



Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 12:51 am BST

PDB ID : 1YRS
Title : Crystal structure of KSP in complex with inhibitor 1
Authors : Cox, C.D.; Breslin, M.J.; Mariano, B.J.; Coleman, P.J.; Buser, C.A.; Walsh, E.S.; Hamilton, K.; Kohl, N.E.; Torrent, M.; Yan, Y.; Kuo, L.C.; Hartman, G.D.
Deposited on : 2005-02-04
Resolution : 2.50 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 following versions of software and data (see [references ⓘ](#)) 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) : 2.11

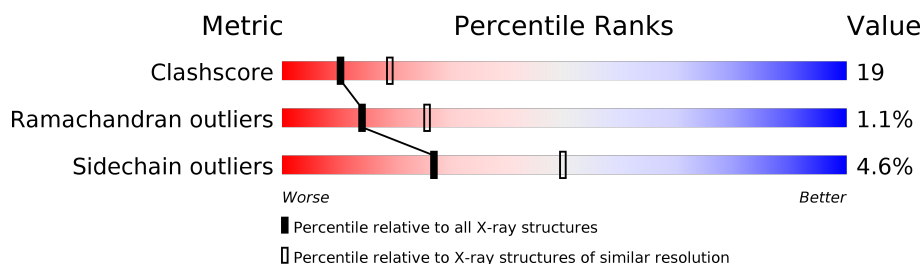
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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 on 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	A	368	
1	B	368	

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
4	L47	A	604	-	X	X	-
4	L47	B	603	-	X	-	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 5454 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.

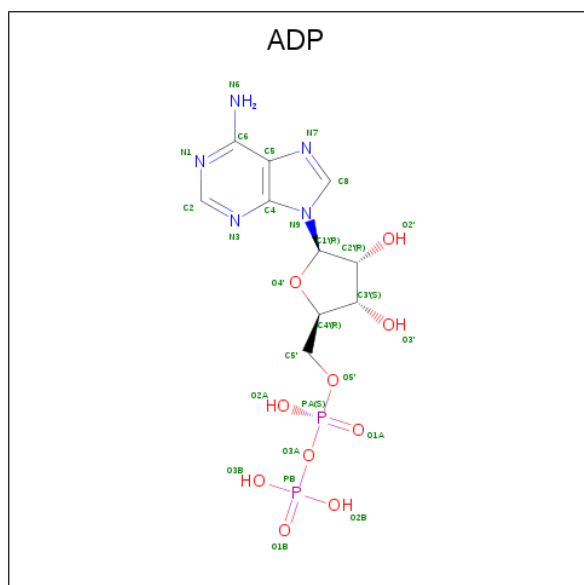
- Molecule 1 is a protein called Kinesin-like protein KIF11.

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

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

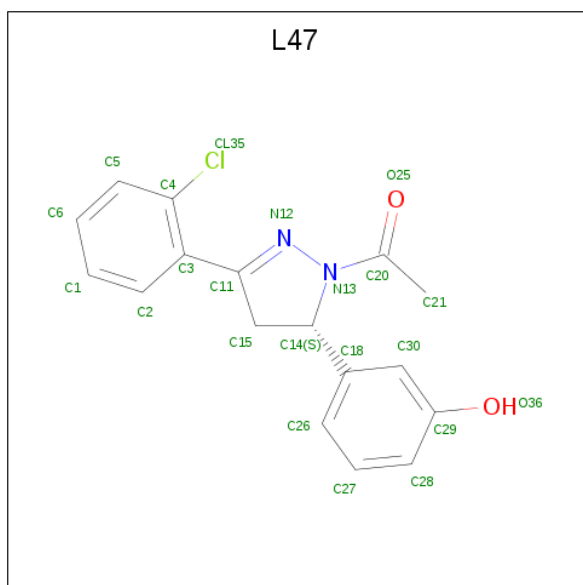
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Mg	0	0
			1	1		
2	A	1	Total	Mg	0	0
			1	1		

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

- Molecule 4 is 3-[(5S)-1-ACETYL-3-(2-CHLOROPHENYL)-4,5-DIHYDRO-1H-PYRAZOL-5-YL]PHENOL (three-letter code: L47) (formula: C₁₇H₁₅ClN₂O₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	Cl	N	O	0	0
			22	17	1	2	2		
4	B	1	Total	C	Cl	N	O	0	0
			22	17	1	2	2		

- Molecule 5 is water.

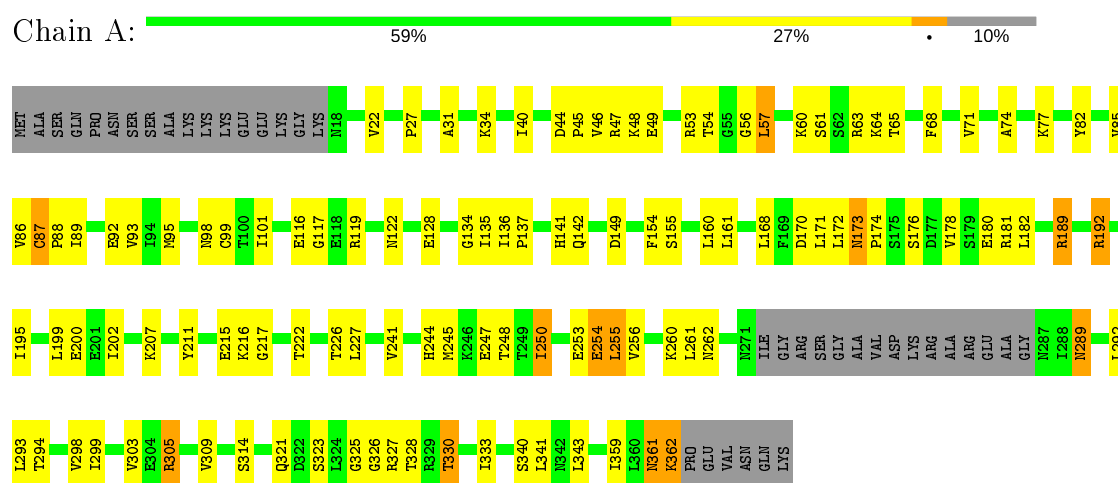
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	97	Total	O	0	0
			97	97		
5	B	69	Total	O	0	0
			69	69		

3 Residue-property plots [i](#)

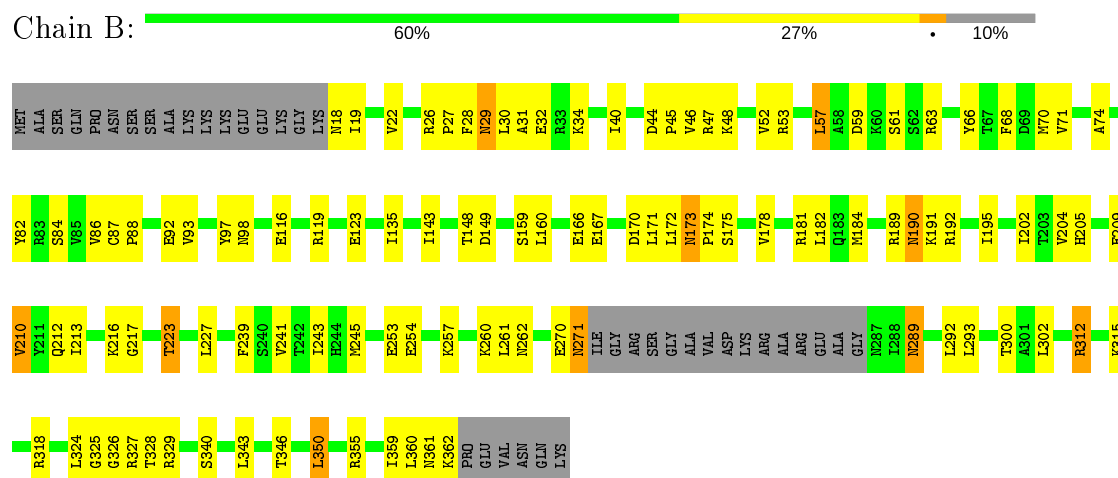
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded according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



- Molecule 1: Kinesin-like protein KIF11



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	69.30 Å 79.80 Å 159.60 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.50	Depositor
% Data completeness (in resolution range)	94.0 (50.00-2.50)	Depositor
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.250 , 0.310	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5454	wwPDB-VP
Average B, all atoms (Å ²)	32.0	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: MG, L47, 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	0.51	0/2632	0.64	0/3559
1	B	0.50	0/2632	0.63	0/3559
All	All	0.51	0/5264	0.64	0/7118

There are no bond length outliers.

There are no bond angle outliers.

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	2594	0	2618	95	0
1	B	2594	0	2617	99	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	22	0	14	8	0
4	B	22	0	14	5	0
5	A	97	0	0	4	0
5	B	69	0	0	1	0
All	All	5454	0	5287	198	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 (198) 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:B:312:ARG:HB3	1:B:312:ARG:HH11	1.29	0.96
1:A:299:ILE:HG23	1:A:359:ILE:HD11	1.52	0.90
1:A:247:GLU:HB3	1:A:255:LEU:HD23	1.57	0.86
1:A:92:GLU:HA	1:A:95:MET:HE3	1.58	0.86
1:A:305:ARG:HH11	1:A:305:ARG:HB3	1.43	0.84
1:A:89:ILE:HD11	1:A:99:CYS:HB3	1.61	0.82
1:A:323:SER:O	1:A:330:THR:HG21	1.78	0.81
4:A:604:L47:H151	4:A:604:L47:CL35	2.20	0.79
1:B:18:ASN:HD22	1:B:361:ASN:H	1.30	0.79
1:A:247:GLU:O	1:A:254:GLU:HA	1.84	0.78
1:A:45:PRO:HA	1:A:71:VAL:HG23	1.66	0.78
1:B:312:ARG:HB3	1:B:312:ARG:NH1	2.00	0.77
1:A:361:ASN:O	1:A:362:LYS:HB2	1.87	0.74
1:B:184:MET:HE3	1:B:318:ARG:HH21	1.55	0.72
1:A:305:ARG:NH1	1:A:305:ARG:HB3	2.05	0.71
1:B:86:VAL:HG21	1:B:135:ILE:HG12	1.73	0.70
1:B:70:MET:CE	1:B:84:SER:HB3	2.21	0.70
1:B:70:MET:HE2	1:B:84:SER:HB3	1.74	0.70
1:A:326:GLY:O	1:A:361:ASN:HB2	1.93	0.69
1:A:27:PRO:HB3	1:A:74:ALA:HB1	1.74	0.69
1:A:250:ILE:HD13	1:A:250:ILE:H	1.58	0.69
1:A:260:LYS:HE2	1:A:262:ASN:HD21	1.56	0.69
1:B:18:ASN:ND2	1:B:361:ASN:H	1.90	0.69
1:B:184:MET:HE3	1:B:318:ARG:NH2	2.08	0.69
1:B:181:ARG:HG2	1:B:181:ARG:HH11	1.57	0.68
4:B:603:L47:H151	4:B:603:L47:CL35	2.31	0.68
1:B:40:ILE:HD13	1:B:340:SER:HA	1.76	0.67
1:A:54:THR:HG21	1:A:64:LYS:HD3	1.76	0.67
1:A:119:ARG:HG2	4:A:604:L47:C21	2.25	0.67
1:A:362:LYS:HA	5:A:606:HOH:O	1.97	0.65
1:B:184:MET:CE	1:B:318:ARG:HE	2.08	0.65
1:B:119:ARG:HG2	4:B:603:L47:H211	1.78	0.65
1:B:31:ALA:O	1:B:34:LYS:HB2	1.97	0.65
1:A:57:LEU:O	1:A:61:SER:HB3	1.97	0.64
1:B:53:ARG:HB2	1:B:63:ARG:NH1	2.13	0.64
1:B:260:LYS:HE2	1:B:262:ASN:HD21	1.64	0.62
1:B:189:ARG:HH21	1:B:195:ILE:HD11	1.64	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:346:THR:O	1:B:350:LEU:HD22	2.00	0.61
1:B:209:GLU:O	1:B:213:ILE:HG13	2.01	0.61
1:A:305:ARG:HH11	1:A:305:ARG:CB	2.13	0.60
1:A:40:ILE:HD12	1:A:343:LEU:HD13	1.83	0.60
1:B:48:LYS:HA	1:B:71:VAL:HG22	1.82	0.60
1:A:172:LEU:O	1:A:174:PRO:HD3	2.02	0.60
1:A:192:ARG:HB2	1:A:321:GLN:OE1	2.01	0.59
1:A:40:ILE:HD13	1:A:340:SER:HA	1.85	0.59
4:A:604:L47:C15	4:A:604:L47:CL35	2.87	0.59
1:A:119:ARG:HG2	4:A:604:L47:H211	1.84	0.59
1:B:326:GLY:O	1:B:362:LYS:N	2.34	0.59
1:B:34:LYS:HA	1:B:34:LYS:HE2	1.85	0.58
1:B:205:HIS:HB2	1:B:209:GLU:OE1	2.03	0.58
1:A:202:ILE:HD12	1:A:202:ILE:N	2.19	0.58
1:A:48:LYS:HA	1:A:71:VAL:HG22	1.84	0.58
1:B:184:MET:HE1	1:B:318:ARG:HE	1.69	0.57
1:B:29:ASN:HD21	1:B:32:GLU:HG3	1.70	0.57
1:A:31:ALA:HA	1:A:34:LYS:HE2	1.86	0.57
1:B:173:ASN:ND2	1:B:175:SER:H	2.02	0.57
1:B:312:ARG:HH11	1:B:312:ARG:CB	2.12	0.56
1:B:325:GLY:H	1:B:361:ASN:ND2	2.04	0.56
1:A:82:TYR:OH	1:A:142:GLN:HG3	2.06	0.56
1:A:77:LYS:HE2	5:A:685:HOH:O	2.05	0.56
1:A:85:VAL:HG23	1:A:86:VAL:N	2.21	0.55
1:A:341:LEU:H	1:A:341:LEU:HD12	1.70	0.55
1:B:87:CYS:HB3	1:B:88:PRO:HD3	1.88	0.55
1:A:56:GLY:O	1:A:57:LEU:HB2	2.07	0.55
1:A:227:LEU:O	1:A:227:LEU:HD12	2.06	0.55
1:B:241:VAL:HG13	1:B:261:LEU:HB3	1.89	0.55
1:A:250:ILE:HD13	1:A:250:ILE:N	2.21	0.55
1:A:98:ASN:O	1:A:328:THR:HG23	2.08	0.54
1:B:171:LEU:O	1:B:217:GLY:HA2	2.07	0.54
1:A:189:ARG:HH21	1:A:195:ILE:HD11	1.72	0.54
1:B:172:LEU:O	1:B:174:PRO:HD3	2.07	0.54
1:A:176:SER:HB2	1:A:180:GLU:OE1	2.08	0.54
1:B:40:ILE:HD12	1:B:343:LEU:HD23	1.90	0.54
1:B:57:LEU:HD12	1:B:57:LEU:H	1.72	0.54
1:A:116:GLU:HB3	4:A:604:L47:CL35	2.45	0.53
1:B:192:ARG:HB3	1:B:192:ARG:NH1	2.24	0.53
1:A:173:ASN:HD22	1:A:174:PRO:N	2.07	0.53
1:A:173:ASN:HD22	1:A:173:ASN:C	2.11	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:27:PRO:HB3	1:B:74:ALA:HB1	1.91	0.53
1:A:361:ASN:C	1:A:361:ASN:HD22	2.13	0.52
1:A:211:TYR:O	1:A:215:GLU:HG3	2.10	0.52
1:B:22:VAL:HG12	1:B:70:MET:HB2	1.92	0.52
1:A:141:HIS:CD2	1:A:207:LYS:HD3	2.45	0.52
1:A:250:ILE:CD1	1:A:250:ILE:H	2.23	0.52
1:A:170:ASP:HB2	1:A:182:LEU:HD11	1.92	0.51
1:A:44:ASP:OD1	1:A:46:VAL:HG12	2.10	0.51
1:B:325:GLY:H	1:B:361:ASN:HD21	1.57	0.51
1:A:155:SER:OG	1:A:244:HIS:HB2	2.10	0.51
1:B:173:ASN:HD21	1:B:175:SER:HB2	1.74	0.51
1:B:318:ARG:HH11	1:B:318:ARG:HG2	1.75	0.51
1:B:44:ASP:O	1:B:48:LYS:N	2.42	0.51
1:B:189:ARG:NH2	1:B:195:ILE:HD11	2.25	0.51
1:B:143:ILE:HD13	1:B:243:ILE:HD11	1.91	0.50
1:A:248:THR:HA	1:A:253:GLU:O	2.11	0.50
1:B:98:ASN:O	1:B:328:THR:HG23	2.10	0.50
1:A:255:LEU:HD22	1:A:255:LEU:N	2.27	0.50
1:A:160:LEU:HB3	1:A:172:LEU:HG	1.93	0.50
1:A:48:LYS:HB3	1:A:68:PHE:O	2.12	0.50
1:B:22:VAL:CG1	1:B:70:MET:HB2	2.42	0.50
1:B:312:ARG:HA	1:B:318:ARG:HG3	1.94	0.50
1:B:57:LEU:HD13	1:B:59:ASP:HB3	1.93	0.50
1:B:166:GLU:OE1	1:B:315:LYS:HG2	2.12	0.49
4:B:603:L47:CL35	4:B:603:L47:C15	2.97	0.49
1:B:82:TYR:CD2	1:B:86:VAL:HB	2.46	0.49
1:B:70:MET:HE1	1:B:84:SER:HB3	1.93	0.49
1:B:57:LEU:O	1:B:61:SER:HB3	2.13	0.49
1:A:119:ARG:HG2	4:A:604:L47:H212	1.94	0.49
1:A:160:LEU:HD22	4:A:604:L47:H6	1.95	0.48
1:B:53:ARG:HB2	1:B:63:ARG:HH12	1.76	0.48
1:B:92:GLU:OE2	1:B:329:ARG:HD2	2.13	0.48
1:B:271:ASN:C	1:B:271:ASN:HD22	2.16	0.48
1:B:270:GLU:O	1:B:270:GLU:HG2	2.13	0.48
1:B:327:ARG:HA	1:B:362:LYS:HA	1.96	0.48
1:B:66:TYR:HB2	1:B:68:PHE:CE1	2.49	0.48
1:B:181:ARG:HG2	1:B:181:ARG:NH1	2.27	0.47
1:B:92:GLU:OE2	1:B:329:ARG:NH1	2.44	0.47
1:B:202:ILE:HD12	1:B:202:ILE:N	2.29	0.47
1:A:255:LEU:H	1:A:255:LEU:HD22	1.79	0.47
1:A:47:ARG:NH2	1:A:49:GLU:OE2	2.48	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:300:THR:HG23	1:B:355:ARG:HH21	1.79	0.47
1:A:241:VAL:HG13	1:A:261:LEU:HB3	1.95	0.47
1:A:154:PHE:HA	1:A:244:HIS:O	2.15	0.47
1:A:128:GLU:HB2	1:A:141:HIS:CD2	2.50	0.47
1:A:173:ASN:ND2	1:A:173:ASN:C	2.68	0.46
1:A:86:VAL:HG21	1:A:135:ILE:HG12	1.98	0.46
1:A:85:VAL:CG2	1:A:86:VAL:N	2.78	0.46
1:B:184:MET:CE	1:B:318:ARG:NE	2.79	0.46
1:A:247:GLU:HB3	1:A:255:LEU:CD2	2.37	0.46
1:A:92:GLU:CA	1:A:95:MET:HE3	2.39	0.46
1:B:184:MET:HE3	1:B:318:ARG:CZ	2.46	0.46
1:B:18:ASN:HB3	1:B:360:LEU:HA	1.97	0.46
1:A:325:GLY:HA2	1:A:361:ASN:HB3	1.98	0.45
1:A:117:GLY:HA3	1:A:134:GLY:N	2.31	0.45
1:A:57:LEU:HB2	1:A:60:LYS:O	2.16	0.45
1:A:289:ASN:HD22	1:A:292:LEU:H	1.65	0.45
1:B:148:THR:HG23	1:B:149:ASP:N	2.31	0.45
1:B:116:GLU:HG2	4:B:603:L47:CL35	2.54	0.45
1:B:47:ARG:HH11	1:B:47:ARG:HG2	1.80	0.45
1:B:212:GLN:HG2	1:B:216:LYS:HZ3	1.80	0.45
1:B:245:MET:HB2	1:B:257:LYS:HB2	1.97	0.45
1:B:119:ARG:HG2	4:B:603:L47:C21	2.44	0.44
1:A:303:VAL:CG1	1:A:359:ILE:HG13	2.46	0.44
1:B:28:PHE:HD2	1:B:32:GLU:OE2	2.00	0.44
1:B:190:ASN:O	1:B:192:ARG:N	2.51	0.44
1:B:19:ILE:HD12	1:B:359:ILE:HB	1.98	0.44
1:B:327:ARG:O	1:B:362:LYS:HG3	2.17	0.44
1:B:184:MET:HE3	1:B:318:ARG:NE	2.32	0.44
1:A:161:LEU:HD21	1:A:168:LEU:HG	2.00	0.43
1:B:289:ASN:HD22	1:B:292:LEU:H	1.66	0.43
1:B:34:LYS:HA	1:B:34:LYS:CE	2.49	0.43
1:B:312:ARG:HA	1:B:318:ARG:CG	2.48	0.43
1:A:202:ILE:HD12	1:A:202:ILE:H	1.83	0.43
1:A:89:ILE:HG21	1:A:101:ILE:HD11	2.00	0.43
1:B:312:ARG:HA	1:B:318:ARG:HD2	2.01	0.43
1:B:302:LEU:HD21	1:B:324:LEU:O	2.19	0.42
1:B:30:LEU:O	1:B:30:LEU:HD12	2.20	0.42
1:A:87:CYS:N	1:A:88:PRO:HD2	2.34	0.42
1:A:85:VAL:O	1:A:88:PRO:HG2	2.19	0.42
1:B:93:VAL:HG21	1:B:261:LEU:HB2	2.01	0.42
1:B:170:ASP:HB2	1:B:182:LEU:HD11	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:205:HIS:CD2	1:B:209:GLU:OE1	2.72	0.42
1:B:223:THR:O	1:B:227:LEU:HD23	2.20	0.42
1:B:52:VAL:HG21	1:B:350:LEU:CD2	2.50	0.42
1:A:245:MET:O	1:A:256:VAL:HA	2.20	0.42
1:A:27:PRO:CB	1:A:74:ALA:HB1	2.45	0.42
1:B:45:PRO:HG2	1:B:46:VAL:H	1.85	0.42
1:A:222:THR:O	1:A:226:THR:HG23	2.20	0.42
1:A:99:CYS:O	1:A:261:LEU:HA	2.19	0.42
1:A:93:VAL:HG21	1:A:261:LEU:HB2	2.02	0.42
1:A:46:VAL:HG13	1:A:47:ARG:N	2.34	0.42
1:B:97:TYR:HE1	1:B:362:LYS:HE3	1.83	0.42
1:A:171:LEU:O	1:A:217:GLY:HA2	2.20	0.42
1:A:22:VAL:HG22	1:A:333:ILE:HG12	2.02	0.42
1:A:53:ARG:HA	1:A:63:ARG:HG2	2.02	0.42
1:A:136:ILE:HB	1:A:137:PRO:CD	2.49	0.41
1:A:160:LEU:CD2	4:A:604:L47:H6	2.50	0.41
1:A:40:ILE:CG2	1:A:340:SER:HB3	2.50	0.41
1:A:173:ASN:HB2	1:A:200:GLU:HG3	2.01	0.41
1:A:294:THR:OG1	1:A:314:SER:HB3	2.20	0.41
1:B:123:GLU:HB2	5:B:631:HOH:O	2.20	0.41
1:B:167:GLU:CD	1:B:181:ARG:HD3	2.40	0.41
1:A:173:ASN:HB2	1:A:200:GLU:CG	2.50	0.41
1:B:160:LEU:HB3	1:B:172:LEU:HG	2.02	0.41
1:B:253:GLU:HG2	1:B:254:GLU:N	2.34	0.41
1:B:98:ASN:C	1:B:328:THR:HG23	2.41	0.41
1:B:82:TYR:HA	1:B:86:VAL:HG23	2.02	0.41
1:A:341:LEU:HD12	1:A:341:LEU:N	2.35	0.41
1:B:29:ASN:ND2	1:B:32:GLU:HG3	2.35	0.41
1:A:298:VAL:HG13	1:A:309:VAL:CG1	2.51	0.41
1:A:89:ILE:CD1	1:A:99:CYS:HB3	2.40	0.41
1:B:159:SER:O	1:B:239:PHE:HA	2.20	0.41
1:B:343:LEU:O	1:B:343:LEU:HD22	2.21	0.41
1:A:56:GLY:O	1:A:57:LEU:CB	2.68	0.41
1:B:47:ARG:NH1	1:B:47:ARG:HG2	2.36	0.41
1:A:86:VAL:HA	1:A:89:ILE:HG22	2.03	0.41
1:A:181:ARG:CZ	5:A:659:HOH:O	2.69	0.41
1:A:327:ARG:HD3	5:A:689:HOH:O	2.21	0.40
1:B:204:VAL:HG11	1:B:210:VAL:HG12	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%)	306 (94%)	16 (5%)	4 (1%)	13	24
1	B	326/368 (89%)	314 (96%)	9 (3%)	3 (1%)	17	31
All	All	652/736 (89%)	620 (95%)	25 (4%)	7 (1%)	14	26

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	191	LYS
1	A	57	LEU
1	A	189	ARG
1	B	178	VAL
1	A	149	ASP
1	B	190	ASN
1	A	178	VAL

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%)	282 (96%)	11 (4%)	33	58
All	All	586/644 (91%)	559 (95%)	27 (5%)	27	50

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	65	THR
1	A	87	CYS
1	A	122	ASN
1	A	173	ASN
1	A	192	ARG
1	A	199	LEU
1	A	216	LYS
1	A	250	ILE
1	A	254	GLU
1	A	255	LEU
1	A	289	ASN
1	A	293	LEU
1	A	305	ARG
1	A	330	THR
1	A	361	ASN
1	A	362	LYS
1	B	26	ARG
1	B	29	ASN
1	B	57	LEU
1	B	173	ASN
1	B	210	VAL
1	B	223	THR
1	B	271	ASN
1	B	289	ASN
1	B	293	LEU
1	B	312	ARG
1	B	350	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	18	ASN
1	A	173	ASN
1	A	229	ASN
1	A	262	ASN
1	A	287	ASN
1	A	289	ASN
1	A	342	ASN
1	B	18	ASN
1	B	106	GLN
1	B	141	HIS
1	B	173	ASN
1	B	183	GLN

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Mol	Chain	Res	Type
1	B	190	ASN
1	B	205	HIS
1	B	262	ASN
1	B	271	ASN
1	B	289	ASN
1	B	290	GLN
1	B	358	ASN
1	B	361	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 carbohydrates 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	L47	A	604	-	24,24,24	2.96	16 (66%)	33,34,34	4.73	16 (48%)
3	ADP	A	601	2	24,29,29	1.46	4 (16%)	29,45,45	1.56	1 (3%)
3	ADP	B	600	2	24,29,29	1.44	3 (12%)	29,45,45	1.62	2 (6%)
4	L47	B	603	-	24,24,24	3.23	17 (70%)	33,34,34	4.91	15 (45%)

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	L47	A	604	-	-	4/12/24/24	0/3/3/3
3	ADP	A	601	2	-	6/12/32/32	0/3/3/3
3	ADP	B	600	2	-	8/12/32/32	0/3/3/3
4	L47	B	603	-	-	4/12/24/24	0/3/3/3

All (40) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	603	L47	C11-N12	6.03	1.38	1.29
4	B	603	L47	C3-C4	5.76	1.47	1.39
4	A	604	L47	C3-C4	5.06	1.46	1.39
4	B	603	L47	C15-C14	-4.77	1.47	1.54
4	A	604	L47	C15-C14	-4.63	1.47	1.54
4	A	604	L47	N13-N12	-4.60	1.29	1.38
4	B	603	L47	N13-N12	-4.17	1.29	1.38
4	B	603	L47	C28-C29	4.07	1.46	1.38
4	B	603	L47	C26-C18	4.06	1.45	1.39
4	A	604	L47	C11-N12	4.06	1.35	1.29
4	A	604	L47	C30-C18	4.03	1.45	1.39
4	B	603	L47	C30-C29	3.89	1.44	1.39
4	A	604	L47	C30-C29	3.78	1.44	1.39
4	A	604	L47	C26-C18	3.74	1.45	1.39
4	B	603	L47	C2-C3	3.62	1.45	1.39
4	B	603	L47	C30-C18	3.49	1.44	1.39
4	A	604	L47	C2-C3	3.37	1.45	1.39
4	A	604	L47	C28-C29	3.27	1.45	1.38
4	B	603	L47	C6-C1	3.20	1.46	1.38
4	A	604	L47	C6-C1	3.09	1.46	1.38
4	B	603	L47	C27-C28	3.04	1.45	1.38
4	A	604	L47	C1-C2	3.04	1.45	1.38
4	A	604	L47	C18-C14	-3.02	1.47	1.51
4	B	603	L47	C27-C26	2.99	1.45	1.38
3	A	601	ADP	C2-N3	2.97	1.36	1.32
3	B	600	ADP	C2-N3	2.91	1.36	1.32
3	A	601	ADP	O4'-C1'	2.90	1.45	1.41
4	A	604	L47	C27-C28	2.78	1.44	1.38
4	A	604	L47	C27-C26	2.76	1.44	1.38
4	A	604	L47	C5-C4	2.74	1.44	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	603	L47	C5-C4	2.63	1.44	1.38
4	B	603	L47	C1-C2	2.53	1.44	1.38
4	B	603	L47	C20-N13	2.51	1.39	1.36
4	B	603	L47	C6-C5	2.48	1.44	1.38
4	A	604	L47	C6-C5	2.43	1.44	1.38
3	A	601	ADP	C5-N7	-2.37	1.31	1.39
3	B	600	ADP	O4'-C1'	2.35	1.44	1.41
4	B	603	L47	C14-N13	2.34	1.50	1.48
3	B	600	ADP	C5-N7	-2.28	1.31	1.39
3	A	601	ADP	C4-N3	2.08	1.38	1.35

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	603	L47	C18-C14-N13	16.71	133.14	112.25
4	A	604	L47	C18-C14-N13	15.22	131.28	112.25
4	A	604	L47	C15-C11-N12	-11.60	99.60	113.69
4	B	603	L47	C15-C11-N12	-11.50	99.73	113.69
4	B	603	L47	C11-N12-N13	11.30	116.02	107.97
4	A	604	L47	C14-C15-C11	9.35	111.41	102.73
4	A	604	L47	C11-N12-N13	9.19	114.52	107.97
4	B	603	L47	C14-C15-C11	9.17	111.24	102.73
4	A	604	L47	C15-C14-C18	7.05	126.75	113.86
4	A	604	L47	C15-C14-N13	-6.91	95.08	100.87
3	B	600	ADP	N3-C2-N1	-6.90	117.90	128.68
4	B	603	L47	C15-C14-C18	6.82	126.34	113.86
3	A	601	ADP	N3-C2-N1	-6.70	118.21	128.68
4	B	603	L47	O25-C20-N13	4.57	126.42	119.34
4	A	604	L47	C29-C30-C18	-4.26	116.45	120.11
4	B	603	L47	C15-C14-N13	-4.16	97.39	100.87
4	B	603	L47	C3-C11-N12	3.90	129.88	121.51
4	B	603	L47	C29-C30-C18	-3.62	117.00	120.11
4	A	604	L47	C3-C4-CL35	-3.17	116.31	121.00
4	A	604	L47	C3-C11-N12	3.02	128.01	121.51
4	A	604	L47	C15-C11-C3	2.97	131.90	125.20
4	B	603	L47	C21-C20-N13	-2.93	114.44	117.30
4	B	603	L47	C28-C27-C26	-2.66	116.47	120.25
4	B	603	L47	C26-C18-C30	2.47	121.61	118.76
4	A	604	L47	C26-C18-C30	2.47	121.60	118.76
4	A	604	L47	C28-C27-C26	-2.36	116.89	120.25
4	B	603	L47	C6-C5-C4	-2.31	115.90	119.39
4	B	603	L47	C14-N13-N12	-2.23	110.86	112.98

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	604	L47	C26-C18-C14	-2.23	116.54	120.76
3	B	600	ADP	C2'-C3'-C4'	2.18	106.87	102.64
4	A	604	L47	C6-C5-C4	-2.16	116.13	119.39
4	A	604	L47	O25-C20-N13	2.12	122.62	119.34
4	A	604	L47	C14-N13-C20	2.08	125.89	124.00
4	B	603	L47	C15-C11-C3	2.01	129.74	125.20

There are no chirality outliers.

All (22) torsion outliers are listed below:

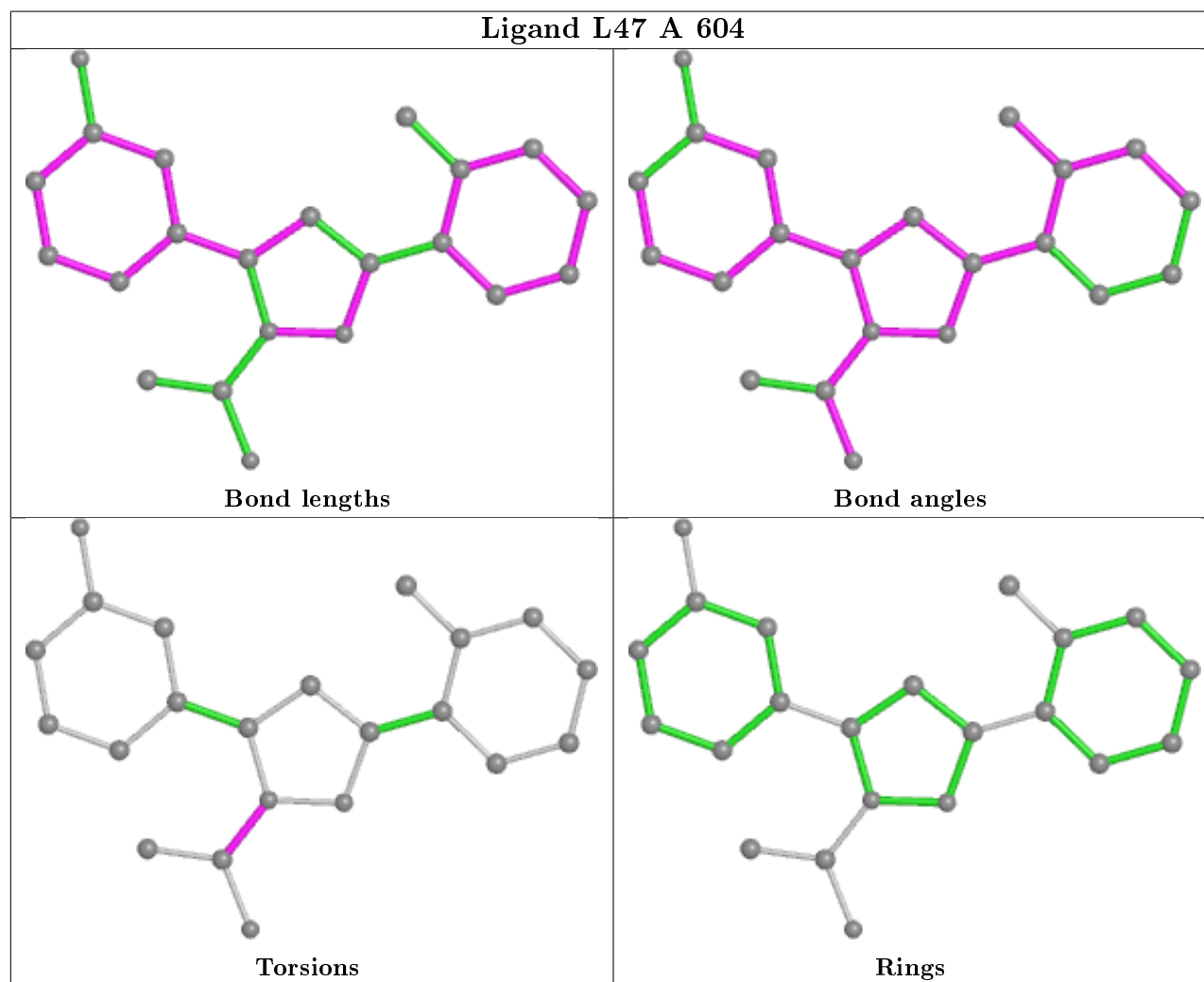
Mol	Chain	Res	Type	Atoms
3	A	601	ADP	C5'-O5'-PA-O1A
3	A	601	ADP	C5'-O5'-PA-O2A
3	A	601	ADP	C5'-O5'-PA-O3A
4	B	603	L47	C21-C20-N13-N12
4	B	603	L47	O25-C20-N13-N12
4	B	603	L47	C21-C20-N13-C14
4	B	603	L47	O25-C20-N13-C14
4	A	604	L47	C21-C20-N13-N12
4	A	604	L47	O25-C20-N13-N12
4	A	604	L47	C21-C20-N13-C14
4	A	604	L47	O25-C20-N13-C14
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	O4'-C4'-C5'-O5'
3	A	601	ADP	C3'-C4'-C5'-O5'
3	B	600	ADP	C3'-C4'-C5'-O5'
3	B	600	ADP	O4'-C4'-C5'-O5'
3	A	601	ADP	PA-O3A-PB-O2B
3	B	600	ADP	PA-O3A-PB-O2B
3	B	600	ADP	PA-O3A-PB-O1B

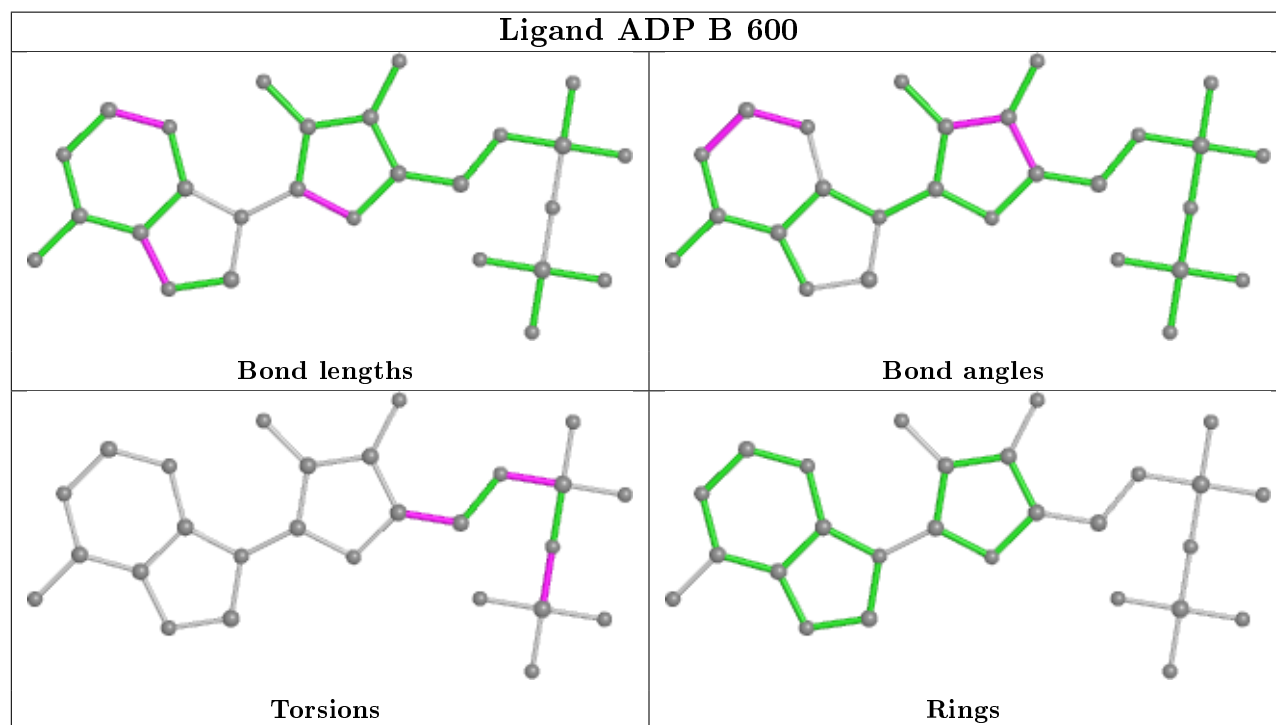
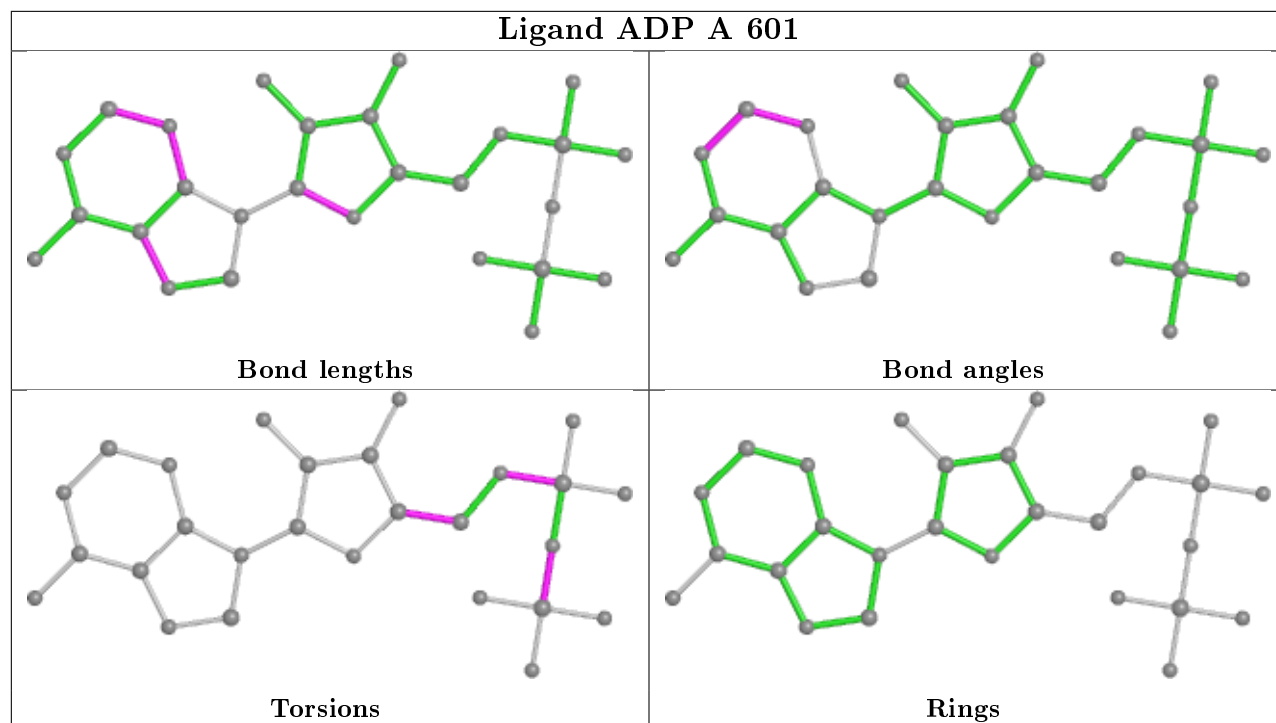
There are no ring outliers.

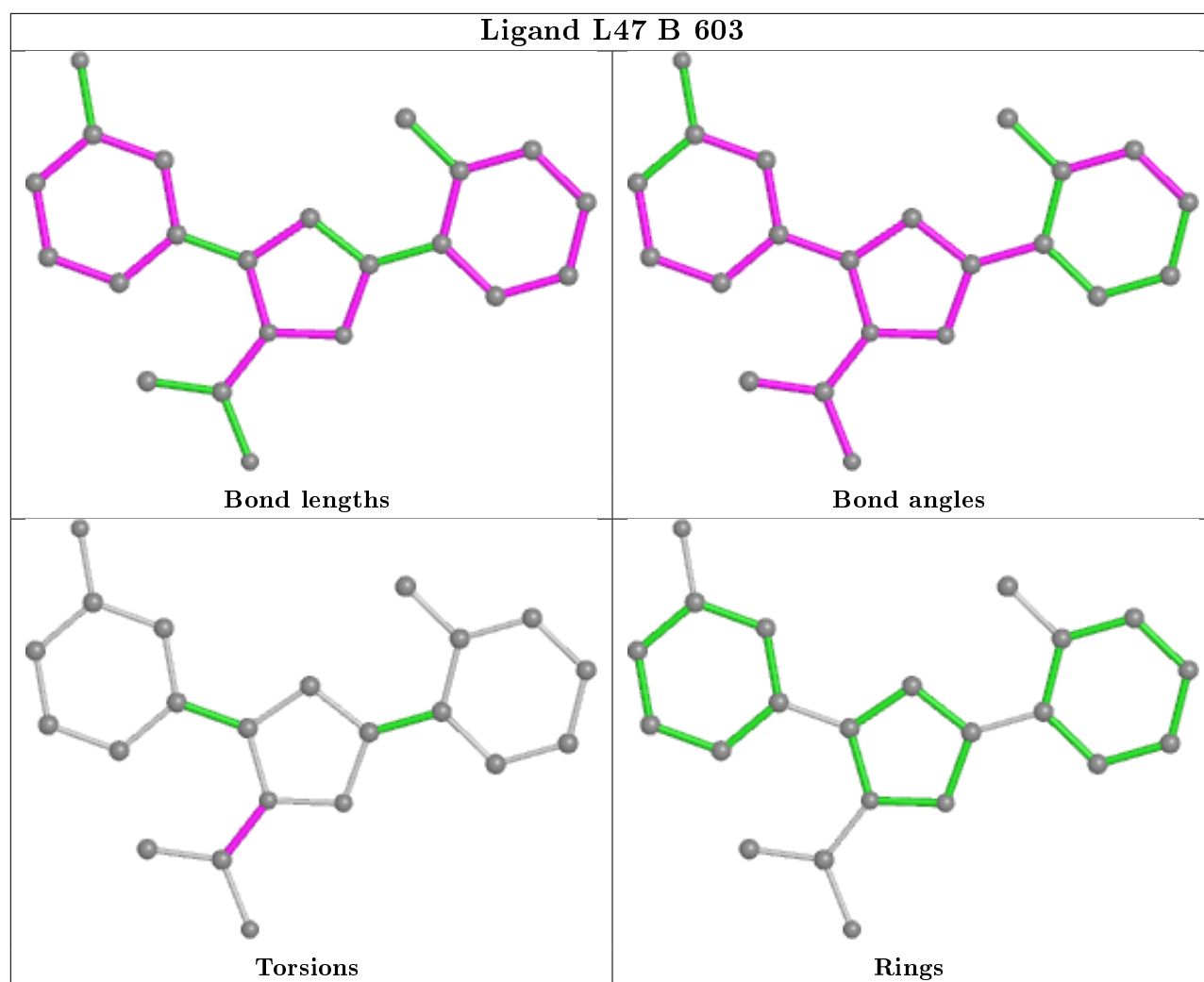
2 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	604	L47	8	0
4	B	603	L47	5	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.