



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 08:18 AM UTC

PDB ID : 1JBP / pdb_00001jbp
Title : Crystal Structure of the Catalytic Subunit of cAMP-dependent Protein Kinase
Complexed with a Substrate Peptide, ADP and Detergent
Authors : Madhusudan; Trafny, E.A.; Xuong, N.H.; Adams, J.A.; Ten Eyck, L.F.; Taylor,
S.S.; Sowadski, J.M.
Deposited on : 2001-06-06
Resolution : 2.20 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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

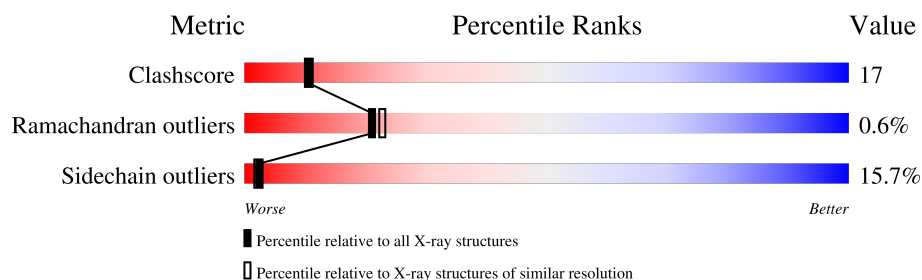
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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))
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	E	350	
2	S	20	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 3069 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 CAMP-DEPENDENT PROTEIN KINASE, ALPHA-CATALYTIC SUBUNIT.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	E	342	Total	C	N	O	P	S	0	0	0
			2711	1756	450	494	3	8			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	10	SEP	SER	modified residue	UNP P05132
E	197	TPO	THR	modified residue	UNP P05132
E	242	GLU	GLN	conflict	UNP P05132
E	338	SEP	SER	modified residue	UNP P05132

- Molecule 2 is a protein called CAMP-DEPENDENT PROTEIN KINASE INHIBITOR, MUSCLE/BRAIN FORM.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	S	20	Total	C	N	O	0	0	0
			155	93	31	31			

There are 2 discrepancies between the modelled and reference sequences:

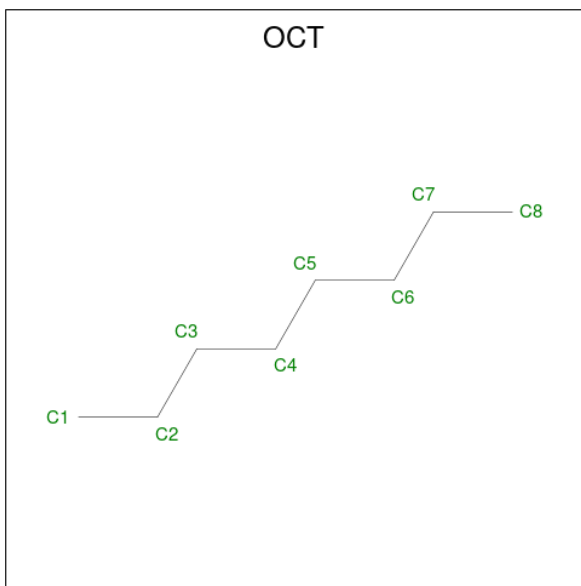
Chain	Residue	Modelled	Actual	Comment	Reference
S	376	ALA	ASN	engineered mutation	UNP P27776
S	377	SER	ALA	engineered mutation	UNP P27776

- Molecule 3 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



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

- Molecule 4 is N-OCTANE (CCD ID: OCT) (formula: C_8H_{18}).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	E	1	Total C 8 8	0	0

- Molecule 5 is water.

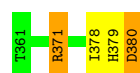
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	E	149	Total 149	O 149	0	0
5	S	19	Total 19	O 19	0	0

Note EDS was not executed.

- Chain E:  57% 31% 9% 3%



Chain S: 80% 10% 10%



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	73.96Å 76.11Å 81.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.20	Depositor
% Data completeness (in resolution range)	(Not available) (30.00-2.20)	Depositor
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	TNT	Depositor
R, R_{free}	0.175 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3069	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: OCT, ADP, TPO, SEP

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	E	1.23	2/2747 (0.1%)	1.68	37/3717 (1.0%)
2	S	1.20	0/157	1.66	1/209 (0.5%)
All	All	1.23	2/2904 (0.1%)	1.68	38/3926 (1.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	121	GLU	CD-OE2	5.15	1.35	1.25
1	E	165	ARG	N-CA	5.02	1.53	1.46

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	164	TYR	CA-C-N	-11.18	102.00	123.13
1	E	164	TYR	C-N-CA	-11.18	102.00	123.13
1	E	12	GLN	N-CA-C	-8.65	92.36	110.80
1	E	54	PHE	CA-CB-CG	-8.53	105.27	113.80
1	E	184	ASP	CA-CB-CG	-8.19	104.41	112.60
1	E	64	GLU	N-CA-C	8.12	120.21	111.36
1	E	163	ILE	CB-CA-C	-7.91	99.42	110.96
1	E	164	TYR	O-C-N	-7.59	114.57	123.22
1	E	255	VAL	CB-CA-C	-6.95	100.12	110.82
1	E	163	ILE	N-CA-CB	6.90	120.15	111.46
1	E	32	THR	CA-C-N	6.73	126.49	119.76
1	E	32	THR	C-N-CA	6.73	126.49	119.76
1	E	324	THR	N-CA-CB	-6.73	100.45	110.61
1	E	124	ALA	N-CA-C	6.52	120.97	112.89
1	E	194	ARG	N-CA-C	6.38	119.97	110.52
1	E	192	LYS	N-CA-C	-6.14	99.24	109.24
1	E	103	LEU	CA-C-N	-6.09	112.44	120.91

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	103	LEU	C-N-CA	-6.09	112.44	120.91
1	E	60	VAL	N-CA-CB	-5.97	104.16	112.52
1	E	192	LYS	CA-C-N	-5.94	111.90	122.00
1	E	192	LYS	C-N-CA	-5.94	111.90	122.00
1	E	75	ASP	CA-CB-CG	5.92	118.52	112.60
1	E	342	LYS	N-CA-C	5.90	123.36	110.80
1	E	73	ILE	N-CA-C	5.70	116.08	107.75
1	E	32	THR	O-C-N	5.54	125.88	121.17
1	E	222	TRP	N-CA-C	-5.52	105.18	111.14
1	E	155	GLU	N-CA-C	-5.50	105.37	111.36
1	E	115	ASN	N-CA-C	5.47	117.29	109.14
2	S	371	ARG	N-CA-C	5.42	116.94	110.44
1	E	302	TRP	N-CA-C	5.38	117.84	111.33
1	E	244	ILE	N-CA-C	5.33	116.06	110.62
1	E	301	ASP	N-CA-CB	5.29	119.30	110.53
1	E	87	HIS	CA-CB-CG	-5.26	108.54	113.80
1	E	174	ILE	CB-CA-C	-5.25	103.30	110.96
1	E	20	ALA	CB-CA-C	-5.13	102.82	110.88
1	E	102	PHE	CA-CB-CG	-5.08	108.72	113.80
1	E	35	GLN	CB-CG-CD	-5.03	104.06	112.60
1	E	85	ILE	N-CA-C	5.01	115.23	110.42

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	E	2711	0	2593	94	0
2	S	155	0	145	4	0
3	E	27	0	12	3	0
4	E	8	0	18	0	0
5	E	149	0	0	10	0
5	S	19	0	0	0	0
All	All	3069	0	2768	96	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 17.

All (96) 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:E:285:LYS:HD3	1:E:286:ASN:H	1.12	1.12
1:E:73:ILE:HD13	1:E:335:ILE:HD11	1.51	0.93
1:E:285:LYS:HD3	1:E:286:ASN:N	1.87	0.88
1:E:268:LEU:HD22	1:E:272:LEU:HD22	1.55	0.88
1:E:210:ILE:HD13	1:E:247:TYR:HB3	1.59	0.84
1:E:11:GLU:C	1:E:12:GLN:O	2.15	0.78
1:E:305:ILE:HD13	1:E:310:VAL:HG21	1.66	0.78
1:E:134:ARG:HD2	5:E:479:HOH:O	1.85	0.76
1:E:236:PRO:HG2	1:E:239:PHE:HB3	1.69	0.73
1:E:142:HIS:CD2	1:E:313:PRO:HG2	2.26	0.71
1:E:104:VAL:HB	5:E:388:HOH:O	1.89	0.71
1:E:131:HIS:O	1:E:135:ILE:HG13	1.89	0.70
1:E:107:GLU:HB3	1:E:108:PHE:CD2	2.28	0.68
1:E:286:ASN:O	1:E:289:ASN:HB2	1.93	0.68
1:E:189:LYS:HG2	1:E:191:VAL:HG13	1.75	0.68
1:E:295:LYS:HD3	1:E:295:LYS:H	1.60	0.67
1:E:284:LEU:HB2	1:E:285:LYS:HD2	1.77	0.67
1:E:189:LYS:HE2	1:E:191:VAL:CG1	2.25	0.66
1:E:50:GLY:HA2	1:E:330:TYR:CE1	2.30	0.66
1:E:305:ILE:CD1	1:E:310:VAL:HG21	2.25	0.65
1:E:12:GLN:O	1:E:14:SER:N	2.28	0.65
1:E:200:GLY:O	2:S:378:ILE:HG13	1.97	0.65
1:E:52:GLY:HA3	3:E:381:ADP:O2B	1.98	0.64
1:E:103:LEU:HD22	1:E:185:PHE:HZ	1.62	0.64
1:E:342:LYS:NZ	5:E:447:HOH:O	2.30	0.63
1:E:305:ILE:HD13	1:E:310:VAL:CG2	2.31	0.61
1:E:47:LYS:HD3	1:E:324:THR:HG21	1.82	0.61
1:E:52:GLY:HA3	3:E:381:ADP:PB	2.41	0.60
1:E:284:LEU:N	1:E:284:LEU:HD12	2.15	0.60
1:E:303:ILE:O	1:E:307:GLN:HG3	1.99	0.60
1:E:247:TYR:O	1:E:251:VAL:HG22	2.01	0.60
1:E:284:LEU:HD12	1:E:284:LEU:H	1.66	0.59
1:E:92:LYS:HE3	1:E:350:PHE:OXT	2.02	0.59
1:E:326:ASN:HB2	5:E:410:HOH:O	2.01	0.59
1:E:262:SER:HB2	5:E:462:HOH:O	2.03	0.57
1:E:9:GLY:N	5:E:514:HOH:O	2.36	0.57
1:E:240:ALA:HB3	1:E:246:ILE:HG13	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:44:ASP:OD1	1:E:63:LYS:HE3	2.06	0.56
1:E:284:LEU:HD13	1:E:290:ASP:OD1	2.05	0.56
1:E:104:VAL:HG13	1:E:121:GLU:OE2	2.05	0.55
1:E:164:TYR:O	1:E:220:ASP:OD1	2.24	0.55
1:E:202:PRO:HA	1:E:205:LEU:HD22	1.88	0.55
1:E:277:LEU:HD12	1:E:277:LEU:N	2.21	0.55
1:E:323:ASP:CG	1:E:325:SER:HG	2.15	0.55
1:E:221:TRP:CH2	1:E:288:VAL:HG23	2.41	0.54
2:S:379:HIS:O	2:S:380:ASP:OXT	2.25	0.54
1:E:85:ILE:O	1:E:89:LEU:HG	2.07	0.54
1:E:295:LYS:H	1:E:295:LYS:CD	2.17	0.54
1:E:216:ASN:HB2	5:E:542:HOH:O	2.08	0.54
1:E:158:HIS:HE1	1:E:220:ASP:OD2	1.92	0.53
1:E:50:GLY:HA2	1:E:330:TYR:CZ	2.45	0.52
1:E:10:SEP:O	1:E:12:GLN:O	2.27	0.51
1:E:203:GLU:OE2	2:S:371:ARG:HD3	2.10	0.51
1:E:323:ASP:OD2	1:E:325:SER:OG	2.30	0.50
1:E:277:LEU:H	1:E:277:LEU:CD1	2.24	0.50
1:E:11:GLU:O	1:E:12:GLN:O	2.30	0.49
1:E:80:VAL:HG22	1:E:85:ILE:HD11	1.94	0.49
1:E:47:LYS:HE3	5:E:384:HOH:O	2.13	0.49
1:E:52:GLY:HA3	3:E:381:ADP:O3B	2.13	0.49
1:E:113:ASN:ND2	1:E:341:GLU:HB2	2.30	0.47
1:E:244:ILE:O	1:E:248:GLU:HG3	2.16	0.46
1:E:175:ASP:OD1	1:E:179:TYR:N	2.45	0.46
1:E:277:LEU:N	1:E:277:LEU:CD1	2.79	0.46
1:E:180:ILE:HG23	1:E:181:GLN:N	2.31	0.45
1:E:228:ILE:O	1:E:229:TYR:C	2.58	0.45
1:E:62:HIS:HD2	1:E:65:SER:OG	1.99	0.45
1:E:111:LYS:NZ	1:E:350:PHE:O	2.45	0.45
1:E:195:THR:CG2	1:E:215:TYR:CE1	3.00	0.45
1:E:335:ILE:O	1:E:335:ILE:HG22	2.15	0.45
1:E:59:LEU:HD22	1:E:60:VAL:N	2.33	0.44
1:E:191:VAL:HG23	1:E:191:VAL:O	2.17	0.44
1:E:157:LEU:O	1:E:162:LEU:HB2	2.16	0.44
1:E:272:LEU:HD12	1:E:272:LEU:HA	1.78	0.44
1:E:107:GLU:N	1:E:119:VAL:O	2.39	0.44
1:E:95:LEU:HD23	1:E:95:LEU:HA	1.77	0.44
1:E:243:PRO:O	1:E:246:ILE:HB	2.18	0.44
1:E:47:LYS:CE	1:E:324:THR:HG21	2.48	0.44
1:E:319:LYS:O	1:E:323:ASP:HB2	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:47:LYS:CD	1:E:324:THR:HG21	2.48	0.43
1:E:122:TYR:CE2	1:E:124:ALA:HB2	2.54	0.43
1:E:197:TPO:HG23	1:E:198:LEU:N	2.32	0.43
1:E:59:LEU:HD23	1:E:59:LEU:HA	1.82	0.43
1:E:195:THR:HG23	1:E:215:TYR:CZ	2.54	0.43
1:E:39:GLN:OE1	1:E:42:GLN:HG3	2.19	0.42
1:E:277:LEU:HD12	1:E:277:LEU:H	1.83	0.42
1:E:285:LYS:N	1:E:285:LYS:CD	2.81	0.42
1:E:338:SEP:C	1:E:340:ASN:H	2.33	0.42
1:E:192:LYS:NZ	5:E:490:HOH:O	2.54	0.41
1:E:25:ASP:N	1:E:25:ASP:OD2	2.53	0.41
1:E:32:THR:N	1:E:33:PRO:HD3	2.36	0.41
1:E:75:ASP:O	1:E:79:VAL:HG23	2.21	0.41
1:E:342:LYS:HD3	1:E:342:LYS:HA	1.76	0.41
1:E:229:TYR:O	1:E:230:GLU:C	2.61	0.40
2:S:371:ARG:HD3	2:S:371:ARG:HH11	1.67	0.40
1:E:273:LEU:HD23	1:E:273:LEU:HA	1.86	0.40
1:E:324:THR:HG23	5:E:384:HOH:O	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	E	337/350 (96%)	317 (94%)	18 (5%)	2 (1%)	21	23
2	S	18/20 (90%)	18 (100%)	0	0	100	100
All	All	355/370 (96%)	335 (94%)	18 (5%)	2 (1%)	21	23

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	12	GLN
1	E	53	SER

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	E	271/302 (90%)	227 (84%)	44 (16%)	2	2
2	S	15/15 (100%)	14 (93%)	1 (7%)	15	17
All	All	286/317 (90%)	241 (84%)	45 (16%)	2	2

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	13	GLU
1	E	19	LEU
1	E	25	ASP
1	E	27	LEU
1	E	36	ASN
1	E	39	GLN
1	E	40	LEU
1	E	54	PHE
1	E	59	LEU
1	E	60	VAL
1	E	61	LYS
1	E	74	LEU
1	E	75	ASP
1	E	78	LYS
1	E	95	LEU
1	E	114	SER
1	E	120	MET
1	E	121	GLU
1	E	133	ARG
1	E	160	LEU
1	E	162	LEU
1	E	163	ILE

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Mol	Chain	Res	Type
1	E	177	GLN
1	E	180	ILE
1	E	198	LEU
1	E	205	LEU
1	E	210	ILE
1	E	213	LYS
1	E	244	ILE
1	E	245	GLN
1	E	255	VAL
1	E	268	LEU
1	E	269	LEU
1	E	272	LEU
1	E	278	THR
1	E	284	LEU
1	E	285	LYS
1	E	288	VAL
1	E	295	LYS
1	E	299	THR
1	E	324	THR
1	E	335	ILE
1	E	343	CYS
1	E	348	THR
2	S	380	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	E	36	ASN
1	E	42	GLN
1	E	62	HIS
1	E	67	ASN
1	E	99	ASN
1	E	113	ASN
1	E	158	HIS
1	E	181	GLN
1	E	245	GLN
1	E	271	ASN
1	E	286	ASN
2	S	379	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

3 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	SEP	E	10	1	8,9,10	0.98	0	7,12,14	2.56	1 (14%)
1	SEP	E	338	1	8,9,10	1.04	0	7,12,14	3.15	3 (42%)
1	TPO	E	197	1	8,10,11	1.04	1 (12%)	10,14,16	1.71	3 (30%)

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	SEP	E	10	1	-	3/6/8/10	-
1	SEP	E	338	1	-	3/6/8/10	-
1	TPO	E	197	1	-	1/9/11/13	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	197	TPO	P-O3P	-2.25	1.46	1.54

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	10	SEP	OG-CB-CA	6.47	114.44	108.14
1	E	338	SEP	O3P-P-O1P	6.05	134.43	110.83
1	E	338	SEP	O3P-P-O2P	-4.57	90.65	107.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	197	TPO	O3P-P-O1P	-3.06	98.90	110.83
1	E	338	SEP	OG-CB-CA	2.99	111.05	108.14
1	E	197	TPO	O3P-P-O2P	2.78	118.21	107.80
1	E	197	TPO	CG2-CB-CA	-2.26	108.85	113.26

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	E	10	SEP	N-CA-CB-OG
1	E	10	SEP	C-CA-CB-OG
1	E	10	SEP	CB-OG-P-O3P
1	E	197	TPO	O-C-CA-CB
1	E	338	SEP	N-CA-CB-OG
1	E	338	SEP	C-CA-CB-OG
1	E	338	SEP	CA-CB-OG-P

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	E	10	SEP	1	0
1	E	338	SEP	1	0
1	E	197	TPO	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 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
4	OCT	E	382	-	7,7,7	0.62	0	6,6,6	0.21	0
3	ADP	E	381	-	28,29,29	0.79	1 (3%)	43,45,45	0.81	1 (2%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	OCT	E	382	-	-	3/5/5/5	-
3	ADP	E	381	-	-	4/16/32/32	0/3/3/3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	381	ADP	PB-O3B	2.33	1.63	1.54

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	381	ADP	O3B-PB-O3A	-2.29	96.95	104.64

There are no chirality outliers.

All (7) torsion outliers are listed below:

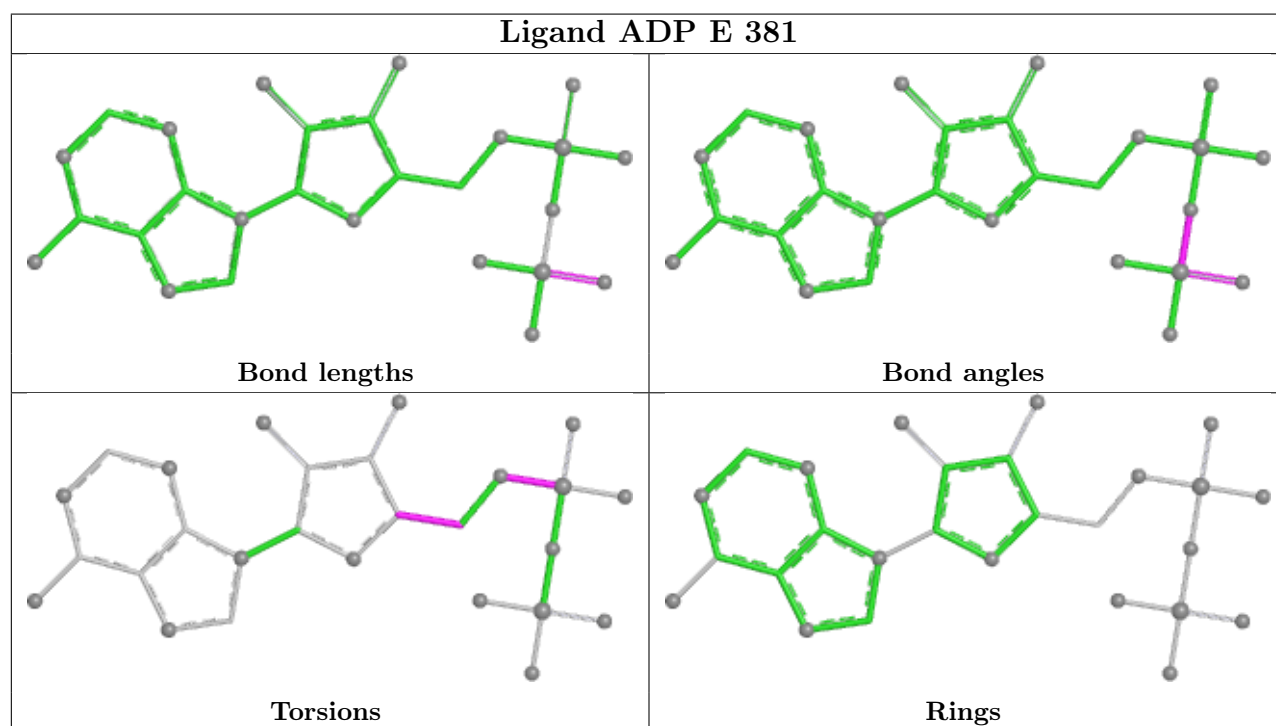
Mol	Chain	Res	Type	Atoms
3	E	381	ADP	C5'-O5'-PA-O2A
4	E	382	OCT	C1-C2-C3-C4
4	E	382	OCT	C2-C3-C4-C5
4	E	382	OCT	C5-C6-C7-C8
3	E	381	ADP	C5'-O5'-PA-O1A
3	E	381	ADP	C5'-O5'-PA-O3A
3	E	381	ADP	O4'-C4'-C5'-O5'

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	E	381	ADP	3	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

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.