



Full wwPDB EM Validation Report ⓘ

Nov 8, 2022 – 01:42 AM EST

PDB ID : 6D00
EMDB ID : EMD-7782
Title : Calcarisporiella thermophila Hsp104
Authors : Zhang, K.; Pintilie, G.
Deposited on : 2018-04-09
Resolution : 4.00 Å(reported)

This is a Full wwPDB EM 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/EMValidationReportHelp>

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:

EMDB validation analysis : 0.0.1.dev43
Mogul : 1.8.5 (274361), CSD as541be (2020)
MolProbity : 4.02b-467
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ : **FAILED**
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.31.2

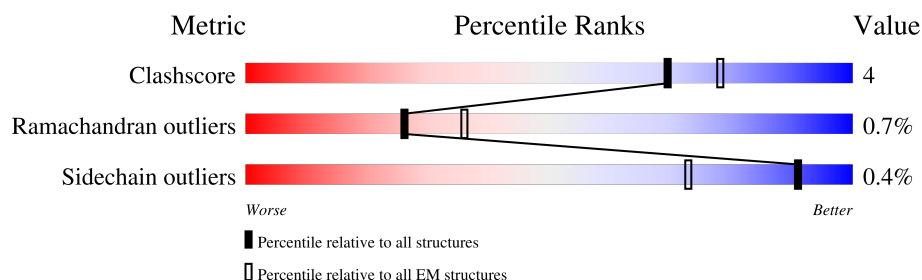
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 4.00 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)	EM structures (#Entries)
Clashscore	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the 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 $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	1	883	81% 10% 9%
1	2	883	80% 11% 9%
1	3	883	80% 11% 9%
1	4	883	81% 10% 9%
1	5	883	80% 11% 9%
1	6	883	81% 10% 9%

2 Entry composition [i](#)

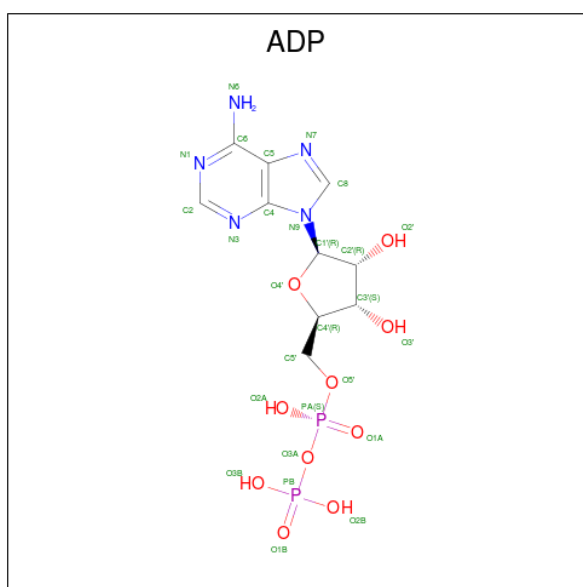
There are 2 unique types of molecules in this entry. The entry contains 38328 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, 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 *Calcarisporiella thermophila* Hsp104.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	1	804	Total	C	N	O	S	0	0
			6334	3965	1150	1194	25		
1	2	804	Total	C	N	O	S	0	0
			6334	3965	1150	1194	25		
1	3	804	Total	C	N	O	S	0	0
			6334	3965	1150	1194	25		
1	4	804	Total	C	N	O	S	0	0
			6334	3965	1150	1194	25		
1	5	804	Total	C	N	O	S	0	0
			6334	3965	1150	1194	25		
1	6	804	Total	C	N	O	S	0	0
			6334	3965	1150	1194	25		

- Molecule 2 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).

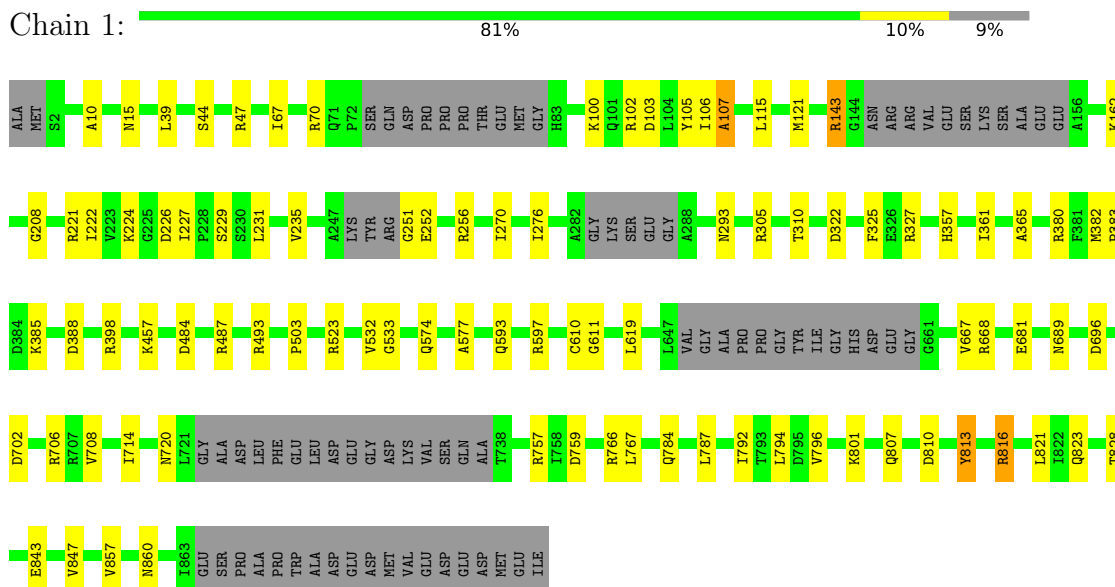


Mol	Chain	Residues	Atoms					AltConf
2	1	1	Total	C	N	O	P	0
			54	20	10	20	4	
2	1	1	Total	C	N	O	P	0
			54	20	10	20	4	
2	2	1	Total	C	N	O	P	0
			54	20	10	20	4	
2	2	1	Total	C	N	O	P	0
			54	20	10	20	4	
2	3	1	Total	C	N	O	P	0
			54	20	10	20	4	
2	3	1	Total	C	N	O	P	0
			54	20	10	20	4	
2	4	1	Total	C	N	O	P	0
			54	20	10	20	4	
2	4	1	Total	C	N	O	P	0
			54	20	10	20	4	
2	5	1	Total	C	N	O	P	0
			54	20	10	20	4	
2	5	1	Total	C	N	O	P	0
			54	20	10	20	4	
2	6	1	Total	C	N	O	P	0
			54	20	10	20	4	
2	6	1	Total	C	N	O	P	0
			54	20	10	20	4	

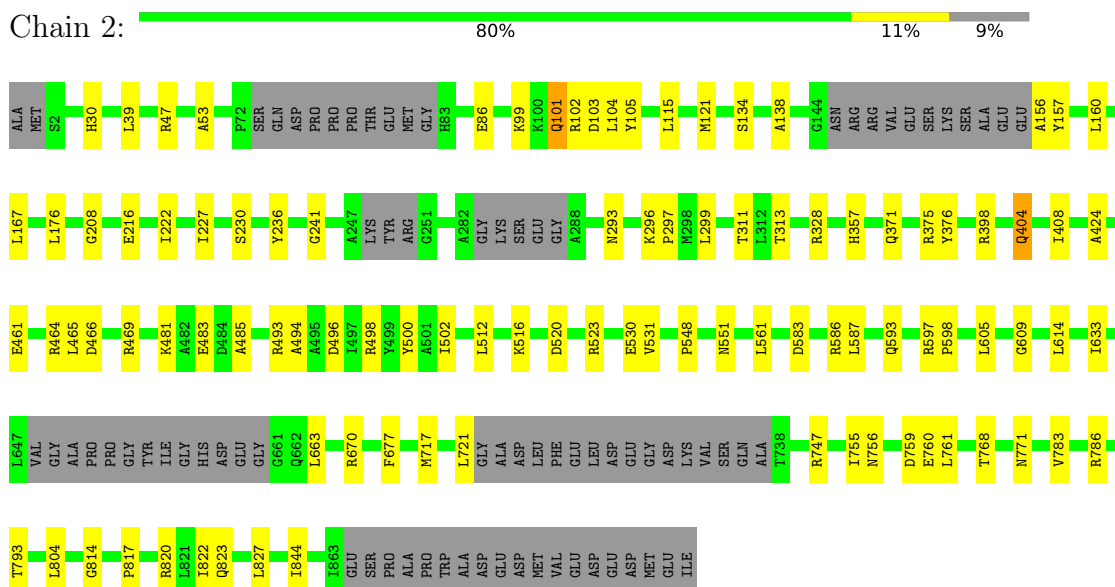
3 Residue-property plots

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.


• Molecule 1: *Calcarisporiella thermophila* Hsp104

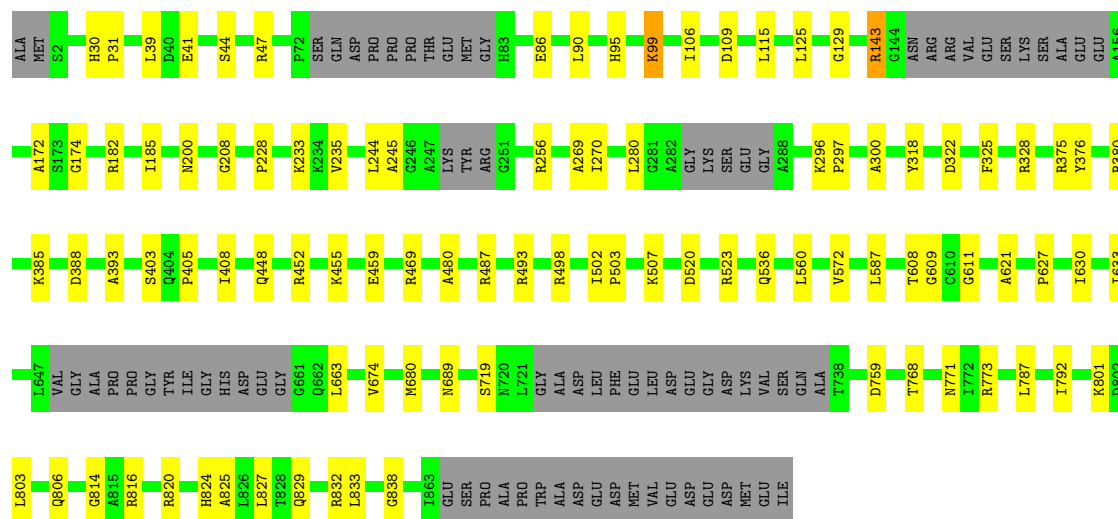


• Molecule 1: *Calcarisporiella thermophila* Hsp104




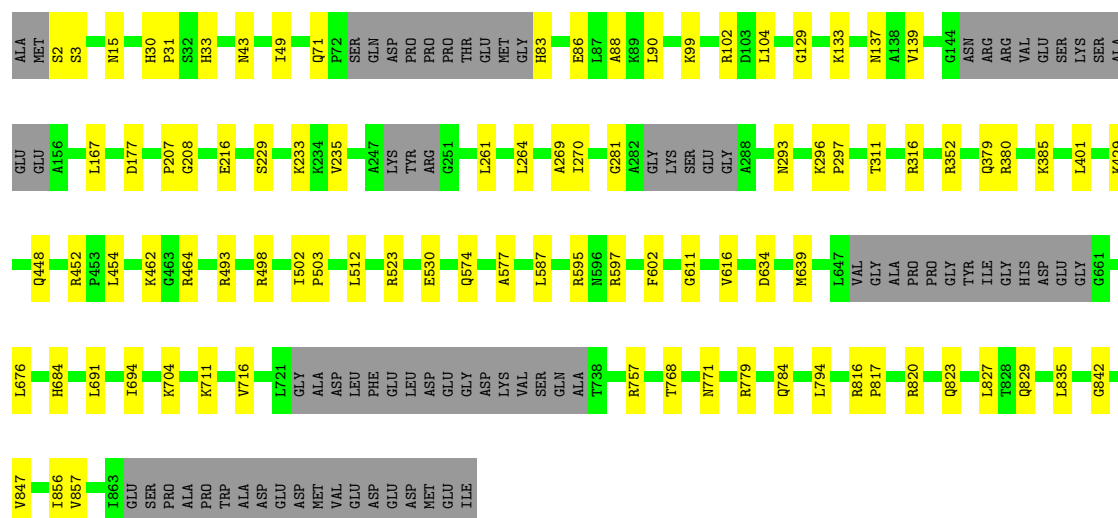
• Molecule 1: *Calcarisporiella thermophila* Hsp104

Chain 3:  80% 11% 9%




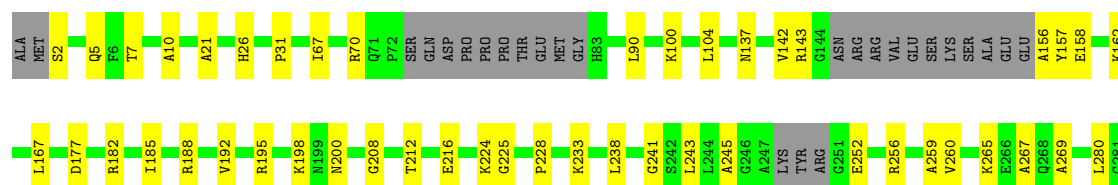
• Molecule 1: *Calcarisporiella thermophila* Hsp104

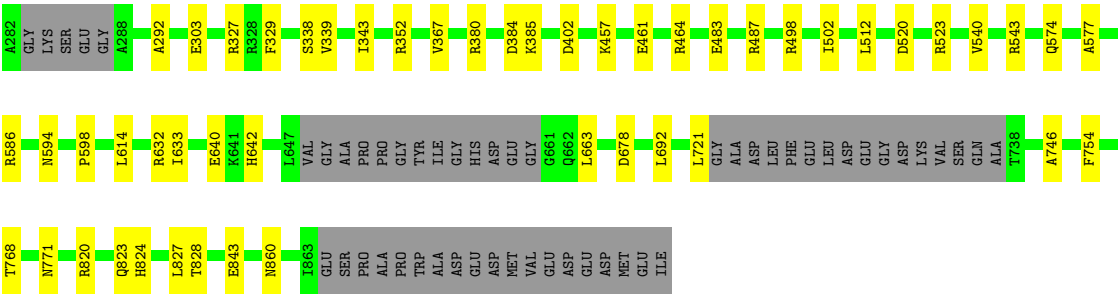
Chain 4:  81% 10% 9%



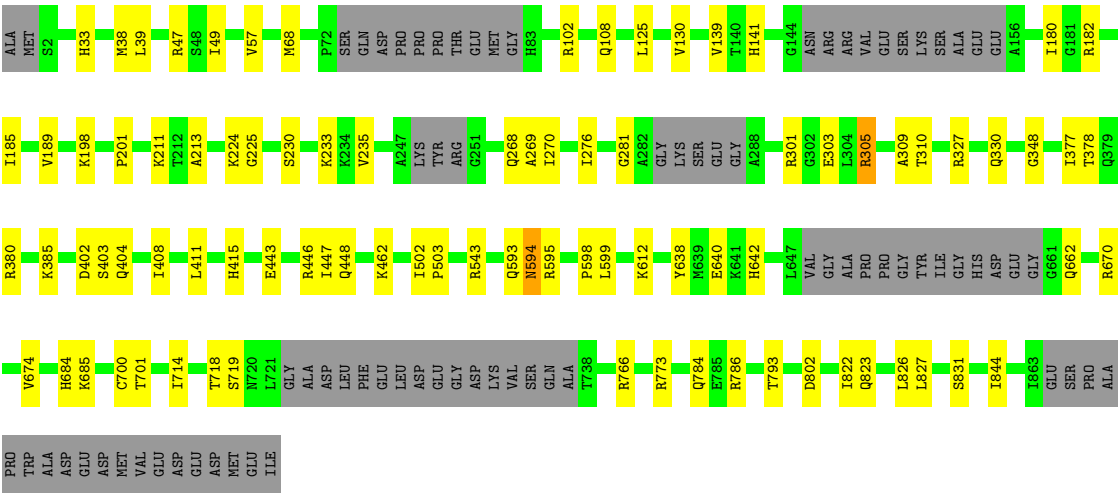
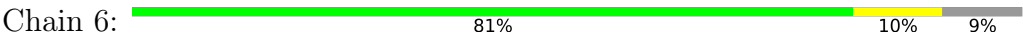
• Molecule 1: *Calcarisporiella thermophila* Hsp104

Chain 5:  80% 11% 9%





● Molecule 1: Calcarisporiella thermophila Hsp104



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	224915	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION; Gctf	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	31.8	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	130000	Depositor
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 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	1	0.22	0/6411	0.38	0/8631
1	2	0.23	0/6411	0.37	0/8631
1	3	0.22	0/6411	0.37	0/8631
1	4	0.22	0/6411	0.37	0/8631
1	5	0.22	0/6411	0.37	0/8631
1	6	0.22	0/6411	0.37	0/8631
All	All	0.22	0/38466	0.37	0/51786

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	1	6334	0	6497	53	0
1	2	6334	0	6497	58	0
1	3	6334	0	6497	56	0
1	4	6334	0	6497	49	0
1	5	6334	0	6497	56	0
1	6	6334	0	6497	50	0
2	1	54	0	24	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	2	54	0	24	3	0
2	3	54	0	24	3	0
2	4	54	0	24	1	0
2	5	54	0	24	2	0
2	6	54	0	24	2	0
All	All	38328	0	39126	287	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 4.

All (287) 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)
1:2:156:ALA:N	1:2:236:TYR:HH	1.72	0.88
1:3:106:ILE:HG13	1:4:102:ARG:HH21	1.58	0.69
1:1:816:ARG:HH21	1:6:303:GLU:HG2	1.61	0.65
1:4:784:GLN:HE22	1:4:794:LEU:H	1.42	0.65
1:6:404:GLN:HE21	1:6:408:ILE:HG22	1.62	0.64
1:2:157:TYR:H	1:2:160:LEU:HB2	1.63	0.62
1:5:594:ASN:HD22	1:6:786:ARG:HH21	1.44	0.62
1:1:102:ARG:HH11	1:2:103:ASP:HA	1.65	0.61
1:1:457:LYS:HG3	1:1:523:ARG:HH21	1.65	0.61
1:3:31:PRO:HB3	1:3:90:LEU:HD22	1.83	0.61
1:4:779:ARG:NH1	2:4:902:ADP:N3	2.49	0.60
1:3:300:ALA:HA	1:3:328:ARG:HE	1.67	0.60
1:2:481:LYS:HB3	1:2:493:ARG:HH12	1.66	0.60
1:1:821:LEU:HD13	1:6:230:SER:HB3	1.84	0.60
1:2:357:HIS:O	1:2:398:ARG:NH2	2.34	0.60
1:1:784:GLN:HE22	1:1:794:LEU:H	1.51	0.58
1:6:443:GLU:OE1	1:6:446:ARG:NH2	2.36	0.58
1:1:668:ARG:HD2	1:1:708:VAL:HG22	1.85	0.58
1:1:357:HIS:O	1:1:398:ARG:NH2	2.36	0.58
1:4:602:PHE:HB2	1:4:716:VAL:HG12	1.84	0.58
1:2:39:LEU:HD22	1:2:47:ARG:HH21	1.69	0.58
1:5:633:ILE:HG13	1:5:663:LEU:HD12	1.84	0.58
1:3:95:HIS:HB3	1:4:99:LYS:HB3	1.86	0.57
1:6:674:VAL:HG22	1:6:714:ILE:HD11	1.86	0.57
1:2:53:ALA:HB1	1:2:134:SER:HB3	1.87	0.57
1:6:38:MET:SD	1:6:108:GLN:NE2	2.75	0.57
1:3:448:GLN:HE22	1:3:452:ARG:HE	1.52	0.57
1:5:7:THR:OG1	1:5:162:LYS:NZ	2.31	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:3:41:GLU:HB2	1:3:44:SER:HB2	1.86	0.57
1:4:380:ARG:HB2	1:4:385:LYS:HB3	1.86	0.56
1:2:593:GLN:HE22	1:2:597:ARG:HH21	1.53	0.56
1:5:192:VAL:HG13	1:5:195:ARG:HH21	1.69	0.56
1:6:213:ALA:HB2	2:6:901:ADP:H2'	1.86	0.56
1:5:245:ALA:O	1:5:256:ARG:NH2	2.39	0.56
1:6:700:CYS:SG	1:6:701:THR:N	2.78	0.56
1:2:404:GLN:HE21	1:2:408:ILE:HG22	1.70	0.56
1:3:816:ARG:NH2	2:3:902:ADP:O2B	2.39	0.56
1:4:464:ARG:HG3	1:4:512:LEU:HD22	1.87	0.56
1:3:109:ASP:OD2	1:3:143:ARG:NH1	2.39	0.55
1:4:823:GLN:HA	1:4:827:LEU:HB2	1.88	0.55
1:5:632:ARG:NH1	1:5:678:ASP:OD2	2.40	0.55
1:5:233:LYS:HB3	1:5:269:ALA:HA	1.89	0.55
1:1:696:ASP:OD2	1:1:757:ARG:NH1	2.40	0.54
1:1:39:LEU:O	1:1:47:ARG:NH2	2.41	0.54
1:5:185:ILE:HG12	1:5:188:ARG:HH21	1.72	0.54
1:2:759:ASP:O	1:3:820:ARG:NH2	2.40	0.54
1:2:167:LEU:HD21	1:2:216:GLU:HA	1.90	0.54
1:3:787:LEU:HB3	1:3:792:ILE:HB	1.90	0.54
1:3:609:GLY:HA3	1:3:814:GLY:HA3	1.90	0.54
1:6:39:LEU:O	1:6:47:ARG:NH2	2.41	0.54
1:6:49:ILE:HD13	1:6:139:VAL:HG12	1.90	0.54
1:1:327:ARG:NH2	2:2:901:ADP:O3A	2.41	0.54
1:2:376:TYR:O	1:2:670:ARG:NH1	2.41	0.54
1:5:21:ALA:HB1	1:5:26:HIS:HB2	1.89	0.53
1:1:235:VAL:HG22	1:1:270:ILE:HD11	1.91	0.53
1:5:457:LYS:NZ	1:5:461:GLU:OE2	2.41	0.53
1:3:39:LEU:HD11	1:3:47:ARG:HE	1.73	0.53
1:5:380:ARG:HH21	1:5:384:ASP:HB3	1.74	0.53
1:4:31:PRO:HB3	1:4:90:LEU:HD22	1.90	0.53
1:2:747:ARG:HH21	1:2:761:LEU:HD22	1.74	0.53
1:6:211:LYS:HD2	1:6:309:ALA:HB1	1.91	0.52
1:5:228:PRO:HG3	1:6:403:SER:HB2	1.90	0.52
1:5:252:GLU:OE2	1:5:256:ARG:NH1	2.43	0.52
1:1:484:ASP:OD1	1:1:487:ARG:NH2	2.38	0.52
1:6:402:ASP:OD1	1:6:462:LYS:NZ	2.42	0.52
1:3:825:ALA:O	1:3:829:GLN:NE2	2.43	0.52
1:3:393:ALA:HB1	1:3:536:GLN:HG3	1.92	0.52
1:3:455:LYS:NZ	1:3:459:GLU:OE2	2.43	0.52
1:3:680:MET:HB3	1:3:719:SER:HB2	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:4:691:LEU:HD13	1:4:694:ILE:HD12	1.92	0.52
1:5:200:ASN:HB2	1:5:329:PHE:HA	1.91	0.52
1:1:224:LYS:HA	1:1:487:ARG:HA	1.92	0.51
1:2:465:LEU:HD12	1:2:516:LYS:HE3	1.92	0.51
1:4:847:VAL:HG22	1:4:857:VAL:HG22	1.93	0.51
1:5:167:LEU:HD21	1:5:216:GLU:HA	1.90	0.51
1:6:233:LYS:HB3	1:6:269:ALA:HA	1.92	0.51
1:1:221:ARG:HB3	1:1:226:ASP:HB2	1.92	0.51
1:1:847:VAL:HG22	1:1:857:VAL:HG22	1.92	0.51
1:5:67:ILE:HG12	1:5:70:ARG:HH11	1.75	0.51
1:5:100:LYS:HD3	1:6:102:ARG:HB3	1.91	0.51
1:1:143:ARG:HD2	1:2:99:LYS:HE3	1.93	0.51
1:3:245:ALA:O	1:3:256:ARG:NH2	2.43	0.51
1:4:177:ASP:OD1	1:4:352:ARG:NE	2.43	0.51
1:6:793:THR:HB	1:6:844:ILE:HG12	1.93	0.51
1:6:773:ARG:NH2	1:6:802:ASP:OD1	2.44	0.51
1:4:49:ILE:HD13	1:4:139:VAL:HG12	1.91	0.51
1:1:231:LEU:HD22	1:2:469:ARG:HH22	1.75	0.51
1:2:614:LEU:HD22	2:2:902:ADP:H2'	1.93	0.51
1:6:198:LYS:HB3	1:6:330:GLN:HB2	1.93	0.50
1:6:380:ARG:HB2	1:6:385:LYS:HB3	1.93	0.50
1:1:807:GLN:NE2	1:6:230:SER:OG	2.44	0.50
1:5:327:ARG:NH2	2:6:901:ADP:O2B	2.44	0.50
1:1:810:ASP:HB3	1:1:813:TYR:HB3	1.93	0.50
1:3:322:ASP:HB2	1:3:325:PHE:HD2	1.76	0.50
1:3:633:ILE:HG13	1:3:663:LEU:HD12	1.92	0.50
1:5:843:GLU:OE2	1:5:860:ASN:ND2	2.45	0.50
1:1:115:LEU:HD12	1:1:121:MET:HG3	1.92	0.50
1:2:520:ASP:OD1	1:2:523:ARG:NH2	2.45	0.50
1:5:614:LEU:HD22	2:5:902:ADP:H2'	1.94	0.50
1:6:182:ARG:HD2	1:6:185:ILE:HD12	1.92	0.50
1:6:640:GLU:HG3	1:6:642:HIS:H	1.74	0.50
1:5:520:ASP:OD1	1:5:523:ARG:NH2	2.45	0.49
1:3:99:LYS:HD3	1:4:104:LEU:HA	1.94	0.49
1:3:689:ASN:ND2	1:4:634:ASP:OD2	2.45	0.49
1:3:773:ARG:HH21	1:3:801:LYS:HE3	1.77	0.49
1:6:189:VAL:HG13	1:6:201:PRO:HG3	1.94	0.49
1:6:784:GLN:HE22	1:6:793:THR:HA	1.76	0.49
1:2:823:GLN:HA	1:2:827:LEU:HB2	1.94	0.49
1:6:612:LYS:HD2	1:6:718:THR:HG23	1.94	0.49
1:4:316:ARG:NH2	1:4:704:LYS:O	2.45	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:5:182:ARG:HD2	1:5:185:ILE:HD12	1.94	0.49
1:6:593:GLN:NE2	1:6:598:PRO:O	2.46	0.49
1:1:787:LEU:HB3	1:1:792:ILE:HB	1.93	0.49
1:2:530:GLU:HG2	1:2:531:VAL:HG23	1.93	0.49
1:3:520:ASP:OD1	1:3:523:ARG:NH2	2.45	0.49
1:3:200:ASN:HD22	1:3:328:ARG:HB3	1.78	0.49
1:3:507:LYS:NZ	1:4:429:LYS:O	2.44	0.49
1:4:233:LYS:HB3	1:4:269:ALA:HA	1.94	0.49
1:3:768:THR:OG1	1:3:771:ASN:OD1	2.31	0.49
1:1:293:ASN:ND2	1:2:241:GLY:O	2.45	0.49
1:2:498:ARG:HD3	1:2:502:ILE:HD12	1.95	0.49
1:5:265:LYS:HD3	1:5:303:GLU:HB3	1.94	0.49
1:5:195:ARG:NH2	1:5:198:LYS:O	2.46	0.48
1:1:102:ARG:HE	1:1:143:ARG:HH21	1.61	0.48
1:1:105:TYR:HH	1:1:251:GLY:N	2.12	0.48
1:1:222:ILE:HG12	1:1:227:ILE:HD13	1.96	0.48
1:4:379:GLN:HG3	1:4:380:ARG:HG2	1.95	0.48
1:4:498:ARG:HD3	1:4:502:ILE:HD12	1.95	0.48
1:1:67:ILE:HG12	1:1:70:ARG:HH11	1.77	0.48
1:2:583:ASP:OD1	1:2:586:ARG:NH2	2.45	0.48
1:5:2:SER:HA	1:6:141:HIS:HB2	1.96	0.48
1:2:817:PRO:HA	1:2:820:ARG:HE	1.77	0.48
1:4:261:LEU:HD13	1:4:264:LEU:HD12	1.94	0.48
1:5:586:ARG:NH1	1:6:831:SER:OG	2.46	0.48
1:3:115:LEU:HB3	1:3:125:LEU:HD11	1.95	0.48
1:1:322:ASP:HB3	1:1:325:PHE:HD2	1.78	0.48
1:2:633:ILE:HG13	1:2:663:LEU:HD12	1.94	0.48
1:4:639:MET:HA	1:4:684:HIS:HB2	1.96	0.48
1:3:233:LYS:HB3	1:3:269:ALA:HA	1.96	0.48
1:3:405:PRO:HD2	1:3:408:ILE:HD12	1.95	0.48
1:2:548:PRO:HD2	1:2:551:ASN:HD22	1.78	0.48
1:2:804:LEU:HD13	1:2:822:ILE:HG12	1.95	0.48
1:5:143:ARG:HG3	1:5:157:TYR:HA	1.96	0.48
1:1:667:VAL:HG23	1:1:708:VAL:HG11	1.96	0.47
1:1:702:ASP:HB3	1:1:706:ARG:H	1.79	0.47
1:1:796:VAL:O	1:1:801:LYS:NZ	2.47	0.47
1:3:182:ARG:HD2	1:3:185:ILE:HD12	1.95	0.47
1:4:448:GLN:HB3	1:4:452:ARG:HE	1.79	0.47
1:2:768:THR:OG1	1:2:771:ASN:OD1	2.31	0.47
1:4:587:LEU:HD11	1:5:827:LEU:HB3	1.95	0.47
1:2:747:ARG:HA	1:2:755:ILE:HG21	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:3:759:ASP:O	1:4:820:ARG:NH2	2.47	0.47
1:4:207:PRO:HD3	1:4:311:THR:HG23	1.96	0.47
1:6:411:LEU:HD22	1:6:447:ILE:HG23	1.96	0.47
1:2:293:ASN:HD21	1:3:244:LEU:HD23	1.78	0.47
1:1:843:GLU:OE1	1:1:860:ASN:ND2	2.48	0.47
1:4:2:SER:OG	1:4:3:SER:N	2.45	0.47
1:4:293:ASN:ND2	1:5:241:GLY:O	2.48	0.47
1:5:498:ARG:HA	1:5:502:ILE:HD12	1.96	0.47
1:6:33:HIS:NE2	1:6:68:MET:SD	2.76	0.47
1:3:380:ARG:NH1	1:3:388:ASP:OD2	2.38	0.46
1:2:230:SER:HA	1:3:469:ARG:HH12	1.81	0.46
1:2:609:GLY:N	2:2:902:ADP:O1B	2.43	0.46
1:4:71:GLN:NE2	1:4:86:GLU:OE1	2.48	0.46
1:1:574:GLN:HB3	1:1:577:ALA:HB3	1.96	0.46
1:2:30:HIS:NE2	1:2:86:GLU:OE1	2.47	0.46
1:2:677:PHE:HB2	1:2:717:MET:HG2	1.97	0.46
1:5:104:LEU:HD23	1:5:259:ALA:HB2	1.96	0.46
1:2:464:ARG:HG3	1:2:512:LEU:HD22	1.97	0.46
1:3:498:ARG:HD3	1:3:502:ILE:HD12	1.98	0.46
1:3:608:THR:OG1	2:3:902:ADP:O2B	2.30	0.46
1:4:574:GLN:HB3	1:4:577:ALA:HB3	1.97	0.46
1:6:235:VAL:HG22	1:6:270:ILE:HD11	1.98	0.46
1:1:252:GLU:OE2	1:1:256:ARG:NE	2.49	0.46
1:6:378:THR:HG21	1:6:670:ARG:HG3	1.98	0.46
1:6:594:ASN:HD22	1:6:595:ARG:H	1.63	0.46
1:1:720:ASN:HB3	1:6:301:ARG:HD3	1.96	0.45
1:2:299:LEU:HB2	1:2:328:ARG:HH11	1.82	0.45
1:3:380:ARG:HB2	1:3:385:LYS:HB3	1.97	0.45
1:4:167:LEU:HD21	1:4:216:GLU:HA	1.97	0.45
1:1:610:CYS:HA	1:1:767:LEU:HD21	1.96	0.45
1:2:587:LEU:HD11	1:3:827:LEU:HB3	1.98	0.45
1:5:31:PRO:HB3	1:5:90:LEU:HD22	1.98	0.45
1:5:158:GLU:HG3	1:5:162:LYS:HE2	1.98	0.45
1:5:343:ILE:HD13	1:5:367:VAL:HG22	1.98	0.45
1:1:10:ALA:HB2	1:1:107:ALA:HA	1.98	0.45
1:5:820:ARG:HH11	1:5:824:HIS:HE1	1.65	0.45
1:6:125:LEU:HB3	1:6:130:VAL:HB	1.98	0.45
1:2:760:GLU:OE2	1:3:824:HIS:NE2	2.50	0.45
1:5:177:ASP:OD1	1:5:352:ARG:NE	2.49	0.45
1:2:485:ALA:HB3	1:2:494:ALA:HB2	1.99	0.45
1:6:612:LYS:NZ	1:6:719:SER:O	2.41	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:44:SER:O	1:1:162:LYS:NZ	2.44	0.45
1:5:768:THR:OG1	1:5:771:ASN:OD1	2.34	0.45
1:6:47:ARG:HG3	1:6:57:VAL:HG11	1.98	0.45
1:1:597:ARG:NH1	1:1:759:ASP:OD1	2.45	0.45
1:2:561:LEU:HA	1:3:832:ARG:HH12	1.82	0.45
1:2:783:VAL:HG22	1:2:786:ARG:HH21	1.82	0.45
1:6:823:GLN:HA	1:6:827:LEU:HB2	1.99	0.45
1:5:380:ARG:HB2	1:5:385:LYS:HB3	1.99	0.45
1:5:157:TYR:HE2	1:5:267:ALA:HB2	1.82	0.44
1:5:692:LEU:HD13	1:5:754:PHE:HD1	1.82	0.44
1:6:684:HIS:CD2	1:6:685:LYS:H	2.36	0.44
1:4:829:GLN:HE22	1:4:856:ILE:HG13	1.82	0.44
1:3:296:LYS:HB3	1:3:297:PRO:HD3	1.99	0.44
1:5:721:LEU:HD23	1:5:746:ALA:HA	2.00	0.44
1:1:100:LYS:HB2	1:2:101:GLN:HB2	2.00	0.44
1:3:587:LEU:HD11	1:4:827:LEU:HB3	1.99	0.44
1:5:5:GLN:NE2	1:5:104:LEU:O	2.51	0.44
1:1:270:ILE:HG22	1:1:305:ARG:HB3	2.00	0.44
1:2:311:THR:HG22	1:2:313:THR:H	1.83	0.44
1:4:83:HIS:HE1	1:4:88:ALA:HB2	1.83	0.44
1:4:133:LYS:NZ	1:4:137:ASN:OD1	2.49	0.44
1:6:377:ILE:HG21	1:6:380:ARG:HH21	1.82	0.44
1:5:280:LEU:HD13	1:5:292:ALA:HA	1.99	0.44
1:2:609:GLY:HA3	1:2:814:GLY:HA3	1.99	0.44
1:6:305:ARG:H	1:6:305:ARG:HD2	1.83	0.44
1:3:630:ILE:HG12	1:3:674:VAL:HB	2.00	0.44
1:2:793:THR:HB	1:2:844:ILE:HG12	2.00	0.43
1:3:621:ALA:HB2	1:3:627:PRO:HG3	2.00	0.43
1:5:823:GLN:O	1:5:828:THR:OG1	2.36	0.43
1:3:803:LEU:HD12	1:3:806:GLN:HE21	1.83	0.43
1:4:296:LYS:HB3	1:4:297:PRO:HD3	1.98	0.43
1:6:822:ILE:HG23	1:6:826:LEU:HD23	2.00	0.43
1:3:560:LEU:HB3	1:4:835:LEU:HD21	1.99	0.43
1:2:222:ILE:HG12	1:2:227:ILE:HD13	2.01	0.43
1:2:461:GLU:HB3	1:2:516:LYS:HE2	2.00	0.43
1:3:280:LEU:HD22	1:3:318:TYR:HB3	2.01	0.43
1:4:768:THR:OG1	1:4:771:ASN:ND2	2.52	0.43
1:5:7:THR:HG23	1:5:10:ALA:H	1.84	0.43
1:4:597:ARG:HH22	1:4:757:ARG:HE	1.67	0.43
1:5:238:LEU:HD13	1:5:260:VAL:HG11	2.01	0.43
1:5:483:GLU:HB3	1:5:487:ARG:HH12	1.84	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:483:GLU:HG3	1:2:498:ARG:HH21	1.84	0.43
1:5:464:ARG:HG2	1:5:512:LEU:HD22	2.01	0.43
1:6:270:ILE:HG22	1:6:305:ARG:HD3	2.01	0.43
1:4:229:SER:H	1:5:402:ASP:HB3	1.84	0.43
1:5:142:VAL:O	1:5:156:ALA:N	2.52	0.43
1:2:176:LEU:HD21	1:2:216:GLU:HB3	2.00	0.42
1:3:228:PRO:HG3	1:4:462:LYS:HD2	2.01	0.42
1:4:502:ILE:HB	1:4:503:PRO:HD3	2.01	0.42
1:5:574:GLN:HB3	1:5:577:ALA:HB3	2.00	0.42
1:3:572:VAL:O	2:3:902:ADP:N6	2.52	0.42
1:3:375:ARG:HH11	1:3:376:TYR:HE2	1.66	0.42
1:1:382:MET:HA	1:1:383:PRO:HA	1.87	0.42
1:3:172:ALA:O	1:3:487:ARG:NH2	2.53	0.42
1:1:593:GLN:HE21	1:1:597:ARG:HD3	1.85	0.42
1:4:401:LEU:HD12	1:4:530:GLU:HG2	2.02	0.42
1:6:233:LYS:HE3	1:6:268:GLN:HG2	2.01	0.42
1:2:53:ALA:HB2	1:2:138:ALA:HB2	2.02	0.42
1:5:338:SER:OG	1:5:339:VAL:N	2.53	0.41
1:4:816:ARG:HB3	1:4:817:PRO:HD3	2.02	0.41
1:1:106:ILE:HD12	1:2:102:ARG:HH21	1.85	0.41
1:6:276:ILE:HG22	1:6:310:THR:HB	2.02	0.41
1:6:638:TYR:OH	1:6:662:GLN:OE1	2.38	0.41
1:3:174:GLY:HA3	1:3:480:ALA:HB1	2.02	0.41
1:4:454:LEU:HD22	1:4:523:ARG:HH12	1.85	0.41
1:5:212:THR:HB	2:5:901:ADP:H3'	2.02	0.41
1:1:361:ILE:HA	1:1:532:VAL:HB	2.02	0.41
1:1:380:ARG:HH22	1:1:388:ASP:HB3	1.86	0.41
1:1:823:GLN:O	1:1:828:THR:OG1	2.35	0.41
1:4:30:HIS:HB2	1:4:33:HIS:HD1	1.85	0.41
1:1:105:TYR:OH	1:1:251:GLY:N	2.54	0.41
1:2:605:LEU:HD22	1:2:721:LEU:HB2	2.01	0.41
1:6:502:ILE:HB	1:6:503:PRO:HD3	2.02	0.41
1:2:496:ASP:HA	1:2:500:TYR:HD2	1.85	0.41
1:3:502:ILE:HB	1:3:503:PRO:HD3	2.02	0.41
1:3:833:LEU:HD22	1:3:838:GLY:HA3	2.03	0.41
1:4:595:ARG:O	1:4:711:LYS:NZ	2.41	0.41
1:4:616:VAL:HG21	1:4:676:LEU:HD13	2.02	0.41
1:6:180:ILE:HD13	1:6:348:GLY:HA3	2.02	0.41
1:1:380:ARG:HB2	1:1:385:LYS:HB3	2.03	0.40
1:2:296:LYS:HB3	1:2:297:PRO:HD3	2.03	0.40
1:4:235:VAL:HG22	1:4:270:ILE:HD11	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:365:ALA:HB2	1:1:533:GLY:HA2	2.02	0.40
1:1:503:PRO:HB3	1:2:424:ALA:HB2	2.03	0.40
1:3:30:HIS:NE2	1:3:86:GLU:OE1	2.55	0.40
1:5:243:LEU:HD22	1:5:256:ARG:HB3	2.02	0.40
1:6:415:HIS:CD2	1:6:448:GLN:HE21	2.39	0.40
1:1:229:SER:OG	1:2:466:ASP:OD2	2.38	0.40
1:1:619:LEU:HD21	1:1:714:ILE:HD13	2.03	0.40
1:2:371:GLN:O	1:2:375:ARG:NE	2.50	0.40
1:3:235:VAL:HG22	1:3:270:ILE:HD11	2.03	0.40
1:5:640:GLU:HG2	1:5:642:HIS:H	1.86	0.40
1:1:276:ILE:HG22	1:1:310:THR:HB	2.02	0.40
1:2:115:LEU:HD12	1:2:121:MET:HG3	2.03	0.40
1:5:540:VAL:HG22	1:5:543:ARG:HH11	1.84	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 PDB entries followed by that with respect to all EM entries.

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	1	790/883 (90%)	716 (91%)	66 (8%)	8 (1%)	15	53
1	2	790/883 (90%)	740 (94%)	44 (6%)	6 (1%)	19	58
1	3	790/883 (90%)	732 (93%)	53 (7%)	5 (1%)	25	63
1	4	790/883 (90%)	747 (95%)	38 (5%)	5 (1%)	25	63
1	5	790/883 (90%)	736 (93%)	50 (6%)	4 (0%)	29	67
1	6	790/883 (90%)	729 (92%)	57 (7%)	4 (0%)	29	67
All	All	4740/5298 (90%)	4400 (93%)	308 (6%)	32 (1%)	26	61

All (32) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	1	103	ASP
1	2	101	GLN
1	2	105	TYR
1	6	224	LYS
1	1	611	GLY
1	2	104	LEU
1	2	404	GLN
1	4	611	GLY
1	5	225	GLY
1	6	225	GLY
1	6	599	LEU
1	1	107	ALA
1	1	143	ARG
1	1	813	TYR
1	2	598	PRO
1	3	403	SER
1	4	129	GLY
1	5	208	GLY
1	1	681	GLU
1	1	816	ARG
1	3	208	GLY
1	5	598	PRO
1	3	143	ARG
1	5	224	LYS
1	1	208	GLY
1	3	611	GLY
1	4	281	GLY
1	6	281	GLY
1	3	129	GLY
1	4	842	GLY
1	4	208	GLY
1	2	208	GLY

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 PDB entries followed by that with respect to all EM entries.

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	1	679/742 (92%)	675 (99%)	4 (1%)	86	92
1	2	679/742 (92%)	678 (100%)	1 (0%)	93	97
1	3	679/742 (92%)	677 (100%)	2 (0%)	92	95
1	4	679/742 (92%)	676 (100%)	3 (0%)	91	94
1	5	679/742 (92%)	678 (100%)	1 (0%)	93	97
1	6	679/742 (92%)	674 (99%)	5 (1%)	84	90
All	All	4074/4452 (92%)	4058 (100%)	16 (0%)	91	94

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	1	15	ASN
1	1	493	ARG
1	1	689	ASN
1	1	766	ARG
1	2	756	ASN
1	3	99	LYS
1	3	493	ARG
1	4	15	ASN
1	4	43	ASN
1	4	493	ARG
1	5	137	ASN
1	6	305	ARG
1	6	327	ARG
1	6	543	ARG
1	6	594	ASN
1	6	766	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (39) such sidechains are listed below:

Mol	Chain	Res	Type
1	1	15	ASN
1	1	71	GLN
1	1	101	GLN
1	1	415	HIS
1	1	551	ASN
1	1	593	GLN
1	1	689	ASN
1	1	784	GLN
1	1	807	GLN

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Mol	Chain	Res	Type
1	1	824	HIS
1	2	293	ASN
1	2	551	ASN
1	2	593	GLN
1	2	625	ASN
1	2	756	ASN
1	2	806	GLN
1	3	108	GLN
1	3	200	ASN
1	3	806	GLN
1	3	829	GLN
1	4	15	ASN
1	4	43	ASN
1	4	83	HIS
1	4	293	ASN
1	4	756	ASN
1	4	771	ASN
1	4	784	GLN
1	5	137	ASN
1	5	293	ASN
1	5	536	GLN
1	5	594	ASN
1	6	26	HIS
1	6	200	ASN
1	6	415	HIS
1	6	536	GLN
1	6	594	ASN
1	6	684	HIS
1	6	720	ASN
1	6	771	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

5.6 Ligand geometry ⓘ

12 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
2	ADP	2	902	-	24,29,29	0.95	1 (4%)	29,45,45	1.46	4 (13%)
2	ADP	6	901	-	24,29,29	0.95	1 (4%)	29,45,45	1.39	4 (13%)
2	ADP	1	902	-	24,29,29	0.96	1 (4%)	29,45,45	1.45	4 (13%)
2	ADP	5	901	-	24,29,29	0.96	1 (4%)	29,45,45	1.42	4 (13%)
2	ADP	4	901	-	24,29,29	0.96	1 (4%)	29,45,45	1.41	4 (13%)
2	ADP	1	901	-	24,29,29	0.95	1 (4%)	29,45,45	1.39	4 (13%)
2	ADP	2	901	-	24,29,29	0.96	1 (4%)	29,45,45	1.40	4 (13%)
2	ADP	3	901	-	24,29,29	0.97	1 (4%)	29,45,45	1.40	4 (13%)
2	ADP	4	902	-	24,29,29	0.96	1 (4%)	29,45,45	1.45	4 (13%)
2	ADP	5	902	-	24,29,29	0.96	1 (4%)	29,45,45	1.45	4 (13%)
2	ADP	6	902	-	24,29,29	0.95	1 (4%)	29,45,45	1.45	4 (13%)
2	ADP	3	902	-	24,29,29	0.96	1 (4%)	29,45,45	1.46	4 (13%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	ADP	2	902	-	-	1/12/32/32	0/3/3/3
2	ADP	6	901	-	-	4/12/32/32	0/3/3/3
2	ADP	1	902	-	-	2/12/32/32	0/3/3/3
2	ADP	5	901	-	-	2/12/32/32	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	ADP	4	901	-	-	3/12/32/32	0/3/3/3
2	ADP	1	901	-	-	5/12/32/32	0/3/3/3
2	ADP	2	901	-	-	5/12/32/32	0/3/3/3
2	ADP	3	901	-	-	5/12/32/32	0/3/3/3
2	ADP	4	902	-	-	0/12/32/32	0/3/3/3
2	ADP	5	902	-	-	1/12/32/32	0/3/3/3
2	ADP	6	902	-	-	2/12/32/32	0/3/3/3
2	ADP	3	902	-	-	1/12/32/32	0/3/3/3

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	3	901	ADP	C5-C4	2.53	1.47	1.40
2	2	901	ADP	C5-C4	2.50	1.47	1.40
2	2	902	ADP	C5-C4	2.50	1.47	1.40
2	1	902	ADP	C5-C4	2.50	1.47	1.40
2	4	901	ADP	C5-C4	2.49	1.47	1.40
2	1	901	ADP	C5-C4	2.49	1.47	1.40
2	5	902	ADP	C5-C4	2.49	1.47	1.40
2	6	902	ADP	C5-C4	2.49	1.47	1.40
2	3	902	ADP	C5-C4	2.48	1.47	1.40
2	5	901	ADP	C5-C4	2.48	1.47	1.40
2	4	902	ADP	C5-C4	2.48	1.47	1.40
2	6	901	ADP	C5-C4	2.47	1.47	1.40

All (48) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	2	902	ADP	PA-O3A-PB	-3.56	120.62	132.83
2	3	902	ADP	PA-O3A-PB	-3.51	120.77	132.83
2	5	902	ADP	PA-O3A-PB	-3.49	120.86	132.83
2	1	902	ADP	PA-O3A-PB	-3.48	120.88	132.83
2	4	902	ADP	PA-O3A-PB	-3.46	120.95	132.83
2	6	902	ADP	PA-O3A-PB	-3.46	120.95	132.83
2	2	901	ADP	C3'-C2'-C1'	3.33	105.99	100.98
2	3	901	ADP	C3'-C2'-C1'	3.32	105.98	100.98
2	2	902	ADP	C3'-C2'-C1'	3.32	105.98	100.98
2	6	901	ADP	C3'-C2'-C1'	3.31	105.97	100.98
2	3	902	ADP	C3'-C2'-C1'	3.30	105.94	100.98
2	6	902	ADP	C3'-C2'-C1'	3.30	105.94	100.98
2	5	902	ADP	C3'-C2'-C1'	3.29	105.93	100.98

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	4	902	ADP	C3'-C2'-C1'	3.28	105.92	100.98
2	4	901	ADP	C3'-C2'-C1'	3.28	105.91	100.98
2	4	901	ADP	PA-O3A-PB	-3.27	121.61	132.83
2	5	901	ADP	C3'-C2'-C1'	3.24	105.86	100.98
2	5	901	ADP	PA-O3A-PB	-3.23	121.76	132.83
2	1	902	ADP	C3'-C2'-C1'	3.23	105.83	100.98
2	1	901	ADP	N3-C2-N1	-3.19	123.69	128.68
2	2	901	ADP	N3-C2-N1	-3.19	123.70	128.68
2	4	901	ADP	N3-C2-N1	-3.18	123.70	128.68
2	1	901	ADP	C3'-C2'-C1'	3.18	105.77	100.98
2	6	901	ADP	N3-C2-N1	-3.18	123.71	128.68
2	5	902	ADP	N3-C2-N1	-3.18	123.71	128.68
2	4	902	ADP	N3-C2-N1	-3.17	123.72	128.68
2	3	902	ADP	N3-C2-N1	-3.17	123.72	128.68
2	2	902	ADP	N3-C2-N1	-3.17	123.73	128.68
2	3	901	ADP	N3-C2-N1	-3.17	123.73	128.68
2	5	901	ADP	N3-C2-N1	-3.17	123.73	128.68
2	1	902	ADP	N3-C2-N1	-3.16	123.73	128.68
2	6	902	ADP	N3-C2-N1	-3.14	123.77	128.68
2	3	901	ADP	PA-O3A-PB	-2.91	122.84	132.83
2	1	901	ADP	PA-O3A-PB	-2.90	122.89	132.83
2	2	901	ADP	PA-O3A-PB	-2.80	123.23	132.83
2	6	901	ADP	PA-O3A-PB	-2.76	123.35	132.83
2	3	902	ADP	C4-C5-N7	-2.72	106.56	109.40
2	1	901	ADP	C4-C5-N7	-2.72	106.56	109.40
2	3	901	ADP	C4-C5-N7	-2.71	106.58	109.40
2	4	902	ADP	C4-C5-N7	-2.69	106.59	109.40
2	6	901	ADP	C4-C5-N7	-2.69	106.59	109.40
2	5	901	ADP	C4-C5-N7	-2.69	106.60	109.40
2	5	902	ADP	C4-C5-N7	-2.69	106.60	109.40
2	2	901	ADP	C4-C5-N7	-2.68	106.61	109.40
2	6	902	ADP	C4-C5-N7	-2.67	106.61	109.40
2	1	902	ADP	C4-C5-N7	-2.67	106.61	109.40
2	4	901	ADP	C4-C5-N7	-2.65	106.64	109.40
2	2	902	ADP	C4-C5-N7	-2.62	106.67	109.40

There are no chirality outliers.

All (31) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	1	901	ADP	PB-O3A-PA-O5'
2	1	901	ADP	C5'-O5'-PA-O1A

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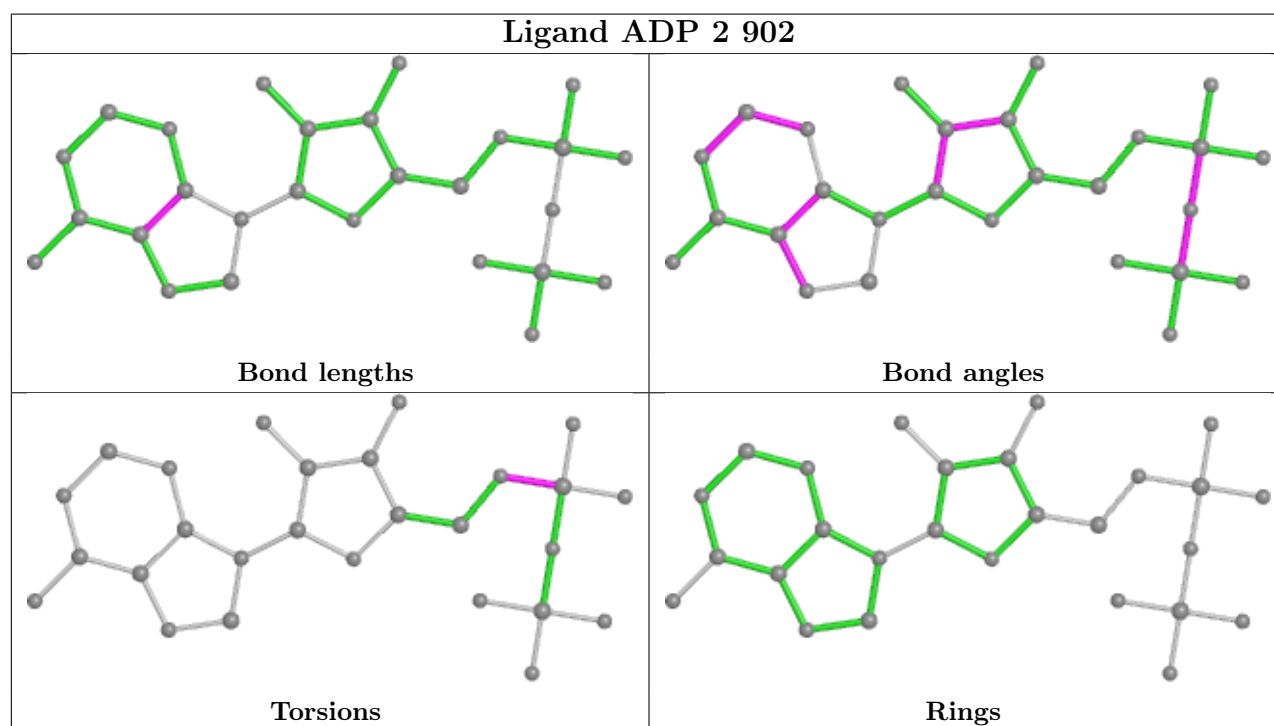
Mol	Chain	Res	Type	Atoms
2	2	901	ADP	C5'-O5'-PA-O1A
2	3	901	ADP	PB-O3A-PA-O5'
2	3	901	ADP	C5'-O5'-PA-O1A
2	3	902	ADP	C5'-O5'-PA-O1A
2	6	902	ADP	C5'-O5'-PA-O1A
2	2	901	ADP	PB-O3A-PA-O5'
2	6	901	ADP	PB-O3A-PA-O5'
2	1	901	ADP	C5'-O5'-PA-O3A
2	2	901	ADP	C5'-O5'-PA-O3A
2	3	901	ADP	C5'-O5'-PA-O3A
2	6	901	ADP	C5'-O5'-PA-O3A
2	6	901	ADP	C5'-O5'-PA-O1A
2	1	901	ADP	O4'-C4'-C5'-O5'
2	5	901	ADP	PB-O3A-PA-O2A
2	2	901	ADP	O4'-C4'-C5'-O5'
2	6	902	ADP	C5'-O5'-PA-O3A
2	1	902	ADP	O4'-C4'-C5'-O5'
2	6	901	ADP	O4'-C4'-C5'-O5'
2	4	901	ADP	PB-O3A-PA-O1A
2	4	901	ADP	PB-O3A-PA-O2A
2	5	901	ADP	PB-O3A-PA-O1A
2	1	901	ADP	C5'-O5'-PA-O2A
2	1	902	ADP	C5'-O5'-PA-O1A
2	2	901	ADP	C5'-O5'-PA-O2A
2	2	902	ADP	C5'-O5'-PA-O1A
2	3	901	ADP	C5'-O5'-PA-O2A
2	4	901	ADP	C5'-O5'-PA-O1A
2	5	902	ADP	C5'-O5'-PA-O1A
2	3	901	ADP	O4'-C4'-C5'-O5'

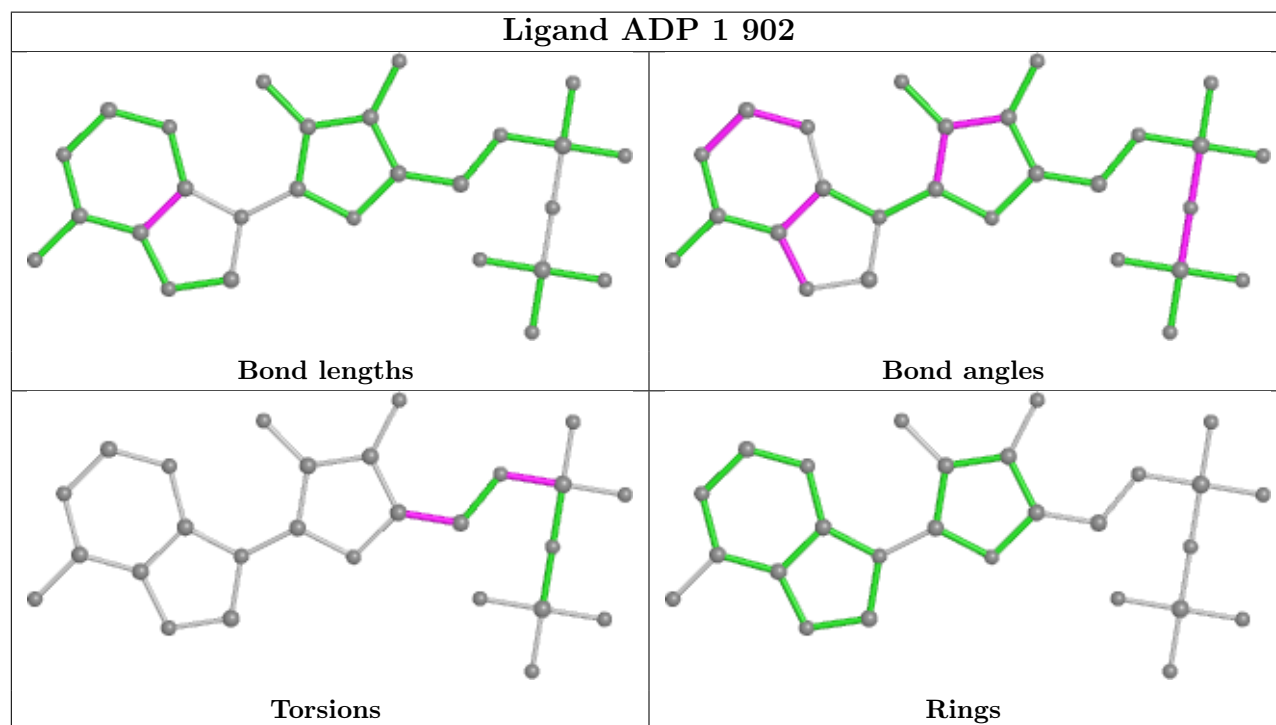
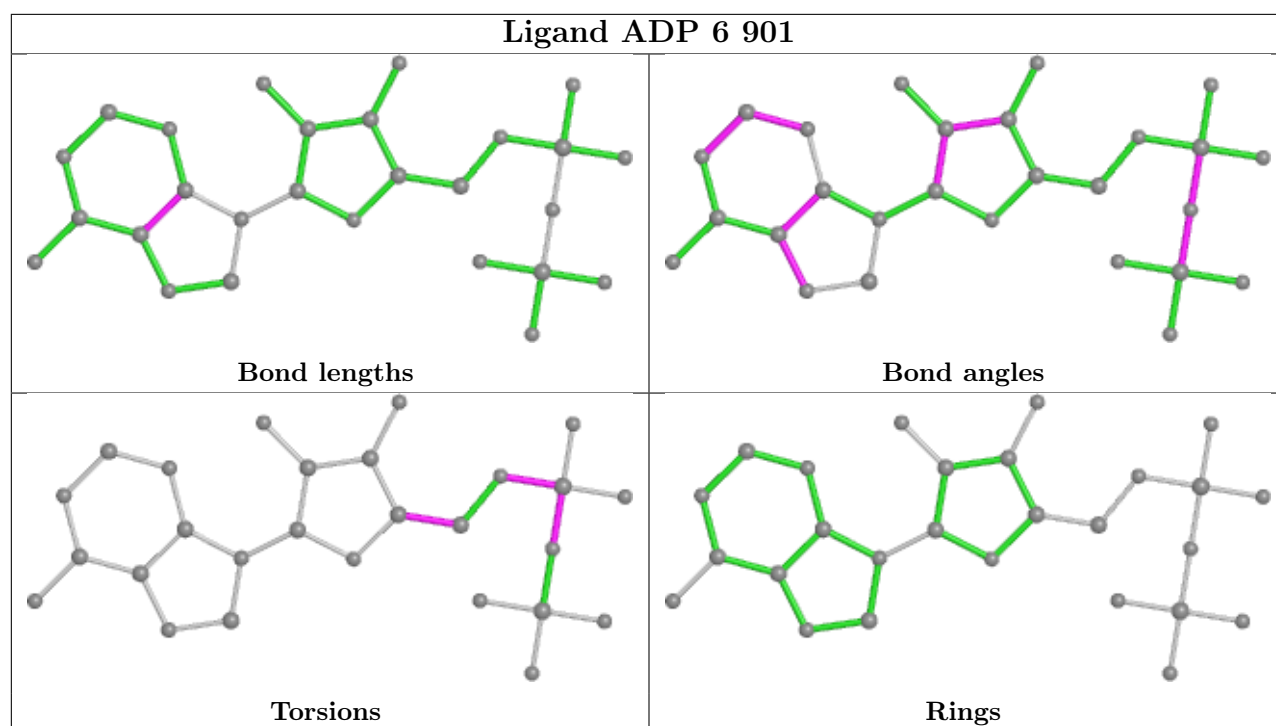
There are no ring outliers.

7 monomers are involved in 11 short contacts:

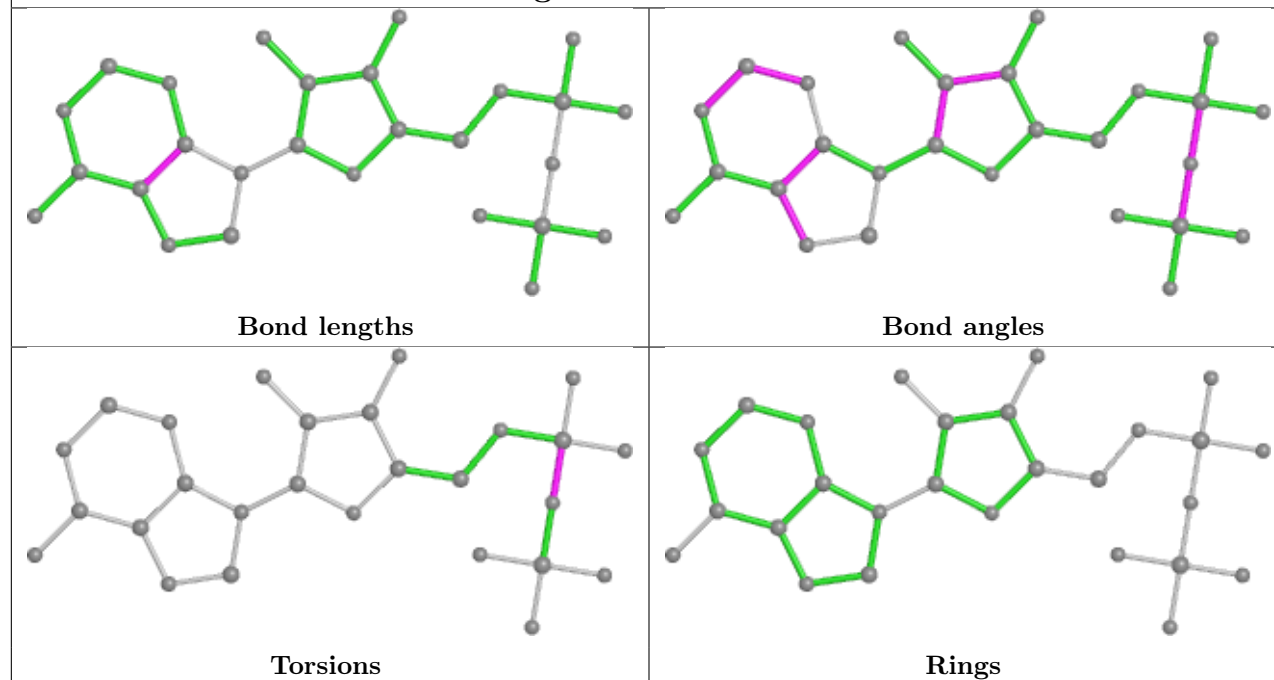
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	2	902	ADP	2	0
2	6	901	ADP	2	0
2	5	901	ADP	1	0
2	2	901	ADP	1	0
2	4	902	ADP	1	0
2	5	902	ADP	1	0
2	3	902	ADP	3	0

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.

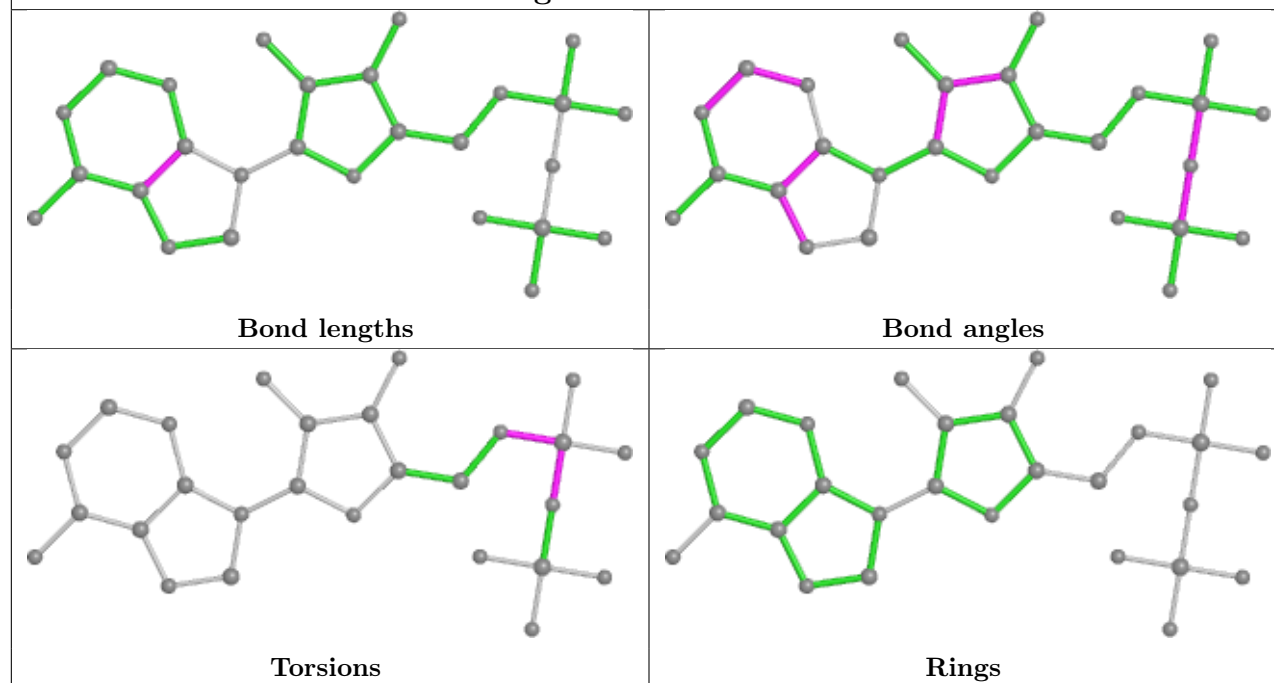


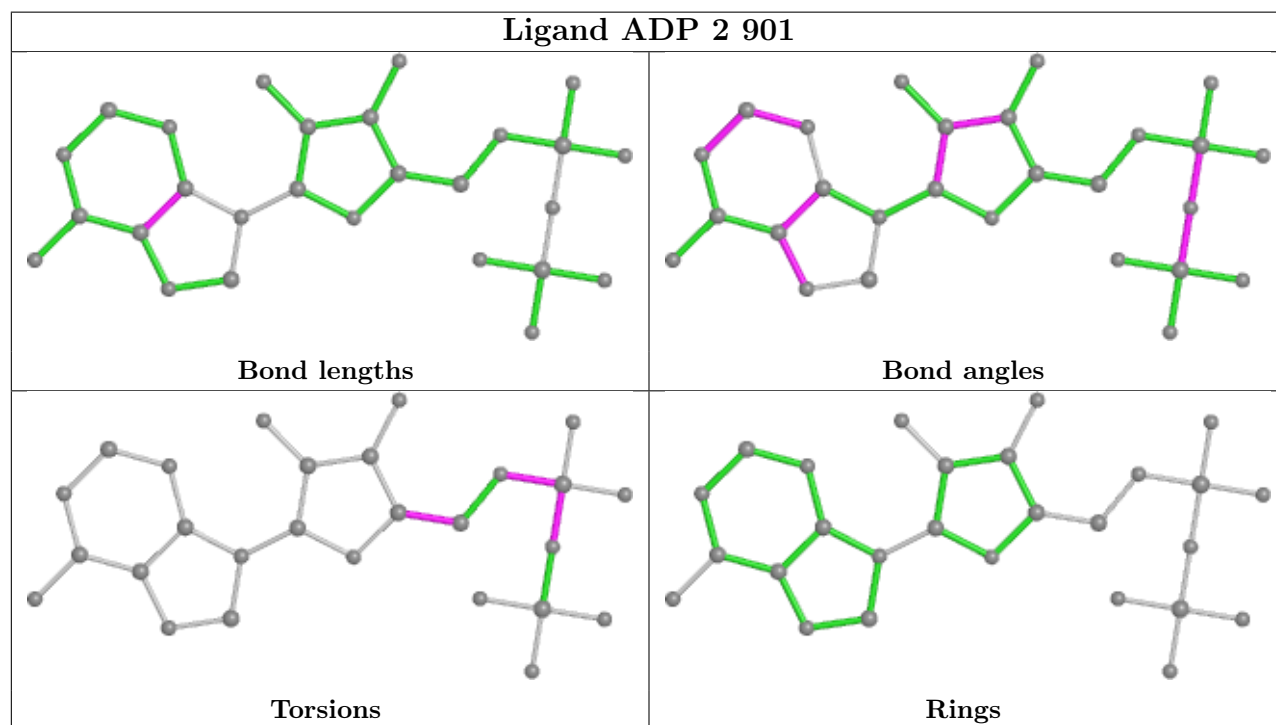
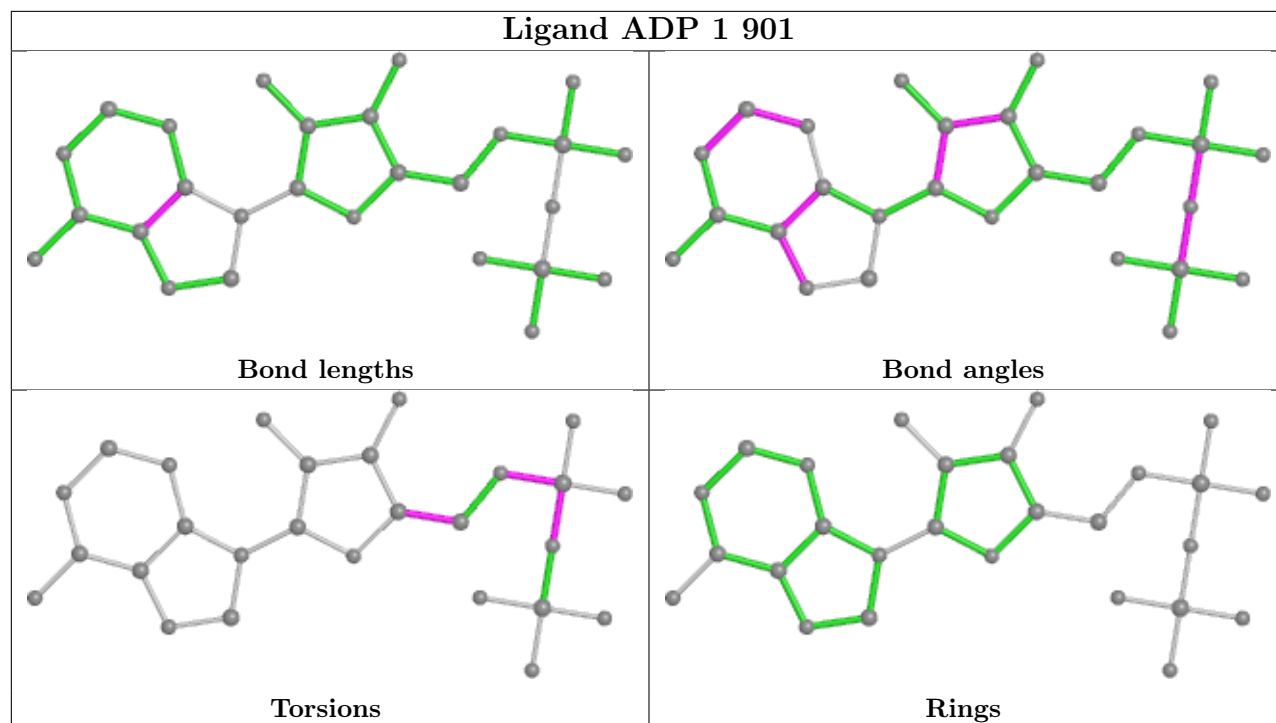


Ligand ADP 5 901

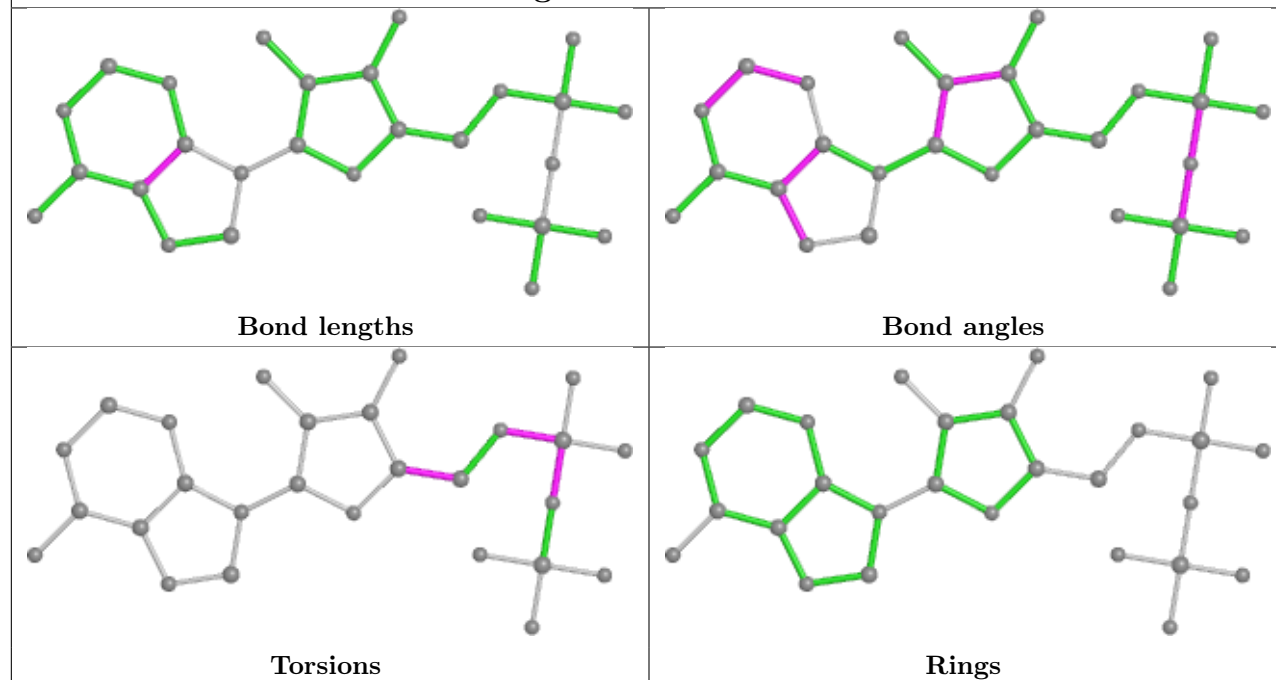


Ligand ADP 4 901

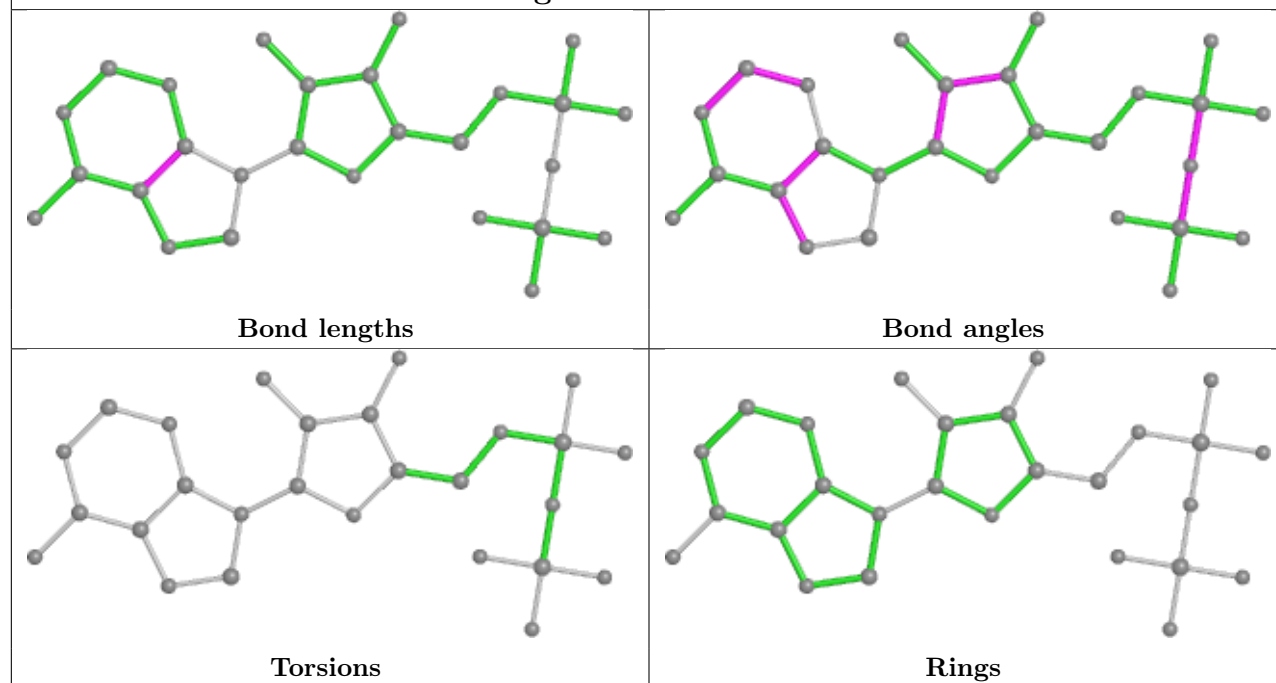


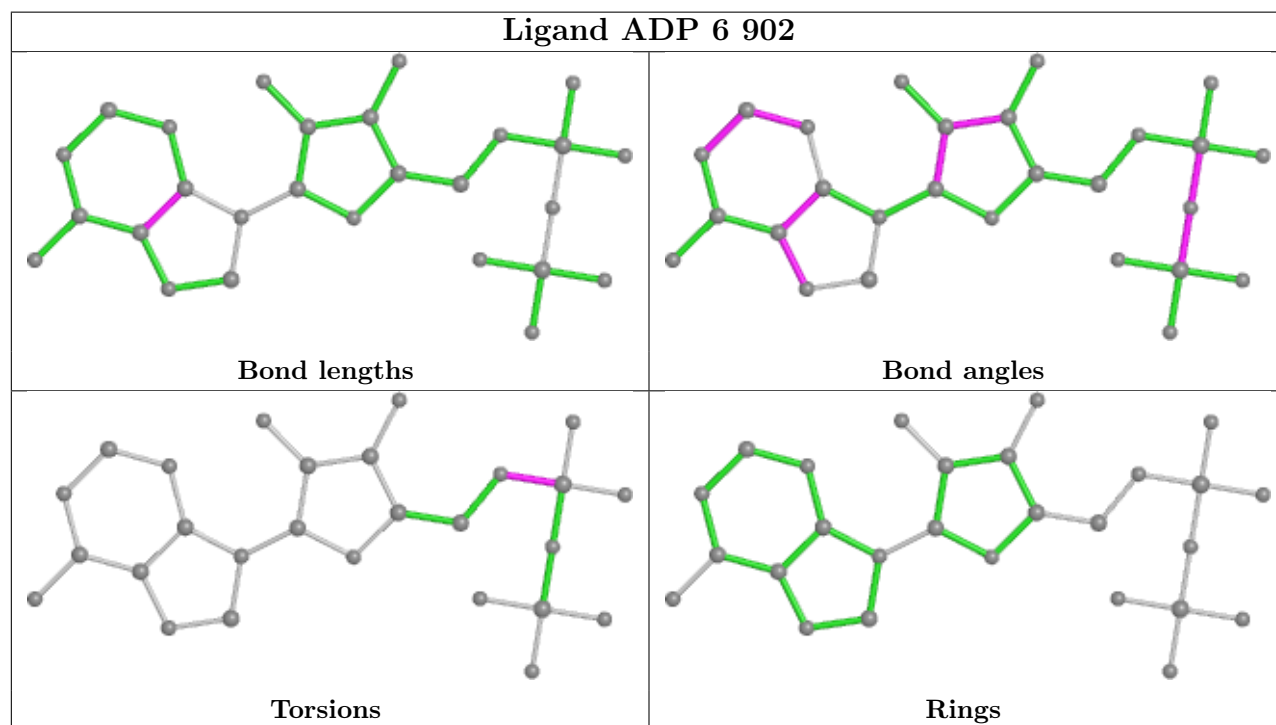
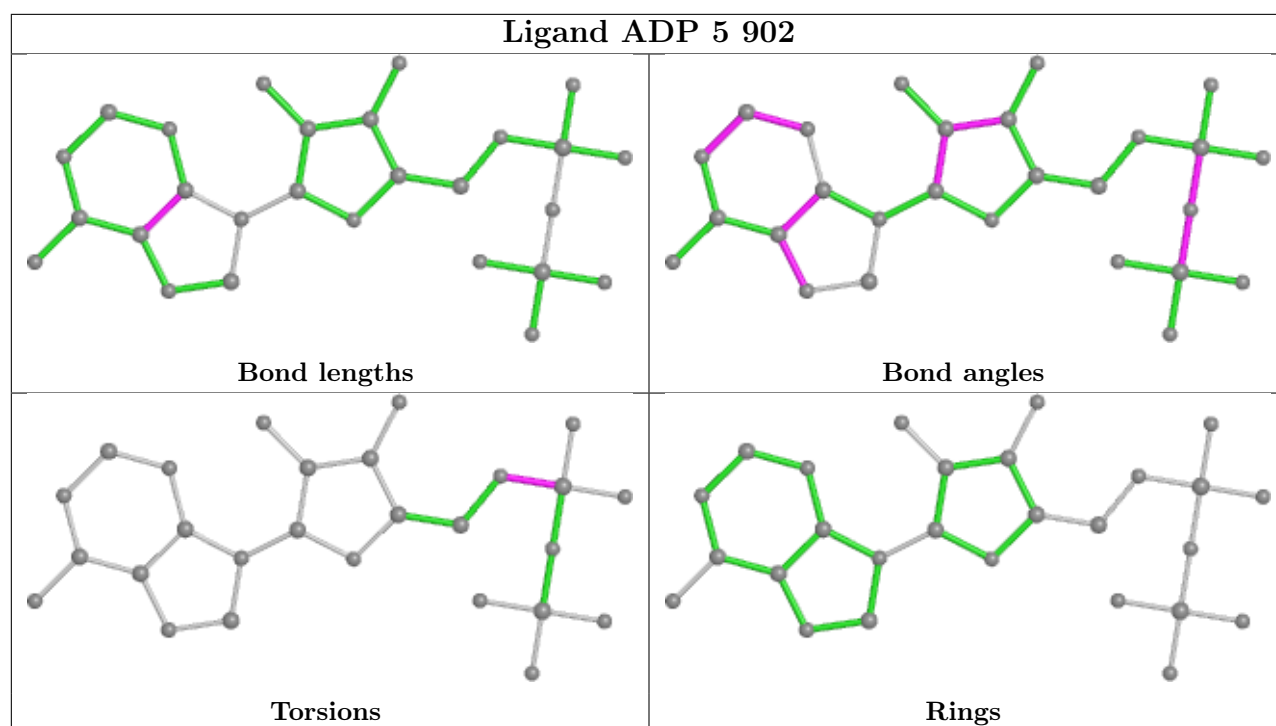


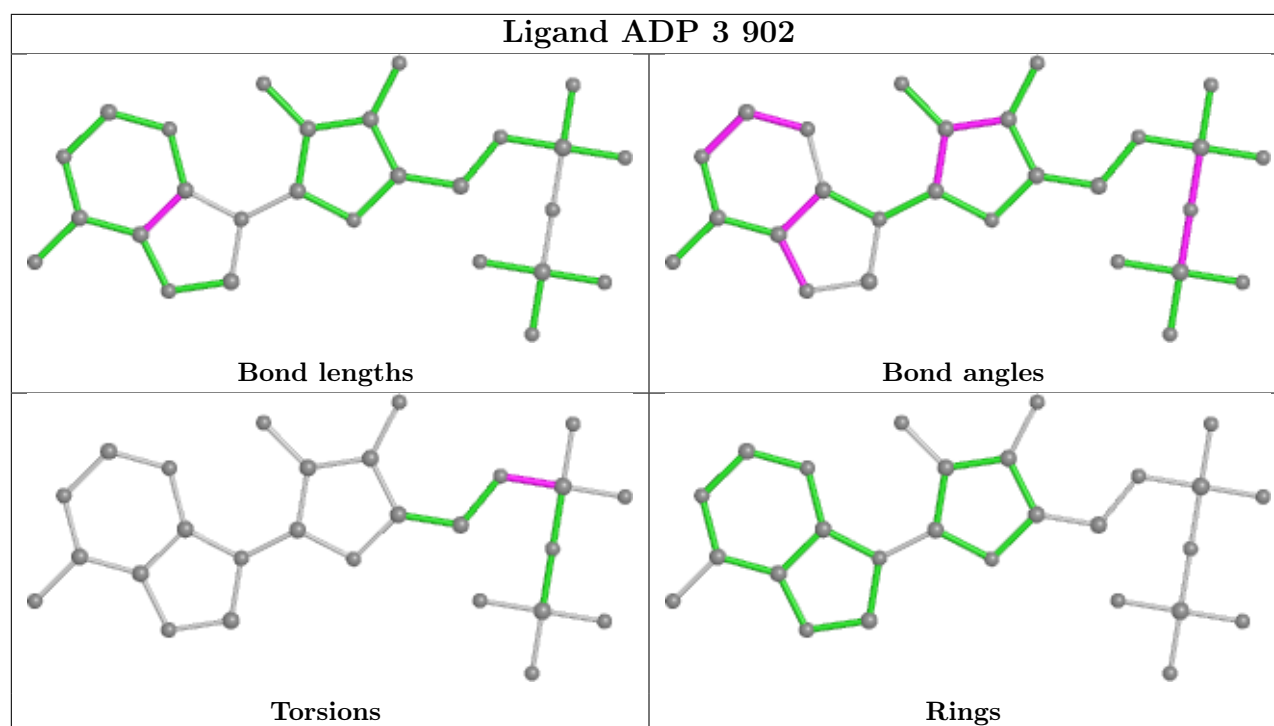
Ligand ADP 3 901



Ligand ADP 4 902







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 Map visualisation

This section contains visualisations of the EMDB entry EMD-7782. These allow visual inspection of the internal detail of the map and identification of artifacts.

No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections

This section was not generated.

6.2 Central slices

This section was not generated.

6.3 Largest variance slices

This section was not generated.

6.4 Orthogonal surface views

This section was not generated.

6.5 Mask visualisation

This section was not generated. No masks/segmentation were deposited.

7 Map analysis ⓘ

This section contains the results of statistical analysis of the map.

7.1 Map-value distribution ⓘ

This section was not generated.

7.2 Volume estimate versus contour level ⓘ

This section was not generated.

7.3 Rotationally averaged power spectrum ⓘ

This section was not generated. The rotationally averaged power spectrum had issues being displayed.

8 Fourier-Shell correlation ⓘ

This section was not generated. No FSC curve or half-maps provided.

9 Map-model fit

This section was not generated.