:344:SER:OG	2.46	0.44
1:A:170:ALA:O	1:A:172:PRO:HD3	2.18	0.44
1:A:18:LYS:NZ	4:A:377:ATP:O2B	2.51	0.43
1:A:38:PRO:CD	1:A:47:MET:HE3	2.48	0.43
1:A:366:GLY:O	1:A:369:ILE:HG22	2.18	0.43
2:B:214:LEU:HD11	2:B:217:THR:CG2	2.49	0.43
2:B:283:VAL:HG13	2:B:289:MET:HB3	2.00	0.43
1:A:202:THR:HG22	1:A:205:GLU:H	1.83	0.42
1:A:189:LEU:HD23	1:A:209:VAL:CG1	2.49	0.42
1:A:192:ILE:HD12	1:A:253:GLU:HG3	2.01	0.42
1:A:329:ILE:N	1:A:329:ILE:CD1	2.83	0.42
1:A:159:VAL:HG21	1:A:161:HIS:CE1	2.55	0.42
2:B:283:VAL:HG13	2:B:289:MET:HB2	2.01	0.42
1:A:283:MET:HE1	1:A:290:ARG:NH1	2.34	0.42
1:A:363:ASP:O	1:A:365:ALA:HB2	2.19	0.41
2:B:194:SER:HA	2:B:289:MET:CE	2.49	0.41
1:A:189:LEU:HD22	1:A:213:LYS:HB2	2.02	0.41
2:B:284:GLU:OE2	2:B:289:MET:O	2.39	0.41
1:A:365:ALA:HB1	1:A:366:GLY:HA2	2.01	0.41
1:A:148:THR:OG1	2:B:296:GLU:OE1	2.26	0.41
1:A:8:LEU:HD11	1:A:96:VAL:HG21	2.03	0.41
1:A:47:MET:HE1	1:A:65:LEU:CD1	2.49	0.41
1:A:113:LYS:HB3	1:A:371:HIS:HE2	1.86	0.41
1:A:180:LEU:HD22	1:A:269:MET:HE3	2.02	0.41
1:A:151:ILE:HG23	1:A:151:ILE:O	2.21	0.41
1:A:286:ASP:HB2	1:A:289:ILE:CG2	2.39	0.40
1:A:144:ALA:CB	2:B:270:MET:CE	2.99	0.40
1:A:375:PHE:O	2:B:176:GLN:NE2	2.55	0.40
2:B:257:TYR:HB3	2:B:296:GLU:HG2	2.04	0.40
2:B:308:LEU:HD12	2:B:308:LEU:HA	1.96	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	370/377 (98%)	345 (93%)	24 (6%)	1 (0%)	36	45
2	B	138/164 (84%)	130 (94%)	7 (5%)	1 (1%)	18	25
All	All	508/541 (94%)	475 (94%)	31 (6%)	2 (0%)	30	39

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	369	ILE
2	B	246	GLU

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	313/320 (98%)	285 (91%)	28 (9%)	9	12
2	B	130/153 (85%)	117 (90%)	13 (10%)	7	9
All	All	443/473 (94%)	402 (91%)	41 (9%)	8	10

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	THR
1	A	10	CYS
1	A	18	LYS
1	A	28	ARG

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Mol	Chain	Res	Type
1	A	34	ILE
1	A	43	VAL
1	A	44	MET
1	A	65	LEU
1	A	84	LYS
1	A	95	ARG
1	A	120	THR
1	A	125	GLU
1	A	167	GLU
1	A	180	LEU
1	A	183	ARG
1	A	201	VAL
1	A	202	THR
1	A	207	GLU
1	A	216	LEU
1	A	278	THR
1	A	317	ILE
1	A	326	LYS
1	A	329	ILE
1	A	334	GLU
1	A	346	LEU
1	A	350	SER
1	A	369	ILE
1	A	370	VAL
2	B	184	ARG
2	B	190	LEU
2	B	203	LEU
2	B	207	ILE
2	B	211	THR
2	B	219	ASN
2	B	242	LYS
2	B	252	SER
2	B	281	GLU
2	B	283	VAL
2	B	289	MET
2	B	292	ILE
2	B	294	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	128	ASN

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Mol	Chain	Res	Type
1	A	161	HIS
2	B	176	GLN
2	B	188	GLN
2	B	245	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 1 is 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$
4	ATP	A	377	3	32,33,33	1.47	6 (18%)	48,52,52	1.87	13 (27%)

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	ATP	A	377	3	-	3/22/38/38	0/3/3/3

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	377	ATP	C5-C4	4.40	1.46	1.39
4	A	377	ATP	PB-O3A	2.80	1.62	1.59
4	A	377	ATP	C4-N9	-2.73	1.32	1.37
4	A	377	ATP	C5-C6	2.57	1.48	1.41
4	A	377	ATP	PB-O3B	2.42	1.62	1.59
4	A	377	ATP	C8-N7	2.34	1.36	1.31

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	377	ATP	C5-C4-N3	-4.82	120.09	126.72
4	A	377	ATP	N3-C4-N9	4.33	134.53	127.17
4	A	377	ATP	C4-N9-C8	4.30	110.25	105.74
4	A	377	ATP	N3-C2-N1	-3.91	122.66	128.58
4	A	377	ATP	C2-N3-C4	3.62	120.67	111.83
4	A	377	ATP	O2B-PB-O3B	3.08	115.59	107.27
4	A	377	ATP	C4-C5-N7	-2.82	107.36	110.58
4	A	377	ATP	N9-C8-N7	-2.64	110.20	113.94
4	A	377	ATP	C6-C5-N7	2.49	136.90	132.09
4	A	377	ATP	C2'-C1'-N9	-2.28	107.65	113.30
4	A	377	ATP	C2-N1-C6	2.21	122.36	118.73
4	A	377	ATP	C5-N7-C8	2.14	106.82	103.45
4	A	377	ATP	O3G-PG-O2G	2.14	115.81	107.80

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	377	ATP	PB-O3B-PG-O1G
4	A	377	ATP	PG-O3B-PB-O1B
4	A	377	ATP	PG-O3B-PB-O2B

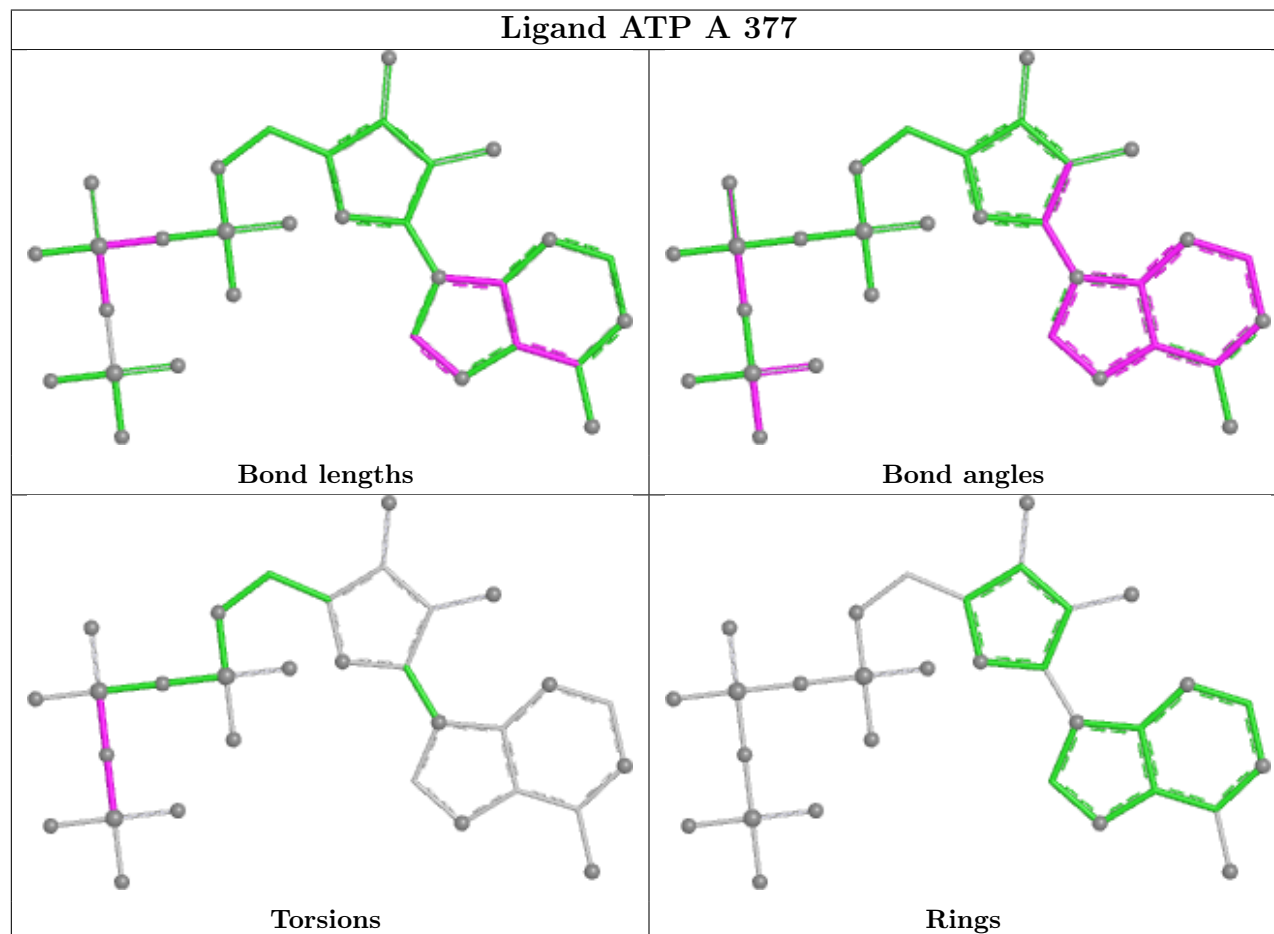
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	377	ATP	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	372/377 (98%)	0.10	8 (2%) 62 63	30, 47, 73, 77	0
2	B	140/164 (85%)	-0.08	0 100 100	33, 43, 52, 62	0
All	All	512/541 (94%)	0.05	8 (1%) 70 71	30, 45, 70, 77	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	45	VAL	3.9
1	A	46	GLY	3.4
1	A	42	GLY	3.4
1	A	4	GLU	2.4
1	A	250	ILE	2.3
1	A	201	VAL	2.1
1	A	47	MET	2.1
1	A	198	TYR	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.

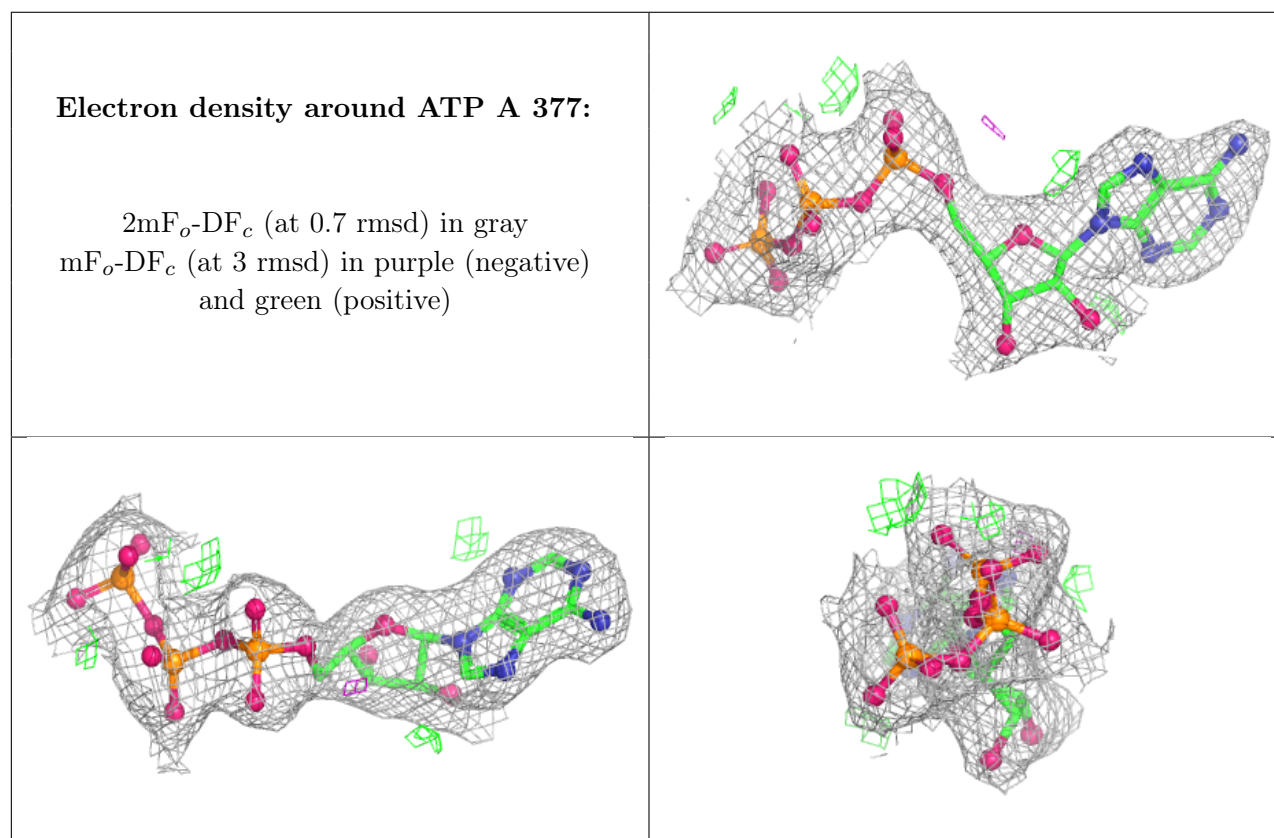
6.4 Ligands [i](#)

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(Å ²)	Q<0.9
3	CA	A	376	1/1	0.98	0.04	33,33,33,33	0
4	ATP	A	377	31/31	0.98	0.06	27,34,36,39	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.