



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:16 PM UTC

PDB ID : 1OL7 / pdb_00001ol7
Title : Structure of Human Aurora-A 122-403 phosphorylated on Thr287, Thr288
Authors : Bayliss, R.; Conti, E.
Deposited on : 2003-08-06
Resolution : 2.75 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

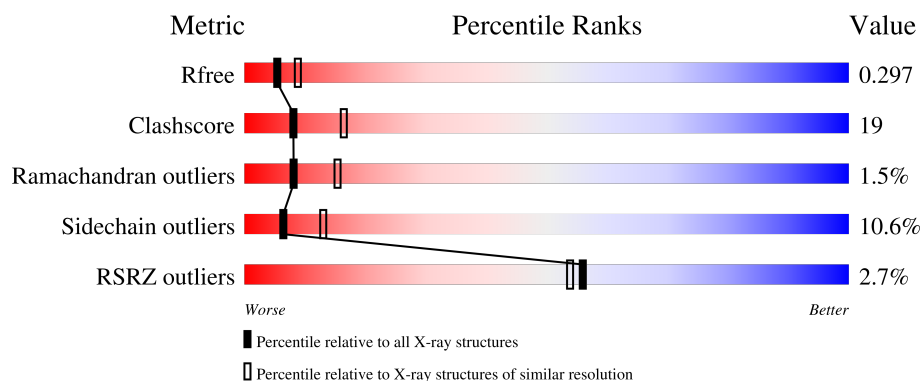
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.75 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	1009 (2.76-2.76)
Clashscore	190562	1044 (2.76-2.76)
Ramachandran outliers	187476	1024 (2.76-2.76)
Sidechain outliers	187428	1024 (2.76-2.76)
RSRZ outliers	180081	1009 (2.76-2.76)

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	282	

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
1	TPO	A	287	-	X	-	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
1	TPO	A	288	-	X	-	-

2 Entry composition [i](#)

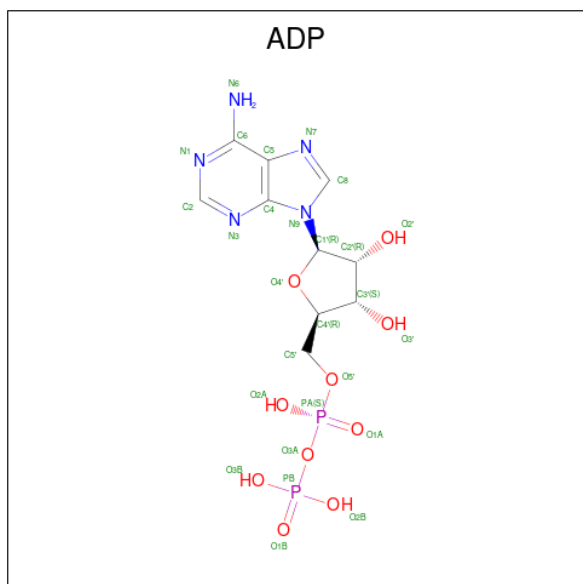
There are 4 unique types of molecules in this entry. The entry contains 2155 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 SERINE/THREONINE KINASE 6.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	P	S			
1	A	264	2117	1356	368	387	2	4	0	0	1

- 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
			Total	C	N	O	P		
2	A	1	27	10	5	10	2	0	0

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

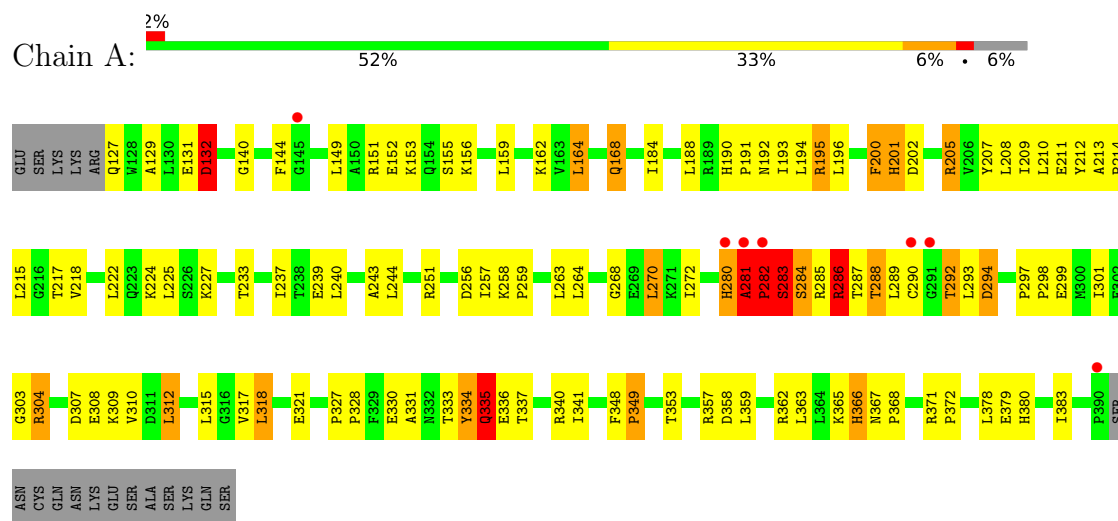
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	9	Total	O	0	0
			9	9		

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: SERINE/THREONINE KINASE 6



4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	81.18Å 81.18Å 169.62Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.75 20.00 – 2.75	Depositor EDS
% Data completeness (in resolution range)	100.0 (20.00-2.75) 99.6 (20.00-2.75)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.97 (at 2.73Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.257 , 0.296 0.248 , 0.297	Depositor DCC
R_{free} test set	440 reflections (4.79%)	wwPDB-VP
Wilson B-factor (Å ²)	61.6	Xtriage
Anisotropy	0.139	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 54.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	2155	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.20% 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, TPO, MG

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	1.03	8/2148 (0.4%)	1.46	32/2913 (1.1%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	200	PHE	CA-C	-8.18	1.42	1.53
1	A	282	PRO	CG-CD	6.43	1.72	1.50
1	A	290	CYS	C-O	6.31	1.31	1.23
1	A	290	CYS	CA-CB	-6.19	1.43	1.53
1	A	286	ARG	C-O	5.96	1.35	1.23
1	A	367	ASN	N-CA	-5.60	1.38	1.46
1	A	292	THR	N-CA	-5.54	1.39	1.46
1	A	367	ASN	CA-C	-5.26	1.47	1.52

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	281	ALA	CA-C-N	24.29	150.20	119.84
1	A	281	ALA	C-N-CA	24.29	150.20	119.84
1	A	290	CYS	N-CA-C	-14.88	85.36	109.76
1	A	282	PRO	N-CA-C	-10.55	90.74	112.47
1	A	281	ALA	C-N-CD	-10.37	82.47	125.00
1	A	335	GLN	N-CA-C	-9.36	101.11	114.12
1	A	348	PHE	CA-C-N	8.97	128.82	120.21
1	A	348	PHE	C-N-CA	8.97	128.82	120.21
1	A	292	THR	N-CA-C	-8.16	93.63	107.99
1	A	336	GLU	N-CA-C	-8.11	100.36	112.04
1	A	201	HIS	N-CA-C	8.04	119.23	108.38
1	A	283	SER	N-CA-C	7.99	120.87	108.96
1	A	153	LYS	N-CA-C	7.22	120.06	111.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	149	LEU	N-CA-C	-6.92	101.61	110.53
1	A	303	GLY	N-CA-C	6.75	121.69	113.79
1	A	281	ALA	N-CA-C	6.68	124.58	109.81
1	A	153	LYS	CA-C-N	-6.54	110.39	122.09
1	A	153	LYS	C-N-CA	-6.54	110.39	122.09
1	A	282	PRO	CA-C-N	-6.53	111.17	121.87
1	A	282	PRO	C-N-CA	-6.53	111.17	121.87
1	A	335	GLN	N-CA-CB	-6.07	102.50	110.88
1	A	284	SER	N-CA-C	-5.94	101.40	109.95
1	A	282	PRO	CA-N-CD	-5.75	103.96	112.00
1	A	366	HIS	N-CA-C	-5.71	106.08	113.16
1	A	213	ALA	CA-C-N	5.54	125.00	119.24
1	A	213	ALA	C-N-CA	5.54	125.00	119.24
1	A	218	VAL	N-CA-C	-5.46	104.91	112.50
1	A	349	PRO	N-CA-C	-5.46	103.58	111.22
1	A	132	ASP	N-CA-C	-5.42	106.84	113.50
1	A	312	LEU	N-CA-C	-5.26	105.21	111.69
1	A	327	PRO	CA-C-N	5.09	124.70	119.56
1	A	327	PRO	C-N-CA	5.09	124.70	119.56

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	2117	0	2056	82	0
2	A	27	0	12	1	0
3	A	2	0	0	0	0
4	A	9	0	0	0	0
All	All	2155	0	2068	82	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 (82) 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:334:TYR:O	1:A:334:TYR:CD2	2.21	0.94
1:A:366:HIS:O	1:A:366:HIS:ND1	2.09	0.86
1:A:292:THR:O	1:A:292:THR:HG23	1.79	0.82
1:A:196:LEU:CD1	1:A:210:LEU:HD23	2.13	0.79
1:A:288:TPO:HG23	1:A:289:LEU:H	1.46	0.79
1:A:301:ILE:HD12	1:A:341:ILE:HG21	1.70	0.73
1:A:196:LEU:HD12	1:A:210:LEU:HD23	1.73	0.70
1:A:317:VAL:HG13	1:A:328:PRO:HD2	1.73	0.70
1:A:280:HIS:O	1:A:281:ALA:HB2	1.91	0.68
1:A:292:THR:O	1:A:292:THR:CG2	2.41	0.68
1:A:155:SER:O	1:A:156:LYS:HB2	1.95	0.66
1:A:297:PRO:HG3	1:A:310:VAL:HB	1.77	0.65
1:A:188:LEU:HD23	1:A:193:ILE:HG21	1.80	0.64
1:A:309:LYS:HD3	1:A:312:LEU:HD12	1.81	0.62
1:A:196:LEU:HD13	1:A:210:LEU:HD23	1.80	0.62
1:A:301:ILE:CD1	1:A:341:ILE:HG21	2.29	0.62
1:A:129:ALA:O	1:A:132:ASP:HB2	2.01	0.61
1:A:259:PRO:HD3	1:A:318:LEU:HD11	1.81	0.61
1:A:307:ASP:O	1:A:310:VAL:HG13	2.01	0.60
1:A:217:THR:HG22	1:A:263:LEU:HD23	1.82	0.59
1:A:200:PHE:O	1:A:201:HIS:HB3	2.01	0.59
1:A:196:LEU:HD13	1:A:210:LEU:CD2	2.32	0.59
1:A:299:GLU:HA	1:A:304:ARG:HE	1.67	0.59
1:A:164:LEU:O	1:A:205:ARG:HB2	2.02	0.59
1:A:286:ARG:N	1:A:286:ARG:HD2	2.16	0.58
1:A:190:HIS:ND1	1:A:191:PRO:HD2	2.18	0.58
1:A:215:LEU:HD12	1:A:268:GLY:HA2	1.86	0.58
1:A:194:LEU:HD11	1:A:211:GLU:O	2.03	0.58
1:A:192:ASN:HB2	1:A:243:ALA:HB2	1.86	0.58
1:A:307:ASP:OD1	1:A:309:LYS:HB2	2.04	0.58
1:A:283:SER:OG	1:A:284:SER:O	2.23	0.57
1:A:285:ARG:C	1:A:286:ARG:HD2	2.30	0.56
1:A:334:TYR:CD2	1:A:334:TYR:C	2.83	0.56
1:A:217:THR:HG22	1:A:263:LEU:CD2	2.36	0.55
1:A:244:LEU:HD21	1:A:257:ILE:HD11	1.89	0.55
1:A:184:ILE:O	1:A:188:LEU:HD13	2.06	0.55
1:A:162:LYS:HB3	1:A:208:LEU:HB2	1.88	0.54
1:A:299:GLU:HA	1:A:304:ARG:HH21	1.75	0.52
1:A:353:THR:O	1:A:357:ARG:HG3	2.10	0.52
1:A:144:PHE:O	1:A:164:LEU:HA	2.09	0.52
1:A:283:SER:OG	1:A:284:SER:N	2.41	0.52
1:A:129:ALA:HB1	1:A:131:GLU:OE1	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:288:TPO:HG23	1:A:289:LEU:N	2.21	0.51
1:A:288:TPO:CG2	1:A:289:LEU:H	2.18	0.51
1:A:286:ARG:N	1:A:286:ARG:CD	2.70	0.49
1:A:140:GLY:HA3	2:A:531:ADP:H1'	1.95	0.49
1:A:196:LEU:CD1	1:A:210:LEU:CD2	2.85	0.49
1:A:225:LEU:O	1:A:227:LYS:N	2.43	0.49
1:A:379:GLU:O	1:A:380:HIS:C	2.55	0.48
1:A:359:LEU:HD22	1:A:383:ILE:HG12	1.94	0.48
1:A:240:LEU:HD13	1:A:270:LEU:CD2	2.44	0.48
1:A:358:ASP:O	1:A:362:ARG:HG3	2.14	0.48
1:A:202:ASP:OD2	1:A:207:TYR:OH	2.30	0.47
1:A:334:TYR:O	1:A:334:TYR:HD2	1.88	0.47
1:A:363:LEU:O	1:A:371:ARG:HG2	2.15	0.47
1:A:240:LEU:HD13	1:A:270:LEU:HD21	1.96	0.47
1:A:151:ARG:HG2	1:A:156:LYS:HA	1.97	0.46
1:A:215:LEU:HB2	1:A:264:LEU:HB2	1.98	0.45
1:A:222:LEU:HD21	1:A:321:GLU:HG2	1.98	0.45
1:A:164:LEU:O	1:A:205:ARG:CB	2.65	0.44
1:A:192:ASN:ND2	1:A:239:GLU:HB3	2.32	0.44
1:A:233:THR:O	1:A:237:ILE:HG12	2.18	0.44
1:A:365:LYS:HB2	1:A:371:ARG:HG3	1.98	0.44
1:A:288:TPO:CG2	1:A:289:LEU:N	2.81	0.44
1:A:195:ARG:HE	1:A:195:ARG:HB3	1.37	0.43
1:A:292:THR:O	1:A:293:LEU:C	2.56	0.43
1:A:200:PHE:HD2	1:A:209:ILE:HD11	1.84	0.43
1:A:240:LEU:HD12	1:A:272:ILE:HD11	2.01	0.43
1:A:334:TYR:C	1:A:335:GLN:HG3	2.42	0.43
1:A:334:TYR:O	1:A:334:TYR:CG	2.69	0.43
1:A:366:HIS:O	1:A:366:HIS:CG	2.70	0.43
1:A:330:GLU:O	1:A:340:ARG:NH1	2.53	0.42
1:A:240:LEU:HD23	1:A:315:LEU:CD1	2.50	0.42
1:A:190:HIS:HB3	1:A:193:ILE:HD12	2.01	0.41
1:A:256:ASP:OD1	1:A:258:LYS:HE2	2.21	0.41
1:A:144:PHE:HA	1:A:168:GLN:NE2	2.36	0.41
1:A:200:PHE:CD2	1:A:209:ILE:HD11	2.56	0.41
1:A:293:LEU:O	1:A:294:ASP:C	2.62	0.41
1:A:298:PRO:C	1:A:304:ARG:HH21	2.29	0.41
1:A:331:ALA:HB3	1:A:337:THR:OG1	2.21	0.41
1:A:312:LEU:HD13	1:A:372:PRO:O	2.21	0.41
1:A:212:TYR:O	1:A:214:PRO:HD3	2.21	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	260/282 (92%)	240 (92%)	16 (6%)	4 (2%)	8	15

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	282	PRO
1	A	304	ARG
1	A	281	ALA
1	A	294	ASP

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	218/248 (88%)	195 (89%)	23 (11%)	6	13

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	127	GLN
1	A	132	ASP
1	A	152	GLU
1	A	159	LEU
1	A	164	LEU
1	A	168	GLN
1	A	195	ARG

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Mol	Chain	Res	Type
1	A	205	ARG
1	A	224	LYS
1	A	251	ARG
1	A	270	LEU
1	A	280	HIS
1	A	282	PRO
1	A	283	SER
1	A	286	ARG
1	A	308	GLU
1	A	318	LEU
1	A	333	THR
1	A	334	TYR
1	A	335	GLN
1	A	349	PRO
1	A	368	PRO
1	A	378	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	127	GLN
1	A	168	GLN
1	A	177	GLN
1	A	201	HIS
1	A	242	ASN
1	A	335	GLN
1	A	370	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

2 non-standard protein/DNA/RNA residues 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
1	TPO	A	288	1	8,10,11	4.43	7 (87%)	10,14,16	2.56	4 (40%)
1	TPO	A	287	1	8,10,11	4.12	7 (87%)	10,14,16	2.47	4 (40%)

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
1	TPO	A	288	1	-	4/9/11/13	-
1	TPO	A	287	1	-	3/9/11/13	-

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	288	TPO	P-OG1	8.71	1.74	1.59
1	A	287	TPO	P-OG1	7.66	1.72	1.59
1	A	287	TPO	P-O1P	4.88	1.65	1.50
1	A	288	TPO	P-O1P	4.76	1.65	1.50
1	A	287	TPO	P-O2P	4.30	1.70	1.54
1	A	288	TPO	O-C	4.03	1.35	1.20
1	A	287	TPO	CG2-CB	3.95	1.61	1.51
1	A	288	TPO	P-O2P	3.89	1.69	1.54
1	A	288	TPO	CG2-CB	3.25	1.59	1.51
1	A	287	TPO	P-O3P	2.94	1.65	1.54
1	A	288	TPO	P-O3P	2.88	1.65	1.54
1	A	287	TPO	O-C	2.52	1.29	1.20
1	A	288	TPO	CA-N	-2.35	1.40	1.47
1	A	287	TPO	CB-CA	2.03	1.58	1.53

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	287	TPO	O3P-P-O2P	5.63	128.93	107.80
1	A	288	TPO	O3P-P-O2P	5.32	127.75	107.80
1	A	288	TPO	CG2-CB-CA	4.66	122.35	113.26
1	A	287	TPO	CG2-CB-CA	3.32	119.72	113.26
1	A	287	TPO	O3P-P-O1P	-2.60	100.72	110.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	288	TPO	P-OG1-CB	2.54	130.22	123.33
1	A	288	TPO	O3P-P-O1P	-2.25	102.06	110.83
1	A	287	TPO	P-OG1-CB	2.21	129.32	123.33

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	287	TPO	N-CA-CB-OG1
1	A	287	TPO	O-C-CA-CB
1	A	288	TPO	N-CA-CB-OG1
1	A	288	TPO	CG2-CB-OG1-P
1	A	287	TPO	CG2-CB-OG1-P
1	A	288	TPO	CA-CB-OG1-P
1	A	288	TPO	N-CA-CB-CG2

There are no ring outliers.

1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	288	TPO	4	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 2 are 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	531	3	28,29,29	1.44	4 (14%)	43,45,45	1.08	2 (4%)

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	531	3	-	6/16/32/32	0/3/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	531	ADP	C2-N1	3.88	1.40	1.33
2	A	531	ADP	C8-N7	2.98	1.37	1.31
2	A	531	ADP	C5-C6	2.36	1.47	1.41
2	A	531	ADP	C6-N1	2.19	1.41	1.35

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	531	ADP	N3-C2-N1	-3.02	124.01	128.58
2	A	531	ADP	C2-N1-C6	2.69	123.16	118.73

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	531	ADP	O4'-C4'-C5'-O5'
2	A	531	ADP	C3'-C4'-C5'-O5'
2	A	531	ADP	PA-O3A-PB-O1B
2	A	531	ADP	C4'-C5'-O5'-PA
2	A	531	ADP	PA-O3A-PB-O2B
2	A	531	ADP	PA-O3A-PB-O3B

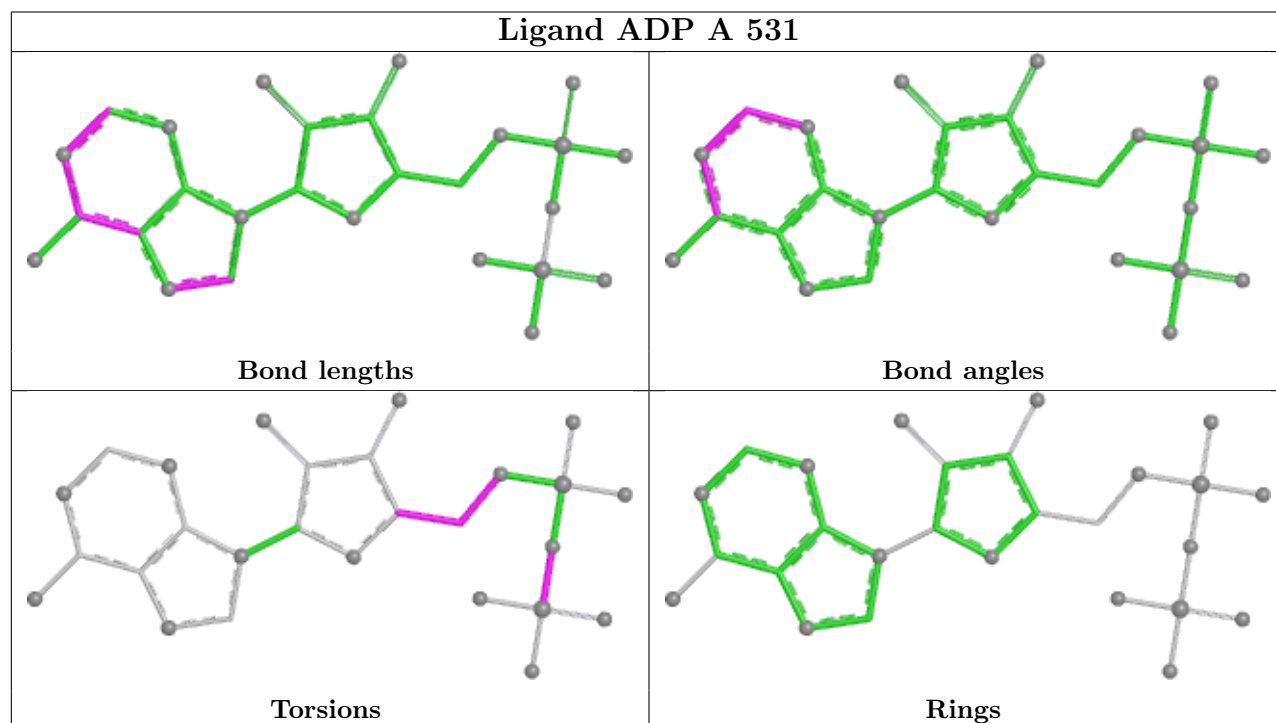
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	531	ADP	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is

within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	262/282 (92%)	0.21	7 (2%) 56 54	32, 53, 72, 88	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	290	CYS	3.4
1	A	390	PRO	2.8
1	A	145	GLY	2.5
1	A	291	GLY	2.3
1	A	281	ALA	2.3
1	A	282	PRO	2.2
1	A	280	HIS	2.2

6.2 Non-standard residues in protein, DNA, RNA chains

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
1	TPO	A	287	11/12	0.68	0.18	78,84,103,105	0
1	TPO	A	288	11/12	0.69	0.18	82,87,93,95	0

6.3 Carbohydrates

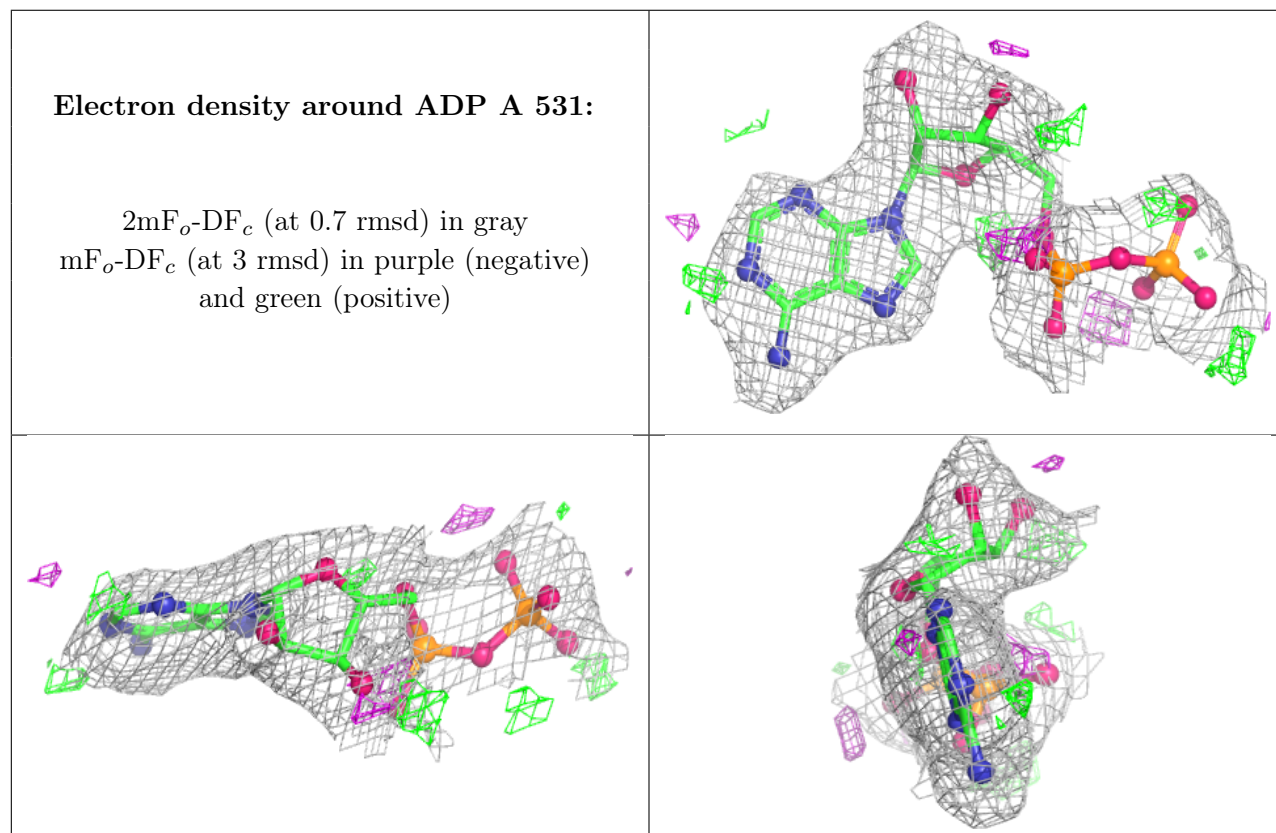
There are no oligosaccharides in this entry.

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
3	MG	A	1392	1/1	0.82	0.18	47,47,47,47	0
3	MG	A	1393	1/1	0.84	0.16	52,52,52,52	0
2	ADP	A	531	27/27	0.92	0.10	22,46,54,62	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.