Full wwPDB X-ray Structure Validation Report (i)

Feb 6, 2024 – 08:03 PM EST

PDB ID	:	2G1Q
Title	:	crystal structure of KSP in complex with inhibitor 9h
Authors	:	Yan, Y.
Deposited on	:	2006-02-14
Resolution	:	2.51 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 (i) 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 (1)) were used in the production of this report:

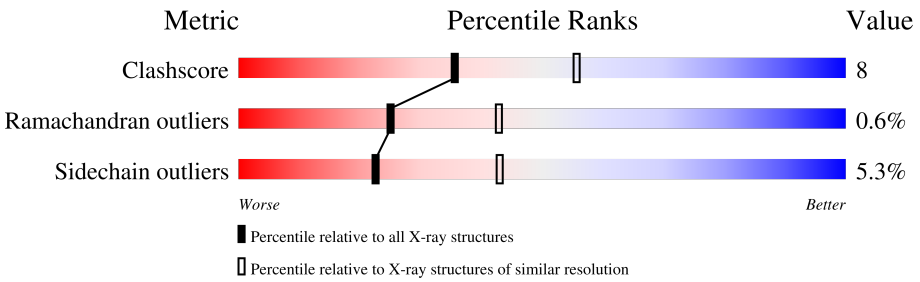
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.51 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	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)

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 <=5%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	368	67% 19% • 10%
1	B	368	70% 18% •• 10%

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 5469 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.

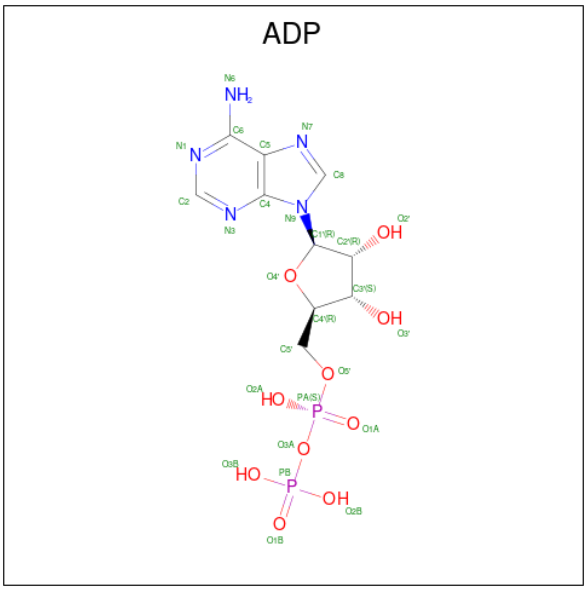
- Molecule 1 is a protein called Kinesin-like protein KIF11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	330	2594	1624	452	508	10	0	0	0
1	B	330	2594	1624	452	508	10	0	0	0

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

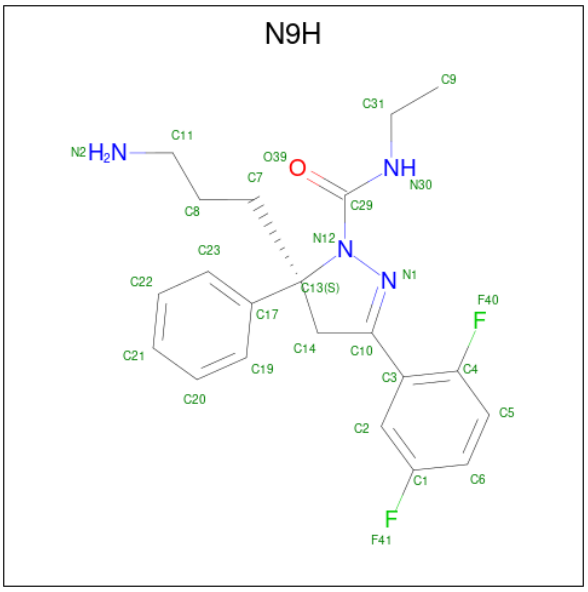
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Mg		
2	A	1	1	1	0	0
2	B	1	1	1	0	0

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



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

- Molecule 4 is (5S)-5-(3-AMINOPROPYL)-3-(2,5-DIFLUOROPHENYL)-N-ETHYL-5-PHENYL-4,5-DIHYDRO-1H-PYRAZOLE-1-CARBOXAMIDE (three-letter code: N9H) (formula: $C_{21}H_{24}F_2N_4O$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	F	N	O		
4	A	1	28	21	2	4	1	0	0
			Total	C	F	N	O		
4	B	1	28	21	2	4	1	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	A	90	90	90	0	0
			Total	O		
5	B	79	79	79	0	0

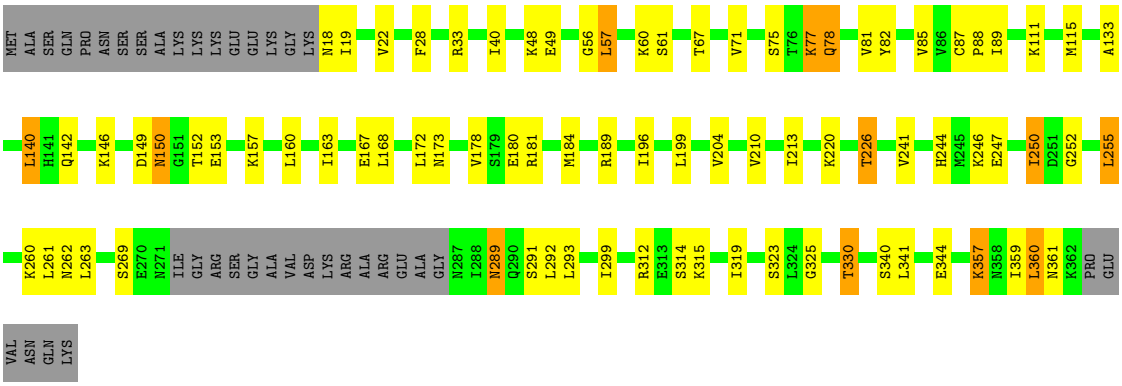
3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. 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. 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.

Note EDS was not executed.

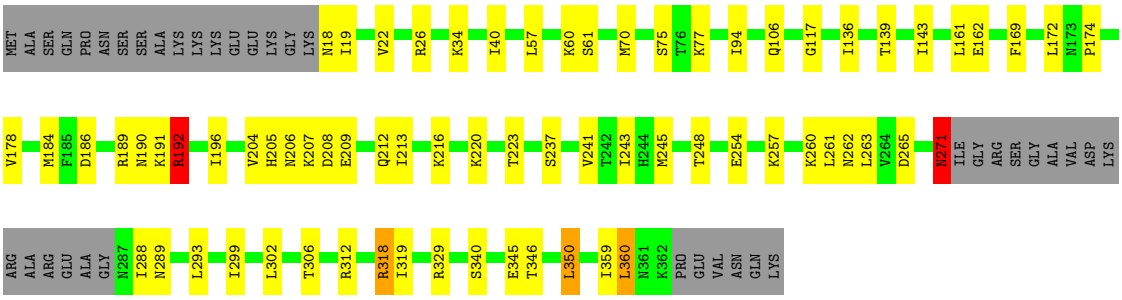
• Molecule 1: Kinesin-like protein KIF11

Chain A: 67% 19% • 10%



• Molecule 1: Kinesin-like protein KIF11

Chain B: 70% 18% •• 10%



4 Data and refinement statistics (i)

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	68.68Å 79.64Å 158.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.51	Depositor
% Data completeness (in resolution range)	97.5 (50.00-2.51)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.234 , 0.279	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5469	wwPDB-VP
Average B, all atoms (Å^2)	39.0	wwPDB-VP

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: MG, N9H, 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	0.49	1/2632 (0.0%)	0.71	2/3559 (0.1%)
1	B	0.52	3/2632 (0.1%)	0.72	5/3559 (0.1%)
All	All	0.50	4/5264 (0.1%)	0.71	7/7118 (0.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	344	GLU	CG-CD	5.95	1.60	1.51
1	B	271	ASN	CB-CG	5.68	1.64	1.51
1	B	162	GLU	CG-CD	-5.39	1.43	1.51
1	B	345	GLU	CG-CD	5.06	1.59	1.51

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	192	ARG	CB-CG-CD	8.40	133.44	111.60
1	B	207	LYS	CB-CG-CD	-6.36	95.07	111.60
1	B	34	LYS	CB-CG-CD	6.18	127.68	111.60
1	A	77	LYS	CB-CG-CD	5.46	125.80	111.60
1	B	60	LYS	CB-CG-CD	5.45	125.78	111.60
1	B	318	ARG	CB-CG-CD	5.42	125.69	111.60
1	A	357	LYS	CB-CG-CD	5.09	124.83	111.60

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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	2594	0	2618	50	0
1	B	2594	0	2618	36	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	27	0	12	0	0
3	B	27	0	12	0	0
4	A	28	0	24	0	0
4	B	28	0	24	1	0
5	A	90	0	0	2	0
5	B	79	0	0	0	0
All	All	5469	0	5308	86	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 8.

All (86) 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:78:GLN:HG2	1:A:133:ALA:O	1.81	0.80
1:B:248:THR:HG22	1:B:254:GLU:HG2	1.65	0.79
1:A:204:VAL:HG11	1:A:210:VAL:HG12	1.69	0.73
1:A:299:ILE:HG23	1:A:359:ILE:HD11	1.71	0.70
1:A:250:ILE:H	1:A:250:ILE:HD13	1.59	0.68
1:A:167:GLU:CD	1:A:181:ARG:HE	1.96	0.68
1:B:18:ASN:HB3	1:B:360:LEU:HA	1.76	0.68
1:A:167:GLU:HG2	1:A:181:ARG:HH21	1.60	0.66
1:A:19:ILE:HG12	1:A:330:THR:HG23	1.81	0.62
1:B:161:LEU:HD21	1:B:319:ILE:HD13	1.81	0.62
1:B:190:ASN:HB2	1:B:192:ARG:HG2	1.82	0.62
1:A:255:LEU:HD23	1:A:255:LEU:H	1.63	0.62
1:A:67:THR:O	1:A:357:LYS:HE3	2.00	0.61
1:B:212:GLN:O	1:B:216:LYS:HG2	2.01	0.60
1:A:85:VAL:O	1:A:89:ILE:HG12	2.02	0.60
1:B:19:ILE:HD12	1:B:359:ILE:HB	1.84	0.59
1:B:346:THR:O	1:B:350:LEU:HD22	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:323:SER:O	1:A:330:THR:HG21	2.02	0.59
1:B:172:LEU:O	1:B:174:PRO:HD3	2.03	0.59
1:A:184:MET:HE1	1:A:315:LYS:HG3	1.83	0.58
1:B:57:LEU:O	1:B:61:SER:HB3	2.04	0.57
1:B:161:LEU:HD22	1:B:196:ILE:HD13	1.86	0.57
1:A:325:GLY:HA2	1:A:361:ASN:HB3	1.85	0.57
1:B:245:MET:HB2	1:B:257:LYS:HG3	1.86	0.57
1:B:206:ASN:HD22	1:B:208:ASP:H	1.53	0.56
1:B:178:VAL:HG12	1:B:220:LYS:HE3	1.88	0.56
1:B:22:VAL:HG12	1:B:70:MET:HB2	1.86	0.56
1:B:260:LYS:HE2	1:B:262:ASN:HD21	1.73	0.54
1:A:312:ARG:HG2	5:A:668:HOH:O	2.08	0.53
1:B:143:ILE:HD13	1:B:243:ILE:HD11	1.90	0.52
1:A:142:GLN:O	1:A:146:LYS:HB2	2.10	0.52
1:A:60:LYS:O	1:A:61:SER:HB3	2.09	0.52
1:B:205:HIS:HB2	1:B:209:GLU:OE2	2.10	0.51
1:A:168:LEU:HD21	1:A:315:LYS:HG2	1.93	0.51
1:A:163:ILE:HG12	1:A:168:LEU:CD1	2.40	0.51
1:A:82:TYR:OH	1:A:142:GLN:HG3	2.10	0.51
1:A:48:LYS:HA	1:A:71:VAL:HG22	1.92	0.51
1:A:204:VAL:HG22	1:A:213:ILE:CD1	2.41	0.50
1:B:245:MET:HB2	1:B:257:LYS:CG	2.42	0.50
1:A:153:GLU:HB3	1:A:246:LYS:HB3	1.95	0.49
1:A:196:ILE:HB	1:A:199:LEU:HD12	1.94	0.48
1:B:248:THR:CG2	1:B:254:GLU:HG2	2.40	0.48
1:B:136:ILE:HG12	1:B:263:LEU:HD13	1.96	0.48
1:B:360:LEU:HD12	1:B:360:LEU:H	1.78	0.48
1:A:157:LYS:HD2	1:A:244:HIS:HE1	1.78	0.48
1:B:271:ASN:C	1:B:271:ASN:HD22	2.17	0.48
1:B:312:ARG:HA	1:B:318:ARG:HD2	1.96	0.47
1:B:94:ILE:O	1:B:257:LYS:HE2	2.14	0.47
1:A:28:PHE:HB2	1:A:33:ARG:NE	2.30	0.47
1:A:81:VAL:O	1:A:85:VAL:HG22	2.15	0.47
1:A:115:MET:CE	1:A:263:LEU:HB3	2.45	0.47
1:B:241:VAL:HG13	1:B:261:LEU:HB3	1.97	0.47
1:B:139:THR:O	1:B:143:ILE:HG13	2.14	0.46
1:B:329:ARG:HG2	1:B:329:ARG:HH11	1.79	0.46
1:A:241:VAL:CG1	1:A:261:LEU:HB3	2.46	0.46
1:B:40:ILE:HD13	1:B:340:SER:HA	1.98	0.46
1:B:299:ILE:HG23	1:B:359:ILE:HD11	1.98	0.46
1:A:260:LYS:HE2	1:A:262:ASN:HD21	1.80	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:288:ILE:O	1:B:288:ILE:HG13	2.16	0.45
1:A:111:LYS:HE3	1:A:111:LYS:HB2	1.53	0.45
1:A:75:SER:O	1:A:77:LYS:HE3	2.16	0.45
1:A:315:LYS:O	1:A:319:ILE:HG13	2.17	0.45
1:B:237:SER:OG	1:B:265:ASP:HB3	2.18	0.44
1:B:169:PHE:HD2	1:B:178:VAL:O	2.01	0.43
1:A:160:LEU:HB3	1:A:172:LEU:HG	2.00	0.43
1:A:330:THR:HG22	1:A:361:ASN:HD21	1.82	0.43
1:A:152:THR:HB	1:A:247:GLU:HA	2.00	0.43
1:A:18:ASN:HB2	1:A:360:LEU:HD13	2.00	0.43
1:A:241:VAL:O	1:A:241:VAL:HG13	2.18	0.43
1:A:226:THR:HG22	5:A:663:HOH:O	2.19	0.42
1:A:87:CYS:HB3	1:A:88:PRO:HD3	2.02	0.42
1:A:150:ASN:HD22	1:A:150:ASN:HA	1.57	0.42
1:A:204:VAL:CG1	1:A:210:VAL:HG12	2.43	0.42
1:B:243:ILE:HG22	1:B:245:MET:HG3	2.01	0.42
1:A:78:GLN:CG	1:A:133:ALA:O	2.62	0.41
1:A:140:LEU:HD12	1:A:140:LEU:HA	1.90	0.41
1:A:178:VAL:HG13	1:A:220:LYS:NZ	2.35	0.41
1:B:184:MET:HE3	1:B:318:ARG:NH2	2.34	0.41
1:A:56:GLY:O	1:A:57:LEU:HB2	2.20	0.41
1:B:117:GLY:O	4:B:605:N9H:H111	2.20	0.41
1:A:269:SER:HA	1:A:292:LEU:HD21	2.03	0.41
1:B:204:VAL:HG22	1:B:213:ILE:CD1	2.51	0.41
1:A:167:GLU:HG2	1:A:181:ARG:NH2	2.33	0.41
1:A:289:ASN:ND2	1:A:292:LEU:H	2.18	0.41
1:A:291:SER:HA	1:A:314:SER:HB2	2.02	0.40
1:A:40:ILE:HD13	1:A:340:SER:HA	2.03	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	326/368 (89%)	305 (94%)	18 (6%)	3 (1%)	17	31
1	B	326/368 (89%)	316 (97%)	9 (3%)	1 (0%)	41	61
All	All	652/736 (89%)	621 (95%)	27 (4%)	4 (1%)	25	43

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	189	ARG
1	A	252	GLY
1	B	189	ARG
1	A	57	LEU

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	293/322 (91%)	277 (94%)	16 (6%)	21	41
1	B	293/322 (91%)	278 (95%)	15 (5%)	24	45
All	All	586/644 (91%)	555 (95%)	31 (5%)	22	43

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	VAL
1	A	49	GLU
1	A	78	GLN
1	A	140	LEU
1	A	149	ASP
1	A	150	ASN
1	A	173	ASN
1	A	180	GLU
1	A	226	THR
1	A	250	ILE
1	A	255	LEU
1	A	289	ASN

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Mol	Chain	Res	Type
1	A	293	LEU
1	A	330	THR
1	A	341	LEU
1	A	360	LEU
1	B	26	ARG
1	B	75	SER
1	B	77	LYS
1	B	106	GLN
1	B	186	ASP
1	B	191	LYS
1	B	192	ARG
1	B	223	THR
1	B	271	ASN
1	B	289	ASN
1	B	293	LEU
1	B	302	LEU
1	B	306	THR
1	B	350	LEU
1	B	360	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	122	ASN
1	A	150	ASN
1	A	229	ASN
1	A	244	HIS
1	A	262	ASN
1	A	271	ASN
1	A	289	ASN
1	A	321	GLN
1	A	361	ASN
1	B	18	ASN
1	B	205	HIS
1	B	206	ASN
1	B	262	ASN
1	B	271	ASN
1	B	287	ASN
1	B	289	ASN
1	B	290	GLN
1	B	321	GLN
1	B	342	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 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	$\#\|Z\| > 2$
4	N9H	B	605	-	29,30,30	2.39	14 (48%)	32,42,42	4.17	11 (34%)
3	ADP	B	600	2	24,29,29	1.42	4 (16%)	29,45,45	1.62	3 (10%)
4	N9H	A	604	-	29,30,30	3.06	17 (58%)	32,42,42	4.33	9 (28%)
3	ADP	A	601	2	24,29,29	1.53	5 (20%)	29,45,45	1.56	2 (6%)

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	N9H	B	605	-	-	3/21/37/37	0/3/3/3
3	ADP	B	600	2	-	8/12/32/32	0/3/3/3
4	N9H	A	604	-	-	3/21/37/37	0/3/3/3
3	ADP	A	601	2	-	8/12/32/32	0/3/3/3

All (40) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	604	N9H	C7-C13	6.99	1.59	1.53
4	B	605	N9H	C3-C4	5.80	1.46	1.38
4	A	604	N9H	C3-C4	4.86	1.45	1.38
4	B	605	N9H	C23-C17	4.85	1.47	1.39
4	A	604	N9H	C19-C17	4.79	1.46	1.39
4	A	604	N9H	C23-C17	4.71	1.46	1.39
4	A	604	N9H	C20-C19	4.50	1.48	1.38
4	A	604	N9H	C22-C23	4.22	1.47	1.38
4	B	605	N9H	C6-C1	4.02	1.45	1.37
4	B	605	N9H	C19-C17	3.91	1.45	1.39
4	A	604	N9H	C2-C1	3.55	1.43	1.37
4	A	604	N9H	C21-C20	3.45	1.47	1.38
4	A	604	N9H	C10-N1	3.36	1.34	1.29
4	A	604	N9H	C6-C1	3.30	1.43	1.37
4	B	605	N9H	C22-C23	3.16	1.45	1.38
3	A	601	ADP	C2-N3	3.00	1.36	1.32
4	A	604	N9H	C2-C3	2.91	1.44	1.39
4	A	604	N9H	C3-C10	2.90	1.50	1.47
4	A	604	N9H	F41-C1	-2.87	1.29	1.36
3	B	600	ADP	O4'-C1'	2.81	1.45	1.41
4	B	605	N9H	C2-C1	2.81	1.42	1.37
4	B	605	N9H	C10-N1	2.79	1.33	1.29
4	A	604	N9H	C5-C4	2.74	1.43	1.37
4	A	604	N9H	C22-C21	2.68	1.45	1.38
4	B	605	N9H	C21-C20	2.68	1.45	1.38
4	B	605	N9H	C22-C21	2.63	1.45	1.38
3	A	601	ADP	O4'-C1'	2.63	1.44	1.41
3	A	601	ADP	C4-N3	2.58	1.39	1.35
3	B	600	ADP	C2-N3	2.50	1.36	1.32
4	B	605	N9H	C20-C19	2.48	1.44	1.38
4	B	605	N9H	N12-N1	-2.41	1.33	1.39
4	A	604	N9H	C14-C13	-2.36	1.52	1.54
4	B	605	N9H	C6-C5	2.26	1.42	1.38
3	B	600	ADP	C5-N7	-2.25	1.31	1.39
3	B	600	ADP	C4-N3	2.24	1.38	1.35
4	A	604	N9H	C7-C8	2.17	1.60	1.52
4	B	605	N9H	C5-C4	2.12	1.42	1.37
4	B	605	N9H	C2-C3	2.09	1.43	1.39
3	A	601	ADP	C2-N1	2.07	1.37	1.33
3	A	601	ADP	C5-N7	-2.01	1.32	1.39

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	B	605	N9H	C10-N1-N12	17.51	115.06	107.99
4	A	604	N9H	C14-C10-N1	-15.43	104.59	113.83
4	A	604	N9H	C10-N1-N12	13.57	113.47	107.99
4	B	605	N9H	C14-C10-N1	-11.13	107.16	113.83
4	A	604	N9H	C7-C8-C11	7.63	126.06	112.35
4	A	604	N9H	N30-C29-N12	7.54	120.45	114.17
3	A	601	ADP	N3-C2-N1	-6.96	117.80	128.68
3	B	600	ADP	N3-C2-N1	-6.91	117.88	128.68
4	B	605	N9H	N30-C29-N12	6.14	119.29	114.17
4	A	604	N9H	O39-C29-N12	-4.65	114.64	121.03
4	B	605	N9H	C6-C5-C4	-4.14	114.04	119.05
4	B	605	N9H	O39-C29-N12	-3.14	116.72	121.03
4	B	605	N9H	C2-C3-C4	2.98	120.29	116.66
4	A	604	N9H	C31-N30-C29	2.72	126.17	120.77
4	B	605	N9H	C5-C6-C1	2.44	120.89	118.36
4	A	604	N9H	C3-C10-N1	2.40	126.68	121.51
4	B	605	N9H	C3-C2-C1	-2.36	114.07	117.98
4	B	605	N9H	C14-C10-C3	2.34	127.39	124.63
3	B	600	ADP	C2'-C3'-C4'	2.31	107.13	102.64
4	A	604	N9H	C6-C5-C4	-2.21	116.37	119.05
4	A	604	N9H	C5-C4-C3	2.20	125.52	123.11
4	B	605	N9H	F40-C4-C3	-2.20	115.93	119.67
4	B	605	N9H	C21-C22-C23	-2.13	116.95	120.19
3	B	600	ADP	C2-N1-C6	2.10	122.35	118.75
3	A	601	ADP	C2-N1-C6	2.05	122.27	118.75

There are no chirality outliers.

All (22) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	601	ADP	PA-O3A-PB-O3B
3	A	601	ADP	C5'-O5'-PA-O1A
3	A	601	ADP	C5'-O5'-PA-O2A
3	A	601	ADP	C5'-O5'-PA-O3A
3	B	600	ADP	PA-O3A-PB-O3B
3	B	600	ADP	C5'-O5'-PA-O1A
3	B	600	ADP	C5'-O5'-PA-O2A
3	B	600	ADP	C5'-O5'-PA-O3A
3	A	601	ADP	C3'-C4'-C5'-O5'
3	B	600	ADP	O4'-C4'-C5'-O5'
3	B	600	ADP	C3'-C4'-C5'-O5'
4	B	605	N9H	C13-C7-C8-C11
3	A	601	ADP	O4'-C4'-C5'-O5'

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Mol	Chain	Res	Type	Atoms
3	A	601	ADP	PA-O3A-PB-O2B
3	B	600	ADP	PA-O3A-PB-O2B
4	A	604	N9H	N12-C13-C17-C19
4	B	605	N9H	N12-C13-C17-C19
3	A	601	ADP	PA-O3A-PB-O1B
3	B	600	ADP	PA-O3A-PB-O1B
4	A	604	N9H	C7-C13-C17-C19
4	B	605	N9H	N12-C13-C17-C23
4	A	604	N9H	N30-C29-N12-N1

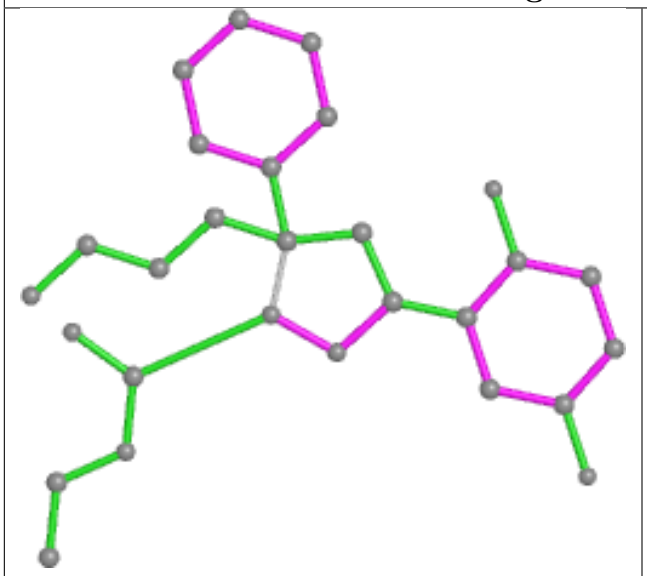
There are no ring outliers.

1 monomer is involved in 1 short contact:

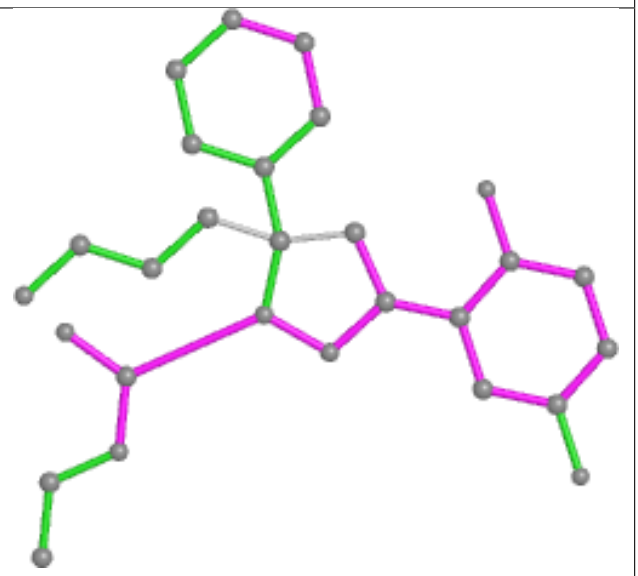
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	605	N9H	1	0

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 then 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.

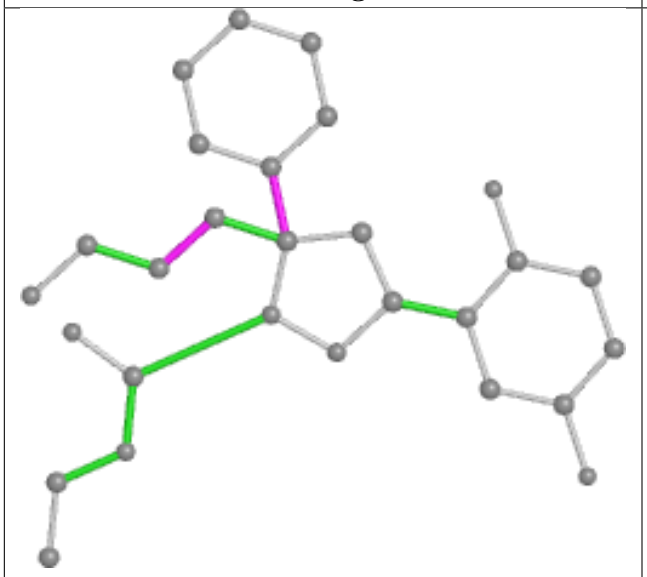
Ligand N9H B 605



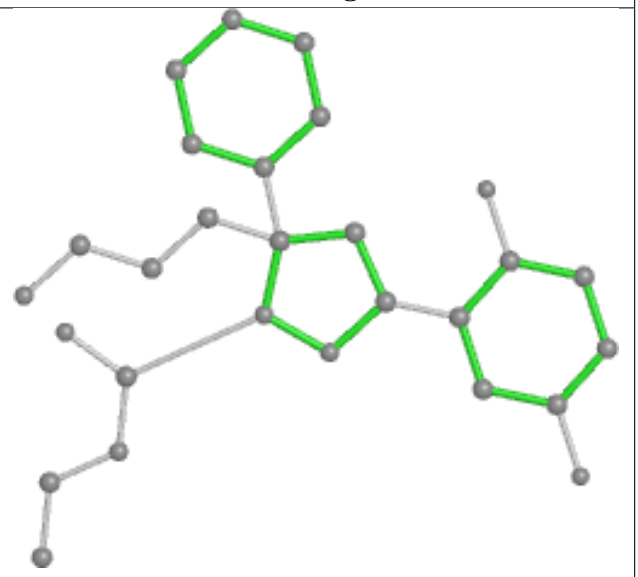
Bond lengths



Bond angles

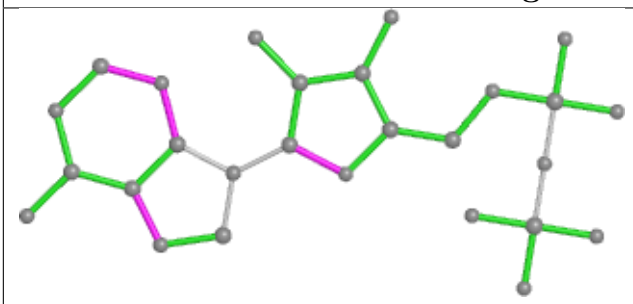


Torsions

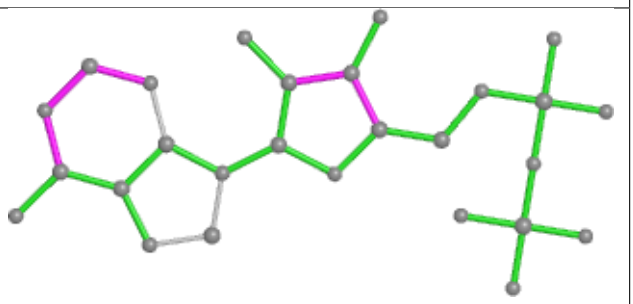


Rings

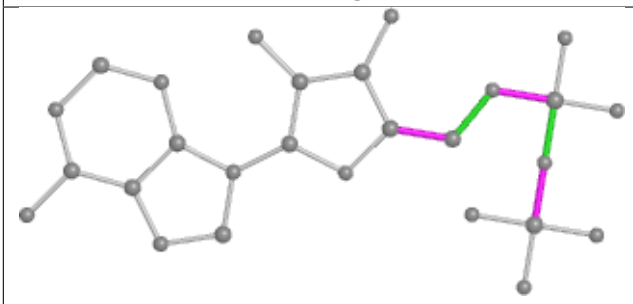
Ligand ADP B 600



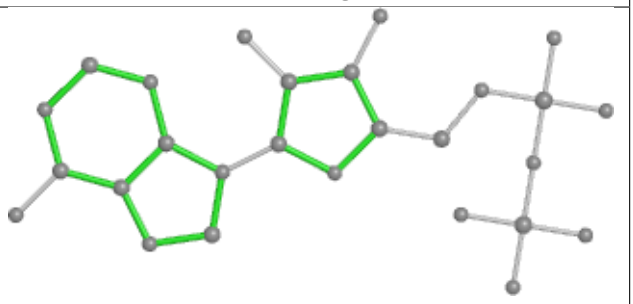
Bond lengths



Bond angles

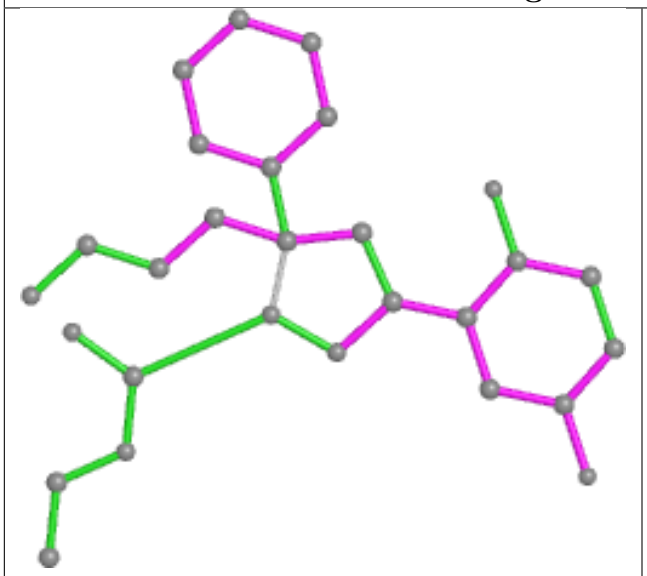


Torsions

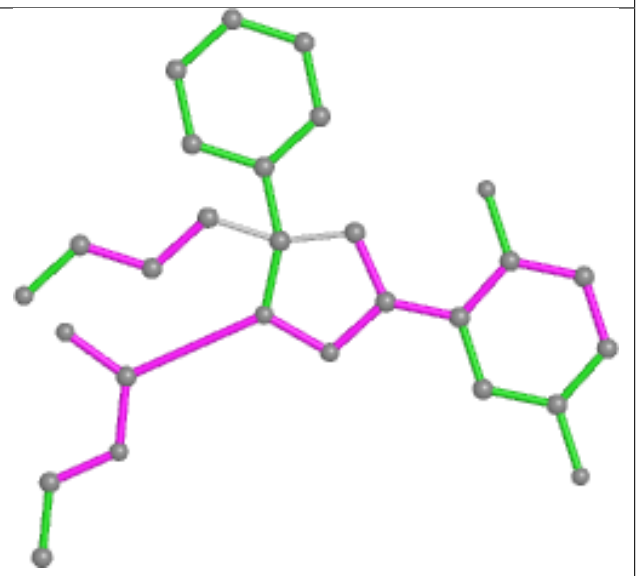


Rings

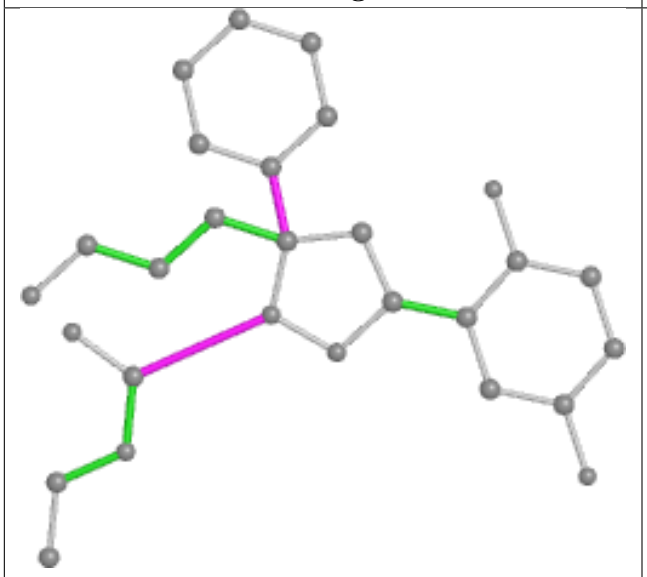
Ligand N9H A 604



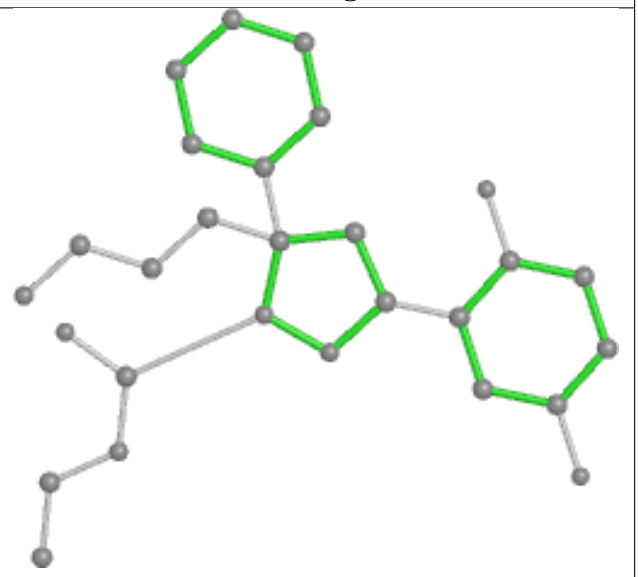
Bond lengths



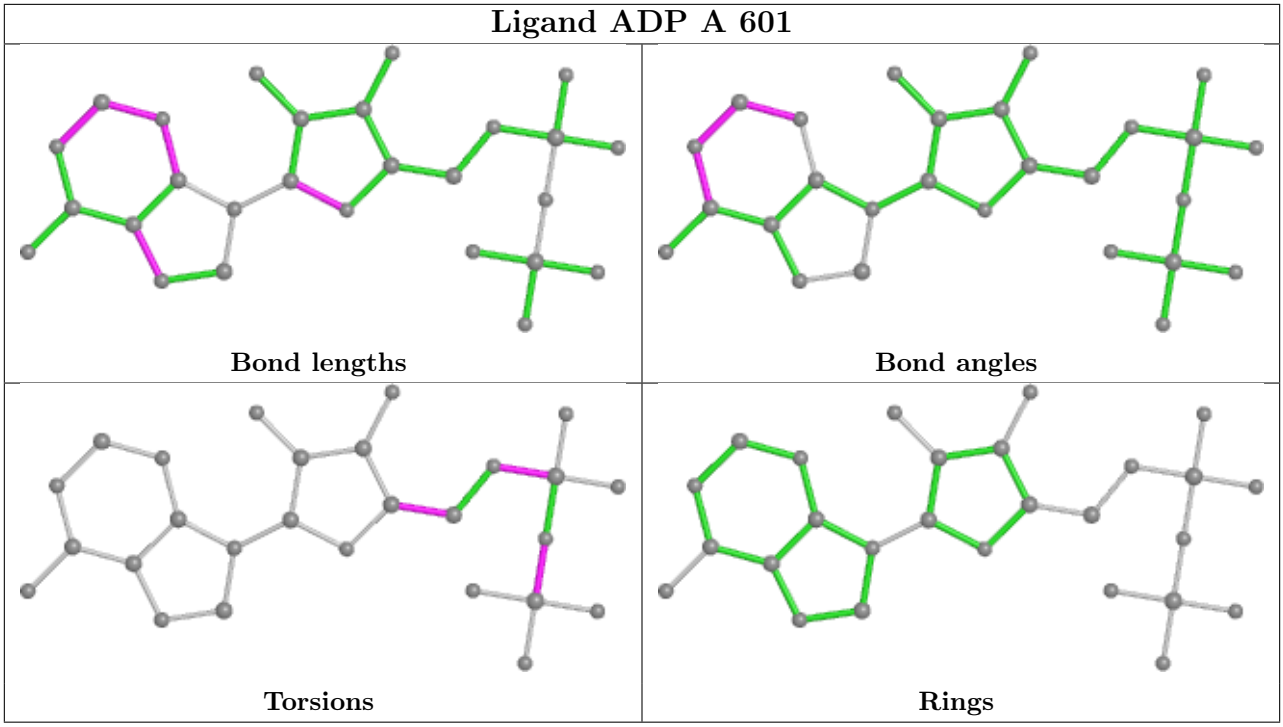
Bond angles



Torsions



Rings



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)

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains (i)

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.