



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 03:00 PM UTC

PDB ID : 2ZJ5 / pdb_00002zj5
Title : Archaeal DNA helicase H_{jm} complexed with ADP in form 1
Authors : Oyama, T.; Oka, H.; Fujikane, R.; Ishino, Y.; Morikawa, K.
Deposited on : 2008-02-29
Resolution : 2.40 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

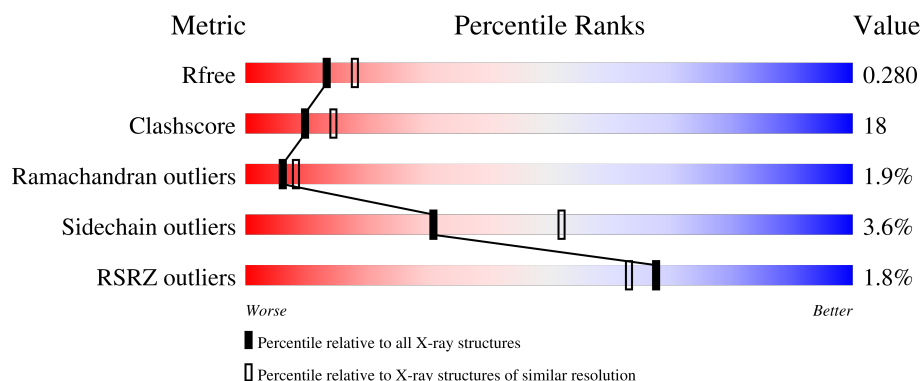
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

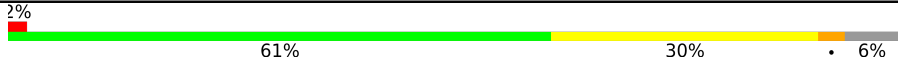
The reported resolution of this entry is 2.40 Å.

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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	720	

2 Entry composition [i](#)

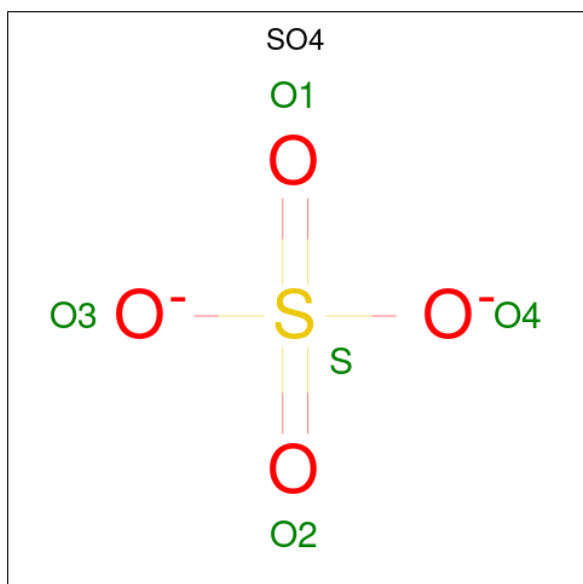
There are 4 unique types of molecules in this entry. The entry contains 5538 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Putative ski2-type helicase.

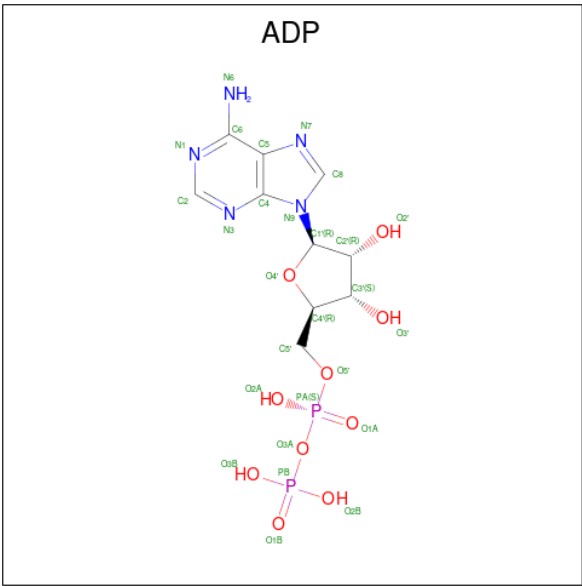
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	680	5373	3473	906	983	11	0	0	0

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

- 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 water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	113	Total	O	0	0
			113	113		

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	117.20Å 83.53Å 96.04Å 90.00° 121.22° 90.00°	Depositor
Resolution (Å)	43.98 – 2.40 43.98 – 2.40	Depositor EDS
% Data completeness (in resolution range)	95.6 (43.98-2.40) 95.7 (43.98-2.40)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.92 (at 2.39Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.216 , 0.290 0.209 , 0.280	Depositor DCC
R_{free} test set	1528 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	34.4	Xtriage
Anisotropy	0.317	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 56.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5538	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 5.88% 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: SO4, ADP

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	0.50	1/5480 (0.0%)	0.99	23/7413 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	365	MET	SD-CE	-5.64	1.65	1.79

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	474	GLU	N-CA-C	-7.70	103.71	113.72
1	A	388	ASP	CA-C-N	6.61	126.85	119.32
1	A	388	ASP	C-N-CA	6.61	126.85	119.32
1	A	526	THR	CA-C-N	6.47	126.45	119.78
1	A	526	THR	C-N-CA	6.47	126.45	119.78
1	A	75	VAL	CA-C-N	6.29	125.70	118.85
1	A	75	VAL	C-N-CA	6.29	125.70	118.85
1	A	12	ILE	N-CA-C	-6.16	105.72	111.45
1	A	468	PHE	N-CA-C	-6.03	105.00	112.90
1	A	371	ARG	CA-C-N	5.97	127.30	119.84
1	A	371	ARG	C-N-CA	5.97	127.30	119.84
1	A	190	ASN	N-CA-C	-5.95	103.70	111.74
1	A	121	THR	N-CA-C	-5.76	102.24	110.59
1	A	381	ILE	N-CA-C	5.73	116.12	107.75
1	A	105	TYR	N-CA-C	5.68	119.02	111.75
1	A	643	GLU	N-CA-C	5.57	118.07	111.33
1	A	91	TRP	N-CA-C	-5.38	106.56	113.23
1	A	192	GLU	N-CA-C	-5.35	101.75	110.20
1	A	351	ASP	N-CA-C	-5.31	106.64	113.01
1	A	327	THR	N-CA-C	-5.28	103.29	110.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	169	LYS	N-CA-C	-5.08	108.11	114.56
1	A	396	TYR	N-CA-C	5.08	119.75	113.50
1	A	245	ASN	N-CA-C	5.01	119.03	113.02

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	5373	0	5365	190	0
2	A	25	0	0	2	0
3	A	27	0	12	2	0
4	A	113	0	0	6	0
All	All	5538	0	5377	190	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 18.

All (190) 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:652:PRO:HG3	1:A:698:PHE:HB2	1.48	0.95
1:A:495:THR:HG22	1:A:523:PRO:HD2	1.45	0.94
1:A:76:PRO:HG2	1:A:146:GLU:HG3	1.53	0.90
1:A:53:THR:HB	3:A:801:ADP:O1A	1.71	0.90
1:A:492:ASP:HB3	1:A:495:THR:HG23	1.61	0.82
1:A:495:THR:HG21	1:A:524:ASP:OD2	1.80	0.81
1:A:485:ARG:HG2	1:A:620:VAL:HG21	1.64	0.80
1:A:201:VAL:HG11	1:A:370:GLY:O	1.82	0.79
1:A:97:ARG:HH11	1:A:97:ARG:HG2	1.50	0.77
1:A:663:TYR:HA	1:A:667:PHE:O	1.85	0.77
1:A:147:ILE:HG22	1:A:176:SER:HB3	1.66	0.76
1:A:657:ARG:HD3	1:A:657:ARG:H	1.51	0.74
1:A:28:GLN:HA	1:A:55:ILE:HD11	1.70	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:177:ALA:HB2	1:A:364:GLN:NE2	2.03	0.73
1:A:66:THR:HG22	1:A:67:GLN:HG2	1.71	0.71
1:A:528:PHE:CE2	1:A:603:ILE:HG13	2.26	0.70
1:A:155:ARG:HH11	1:A:155:ARG:HB2	1.57	0.69
1:A:147:ILE:CG2	1:A:176:SER:HB3	2.21	0.69
1:A:385:THR:HG21	2:A:901:SO4:O3	1.92	0.68
1:A:455:LYS:O	1:A:459:ILE:HG13	1.94	0.68
1:A:245:ASN:ND2	1:A:347:TRP:O	2.25	0.67
1:A:306:ARG:O	1:A:310:VAL:HG23	1.94	0.67
1:A:83:GLU:O	1:A:87:GLU:HG2	1.95	0.67
1:A:306:ARG:NH1	1:A:309:ARG:HD2	2.11	0.66
1:A:66:THR:HG22	1:A:67:GLN:CG	2.25	0.65
1:A:243:PHE:CG	1:A:365:MET:HE2	2.33	0.64
1:A:246:MET:HE3	2:A:902:SO4:O3	1.99	0.63
1:A:3:VAL:HG23	1:A:22:GLU:C	2.23	0.63
1:A:652:PRO:HG2	1:A:695:ILE:HA	1.81	0.63
1:A:423:ALA:HB1	1:A:484:ILE:HA	1.79	0.62
1:A:46:ILE:HG22	1:A:196:SER:HB3	1.84	0.59
1:A:219:SER:O	1:A:220:ILE:HD13	2.01	0.59
1:A:97:ARG:HG2	1:A:97:ARG:NH1	2.13	0.59
1:A:222:ARG:HH11	1:A:222:ARG:HB2	1.68	0.58
1:A:276:LEU:CD2	1:A:311:LEU:HD21	2.34	0.58
1:A:300:HIS:CD2	1:A:300:HIS:C	2.81	0.58
1:A:429:THR:HG21	4:A:1034:HOH:O	2.03	0.58
1:A:655:GLY:HA3	1:A:657:ARG:NH1	2.19	0.58
1:A:211:GLY:HA2	1:A:223:PHE:O	2.04	0.58
1:A:310:VAL:O	1:A:314:GLU:HG2	2.04	0.58
1:A:520:SER:HA	1:A:525:ILE:HG21	1.85	0.58
1:A:652:PRO:CG	1:A:698:PHE:HB2	2.29	0.57
1:A:677:ARG:O	1:A:680:GLU:N	2.34	0.57
1:A:429:THR:HG22	1:A:431:GLU:N	2.19	0.57
1:A:209:TYR:CG	1:A:389:PRO:HG3	2.39	0.56
1:A:92:GLU:HG2	1:A:96:LEU:O	2.04	0.56
1:A:16:LEU:HD11	1:A:54:LEU:HD11	1.86	0.56
1:A:390:ARG:HB3	1:A:390:ARG:NH1	2.20	0.56
1:A:3:VAL:HG23	1:A:22:GLU:O	2.04	0.56
1:A:39:GLY:O	1:A:171:GLN:HG3	2.06	0.56
1:A:586:GLU:OE1	1:A:601:TYR:OH	2.23	0.55
1:A:76:PRO:CG	1:A:146:GLU:HG3	2.31	0.55
1:A:243:PHE:CD2	1:A:365:MET:HE2	2.41	0.55
1:A:276:LEU:HD11	1:A:315:ASN:ND2	2.22	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:311:LEU:HD12	1:A:315:ASN:ND2	2.22	0.55
1:A:590:VAL:HA	1:A:595:VAL:O	2.06	0.55
1:A:492:ASP:HB3	1:A:495:THR:CG2	2.35	0.55
1:A:635:VAL:HG11	1:A:646:ILE:HG23	1.89	0.55
1:A:99:ALA:HB2	1:A:115:TYR:CD2	2.42	0.54
1:A:276:LEU:HD11	1:A:315:ASN:CG	2.31	0.54
1:A:266:THR:O	1:A:270:ILE:HG13	2.08	0.54
1:A:668:ARG:HG3	1:A:668:ARG:HH11	1.72	0.54
1:A:6:LEU:HD23	1:A:8:VAL:CG2	2.39	0.53
1:A:534:GLU:C	1:A:536:GLU:N	2.65	0.53
1:A:222:ARG:HH11	1:A:222:ARG:CB	2.21	0.53
1:A:625:GLU:HG2	4:A:1060:HOH:O	2.09	0.53
1:A:155:ARG:HH11	1:A:155:ARG:CB	2.22	0.52
1:A:209:TYR:CD2	1:A:389:PRO:HG2	2.43	0.52
1:A:222:ARG:HB2	1:A:222:ARG:NH1	2.24	0.52
1:A:663:TYR:CD1	1:A:668:ARG:HG2	2.45	0.52
1:A:223:PHE:CE2	1:A:229:LEU:HD23	2.45	0.52
1:A:652:PRO:HG3	1:A:698:PHE:CB	2.33	0.52
1:A:280:LEU:HD21	1:A:311:LEU:HD23	1.91	0.51
1:A:520:SER:HA	1:A:525:ILE:HD13	1.91	0.51
1:A:180:GLY:C	1:A:182:PRO:HD3	2.35	0.51
1:A:75:VAL:HG11	1:A:81:ALA:HB2	1.93	0.51
1:A:205:ARG:HD3	1:A:216:GLU:OE2	2.10	0.51
1:A:583:GLU:OE2	1:A:660:ARG:HD3	2.10	0.51
1:A:246:MET:O	1:A:249:LYS:HB3	2.10	0.51
1:A:681:LEU:C	1:A:683:LYS:N	2.67	0.51
1:A:429:THR:HG22	1:A:431:GLU:H	1.74	0.51
1:A:534:GLU:C	1:A:536:GLU:H	2.18	0.51
1:A:242:ILE:HD12	1:A:242:ILE:N	2.26	0.51
1:A:161:VAL:HG21	1:A:417:GLN:HG2	1.91	0.51
1:A:225:SER:O	1:A:228:GLU:HB2	2.10	0.51
1:A:616:GLU:O	1:A:620:VAL:HG23	2.11	0.51
1:A:8:VAL:HB	1:A:12:ILE:HD12	1.93	0.50
1:A:108:LYS:HG3	1:A:110:GLU:HG2	1.93	0.50
1:A:276:LEU:HD21	1:A:311:LEU:HD11	1.94	0.50
1:A:578:LEU:O	1:A:582:ASN:ND2	2.44	0.50
1:A:528:PHE:CZ	1:A:603:ILE:HG13	2.46	0.50
1:A:485:ARG:HG2	1:A:620:VAL:CG2	2.38	0.50
1:A:266:THR:OG1	1:A:268:PRO:HD2	2.12	0.50
1:A:362:VAL:HG21	1:A:396:TYR:CE2	2.47	0.50
1:A:466:ASN:OD1	1:A:497:LYS:HE2	2.11	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:466:ASN:C	1:A:467:GLU:HG3	2.37	0.49
1:A:462:PHE:CE1	1:A:497:LYS:HD2	2.47	0.49
1:A:604:VAL:HG13	1:A:637:VAL:O	2.13	0.49
1:A:554:PRO:HG3	1:A:567:PHE:CE2	2.47	0.48
1:A:123:GLU:OE2	1:A:155:ARG:HD2	2.13	0.48
1:A:89:GLN:C	1:A:91:TRP:H	2.22	0.48
1:A:403:LYS:HD3	1:A:405:PHE:CZ	2.49	0.48
1:A:577:LEU:O	1:A:580:TRP:HB3	2.14	0.48
1:A:682:LEU:C	1:A:684:ILE:H	2.21	0.48
1:A:134:SER:O	1:A:137:LYS:HG2	2.14	0.48
1:A:12:ILE:O	1:A:16:LEU:HD23	2.14	0.47
1:A:148:HIS:HA	1:A:404:LEU:HD11	1.95	0.47
1:A:42:ALA:HA	1:A:191:ALA:HB1	1.95	0.47
1:A:481:SER:HB2	4:A:1097:HOH:O	2.13	0.47
1:A:6:LEU:HD23	1:A:8:VAL:HG21	1.97	0.47
1:A:108:LYS:HB3	1:A:131:HIS:HB3	1.95	0.47
1:A:311:LEU:CD1	1:A:315:ASN:ND2	2.78	0.47
1:A:374:TYR:CD1	1:A:374:TYR:N	2.83	0.47
1:A:478:ARG:HD2	4:A:1076:HOH:O	2.15	0.47
1:A:578:LEU:CD1	1:A:582:ASN:HD21	2.28	0.47
1:A:148:HIS:C	1:A:150:ILE:H	2.23	0.46
1:A:677:ARG:C	1:A:679:GLU:N	2.72	0.46
1:A:681:LEU:C	1:A:683:LYS:H	2.21	0.46
1:A:353:GLY:O	1:A:354:MET:C	2.57	0.46
1:A:530:TYR:HA	1:A:534:GLU:OE2	2.15	0.46
1:A:75:VAL:HG22	1:A:76:PRO:HD2	1.98	0.46
1:A:182:PRO:HG3	4:A:1001:HOH:O	2.15	0.46
1:A:527:PRO:HB2	1:A:569:ARG:HB3	1.97	0.46
1:A:622:GLY:O	1:A:624:TYR:HD1	1.99	0.45
1:A:28:GLN:HA	1:A:55:ILE:CD1	2.44	0.45
1:A:209:TYR:CD1	1:A:389:PRO:HG3	2.51	0.45
1:A:622:GLY:HA2	1:A:624:TYR:CE1	2.52	0.45
1:A:474:GLU:O	1:A:475:ASP:HB2	2.16	0.45
1:A:26:PRO:HB2	1:A:27:PRO:HD3	1.99	0.45
1:A:528:PHE:O	1:A:569:ARG:HD2	2.17	0.45
1:A:469:ILE:HD12	1:A:477:ILE:CG2	2.47	0.45
1:A:299:PHE:HA	1:A:325:VAL:O	2.17	0.45
1:A:343:ILE:HD11	1:A:365:MET:HB2	1.97	0.45
1:A:160:GLU:OE2	1:A:440:THR:HB	2.17	0.45
1:A:497:LYS:NZ	1:A:501:ASP:OD2	2.50	0.45
1:A:155:ARG:HB2	1:A:155:ARG:NH1	2.28	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:229:LEU:HD13	1:A:381:ILE:CD1	2.47	0.44
1:A:519:ILE:O	1:A:525:ILE:HG21	2.18	0.44
1:A:219:SER:C	1:A:220:ILE:HD13	2.43	0.44
1:A:56:ALA:O	1:A:60:MET:HG3	2.17	0.44
1:A:300:HIS:CD2	1:A:309:ARG:HG2	2.53	0.44
1:A:586:GLU:CD	1:A:601:TYR:HH	2.26	0.44
1:A:692:VAL:O	1:A:693:GLU:C	2.60	0.44
1:A:97:ARG:HH11	1:A:97:ARG:CG	2.26	0.43
1:A:11:ARG:HD2	1:A:94:ILE:HB	2.01	0.43
1:A:148:HIS:C	1:A:150:ILE:N	2.77	0.43
1:A:267:LYS:N	1:A:268:PRO:CD	2.81	0.43
1:A:427:TYR:HD2	1:A:432:GLU:HB3	1.84	0.43
1:A:48:THR:O	3:A:801:ADP:O1B	2.36	0.43
1:A:442:TYR:OH	1:A:448:ASP:O	2.26	0.43
1:A:75:VAL:HA	1:A:76:PRO:HD3	1.88	0.43
1:A:429:THR:HG22	1:A:430:VAL:N	2.33	0.43
1:A:25:TYR:HB3	1:A:27:PRO:HD2	2.00	0.42
1:A:276:LEU:HD21	1:A:311:LEU:CG	2.49	0.42
1:A:694:ALA:O	1:A:698:PHE:N	2.51	0.42
1:A:306:ARG:HH11	1:A:309:ARG:HD2	1.83	0.42
1:A:145:ASP:O	1:A:146:GLU:HB2	2.20	0.42
1:A:181:ASN:OD1	1:A:181:ASN:N	2.52	0.42
1:A:225:SER:HB2	1:A:227:GLU:OE2	2.19	0.42
1:A:152:SER:C	1:A:407:GLN:HG3	2.44	0.42
1:A:255:LEU:O	1:A:259:LYS:HG3	2.20	0.42
1:A:316:PHE:CD1	1:A:324:VAL:HG23	2.54	0.42
1:A:340:ARG:HG2	1:A:379:GLU:HB2	2.02	0.42
1:A:429:THR:CG2	1:A:430:VAL:N	2.83	0.42
1:A:628:ASP:HB2	4:A:1044:HOH:O	2.19	0.42
1:A:654:VAL:O	1:A:657:ARG:CZ	2.68	0.42
1:A:695:ILE:HG22	1:A:696:PHE:N	2.34	0.42
1:A:47:PRO:HD3	1:A:196:SER:O	2.20	0.42
1:A:272:ALA:O	1:A:276:LEU:HB2	2.20	0.42
1:A:553:ASP:HA	1:A:554:PRO:HD3	1.88	0.41
1:A:177:ALA:HB2	1:A:364:GLN:HE21	1.81	0.41
1:A:241:LEU:HD21	1:A:243:PHE:CZ	2.55	0.41
1:A:42:ALA:CA	1:A:191:ALA:HB1	2.51	0.41
1:A:142:LEU:C	1:A:142:LEU:HD23	2.45	0.41
1:A:276:LEU:O	1:A:276:LEU:HD23	2.20	0.41
1:A:276:LEU:HD23	1:A:311:LEU:HD21	2.02	0.41
1:A:409:SER:O	1:A:410:ASN:C	2.64	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:513:ILE:HD13	1:A:513:ILE:HA	1.82	0.41
1:A:624:TYR:HA	1:A:627:VAL:HG23	2.01	0.41
1:A:495:THR:HB	1:A:522:THR:HB	2.01	0.41
1:A:183:GLU:CD	1:A:183:GLU:H	2.29	0.41
1:A:229:LEU:HD13	1:A:381:ILE:HD13	2.03	0.41
1:A:103:GLY:O	1:A:124:LYS:NZ	2.40	0.41
1:A:276:LEU:HD21	1:A:311:LEU:HD21	2.00	0.41
1:A:658:ARG:O	1:A:659:ALA:C	2.64	0.40
1:A:89:GLN:C	1:A:91:TRP:N	2.79	0.40
1:A:255:LEU:O	1:A:255:LEU:HD12	2.21	0.40
1:A:6:LEU:HD23	1:A:8:VAL:HG22	2.02	0.40
1:A:75:VAL:O	1:A:121:THR:HA	2.21	0.40
1:A:257:LEU:O	1:A:261:VAL:HG23	2.21	0.40

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	672/720 (93%)	616 (92%)	43 (6%)	13 (2%)	6 8

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	448	ASP
1	A	654	VAL
1	A	667	PHE
1	A	684	ILE
1	A	354	MET
1	A	592	LYS
1	A	90	ASP
1	A	536	GLU

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Mol	Chain	Res	Type
1	A	546	LYS
1	A	283	ASN
1	A	328	PRO
1	A	372	PRO
1	A	622	GLY

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	553/626 (88%)	533 (96%)	20 (4%)	31 52

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	75	VAL
1	A	107	SER
1	A	154	ASP
1	A	155	ARG
1	A	205	ARG
1	A	222	ARG
1	A	227	GLU
1	A	356	ARG
1	A	372	PRO
1	A	387	ASP
1	A	448	ASP
1	A	449	THR
1	A	495	THR
1	A	513	ILE
1	A	547	ASP
1	A	594	SER
1	A	644	GLU
1	A	656	ARG
1	A	657	ARG
1	A	678	PRO

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	62	HIS
1	A	67	GLN
1	A	171	GLN
1	A	190	ASN
1	A	300	HIS
1	A	364	GLN
1	A	407	GLN
1	A	582	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](#)

6 ligands are modelled in this entry.

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	$\# Z > 2$
2	SO4	A	905	-	4,4,4	0.41	0	6,6,6	0.15	0
2	SO4	A	901	-	4,4,4	1.91	2 (50%)	6,6,6	0.93	0
2	SO4	A	903	-	4,4,4	0.52	0	6,6,6	0.18	0
2	SO4	A	902	-	4,4,4	0.50	0	6,6,6	0.13	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	ADP	A	801	-	28,29,29	1.54	5 (17%)	43,45,45	1.17	4 (9%)
2	SO4	A	904	-	4,4,4	0.54	0	6,6,6	0.18	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	801	-	-	3/16/32/32	0/3/3/3

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	801	ADP	C1'-N9	3.56	1.56	1.46
2	A	901	SO4	O1-S	3.05	1.63	1.44
3	A	801	ADP	C2'-C3'	2.94	1.61	1.53
3	A	801	ADP	PA-O3A	2.87	1.62	1.59
3	A	801	ADP	C8-N9	-2.64	1.33	1.37
2	A	901	SO4	O3-S	-2.23	1.29	1.48
3	A	801	ADP	O2'-C2'	2.10	1.48	1.43

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	801	ADP	O5'-C5'-C4'	2.76	118.38	108.99
3	A	801	ADP	O4'-C4'-C3'	-2.39	100.41	105.15
3	A	801	ADP	O4'-C1'-N9	2.16	112.24	108.09
3	A	801	ADP	C5'-C4'-C3'	2.05	122.58	115.21

There are no chirality outliers.

All (3) torsion outliers are listed below:

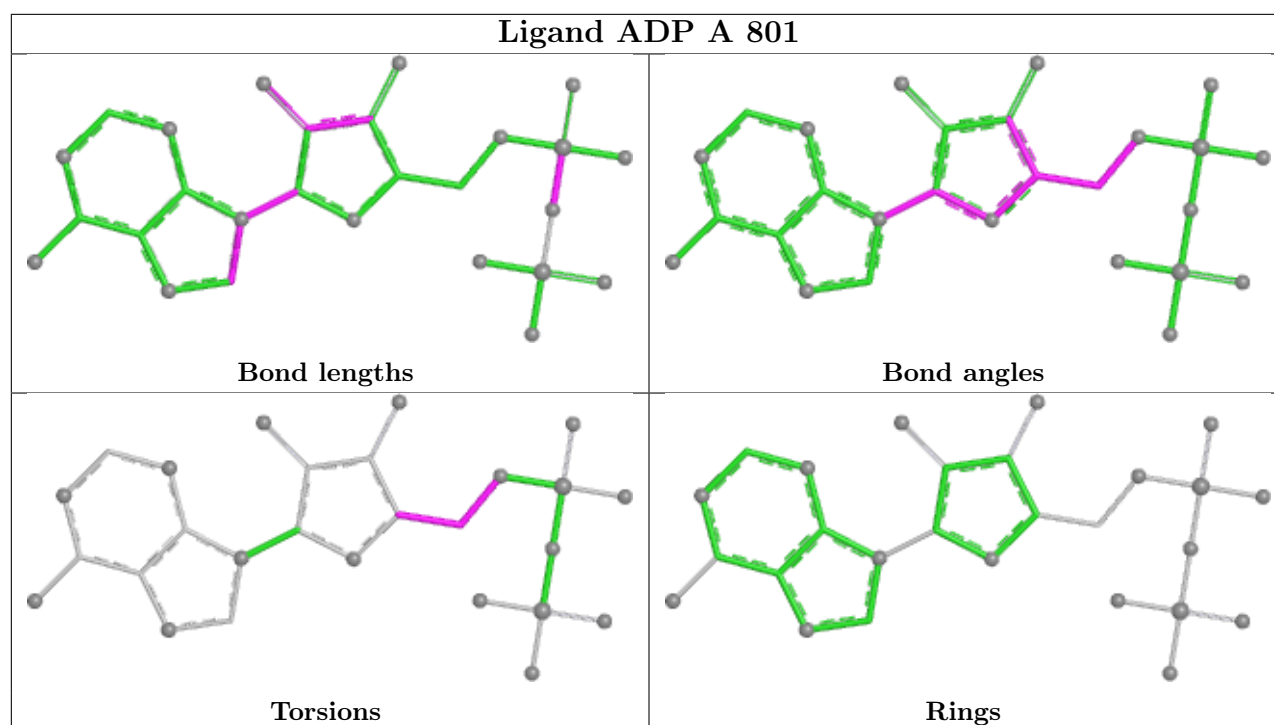
Mol	Chain	Res	Type	Atoms
3	A	801	ADP	C3'-C4'-C5'-O5'
3	A	801	ADP	O4'-C4'-C5'-O5'
3	A	801	ADP	C4'-C5'-O5'-PA

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	901	SO4	1	0
2	A	902	SO4	1	0
3	A	801	ADP	2	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	680/720 (94%)	0.02	12 (1%) 67 63	20, 37, 58, 75	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	673	ILE	4.3
1	A	654	VAL	4.0
1	A	351	ASP	3.7
1	A	653	LEU	3.0
1	A	679	GLU	2.9
1	A	676	ALA	2.7
1	A	680	GLU	2.5
1	A	448	ASP	2.5
1	A	675	GLN	2.4
1	A	699	LEU	2.3
1	A	667	PHE	2.2
1	A	652	PRO	2.1

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.

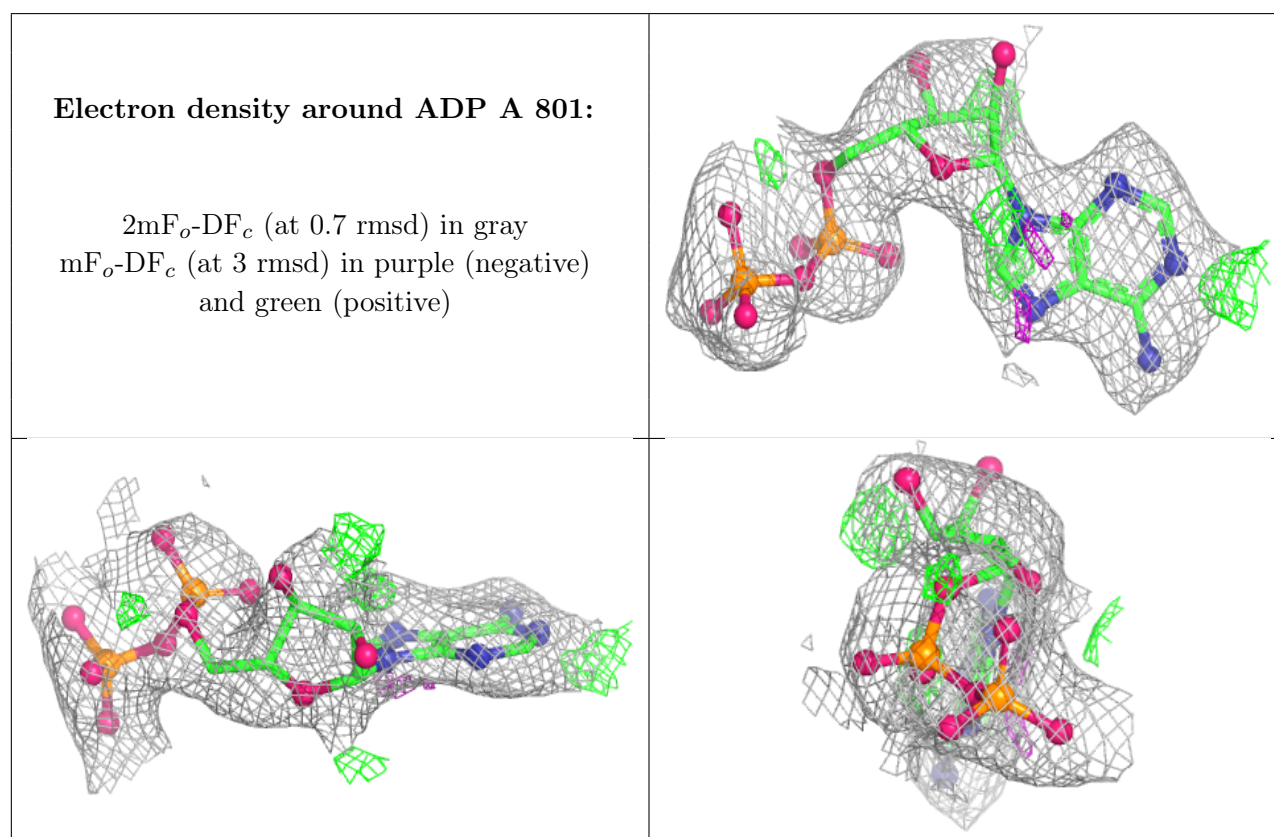
6.4 Ligands [i](#)

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	ADP	A	801	27/27	0.88	0.12	28,47,99,112	0
2	SO4	A	902	5/5	0.90	0.12	38,55,75,79	0
2	SO4	A	903	5/5	0.90	0.14	49,50,65,72	0
2	SO4	A	901	5/5	0.90	0.13	45,48,70,79	5
2	SO4	A	904	5/5	0.91	0.15	49,49,60,65	0
2	SO4	A	905	5/5	0.95	0.09	41,46,63,68	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.