



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 07:21 AM UTC

PDB ID : 3FCC / pdb_00003fcc
Title : CRYSTAL STRUCTURE OF DLTA PROTEIN IN COMPLEX WITH ATP
and MAGNESIUM
Authors : Osman, K.T.; Du, L.; He, Y.; Luo, Y.
Deposited on : 2008-11-21
Resolution : 2.32 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)
Xtrriage (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

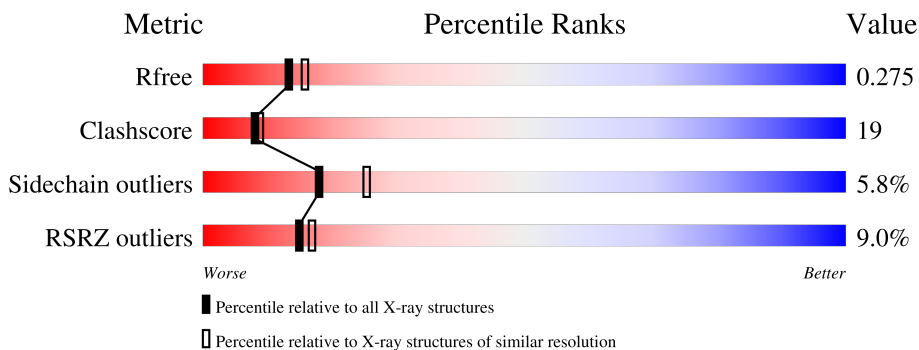
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.32 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	7754 (2.34-2.30)
Clashscore	190562	8383 (2.34-2.30)
Sidechain outliers	187428	8303 (2.34-2.30)
RSRZ outliers	180081	7760 (2.34-2.30)

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	512	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4108 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 D-alanine--poly(phosphoribitol) ligase subunit 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	500	3943	2531	640	749	23	0	0	0

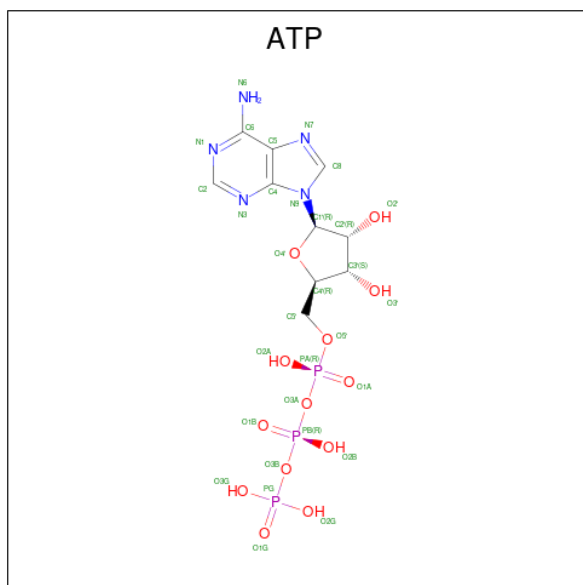
There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	505	LEU	-	expression tag	UNP Q81G39
A	506	GLU	-	expression tag	UNP Q81G39
A	507	HIS	-	expression tag	UNP Q81G39
A	508	HIS	-	expression tag	UNP Q81G39
A	509	HIS	-	expression tag	UNP Q81G39
A	510	HIS	-	expression tag	UNP Q81G39
A	511	HIS	-	expression tag	UNP Q81G39
A	512	HIS	-	expression tag	UNP Q81G39

- 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'-TRIPHOSPHATE (CCD ID: ATP) (formula: C₁₀H₁₆N₅O₁₃P₃).



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

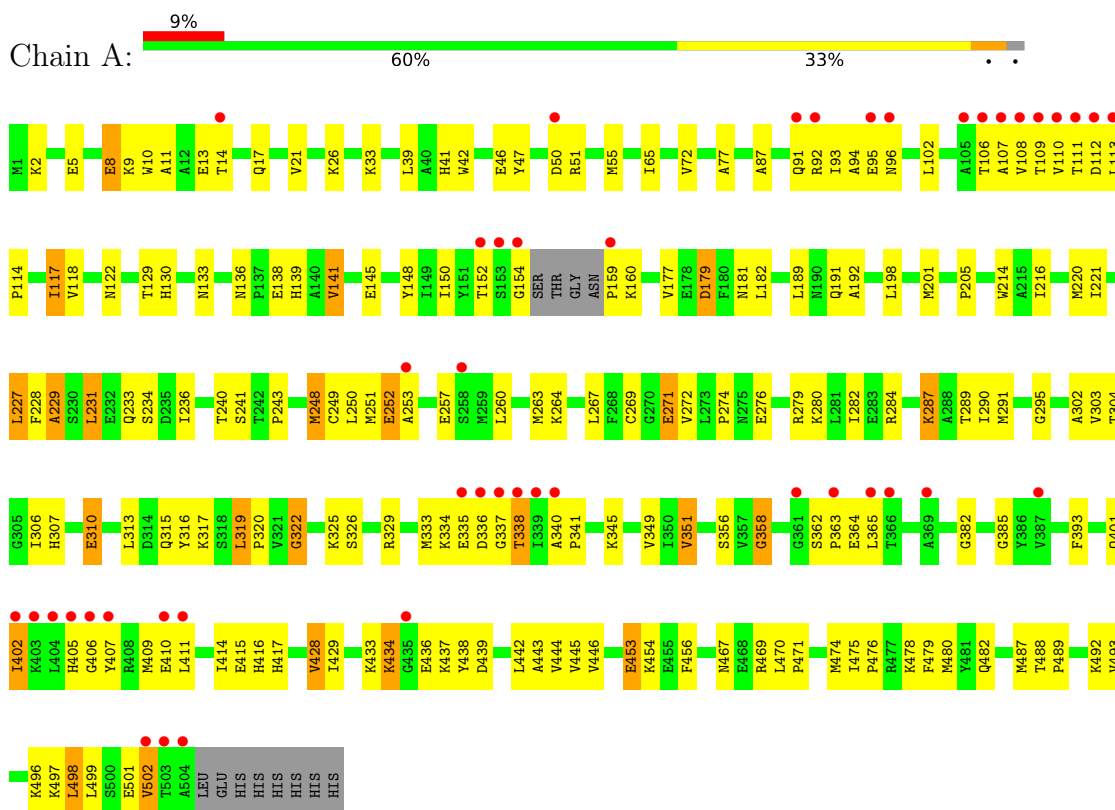
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	133	Total	O	0	0
			133	133		

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: D-alanine--poly(phosphoribitol) ligase subunit 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	61.60Å 87.00Å 60.10Å 90.00° 115.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.32 20.00 – 2.32	Depositor EDS
% Data completeness (in resolution range)	95.4 (20.00-2.32) 98.7 (20.00-2.32)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.32 (at 2.23Å)	Xtrriage
Refinement program	CNS	Depositor
R, R_{free}	0.241 , 0.286 0.231 , 0.275	Depositor DCC
R_{free} test set	1291 reflections (4.84%)	wwPDB-VP
Wilson B-factor (Å ²)	29.5	Xtrriage
Anisotropy	0.305	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 40.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.026 for l,-k,h	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4108	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.78% 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, ATP

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	0.43	0/4034	0.94	16/5469 (0.3%)

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	322	GLY	N-CA-C	7.39	121.17	111.70
1	A	326	SER	N-CA-C	7.14	121.09	112.38
1	A	502	VAL	N-CA-C	-6.44	103.27	112.35
1	A	382	GLY	N-CA-C	-6.17	106.72	115.30
1	A	358	GLY	N-CA-C	6.16	117.80	111.95
1	A	117	ILE	N-CA-C	5.56	115.80	107.51
1	A	446	VAL	N-CA-C	-5.53	102.07	107.55
1	A	319	LEU	N-CA-C	5.52	116.84	109.83
1	A	192	ALA	N-CA-C	5.34	117.35	109.50
1	A	109	THR	N-CA-C	5.33	116.89	111.14
1	A	253	ALA	N-CA-C	5.29	117.46	111.11
1	A	11	ALA	N-CA-C	-5.22	105.76	111.82
1	A	77	ALA	N-CA-C	-5.13	103.62	110.55
1	A	229	ALA	N-CA-C	-5.09	105.74	111.28
1	A	179	ASP	N-CA-C	5.03	119.26	113.18
1	A	310	GLU	N-CA-C	-5.02	105.81	111.28

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	3943	0	3922	154	0
2	A	1	0	0	0	0
3	A	31	0	12	0	0
4	A	133	0	0	7	0
All	All	4108	0	3934	154	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 19.

All (154) 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:152:THR:HG22	1:A:160:LYS:H	1.15	1.11
1:A:333:MET:HE2	1:A:349:VAL:HG21	1.42	0.98
1:A:201:MET:HE3	1:A:267:LEU:HB3	1.49	0.92
1:A:136:ASN:HB3	1:A:139:HIS:CD2	2.17	0.80
1:A:92:ARG:NH2	1:A:96:ASN:HD22	1.81	0.78
1:A:2:LYS:HB2	1:A:5:GLU:HG3	1.66	0.77
1:A:191:GLN:HE21	1:A:241:SER:HB2	1.49	0.75
1:A:152:THR:HG22	1:A:160:LYS:N	1.99	0.74
1:A:480:MET:HE3	1:A:482:GLN:HE21	1.53	0.73
1:A:402:ILE:H	1:A:402:ILE:HD13	1.52	0.72
1:A:152:THR:HG21	1:A:160:LYS:CE	2.19	0.72
1:A:336:ASP:CG	1:A:337:GLY:H	1.98	0.72
1:A:444:VAL:HG22	1:A:480:MET:HE2	1.71	0.71
1:A:497:LYS:O	1:A:501:GLU:HG2	1.90	0.70
1:A:152:THR:HG21	1:A:160:LYS:HE2	1.74	0.70
1:A:336:ASP:C	1:A:338:THR:H	1.97	0.70
1:A:228:PHE:CE2	1:A:252:GLU:HG3	2.28	0.69
1:A:401:GLN:HG2	1:A:410:GLU:HG3	1.75	0.69
1:A:233:GLN:HG2	4:A:849:HOH:O	1.93	0.68
1:A:152:THR:CG2	1:A:160:LYS:H	2.02	0.68
1:A:428:VAL:HG11	1:A:498:LEU:HD21	1.75	0.68
1:A:454:LYS:HD3	1:A:456:PHE:HE1	1.58	0.67
1:A:471:PRO:HD2	1:A:474:MET:SD	2.35	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:334:LYS:HG3	1:A:340:ALA:HB2	1.77	0.66
1:A:201:MET:HE3	1:A:267:LEU:CB	2.23	0.65
1:A:498:LEU:O	1:A:502:VAL:HG13	1.97	0.64
1:A:433:LYS:HB3	1:A:438:TYR:CD2	2.32	0.64
1:A:480:MET:HE3	1:A:482:GLN:NE2	2.12	0.64
1:A:14:THR:OG1	1:A:17:GLN:HB2	1.97	0.63
1:A:434:LYS:O	1:A:434:LYS:HD2	1.99	0.63
1:A:189:LEU:HB2	1:A:236:ILE:HD13	1.81	0.63
1:A:92:ARG:HH22	1:A:96:ASN:HD22	1.45	0.62
1:A:282:ILE:HG13	1:A:290:ILE:HD12	1.81	0.61
1:A:108:VAL:HG12	4:A:825:HOH:O	1.99	0.61
1:A:201:MET:HE2	1:A:240:THR:HG23	1.83	0.61
1:A:227:LEU:HD22	1:A:231:LEU:HD22	1.82	0.61
1:A:14:THR:HG21	4:A:846:HOH:O	2.02	0.60
1:A:9:LYS:O	1:A:13:GLU:HG2	2.01	0.60
1:A:434:LYS:HD2	1:A:434:LYS:C	2.26	0.60
1:A:433:LYS:HB3	1:A:438:TYR:HD2	1.67	0.59
1:A:50:ASP:O	1:A:51:ARG:NH1	2.35	0.59
1:A:493:VAL:HG22	4:A:874:HOH:O	2.03	0.59
1:A:336:ASP:C	1:A:338:THR:N	2.61	0.59
1:A:428:VAL:HG11	1:A:498:LEU:CD2	2.33	0.59
1:A:41:HIS:CD2	1:A:133:ASN:HB3	2.38	0.58
1:A:442:LEU:HD12	1:A:443:ALA:N	2.19	0.58
1:A:336:ASP:CG	1:A:337:GLY:N	2.62	0.58
1:A:201:MET:CE	1:A:267:LEU:HB3	2.28	0.57
1:A:229:ALA:O	1:A:233:GLN:HG3	2.04	0.57
1:A:108:VAL:HG11	1:A:117:ILE:HG21	1.87	0.57
1:A:417:HIS:ND1	1:A:470:LEU:HD21	2.18	0.57
1:A:499:LEU:O	1:A:502:VAL:HG22	2.05	0.57
1:A:417:HIS:HB3	1:A:469:ARG:HD2	1.86	0.57
1:A:251:MET:HE3	4:A:823:HOH:O	2.05	0.57
1:A:272:VAL:O	1:A:274:PRO:HD3	2.07	0.55
1:A:415:GLU:HG3	1:A:429:ILE:HD12	1.89	0.55
1:A:106:THR:OG1	1:A:107:ALA:N	2.38	0.54
1:A:243:PRO:HD2	1:A:271:GLU:HG3	1.89	0.54
1:A:65:ILE:HD12	1:A:65:ILE:N	2.23	0.54
1:A:402:ILE:HD13	1:A:409:MET:O	2.07	0.54
1:A:95:GLU:HG2	1:A:114:PRO:HG3	1.89	0.54
1:A:152:THR:HG21	1:A:160:LYS:HE3	1.89	0.53
1:A:428:VAL:HG13	1:A:444:VAL:HB	1.89	0.53
1:A:250:LEU:C	1:A:252:GLU:H	2.17	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:453:GLU:H	1:A:453:GLU:CD	2.17	0.53
1:A:434:LYS:HB3	1:A:439:ASP:HB2	1.91	0.53
1:A:333:MET:HE2	1:A:349:VAL:CG2	2.28	0.53
1:A:402:ILE:HD11	1:A:409:MET:HG3	1.89	0.53
1:A:414:ILE:HG22	1:A:429:ILE:HD11	1.91	0.52
1:A:313:LEU:HD23	1:A:319:LEU:HD21	1.91	0.52
1:A:476:PRO:HG2	1:A:479:PHE:CE1	2.45	0.52
1:A:234:SER:HB2	1:A:236:ILE:HG13	1.91	0.52
1:A:289:THR:HG23	1:A:307:HIS:CE1	2.44	0.52
1:A:317:LYS:HG3	1:A:416:HIS:CE1	2.45	0.52
1:A:306:ILE:HB	1:A:322:GLY:HA2	1.91	0.51
1:A:10:TRP:HE3	1:A:14:THR:CG2	2.24	0.50
1:A:14:THR:HG23	1:A:14:THR:O	2.11	0.50
1:A:414:ILE:HG22	1:A:429:ILE:CD1	2.42	0.49
1:A:362:SER:N	1:A:363:PRO:HD3	2.26	0.49
1:A:55:MET:HE3	1:A:93:ILE:HG22	1.94	0.49
1:A:154:GLY:HA2	1:A:159:PRO:N	2.28	0.49
1:A:287:LYS:HE3	1:A:287:LYS:HA	1.95	0.49
1:A:329:ARG:HG2	1:A:351:VAL:HG22	1.96	0.48
1:A:433:LYS:HB2	1:A:437:LYS:O	2.13	0.48
1:A:129:THR:HG22	1:A:130:HIS:ND1	2.27	0.48
1:A:303:VAL:HG23	1:A:304:THR:HG23	1.94	0.48
1:A:250:LEU:C	1:A:252:GLU:N	2.72	0.48
1:A:362:SER:C	1:A:364:GLU:N	2.71	0.48
1:A:405:HIS:C	1:A:407:TYR:N	2.70	0.48
1:A:42:TRP:HD1	1:A:130:HIS:CE1	2.31	0.48
1:A:177:VAL:HA	1:A:182:LEU:HD12	1.96	0.48
1:A:362:SER:HB3	1:A:365:LEU:HG	1.96	0.48
1:A:118:VAL:HG13	1:A:122:ASN:HB3	1.96	0.48
1:A:405:HIS:O	1:A:407:TYR:N	2.47	0.48
1:A:92:ARG:HH21	1:A:92:ARG:HG3	1.79	0.47
1:A:248:MET:HG3	1:A:249:CYS:N	2.30	0.47
1:A:362:SER:C	1:A:364:GLU:H	2.22	0.47
1:A:111:THR:O	1:A:112:ASP:C	2.58	0.47
1:A:138:GLU:O	1:A:138:GLU:HG2	2.14	0.47
1:A:179:ASP:HB3	4:A:920:HOH:O	2.14	0.47
1:A:130:HIS:HA	1:A:133:ASN:ND2	2.30	0.46
1:A:189:LEU:CD1	1:A:216:ILE:HD13	2.45	0.46
1:A:214:TRP:CZ3	1:A:234:SER:HA	2.50	0.46
1:A:320:PRO:HB3	1:A:393:PHE:CE2	2.50	0.46
1:A:95:GLU:CG	1:A:114:PRO:HG3	2.45	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:87:ALA:HB1	1:A:110:VAL:HG13	1.98	0.46
1:A:9:LYS:HZ3	1:A:13:GLU:HG3	1.81	0.46
1:A:267:LEU:HD23	1:A:291:MET:HB2	1.97	0.46
1:A:453:GLU:OE1	1:A:453:GLU:N	2.40	0.46
1:A:42:TRP:CZ2	1:A:46:GLU:HG3	2.51	0.45
1:A:72:VAL:HB	1:A:141:VAL:CG1	2.47	0.45
1:A:191:GLN:HE21	1:A:241:SER:CB	2.24	0.45
1:A:220:MET:HG2	1:A:227:LEU:HA	1.98	0.45
1:A:329:ARG:NH1	1:A:351:VAL:HG21	2.32	0.45
1:A:411:LEU:HD11	1:A:429:ILE:HG22	1.98	0.45
1:A:21:VAL:HG23	1:A:26:LYS:HG2	1.98	0.45
1:A:436:GLU:O	1:A:436:GLU:HG2	2.17	0.44
1:A:201:MET:O	1:A:205:PRO:HG2	2.16	0.44
1:A:102:LEU:C	1:A:102:LEU:HD23	2.42	0.44
1:A:291:MET:HE1	1:A:307:HIS:HB2	1.98	0.44
1:A:405:HIS:O	1:A:405:HIS:ND1	2.50	0.44
1:A:87:ALA:O	1:A:91:GLN:HG3	2.18	0.44
1:A:295:GLY:HA3	1:A:302:ALA:HA	2.00	0.44
1:A:221:ILE:HG23	1:A:248:MET:CE	2.49	0.43
1:A:405:HIS:O	1:A:406:GLY:C	2.60	0.43
1:A:198:LEU:HD13	1:A:269:CYS:HB3	2.00	0.43
1:A:478:LYS:HB3	1:A:478:LYS:HE2	1.78	0.43
1:A:315:GLN:HG3	1:A:316:TYR:CD2	2.54	0.43
1:A:8:GLU:CG	1:A:33:LYS:HE2	2.48	0.43
1:A:488:THR:HB	1:A:489:PRO:CD	2.49	0.43
1:A:414:ILE:HG12	1:A:474:MET:HE2	2.01	0.42
1:A:257:GLU:HG3	4:A:922:HOH:O	2.18	0.42
1:A:260:LEU:HD13	1:A:263:MET:CE	2.49	0.42
1:A:497:LYS:HG2	1:A:501:GLU:OE2	2.19	0.42
1:A:356:SER:C	1:A:358:GLY:H	2.26	0.42
1:A:414:ILE:CG2	1:A:429:ILE:HD13	2.50	0.42
1:A:148:TYR:CD2	1:A:150:ILE:HG13	2.54	0.41
1:A:280:LYS:O	1:A:284:ARG:HG3	2.20	0.41
1:A:329:ARG:HG2	1:A:351:VAL:CG2	2.49	0.41
1:A:362:SER:H	1:A:363:PRO:HD3	1.84	0.41
1:A:487:MET:HA	1:A:492:LYS:O	2.21	0.41
1:A:9:LYS:NZ	1:A:13:GLU:HG3	2.35	0.41
1:A:141:VAL:HG23	1:A:145:GLU:HB2	2.03	0.41
1:A:334:LYS:O	1:A:336:ASP:O	2.38	0.41
1:A:340:ALA:HA	1:A:341:PRO:HD3	1.86	0.41
1:A:214:TRP:CH2	1:A:234:SER:HA	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:334:LYS:O	1:A:335:GLU:C	2.64	0.41
1:A:345:LYS:HA	1:A:385:GLY:O	2.21	0.41
1:A:467:ASN:HA	1:A:475:ILE:HD11	2.02	0.41
1:A:91:GLN:O	1:A:94:ALA:HB3	2.21	0.40
1:A:92:ARG:HH22	1:A:96:ASN:ND2	2.16	0.40
1:A:415:GLU:CG	1:A:429:ILE:HD12	2.52	0.40
1:A:47:TYR:HB3	1:A:50:ASP:HB3	2.03	0.40
1:A:10:TRP:HE3	1:A:14:THR:HG21	1.86	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

There are no protein backbone outliers to report in this entry.

5.3.2 Protein sidechains [i](#)

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	434/446 (97%)	409 (94%)	25 (6%)	18 25

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	GLU
1	A	39	LEU
1	A	113	LEU
1	A	141	VAL
1	A	181	ASN
1	A	227	LEU
1	A	231	LEU
1	A	248	MET

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Mol	Chain	Res	Type
1	A	252	GLU
1	A	264	LYS
1	A	271	GLU
1	A	276	GLU
1	A	279	ARG
1	A	287	LYS
1	A	310	GLU
1	A	325	LYS
1	A	338	THR
1	A	351	VAL
1	A	402	ILE
1	A	428	VAL
1	A	434	LYS
1	A	445	VAL
1	A	453	GLU
1	A	496	LYS
1	A	498	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	41	HIS
1	A	96	ASN
1	A	122	ASN
1	A	130	HIS
1	A	133	ASN
1	A	139	HIS
1	A	146	ASN
1	A	191	GLN
1	A	389	ASN
1	A	395	ASN
1	A	401	GLN
1	A	416	HIS
1	A	467	ASN
1	A	482	GLN

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
3	ATP	A	711	2	32,33,33	2.31	10 (31%)	48,52,52	1.36	6 (12%)

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	ATP	A	711	2	-	4/22/38/38	0/3/3/3

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	711	ATP	C5'-C4'	-5.96	1.33	1.51
3	A	711	ATP	C2-N1	4.69	1.42	1.33
3	A	711	ATP	C4-N3	4.27	1.42	1.34
3	A	711	ATP	O2'-C2'	4.21	1.53	1.43
3	A	711	ATP	C1'-N9	-3.94	1.35	1.46
3	A	711	ATP	PG-O3G	-2.65	1.44	1.54
3	A	711	ATP	O5'-C5'	2.25	1.53	1.44
3	A	711	ATP	PB-O2B	-2.15	1.45	1.55

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	711	ATP	O4'-C4'	2.14	1.49	1.45
3	A	711	ATP	C5-N7	-2.11	1.35	1.39

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	711	ATP	N3-C2-N1	-4.59	121.64	128.58
3	A	711	ATP	C5-N7-C8	3.80	109.43	103.45
3	A	711	ATP	N9-C8-N7	-3.15	109.46	113.94
3	A	711	ATP	O3'-C3'-C2'	2.52	119.90	111.82
3	A	711	ATP	C4-C5-N7	-2.15	108.12	110.58
3	A	711	ATP	C3'-C2'-C1'	2.11	105.46	101.46

There are no chirality outliers.

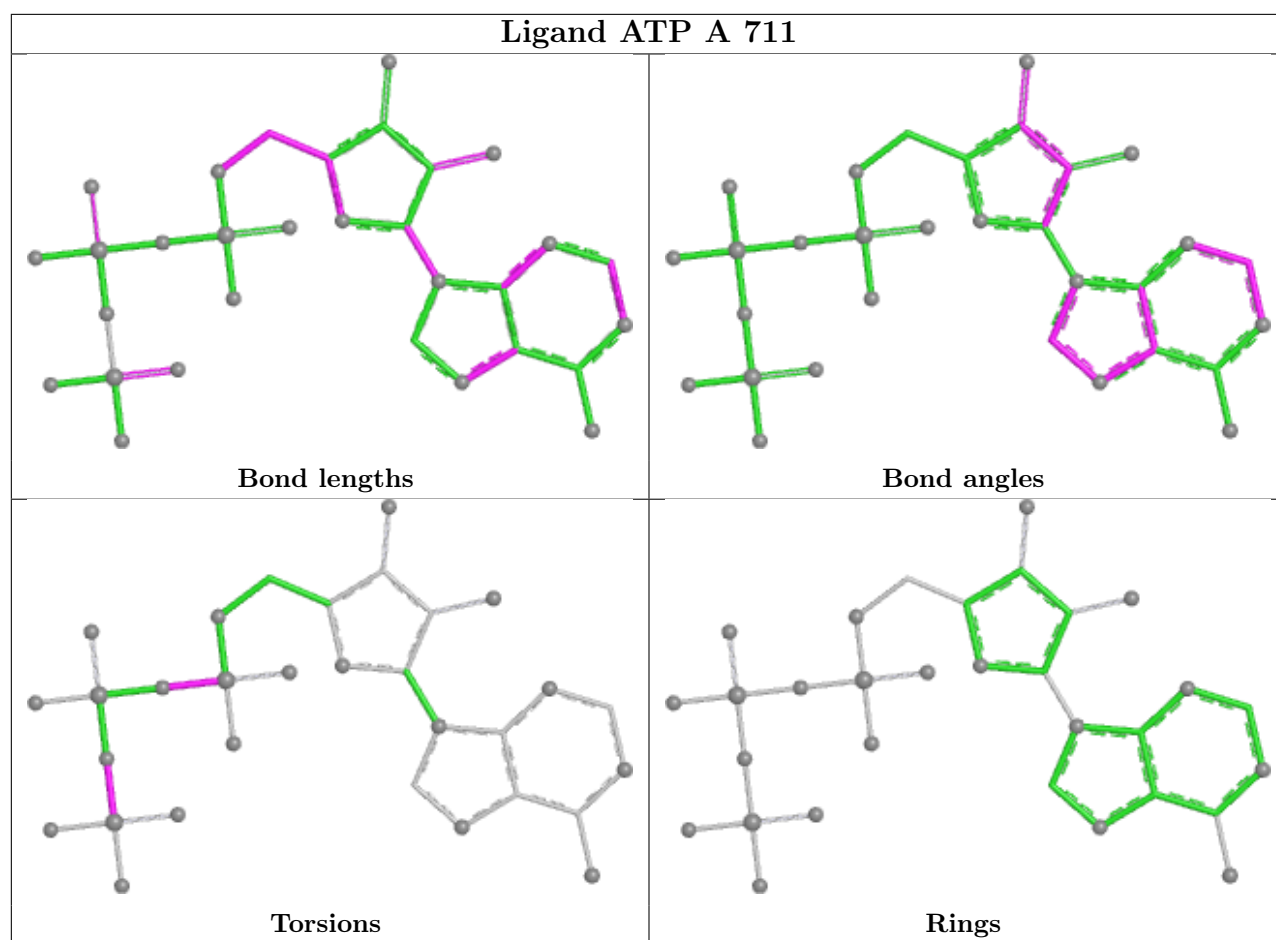
All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	711	ATP	PB-O3A-PA-O5'
3	A	711	ATP	PB-O3B-PG-O1G
3	A	711	ATP	PB-O3B-PG-O2G
3	A	711	ATP	PB-O3B-PG-O3G

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 [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	500/512 (97%)	0.70	45 (9%) 15 17	20, 39, 68, 88	0

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	407	TYR	5.6
1	A	154	GLY	4.5
1	A	153	SER	4.2
1	A	405	HIS	3.9
1	A	337	GLY	3.9
1	A	109	THR	3.8
1	A	111	THR	3.8
1	A	404	LEU	3.4
1	A	503	THR	3.4
1	A	504	ALA	3.4
1	A	502	VAL	3.4
1	A	105	ALA	3.2
1	A	363	PRO	3.1
1	A	110	VAL	3.1
1	A	361	GLY	3.0
1	A	108	VAL	3.0
1	A	338	THR	3.0
1	A	365	LEU	2.9
1	A	112	ASP	2.9
1	A	406	GLY	2.9
1	A	366	THR	2.9
1	A	91	GLN	2.7
1	A	340	ALA	2.7
1	A	335	GLU	2.7
1	A	258	SER	2.7
1	A	336	ASP	2.6
1	A	387	VAL	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	402	ILE	2.5
1	A	152	THR	2.5
1	A	339	ILE	2.4
1	A	159	PRO	2.4
1	A	403	LYS	2.4
1	A	435	GLY	2.3
1	A	369	ALA	2.3
1	A	107	ALA	2.3
1	A	50	ASP	2.3
1	A	410	GLU	2.2
1	A	253	ALA	2.2
1	A	14	THR	2.2
1	A	113	LEU	2.2
1	A	96	ASN	2.1
1	A	95	GLU	2.1
1	A	411	LEU	2.1
1	A	106	THR	2.0
1	A	92	ARG	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.

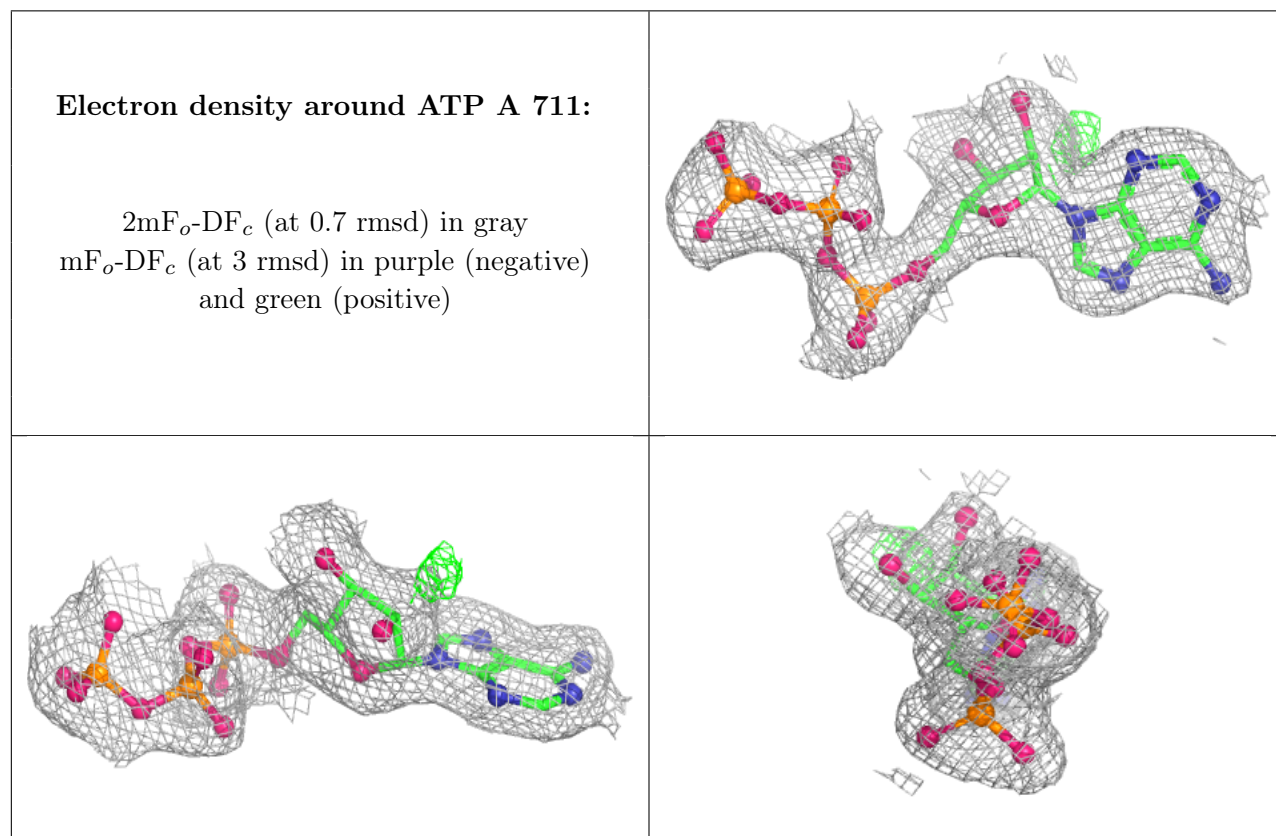
6.4 Ligands [i](#)

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
2	MG	A	710	1/1	0.92	0.07	52,52,52,52	0
3	ATP	A	711	31/31	0.94	0.08	23,30,49,51	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.