



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 02:51 PM UTC

PDB ID : 2FL6 / pdb_00002fl6
Title : crystal structure of KSP in complex with inhibitor 6
Authors : Yan, Y.
Deposited on : 2006-01-05
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 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

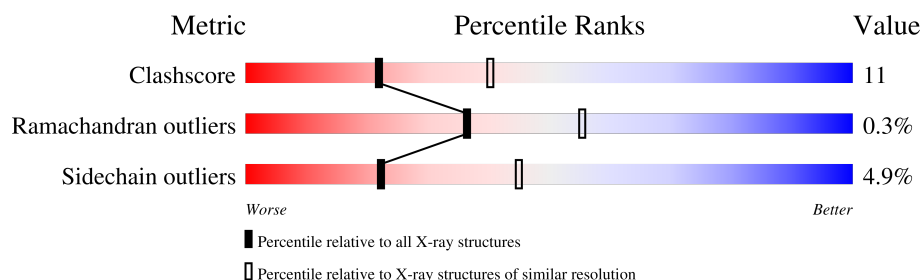
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (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 $\leq 5\%$.

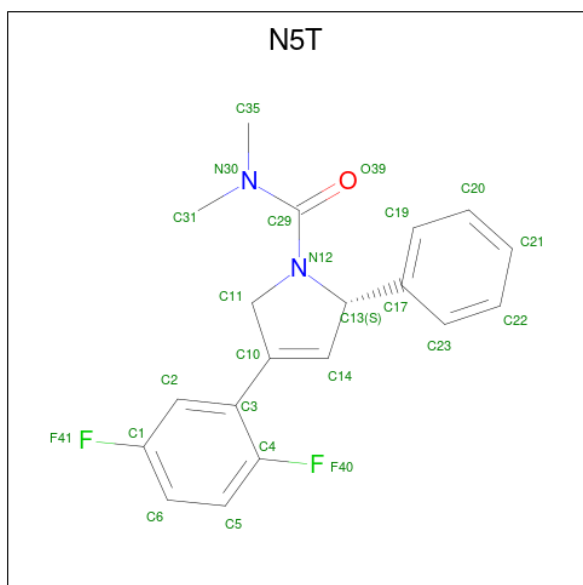
Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	367	
1	B	367	



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 (2S)-4-(2,5-DIFLUOROPHENYL)-N,N-DIMETHYL-2-PHENYL-2,5-DIHYDRO-1H-PYRROLE-1-CARBOXAMIDE (CCD ID: N5T) (formula: C₁₉H₁₈F₂N₂O).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	F	N	O	0	0
			24	19	2	2	1		
4	B	1	Total	C	F	N	O	0	0
			24	19	2	2	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	92	Total O 92 92	0	0
5	B	88	Total O 88 88	0	0

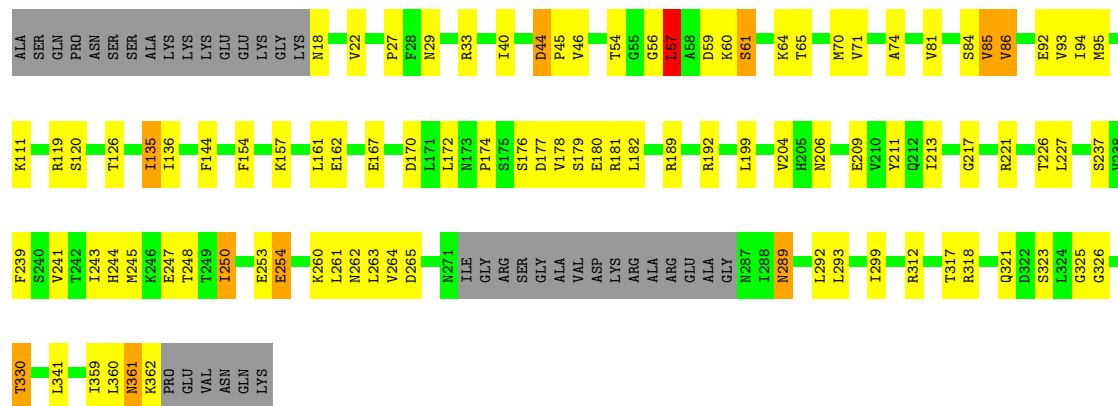
3 Residue-property plots

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. 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.

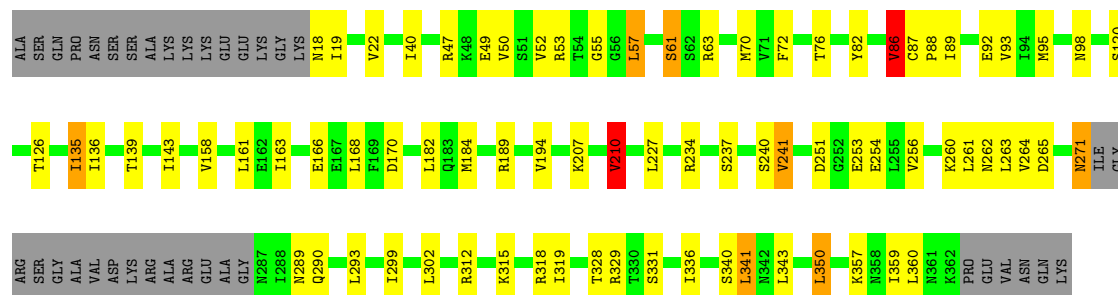
• Molecule 1: Kinesin-like protein KIF11

Chain A: 



• Molecule 1: Kinesin-like protein KIF11

Chain B: 



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	68.90Å 79.40Å 158.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.50	Depositor
% Data completeness (in resolution range)	85.0 (50.00-2.50)	Depositor
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.223 , 0.273	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5472	wwPDB-VP
Average B, all atoms (Å ²)	35.0	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: N5T, ADP, 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	0.59	0/2632	0.97	11/3559 (0.3%)
1	B	0.62	0/2632	0.97	7/3559 (0.2%)
All	All	0.60	0/5264	0.97	18/7118 (0.3%)

There are no bond length outliers.

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	86	VAL	N-CA-C	9.01	119.00	110.53
1	A	126	THR	N-CA-C	-7.42	99.54	110.52
1	B	210	VAL	N-CA-C	6.92	117.67	110.62
1	A	135	ILE	N-CA-C	6.71	117.40	110.36
1	B	135	ILE	N-CA-C	6.63	117.32	110.36
1	B	61	SER	N-CA-C	6.59	117.27	108.38
1	A	86	VAL	N-CA-C	6.49	117.28	110.72
1	B	120	SER	N-CA-C	-6.46	101.44	109.64
1	B	264	VAL	N-CA-C	6.27	117.14	107.99
1	B	126	THR	N-CA-C	-5.80	101.10	110.32
1	A	85	VAL	N-CA-C	5.67	117.05	111.67
1	A	264	VAL	N-CA-C	5.62	115.95	107.80
1	A	120	SER	N-CA-C	-5.61	102.34	110.08
1	A	44	ASP	CA-C-N	5.35	124.80	119.24
1	A	44	ASP	C-N-CA	5.35	124.80	119.24
1	A	211	TYR	N-CA-C	5.16	117.31	111.11
1	A	61	SER	N-CA-C	5.15	115.14	107.88
1	A	40	ILE	N-CA-C	-5.07	107.53	112.29

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	59	0
1	B	2594	0	2618	54	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	24	0	17	1	0
4	B	24	0	17	0	0
5	A	92	0	0	3	0
5	B	88	0	0	4	0
All	All	5472	0	5294	113	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 11.

All (113) 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:40:ILE:HD13	1:B:340:SER:HA	1.48	0.94
1:A:92:GLU:HA	1:A:95:MET:HE3	1.55	0.89
1:A:299:ILE:HG23	1:A:359:ILE:HD11	1.57	0.87
1:B:18:ASN:HB2	1:B:360:LEU:HD12	1.60	0.82
1:A:56:GLY:O	1:A:57:LEU:HB2	1.79	0.80
1:B:341:LEU:O	1:B:341:LEU:HG	1.86	0.76
1:A:326:GLY:O	1:A:361:ASN:HB2	1.86	0.75
1:A:54:THR:HG21	1:A:64:LYS:HD3	1.70	0.73
1:A:323:SER:O	1:A:330:THR:HG21	1.92	0.70
1:A:70:MET:HE1	1:A:84:SER:HB3	1.77	0.67
1:A:18:ASN:HB2	1:A:360:LEU:HD12	1.77	0.66
1:B:92:GLU:OE2	1:B:329:ARG:HD2	1.97	0.64
1:A:176:SER:HB2	1:A:180:GLU:OE1	1.98	0.64
1:A:204:VAL:HG22	1:A:213:ILE:CD1	2.28	0.63
1:A:237:SER:OG	1:A:265:ASP:HB3	1.97	0.63
1:A:260:LYS:HE2	1:A:262:ASN:HD21	1.64	0.61
1:A:170:ASP:HB2	1:A:182:LEU:HD11	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:22:VAL:HG12	1:B:70:MET:HB2	1.83	0.61
1:A:361:ASN:O	1:A:362:LYS:HB2	2.01	0.60
1:A:57:LEU:HB3	1:A:60:LYS:O	2.02	0.59
1:B:57:LEU:HD12	1:B:57:LEU:H	1.66	0.58
1:B:184:MET:HE3	1:B:318:ARG:HH21	1.70	0.57
1:A:29:ASN:O	1:A:33:ARG:HG2	2.05	0.57
1:A:57:LEU:CB	1:A:60:LYS:O	2.54	0.55
1:A:204:VAL:HG22	1:A:213:ILE:HD12	1.88	0.55
1:A:178:VAL:HG23	1:A:227:LEU:HD23	1.89	0.55
1:B:40:ILE:CD1	1:B:340:SER:HA	2.29	0.55
1:A:289:ASN:HD22	1:A:292:LEU:H	1.55	0.54
1:B:92:GLU:HA	1:B:95:MET:HE3	1.89	0.54
1:B:22:VAL:CG1	1:B:70:MET:HB2	2.38	0.54
1:B:312:ARG:HA	1:B:318:ARG:HG3	1.89	0.54
1:B:87:CYS:HB3	1:B:88:PRO:HD3	1.89	0.54
1:A:325:GLY:HA2	1:A:361:ASN:HB3	1.90	0.54
1:A:248:THR:HA	1:A:253:GLU:O	2.08	0.53
1:B:19:ILE:HD12	1:B:359:ILE:HB	1.90	0.53
1:B:189:ARG:HD3	5:B:621:HOH:O	2.07	0.53
1:A:45:PRO:HA	1:A:71:VAL:HG23	1.89	0.53
1:A:136:ILE:HG12	1:A:263:LEU:HD12	1.89	0.53
1:A:93:VAL:HG21	1:A:261:LEU:HB2	1.91	0.53
1:A:227:LEU:HD12	1:A:227:LEU:O	2.09	0.53
1:B:55:GLY:O	1:B:61:SER:HB2	2.07	0.53
1:B:161:LEU:HD21	1:B:319:ILE:HD13	1.91	0.53
1:A:44:ASP:OD1	1:A:46:VAL:HG12	2.08	0.52
1:B:299:ILE:HG23	1:B:359:ILE:HD11	1.90	0.52
1:B:63:ARG:HD2	5:B:611:HOH:O	2.10	0.52
1:B:166:GLU:OE1	1:B:315:LYS:HG2	2.10	0.52
1:B:271:ASN:C	1:B:271:ASN:HD22	2.19	0.51
1:B:76:THR:HA	5:B:630:HOH:O	2.10	0.51
1:A:57:LEU:O	1:A:61:SER:HB3	2.10	0.51
1:A:119:ARG:HG2	4:A:604:N5T:C35	2.41	0.51
1:B:237:SER:OG	1:B:265:ASP:HB3	2.11	0.51
1:B:241:VAL:HG13	1:B:261:LEU:HB3	1.93	0.51
1:B:72:PHE:HD1	1:B:76:THR:HG21	1.77	0.50
1:B:136:ILE:HG12	1:B:263:LEU:HD13	1.93	0.50
1:A:247:GLU:O	1:A:254:GLU:HA	2.12	0.49
1:A:217:GLY:O	1:A:221:ARG:HB2	2.12	0.49
1:B:139:THR:O	1:B:143:ILE:HG13	2.13	0.49
1:B:318:ARG:HH11	1:B:318:ARG:HG2	1.78	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:81:VAL:O	1:A:85:VAL:HG22	2.13	0.48
1:A:192:ARG:HB2	1:A:321:GLN:OE1	2.14	0.48
1:A:312:ARG:HA	1:A:318:ARG:CG	2.44	0.48
1:A:144:PHE:CE1	1:A:206:ASN:HA	2.48	0.48
1:A:172:LEU:O	1:A:174:PRO:HD3	2.13	0.48
1:B:57:LEU:O	1:B:61:SER:HB3	2.14	0.48
1:B:207:LYS:O	1:B:210:VAL:HG13	2.14	0.48
1:A:94:ILE:O	1:A:245:MET:HE1	2.14	0.47
1:A:18:ASN:HB2	1:A:360:LEU:CD1	2.43	0.47
1:A:167:GLU:HG2	1:A:181:ARG:HE	1.78	0.47
1:A:209:GLU:O	1:A:213:ILE:HG13	2.13	0.47
1:A:241:VAL:O	1:A:241:VAL:HG13	2.14	0.47
1:A:341:LEU:HD12	1:A:341:LEU:H	1.80	0.47
1:B:251:ASP:HB2	5:B:644:HOH:O	2.15	0.47
1:B:260:LYS:HE2	1:B:262:ASN:HD21	1.81	0.46
1:B:47:ARG:HG2	1:B:49:GLU:CD	2.41	0.45
1:B:184:MET:HE3	1:B:318:ARG:NH2	2.31	0.45
1:A:18:ASN:CB	1:A:360:LEU:HD12	2.43	0.45
1:A:111:LYS:HE2	5:A:606:HOH:O	2.17	0.45
1:A:177:ASP:OD1	1:A:179:SER:HB2	2.16	0.45
1:A:250:ILE:H	1:A:250:ILE:HD13	1.80	0.45
1:B:82:TYR:CD2	1:B:86:VAL:HB	2.52	0.45
1:B:336:ILE:HG21	1:B:350:LEU:CD1	2.47	0.45
1:B:52:VAL:O	1:B:63:ARG:HA	2.17	0.44
1:A:243:ILE:HG22	1:A:245:MET:HG3	1.99	0.44
1:B:53:ARG:HB2	1:B:63:ARG:NH1	2.33	0.44
1:B:170:ASP:HB2	1:B:182:LEU:HD11	1.99	0.44
1:A:162:GLU:HG2	1:A:237:SER:HA	2.00	0.44
1:A:27:PRO:HB3	1:A:74:ALA:HB1	1.99	0.43
1:B:92:GLU:CD	1:B:329:ARG:HD2	2.43	0.43
1:B:86:VAL:HG21	1:B:135:ILE:HG12	2.00	0.43
1:A:157:LYS:HG3	5:A:665:HOH:O	2.18	0.43
1:A:317:THR:O	1:A:321:GLN:HB3	2.19	0.43
1:B:290:GLN:NE2	1:B:290:GLN:HA	2.33	0.43
1:B:163:ILE:HG12	1:B:168:LEU:HD23	2.00	0.42
1:B:158:VAL:HA	1:B:240:SER:O	2.19	0.42
1:B:184:MET:HG2	1:B:194:VAL:HG21	2.01	0.42
1:B:253:GLU:HG2	1:B:254:GLU:N	2.34	0.42
1:A:86:VAL:HG21	1:A:135:ILE:HG12	2.01	0.42
1:A:239:PHE:CD1	1:A:239:PHE:C	2.98	0.42
1:A:154:PHE:HA	1:A:244:HIS:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:325:GLY:CA	1:A:361:ASN:HB3	2.49	0.41
1:A:85:VAL:HG23	1:A:86:VAL:N	2.35	0.41
1:B:82:TYR:HA	1:B:86:VAL:HG23	2.03	0.41
1:B:98:ASN:C	1:B:328:THR:HG23	2.44	0.41
1:B:184:MET:CE	1:B:318:ARG:HE	2.33	0.41
1:B:312:ARG:HA	1:B:318:ARG:CG	2.50	0.41
1:A:45:PRO:HD2	5:A:647:HOH:O	2.20	0.41
1:B:72:PHE:CD1	1:B:76:THR:HG21	2.56	0.41
1:B:329:ARG:HG2	1:B:329:ARG:HH11	1.86	0.41
1:B:256:VAL:O	1:B:256:VAL:HG23	2.20	0.41
1:B:57:LEU:HD12	1:B:57:LEU:N	2.35	0.41
1:B:89:ILE:O	1:B:93:VAL:HG23	2.21	0.40
1:A:263:LEU:HD23	1:A:263:LEU:HA	1.87	0.40
1:A:57:LEU:HB2	1:A:60:LYS:O	2.22	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	326/367 (89%)	314 (96%)	10 (3%)	2 (1%)	21	38
1	B	326/367 (89%)	320 (98%)	6 (2%)	0	100	100
All	All	652/734 (89%)	634 (97%)	16 (2%)	2 (0%)	36	55

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	57	LEU
1	A	189	ARG

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	293/321 (91%)	280 (96%)	13 (4%)	25	50
1	B	293/321 (91%)	277 (94%)	16 (6%)	19	40
All	All	586/642 (91%)	557 (95%)	29 (5%)	22	45

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	VAL
1	A	57	LEU
1	A	59	ASP
1	A	65	THR
1	A	161	LEU
1	A	199	LEU
1	A	226	THR
1	A	250	ILE
1	A	254	GLU
1	A	289	ASN
1	A	293	LEU
1	A	330	THR
1	A	361	ASN
1	B	50	VAL
1	B	57	LEU
1	B	86	VAL
1	B	210	VAL
1	B	227	LEU
1	B	234	ARG
1	B	241	VAL
1	B	271	ASN
1	B	289	ASN
1	B	293	LEU
1	B	302	LEU
1	B	331	SER
1	B	341	LEU
1	B	343	LEU

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Mol	Chain	Res	Type
1	B	350	LEU
1	B	357	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	173	ASN
1	A	229	ASN
1	A	262	ASN
1	A	287	ASN
1	A	289	ASN
1	A	361	ASN
1	B	18	ASN
1	B	20	GLN
1	B	141	HIS
1	B	212	GLN
1	B	244	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 oligosaccharides 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	N5T	A	604	-	25,26,26	3.04	14 (56%)	29,37,37	3.10	11 (37%)
3	ADP	B	600	2	28,29,29	1.63	5 (17%)	43,45,45	2.02	9 (20%)
3	ADP	A	601	2	28,29,29	1.62	5 (17%)	43,45,45	1.96	9 (20%)
4	N5T	B	605	-	25,26,26	3.14	14 (56%)	29,37,37	3.14	11 (37%)

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	N5T	A	604	-	-	4/16/28/28	0/3/3/3
3	ADP	B	600	2	-	7/16/32/32	0/3/3/3
3	ADP	A	601	2	-	8/16/32/32	0/3/3/3
4	N5T	B	605	-	-	4/16/28/28	0/3/3/3

All (38) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	604	N5T	C14-C10	6.90	1.46	1.34
4	B	605	N5T	C14-C10	6.61	1.46	1.34
4	B	605	N5T	C3-C4	5.13	1.45	1.38
4	B	605	N5T	C23-C17	5.02	1.47	1.39
4	A	604	N5T	C3-C4	4.93	1.45	1.38
4	A	604	N5T	C19-C17	4.88	1.46	1.39
4	B	605	N5T	C19-C17	4.87	1.46	1.39
3	B	600	ADP	C5-N7	-4.56	1.30	1.39
4	B	605	N5T	C22-C23	4.48	1.46	1.38
4	A	604	N5T	C20-C19	4.32	1.46	1.38
3	A	601	ADP	C5-N7	-4.31	1.31	1.39
4	A	604	N5T	C23-C17	4.30	1.45	1.39
4	A	604	N5T	C17-C13	-4.12	1.48	1.52
4	B	605	N5T	C20-C19	4.00	1.45	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	605	N5T	C17-C13	-3.98	1.48	1.52
4	A	604	N5T	C22-C23	3.78	1.45	1.38
4	B	605	N5T	C21-C20	3.63	1.46	1.38
4	A	604	N5T	C21-C20	3.51	1.46	1.38
4	B	605	N5T	C22-C21	3.44	1.45	1.38
4	B	605	N5T	C6-C1	3.32	1.43	1.37
4	A	604	N5T	C6-C1	3.28	1.43	1.37
3	B	600	ADP	C4-N9	-3.23	1.31	1.37
4	A	604	N5T	C5-C4	3.18	1.44	1.37
4	A	604	N5T	C22-C21	2.95	1.44	1.38
4	B	605	N5T	C2-C1	2.77	1.42	1.37
3	A	601	ADP	C4-N9	-2.77	1.31	1.37
4	A	604	N5T	F41-C1	-2.67	1.29	1.36
3	A	601	ADP	C5'-C4'	2.63	1.59	1.51
4	A	604	N5T	C2-C1	2.55	1.41	1.37
4	B	605	N5T	C5-C4	2.54	1.43	1.37
4	B	605	N5T	C2-C3	2.44	1.43	1.39
3	B	600	ADP	C4-N3	2.40	1.38	1.34
4	B	605	N5T	C6-C5	2.36	1.42	1.38
3	A	601	ADP	C4-N3	2.23	1.38	1.34
3	B	600	ADP	C5'-C4'	2.22	1.58	1.51
3	B	600	ADP	C5-C4	2.09	1.42	1.39
3	A	601	ADP	C1'-N9	2.06	1.52	1.46
4	A	604	N5T	C29-N30	-2.01	1.30	1.35

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	605	N5T	C17-C13-N12	11.94	128.11	112.37
4	A	604	N5T	C17-C13-N12	11.17	127.09	112.37
3	A	601	ADP	N3-C2-N1	-7.14	117.78	128.58
3	B	600	ADP	N3-C2-N1	-7.10	117.83	128.58
4	B	605	N5T	C13-N12-C29	7.02	140.65	121.83
4	A	604	N5T	C13-N12-C29	6.37	138.92	121.83
3	B	600	ADP	C5-C4-N3	-4.88	120.00	126.72
3	A	601	ADP	C5-C4-N3	-4.84	120.05	126.72
3	A	601	ADP	C2-N3-C4	4.40	122.57	111.83
3	B	600	ADP	C2-N3-C4	4.37	122.50	111.83
3	B	600	ADP	N3-C4-N9	3.99	133.95	127.17
3	A	601	ADP	N3-C4-N9	3.87	133.75	127.17
4	A	604	N5T	C6-C5-C4	-3.82	114.33	119.04
4	A	604	N5T	C35-N30-C31	3.55	127.21	115.87

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	604	N5T	C31-N30-C29	-3.55	105.80	120.61
4	A	604	N5T	C17-C13-C14	3.50	121.95	112.98
3	B	600	ADP	C5-N7-C8	3.21	108.49	103.45
4	B	605	N5T	C35-N30-C31	3.16	125.97	115.87
4	B	605	N5T	C17-C13-C14	3.14	121.02	112.98
4	B	605	N5T	C23-C17-C19	3.04	122.08	118.30
4	B	605	N5T	C22-C21-C20	2.97	123.93	119.87
3	B	600	ADP	N9-C8-N7	-2.95	109.76	113.94
4	B	605	N5T	C6-C5-C4	-2.95	115.41	119.04
4	A	604	N5T	C5-C4-C3	2.84	126.12	123.07
3	A	601	ADP	C5-N7-C8	2.83	107.89	103.45
3	A	601	ADP	N9-C8-N7	-2.76	110.02	113.94
4	A	604	N5T	C23-C17-C19	2.64	121.57	118.30
3	B	600	ADP	C4-C5-N7	-2.54	107.68	110.58
4	B	605	N5T	C31-N30-C29	-2.52	110.07	120.61
4	A	604	N5T	C6-C1-C2	2.40	126.41	123.23
3	A	601	ADP	C2-N1-C6	2.28	122.48	118.73
4	B	605	N5T	C21-C22-C23	-2.27	117.43	120.24
3	A	601	ADP	C4-C5-N7	-2.27	107.99	110.58
4	A	604	N5T	C3-C10-C14	2.23	131.25	126.89
3	B	600	ADP	C4-N9-C8	2.18	108.03	105.74
4	A	604	N5T	F41-C1-C2	-2.09	115.30	118.28
3	B	600	ADP	C2-N1-C6	2.07	122.14	118.73
3	A	601	ADP	C4-N9-C8	2.06	107.90	105.74
4	B	605	N5T	C23-C17-C13	-2.05	117.45	120.64
4	B	605	N5T	C6-C1-C2	2.01	125.89	123.23

There are no chirality outliers.

All (23) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	601	ADP	PA-O3A-PB-O2B
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-O2B
3	B	600	ADP	C5'-O5'-PA-O1A
3	B	600	ADP	C5'-O5'-PA-O2A
3	B	600	ADP	C5'-O5'-PA-O3A
4	A	604	N5T	N30-C29-N12-C11
4	A	604	N5T	O39-C29-N12-C11

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Mol	Chain	Res	Type	Atoms
4	B	605	N5T	N30-C29-N12-C11
4	B	605	N5T	O39-C29-N12-C11
3	A	601	ADP	C3'-C4'-C5'-O5'
3	B	600	ADP	O4'-C4'-C5'-O5'
3	B	600	ADP	C3'-C4'-C5'-O5'
3	A	601	ADP	O4'-C4'-C5'-O5'
3	B	600	ADP	PA-O3A-PB-O3B
4	B	605	N5T	N30-C29-N12-C13
4	B	605	N5T	O39-C29-N12-C13
4	A	604	N5T	N30-C29-N12-C13
4	A	604	N5T	O39-C29-N12-C13
3	A	601	ADP	PA-O3A-PB-O1B

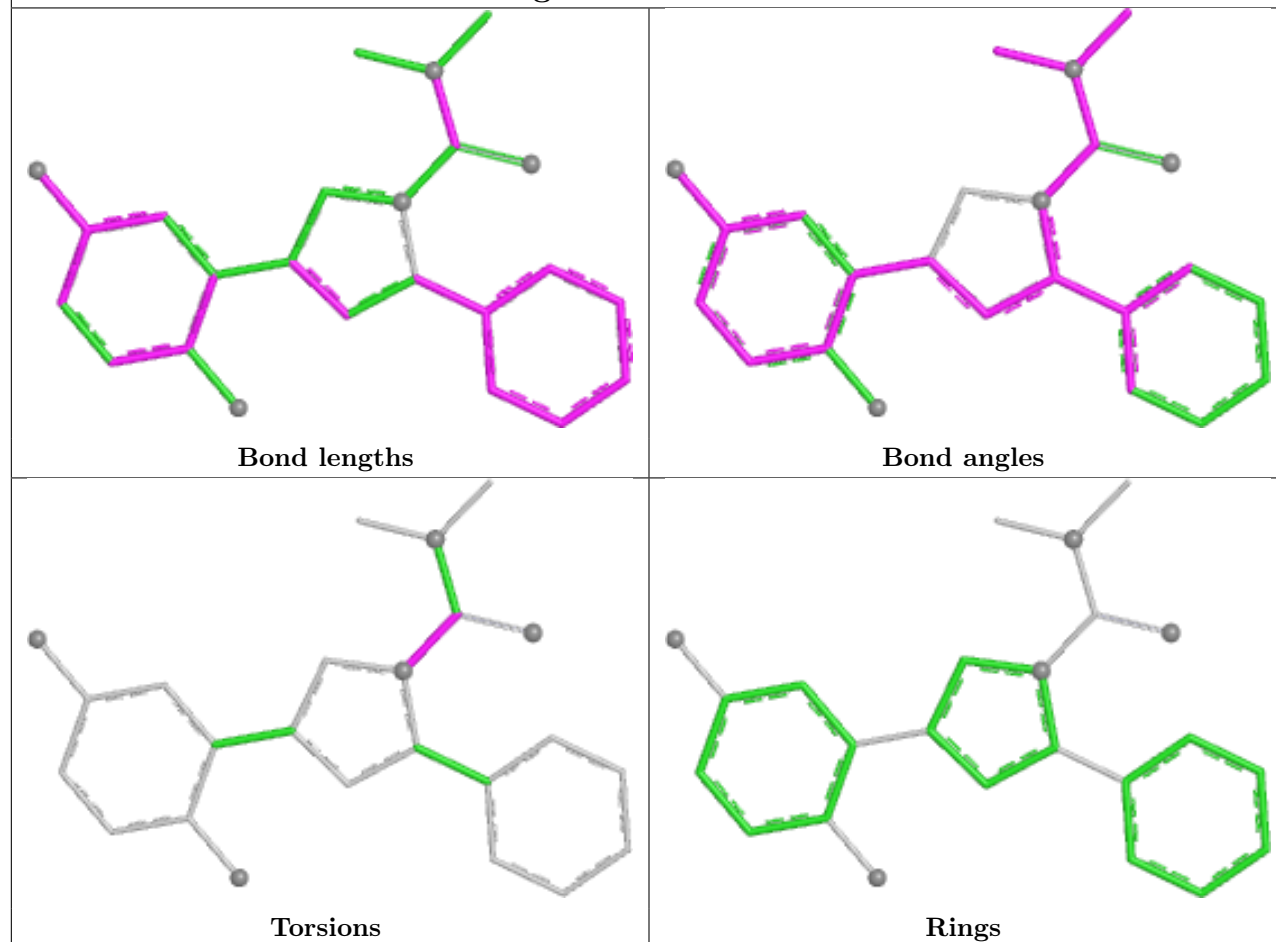
There are no ring outliers.

1 monomer is involved in 1 short contact:

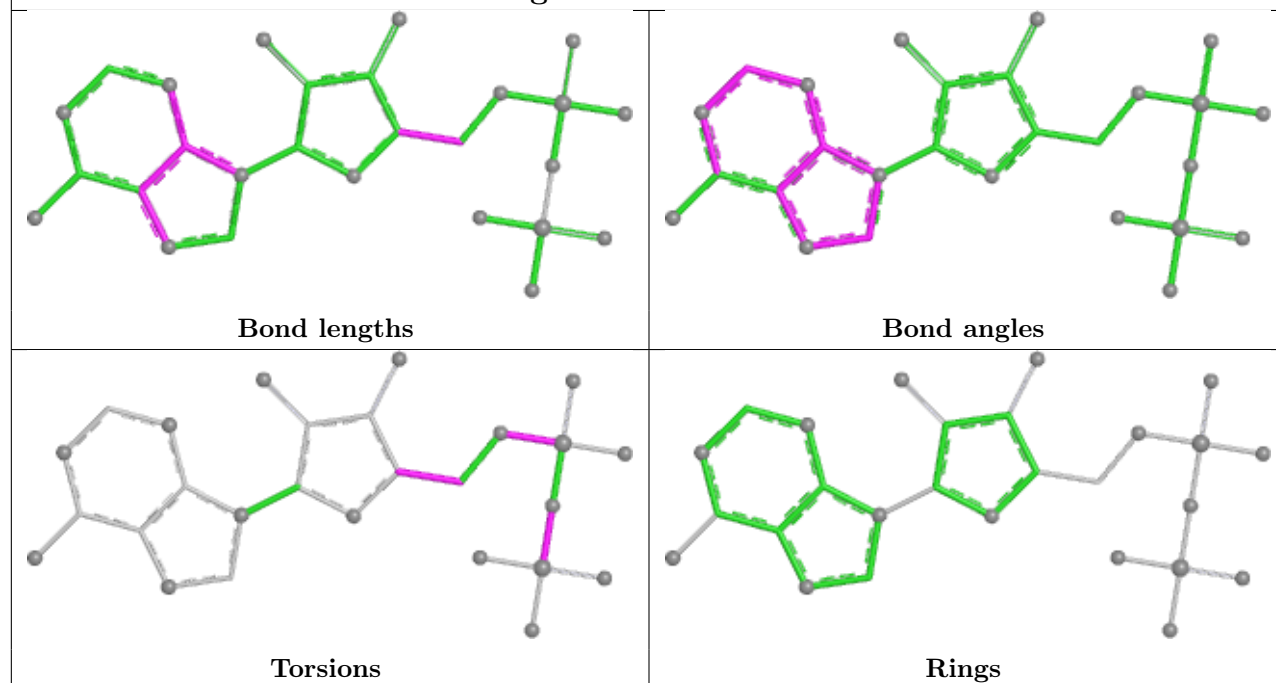
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	604	N5T	1	0

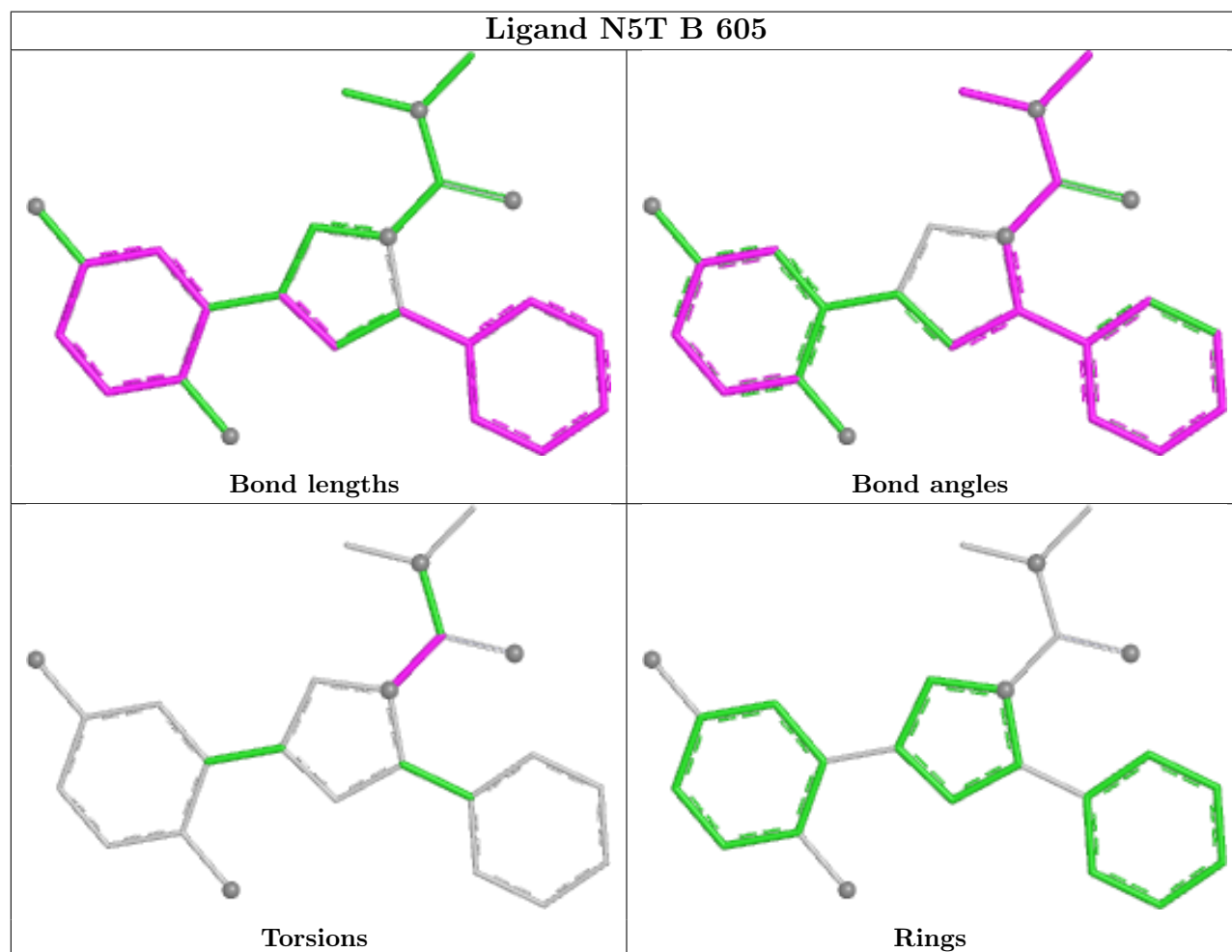
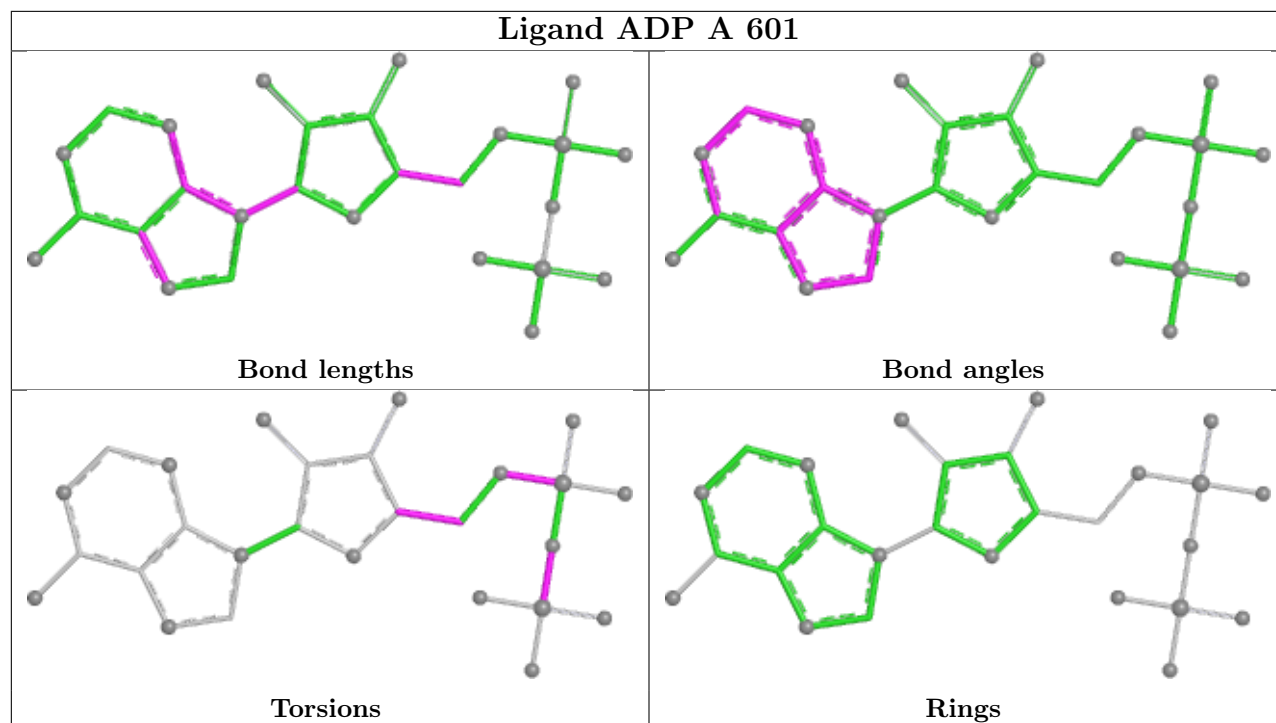
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

Ligand N5T A 604



Ligand ADP B 600





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.