



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 03:18 AM UTC

PDB ID : 3GIF / pdb_00003gif
Title : Crystal structure of DesKC_H188E in complex with ADP
Authors : Trajtenberg, F.; Albanesi, D.; Alzari, P.M.; Buschiazzi, A.; de Mendoza, D.
Deposited on : 2009-03-05
Resolution : 2.70 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

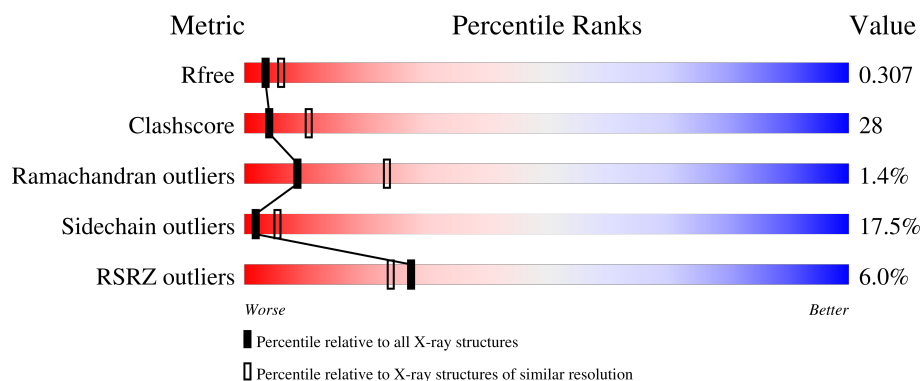
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	218	
1	B	218	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	ADP	A	1301	-	-	X	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 2224 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.

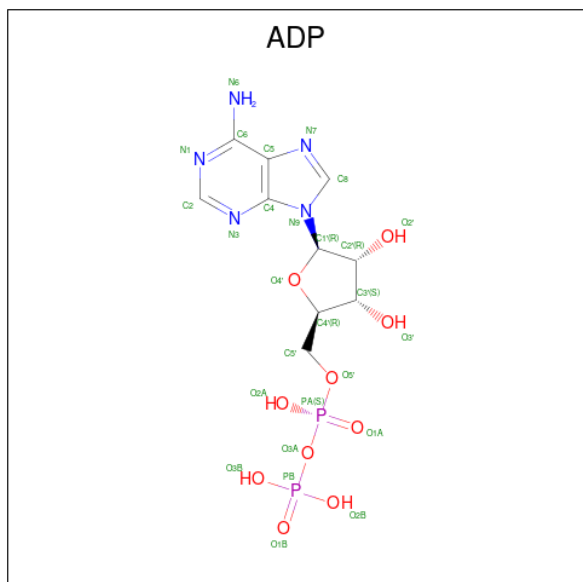
- Molecule 1 is a protein called Sensor histidine kinase desK.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	207	Total	C	N	O	S	0	0	0
			1584	991	276	311	6			
1	B	77	Total	C	N	O		0	0	0
			592	364	110	118				

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	153	GLY	-	expression tag	UNP O34757
A	188	GLU	HIS	engineered mutation	UNP O34757
B	153	GLY	-	expression tag	UNP O34757
B	188	GLU	HIS	engineered mutation	UNP O34757

- Molecule 2 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



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

- 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 water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	15	Total	O	0	0
			15	15		
4	B	5	Total	O	0	0
			5	5		

- Molecule 1: Sensor histidine kinase desK



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	91.58Å 93.48Å 59.42Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.27 – 2.70 36.27 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.7 (36.27-2.70) 99.6 (36.27-2.70)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.64 (at 2.68Å)	Xtriage
Refinement program	PHENIX 2009_02_15_2320_3	Depositor
R, R_{free}	0.254 , 0.317 (Not available) , 0.307	Depositor DCC
R_{free} test set	726 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	58.6	Xtriage
Anisotropy	0.075	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 41.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.008 for k,h,-l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	2224	wwPDB-VP
Average B, all atoms (Å ²)	63.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.69% 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: ADP, 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	1.07	2/1597 (0.1%)	1.39	21/2156 (1.0%)
1	B	1.29	5/593 (0.8%)	1.47	9/797 (1.1%)
All	All	1.14	7/2190 (0.3%)	1.41	30/2953 (1.0%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	270	LYS	C-O	11.83	1.38	1.24
1	B	189	ASP	C-O	10.23	1.36	1.24
1	B	175	VAL	CA-CB	6.86	1.63	1.54
1	B	186	ASP	C-O	6.71	1.31	1.24
1	B	192	GLY	C-O	5.75	1.30	1.23
1	B	188	GLU	C-O	5.32	1.30	1.24
1	A	201	LYS	C-O	-5.16	1.18	1.24

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	192	GLY	N-CA-C	-9.11	101.11	112.77
1	A	156	GLU	N-CA-C	-8.54	103.00	113.50
1	A	311	TRP	N-CA-C	7.13	118.70	111.07
1	A	269	GLU	N-CA-C	7.01	117.53	108.34
1	B	234	VAL	N-CA-C	-7.00	103.70	110.42
1	A	163	LYS	N-CA-C	-6.91	103.80	111.82
1	A	343	ARG	N-CA-C	-6.91	103.36	111.03
1	A	266	TYR	N-CA-C	6.71	118.39	107.32
1	B	212	ASP	CA-C-N	6.43	126.19	119.05
1	B	212	ASP	C-N-CA	6.43	126.19	119.05
1	A	169	GLU	N-CA-C	-6.28	104.52	111.36
1	A	260	ALA	N-CA-C	-6.27	105.91	112.93

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	309	GLN	N-CA-C	6.21	118.09	109.15
1	B	192	GLY	N-CA-C	-6.12	105.24	112.64
1	A	336	GLY	N-CA-C	6.00	119.91	112.64
1	A	244	ILE	N-CA-C	5.88	117.17	109.58
1	B	192	GLY	CA-C-N	-5.85	112.83	120.44
1	B	192	GLY	C-N-CA	-5.85	112.83	120.44
1	A	367	ASN	N-CA-C	5.83	117.31	111.07
1	B	203	ASP	N-CA-C	5.80	117.68	111.36
1	A	270	LYS	O-C-N	5.78	130.28	122.59
1	A	308	GLN	CA-C-N	-5.76	112.67	121.66
1	A	308	GLN	C-N-CA	-5.76	112.67	121.66
1	A	165	GLU	N-CA-C	-5.57	105.13	111.14
1	A	326	GLY	N-CA-C	5.57	126.37	113.18
1	B	208	LEU	N-CA-C	5.56	120.17	113.17
1	A	312	LYS	CA-C-O	5.43	126.04	119.31
1	B	191	LEU	N-CA-C	5.38	116.83	111.07
1	A	280	GLU	N-CA-C	-5.26	105.63	111.36
1	A	293	ASN	N-CA-C	-5.03	105.99	111.82

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	1584	0	1568	94	1
1	B	592	0	603	35	0
2	A	27	0	12	9	0
3	A	1	0	0	0	0
4	A	15	0	0	1	0
4	B	5	0	0	0	0
All	All	2224	0	2183	121	1

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 28.

All (121) 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:297:HIS:CD2	2:A:1301:ADP:H2'	1.95	1.01
1:A:272:PRO:HB2	1:A:275:ILE:HD11	1.45	0.98
1:A:169:GLU:O	1:A:172:ALA:HB3	1.64	0.97
1:A:265:ILE:HD11	1:A:302:THR:HG21	1.58	0.85
1:A:234:VAL:O	1:A:238:VAL:HG13	1.76	0.85
1:A:245:ARG:HD2	1:A:280:GLU:OE1	1.77	0.84
1:A:324:PHE:CD1	1:A:355:THR:HG21	2.13	0.83
1:A:343:ARG:HH11	1:A:343:ARG:CG	1.92	0.82
1:A:341:ARG:O	1:A:345:GLU:HB2	1.79	0.81
1:A:343:ARG:HH11	1:A:343:ARG:HG2	1.44	0.79
1:A:195:LEU:HD12	1:B:231:LEU:HD13	1.66	0.78
1:B:183:ILE:O	1:B:187:LEU:HD23	1.84	0.77
1:B:182:ARG:NH1	1:B:237:ILE:O	2.17	0.77
1:A:233:GLU:O	1:A:237:ILE:HG13	1.88	0.74
1:A:215:GLN:NE2	1:A:218:ARG:HH11	1.85	0.74
1:A:289:GLU:OE2	1:A:336:GLY:HA3	1.90	0.72
1:A:175:VAL:HG22	1:B:175:VAL:HG22	1.72	0.72
1:A:298:SER:C	1:A:299:GLN:HG3	2.14	0.72
1:A:215:GLN:NE2	1:A:218:ARG:NH1	2.38	0.71
1:A:235:ARG:HG2	1:A:235:ARG:HH21	1.54	0.71
1:B:187:LEU:HD22	1:B:237:ILE:HD13	1.73	0.71
1:A:309:GLN:HE22	1:A:367:ASN:HD21	1.38	0.70
1:B:215:GLN:HA	1:B:218:ARG:HH11	1.57	0.68
1:A:300:ALA:HB1	1:A:321:ASP:OD1	1.93	0.68
1:A:297:HIS:CG	2:A:1301:ADP:H2'	2.28	0.67
1:A:203:ASP:OD1	1:A:206:ARG:NH2	2.30	0.65
1:B:225:GLN:HE21	1:B:225:GLN:HA	1.63	0.64
1:A:326:GLY:HA3	2:A:1301:ADP:H4'	1.79	0.64
1:A:184:ALA:HB2	1:B:238:VAL:CG1	2.28	0.64
1:A:351:LEU:O	1:A:352:HIS:CD2	2.50	0.64
1:A:184:ALA:HB2	1:B:238:VAL:HG12	1.80	0.62
1:B:215:GLN:HA	1:B:218:ARG:NH1	2.14	0.62
1:A:245:ARG:HH21	1:A:277:LEU:HD11	1.64	0.62
1:A:257:LEU:HD21	1:A:291:VAL:HG13	1.81	0.60
1:A:309:GLN:HE22	1:A:367:ASN:ND2	1.99	0.60
1:A:336:GLY:O	1:A:340:MET:HG3	2.01	0.60
1:B:187:LEU:CD2	1:B:237:ILE:HG21	2.31	0.60
1:B:191:LEU:O	1:B:192:GLY:C	2.42	0.59
1:A:208:LEU:HD13	1:A:215:GLN:HB3	1.84	0.59
1:A:190:THR:HB	1:A:233:GLU:OE2	2.01	0.59
1:A:204:LEU:O	1:A:207:LYS:HB2	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:215:GLN:HA	1:A:218:ARG:HH11	1.69	0.57
1:B:191:LEU:HD11	1:B:234:VAL:CG2	2.34	0.57
1:A:175:VAL:CG2	1:B:175:VAL:HG22	2.34	0.57
1:A:335:HIS:CE1	2:A:1301:ADP:O2A	2.58	0.56
1:A:347:ALA:CB	1:A:365:ILE:HD11	2.35	0.56
1:A:158:GLU:HA	1:A:161:GLU:HG2	1.88	0.56
1:A:222:SER:O	1:A:226:THR:HG23	2.06	0.55
1:A:265:ILE:HD11	1:A:302:THR:CG2	2.34	0.55
1:A:347:ALA:O	1:A:348:ASN:HB2	2.07	0.55
1:B:166:ASP:O	1:B:167:ALA:C	2.51	0.54
1:B:182:ARG:O	1:B:183:ILE:C	2.51	0.54
1:A:244:ILE:HG13	1:A:245:ARG:H	1.72	0.53
1:A:298:SER:O	1:A:299:GLN:HG3	2.08	0.53
1:A:244:ILE:HG13	1:A:245:ARG:N	2.24	0.53
1:A:215:GLN:HE21	1:A:218:ARG:NH1	2.06	0.52
1:A:246:LEU:HB2	1:A:280:GLU:HG3	1.92	0.52
1:A:265:ILE:HB	1:A:304:ARG:HB2	1.92	0.52
1:A:343:ARG:CG	1:A:343:ARG:NH1	2.63	0.52
1:B:208:LEU:HB2	1:B:216:ALA:HB2	1.90	0.52
1:A:175:VAL:HG22	1:B:175:VAL:CG2	2.39	0.51
1:A:324:PHE:CD1	1:A:355:THR:CG2	2.91	0.51
1:B:166:ASP:O	1:B:169:GLU:N	2.43	0.50
1:A:335:HIS:ND1	2:A:1301:ADP:O2A	2.44	0.50
1:A:245:ARG:NH2	1:A:277:LEU:HD11	2.27	0.50
1:B:187:LEU:O	1:B:188:GLU:C	2.54	0.50
1:A:235:ARG:HG2	1:A:235:ARG:NH2	2.23	0.50
1:A:335:HIS:C	1:A:335:HIS:CD2	2.90	0.50
1:B:179:GLU:HA	1:B:179:GLU:OE2	2.12	0.49
1:A:188:GLU:OE2	1:A:188:GLU:HA	2.12	0.49
1:A:215:GLN:HA	1:A:218:ARG:NH1	2.26	0.49
1:A:156:GLU:C	1:A:158:GLU:N	2.68	0.49
1:A:317:THR:HA	1:A:361:LEU:O	2.13	0.48
1:B:214:GLU:O	1:B:218:ARG:HG3	2.13	0.48
1:A:324:PHE:HD1	1:A:355:THR:HG21	1.74	0.48
1:A:297:HIS:CD2	2:A:1301:ADP:C2'	2.83	0.48
1:A:173:GLU:O	1:A:177:LEU:HG	2.14	0.47
1:A:159:ARG:O	1:A:163:LYS:HG3	2.13	0.47
1:A:306:ASP:HB3	1:A:308:GLN:NE2	2.29	0.47
1:A:291:VAL:O	1:A:295:VAL:HG23	2.13	0.47
1:A:209:ILE:HD12	1:B:217:ALA:HB2	1.96	0.47
1:A:338:LEU:HD22	1:A:338:LEU:O	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:343:ARG:O	1:A:346:PHE:HB3	2.13	0.47
1:B:166:ASP:OD1	1:B:170:ARG:NH1	2.47	0.47
1:B:238:VAL:HG12	1:B:238:VAL:O	2.15	0.47
1:B:177:LEU:HD12	1:B:181:GLN:CD	2.40	0.46
1:B:190:THR:HG21	1:B:233:GLU:OE1	2.15	0.46
1:B:201:LYS:HE3	1:B:222:SER:OG	2.15	0.46
1:A:337:LEU:HD12	2:A:1301:ADP:H5'1	1.99	0.45
1:A:267:GLU:O	1:A:267:GLU:CD	2.60	0.45
1:A:365:ILE:HG12	1:A:366:PRO:HD2	1.99	0.44
1:A:164:LEU:HA	1:A:164:LEU:HD23	1.60	0.44
1:A:277:LEU:HD12	1:A:277:LEU:HA	1.65	0.44
1:A:250:LEU:O	1:A:253:ILE:HG13	2.17	0.44
1:A:337:LEU:HA	1:A:337:LEU:HD23	1.78	0.44
1:B:208:LEU:O	1:B:209:ILE:C	2.58	0.44
1:A:187:LEU:O	1:A:188:GLU:C	2.59	0.43
1:A:203:ASP:O	1:A:207:LYS:HG2	2.18	0.43
1:A:269:GLU:OE2	1:A:308:GLN:HG2	2.18	0.43
1:A:326:GLY:HA3	2:A:1301:ADP:C4'	2.46	0.43
1:A:340:MET:HE1	1:A:361:LEU:CD1	2.49	0.43
1:A:184:ALA:O	1:A:188:GLU:HG2	2.18	0.43
1:A:188:GLU:OE1	1:B:235:ARG:HD2	2.19	0.43
1:A:232:ASN:O	1:A:233:GLU:C	2.58	0.43
1:B:203:ASP:O	1:B:207:LYS:HG2	2.18	0.43
1:A:365:ILE:CG1	1:A:366:PRO:HD2	2.49	0.43
1:B:192:GLY:O	1:B:193:GLN:C	2.61	0.42
1:A:245:ARG:HE	1:A:277:LEU:HD11	1.84	0.42
1:A:309:GLN:N	4:A:2:HOH:O	2.52	0.42
1:A:351:LEU:O	1:A:352:HIS:HD2	2.00	0.42
1:A:179:GLU:HA	1:A:179:GLU:OE1	2.19	0.42
1:B:215:GLN:NE2	1:B:218:ARG:NH1	2.68	0.42
1:B:168:ASN:O	1:B:169:GLU:C	2.63	0.41
1:A:271:TRP:HA	1:A:272:PRO:HD3	1.79	0.41
1:A:215:GLN:O	1:A:219:GLU:HG2	2.21	0.41
1:A:187:LEU:HD23	1:A:187:LEU:HA	1.74	0.41
1:A:347:ALA:HB1	1:A:365:ILE:HD11	2.03	0.41
1:A:182:ARG:HG2	1:A:182:ARG:NH1	2.37	0.40
1:A:247:LYS:HG2	1:A:271:TRP:CZ2	2.56	0.40
1:A:298:SER:HB3	2:A:1301:ADP:N1	2.36	0.40
1:B:187:LEU:HD22	1:B:237:ILE:CD1	2.48	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:311:TRP:O	1:A:311:TRP:O[2_665]	2.05	0.15

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	203/218 (93%)	186 (92%)	15 (7%)	2 (1%)	12	32
1	B	75/218 (34%)	70 (93%)	3 (4%)	2 (3%)	4	10
All	All	278/436 (64%)	256 (92%)	18 (6%)	4 (1%)	9	23

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	270	LYS
1	A	326	GLY
1	B	189	ASP
1	B	169	GLU

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	166/194 (86%)	139 (84%)	27 (16%)	2	6
1	B	62/194 (32%)	49 (79%)	13 (21%)	1	3
All	All	228/388 (59%)	188 (82%)	40 (18%)	2	5

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	158	GLU
1	A	175	VAL
1	A	226	THR
1	A	228	ARG
1	A	231	LEU
1	A	238	VAL
1	A	239	SER
1	A	244	ILE
1	A	245	ARG
1	A	250	LEU
1	A	253	ILE
1	A	261	ASP
1	A	262	ILE
1	A	277	LEU
1	A	284	SER
1	A	299	GLN
1	A	302	THR
1	A	311	TRP
1	A	313	GLU
1	A	319	SER
1	A	335	HIS
1	A	338	LEU
1	A	343	ARG
1	A	351	LEU
1	A	353	ILE
1	A	359	THR
1	A	361	LEU
1	B	166	ASP
1	B	168	ASN
1	B	175	VAL
1	B	177	LEU
1	B	190	THR
1	B	196	SER
1	B	201	LYS
1	B	225	GLN
1	B	230	SER
1	B	231	LEU
1	B	233	GLU
1	B	235	ARG
1	B	236	LYS

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	215	GLN
1	A	224	GLN
1	A	232	ASN
1	A	279	ASN
1	A	281	ASN
1	A	348	ASN
1	A	352	HIS
1	A	367	ASN
1	B	193	GLN
1	B	215	GLN
1	B	225	GLN
1	B	232	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 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	$\# Z > 2$
2	ADP	A	1301	3	28,29,29	1.47	4 (14%)	43,45,45	2.24	14 (32%)

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
2	ADP	A	1301	3	-	2/16/32/32	0/3/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1301	ADP	C5-C4	4.36	1.46	1.39
2	A	1301	ADP	C5-C6	3.23	1.50	1.41
2	A	1301	ADP	PA-O3A	3.22	1.63	1.59
2	A	1301	ADP	C8-N7	2.05	1.35	1.31

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1301	ADP	C5-C4-N3	-5.96	118.52	126.72
2	A	1301	ADP	N3-C4-N9	4.72	135.20	127.17
2	A	1301	ADP	C2-N3-C4	4.47	122.75	111.83
2	A	1301	ADP	C4-C5-N7	-4.31	105.65	110.58
2	A	1301	ADP	N3-C2-N1	-4.31	122.06	128.58
2	A	1301	ADP	C2'-C1'-N9	-4.23	102.79	113.30
2	A	1301	ADP	C4-N9-C8	3.44	109.35	105.74
2	A	1301	ADP	C5-N7-C8	3.31	108.66	103.45
2	A	1301	ADP	O4'-C1'-N9	2.94	113.74	108.09
2	A	1301	ADP	C6-C5-N7	2.78	137.44	132.09
2	A	1301	ADP	N9-C8-N7	-2.74	110.05	113.94
2	A	1301	ADP	O3B-PB-O2B	2.40	116.80	107.80
2	A	1301	ADP	C5'-C4'-C3'	-2.22	107.21	115.21
2	A	1301	ADP	C2-N1-C6	2.15	122.27	118.73

There are no chirality outliers.

All (2) torsion outliers are listed below:

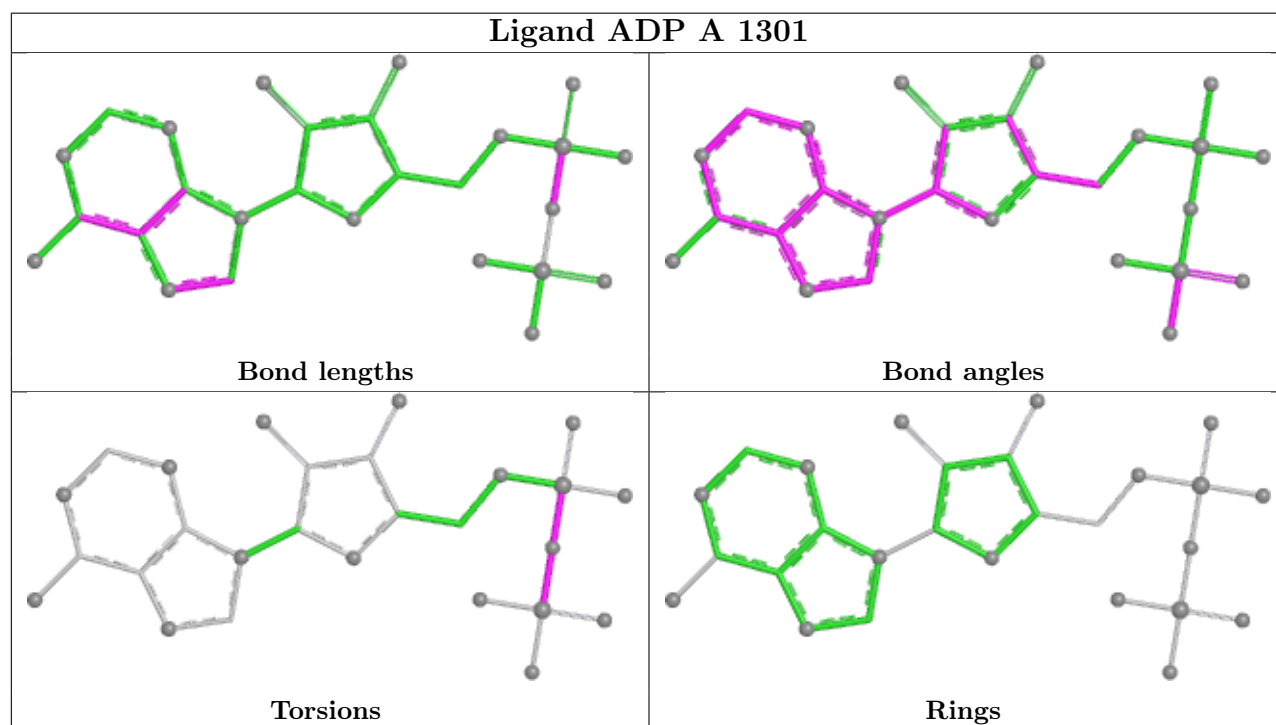
Mol	Chain	Res	Type	Atoms
2	A	1301	ADP	PB-O3A-PA-O5'
2	A	1301	ADP	PA-O3A-PB-O1B

There are no ring outliers.

1 monomer is involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1301	ADP	9	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	207/218 (94%)	0.52	12 (5%) 29 25	46, 62, 82, 89	0
1	B	77/218 (35%)	0.34	5 (6%) 25 22	45, 56, 87, 101	0
All	All	284/436 (65%)	0.47	17 (5%) 27 24	45, 60, 83, 101	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	270	LYS	6.9
1	B	165	GLU	4.1
1	B	167	ALA	3.5
1	A	353	ILE	3.4
1	A	327	GLU	3.2
1	A	301	LYS	2.9
1	B	168	ASN	2.7
1	A	264	PHE	2.7
1	A	263	MET	2.5
1	A	318	VAL	2.3
1	B	170	ARG	2.3
1	A	155	LYS	2.2
1	A	326	GLY	2.2
1	B	241	MET	2.1
1	A	346	PHE	2.1
1	A	157	ARG	2.1
1	A	323	THR	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](#)

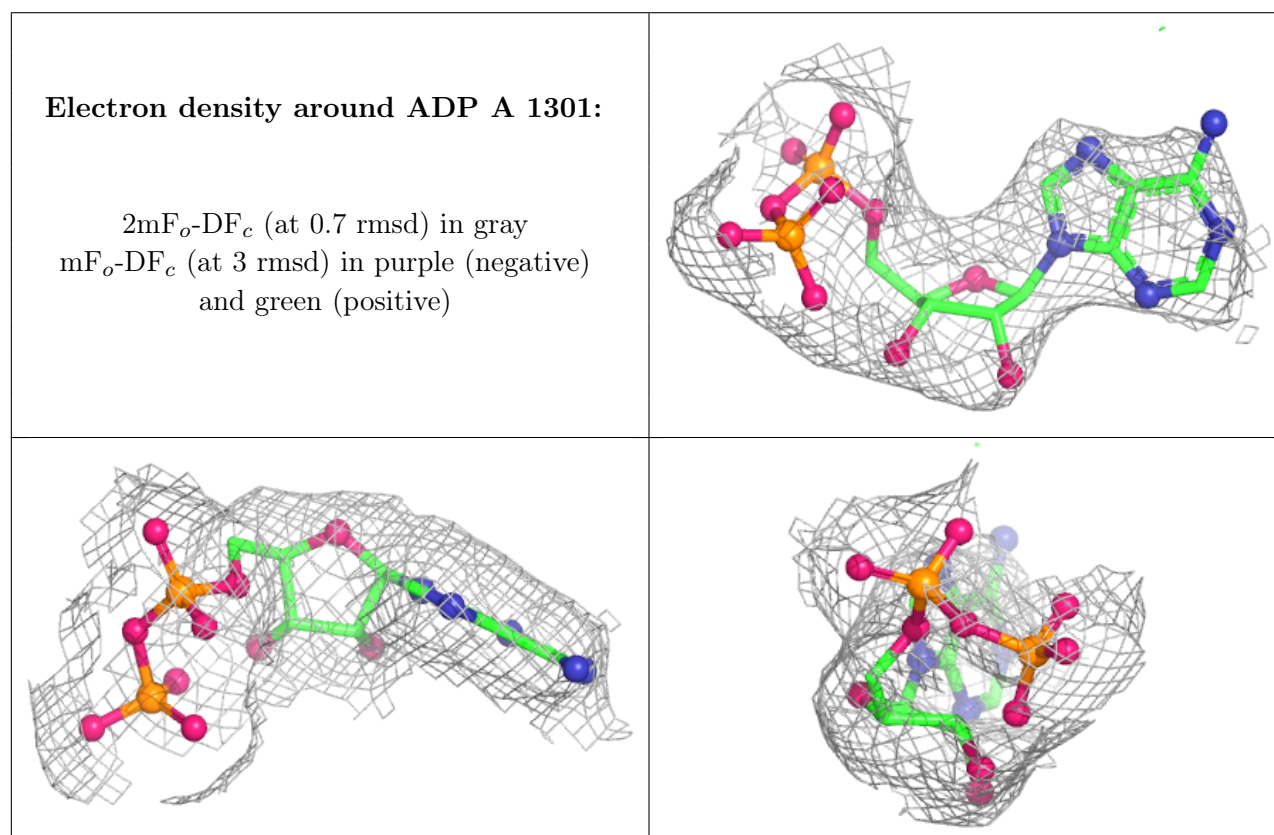
There are no oligosaccharides in this entry.

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
2	ADP	A	1301	27/27	0.94	0.09	69,76,81,87	0
3	CA	A	1303	1/1	0.96	0.07	66,66,66,66	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.