



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 13, 2023 – 03:43 AM EDT

PDB ID : 4NB4
Title : Pantothenamide-bound Pantothenate kinase from *Staphylococcus aureus*
Authors : Hughes, S.J.; Antoshchenko, T.; Smil, D.; Park, H.W.
Deposited on : 2013-10-22
Resolution : 2.25 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.35.1
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.35.1

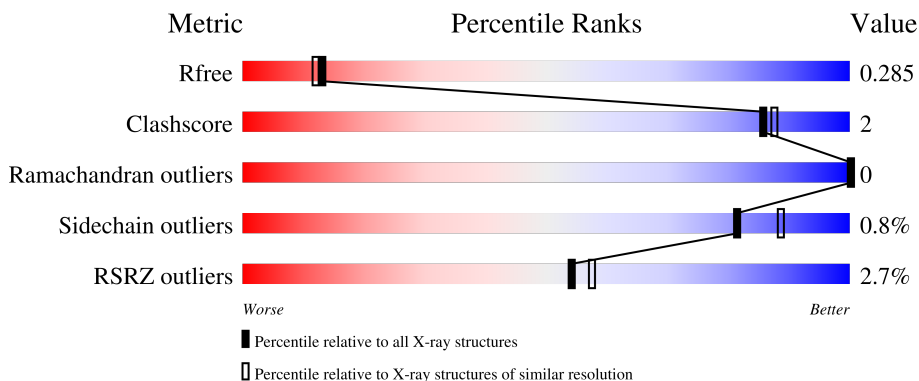
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 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))
R_{free}	130704	1377 (2.26-2.26)
Clashscore	141614	1487 (2.26-2.26)
Ramachandran outliers	138981	1449 (2.26-2.26)
Sidechain outliers	138945	1450 (2.26-2.26)
RSRZ outliers	127900	1356 (2.26-2.26)

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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	285	
1	B	285	
1	C	285	
1	D	285	
1	E	285	

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Mol	Chain	Length	Quality of chain
1	F	285	
1	G	285	
1	H	285	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 17249 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 Type II pantothenate kinase.

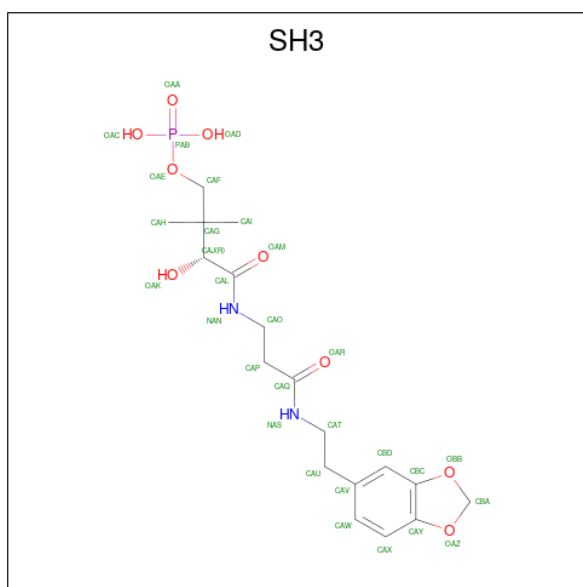
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	265	Total 2067	C 1308	N 358	O 395	S 6	0	3	0
1	B	265	Total 2061	C 1302	N 356	O 397	S 6	0	3	0
1	C	265	Total 2034	C 1289	N 350	O 389	S 6	0	0	0
1	D	265	Total 2053	C 1301	N 351	O 395	S 6	0	2	0
1	E	259	Total 1988	C 1258	N 342	O 382	S 6	0	1	0
1	F	262	Total 2016	C 1276	N 348	O 386	S 6	0	1	0
1	G	265	Total 2052	C 1299	N 354	O 393	S 6	0	2	0
1	H	267	Total 2053	C 1299	N 352	O 395	S 7	0	1	0

- Molecule 2 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		
2	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
2	B	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
2	C	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
2	D	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
2	E	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
2	F	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
2	G	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
2	H	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

- Molecule 3 is N-[2-(1,3-benzodioxol-5-yl)ethyl]-N 3 -[(2R)-2-hydroxy-3,3-dimethyl-4-(phosphonoxy)butanoyl]-beta-alaninamide (three-letter code: SH3) (formula: C₁₈H₂₇N₂O₉P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
3	A	1	Total	C	N	O	P	0	0
			30	18	2	9	1		
3	A	1	Total	C	N	O	P	0	0
			30	18	2	9	1		
3	C	1	Total	C	N	O	P	0	0
			30	18	2	9	1		
3	C	1	Total	C	N	O	P	0	0
			30	18	2	9	1		
3	F	1	Total	C	N	O	P	0	0
			30	18	2	9	1		
3	F	1	Total	C	N	O	P	0	0
			30	18	2	9	1		
3	G	1	Total	C	N	O	P	0	0
			30	18	2	9	1		
3	H	1	Total	C	N	O	P	0	0
			30	18	2	9	1		

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Mg	0	0
			1	1		
4	B	1	Total	Mg	0	0
			1	1		
4	C	1	Total	Mg	0	0
			1	1		
4	D	1	Total	Mg	0	0
			1	1		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	E	1	Total Mg 1 1	0	0
4	F	1	Total Mg 1 1	0	0
4	G	1	Total Mg 1 1	0	0
4	H	1	Total Mg 1 1	0	0

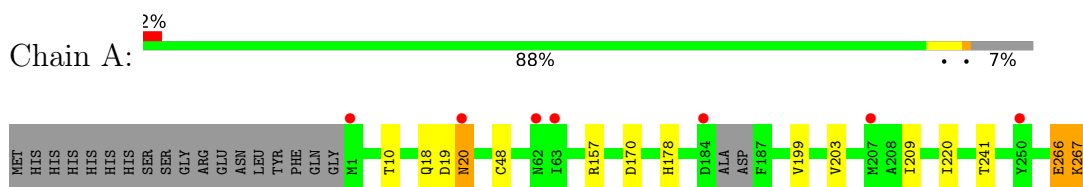
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	58	Total O 58 58	0	0
5	B	83	Total O 83 83	0	0
5	C	58	Total O 58 58	0	0
5	D	42	Total O 42 42	0	0
5	E	40	Total O 40 40	0	0
5	F	31	Total O 31 31	0	0
5	G	71	Total O 71 71	0	0
5	H	78	Total O 78 78	0	0

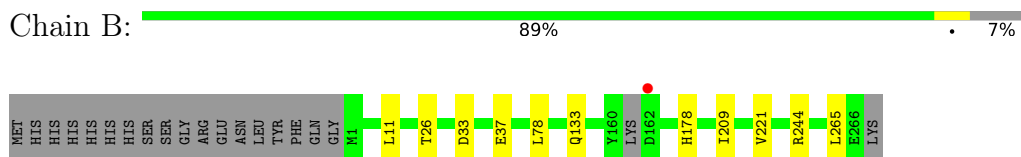
3 Residue-property plots [i](#)

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 and electron density. 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. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). 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.

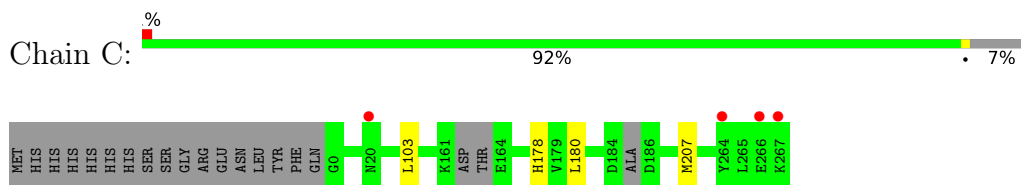
- Molecule 1: Type II pantothenate kinase



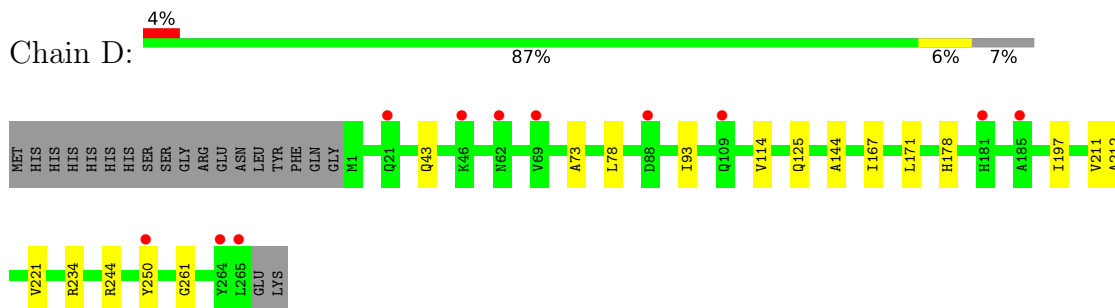
- Molecule 1: Type II pantothenate kinase



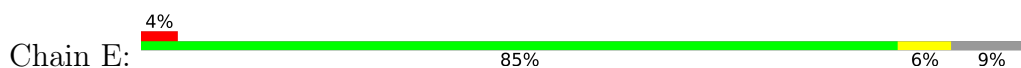
- Molecule 1: Type II pantothenate kinase

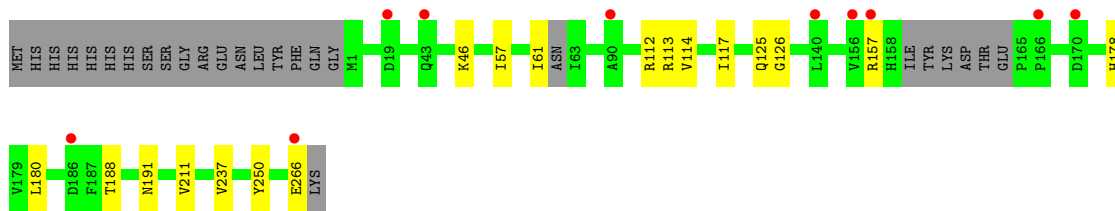


- Molecule 1: Type II pantothenate kinase

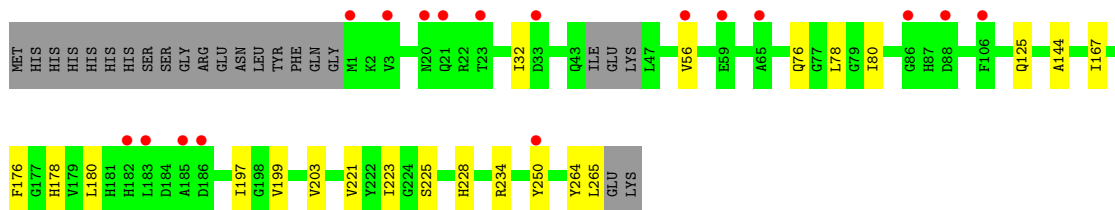
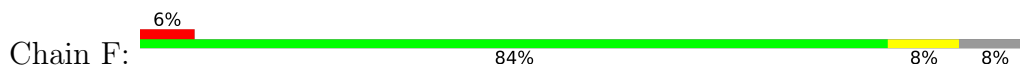


- Molecule 1: Type II pantothenate kinase

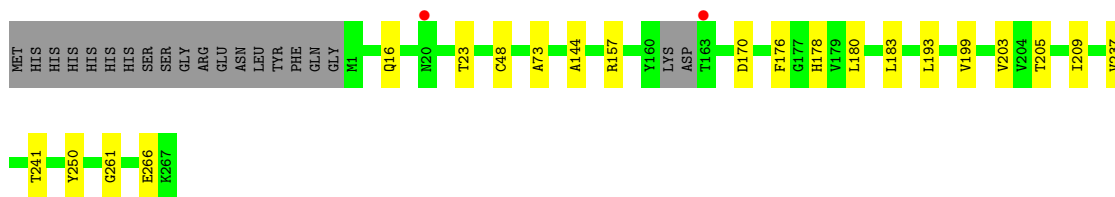
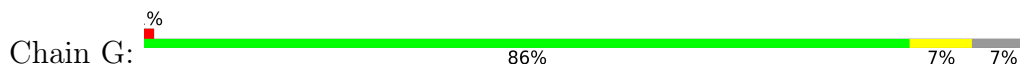




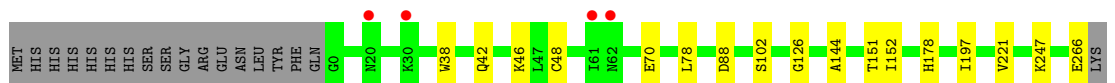
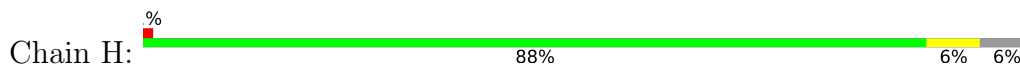
● Molecule 1: Type II pantothenate kinase



● Molecule 1: Type II pantothenate kinase



● Molecule 1: Type II pantothenate kinase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	103.01Å 70.74Å 143.12Å 90.00° 90.64° 90.00°	Depositor
Resolution (Å)	39.91 – 2.25 39.88 – 2.25	Depositor EDS
% Data completeness (in resolution range)	99.3 (39.91-2.25) 96.3 (39.88-2.25)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.83 (at 2.24Å)	Xtrriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.230 , 0.284 0.233 , 0.285	Depositor DCC
R_{free} test set	4896 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	20.6	Xtrriage
Anisotropy	0.049	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 5.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.41$, $\langle L^2 \rangle = 0.23$	Xtrriage
Estimated twinning fraction	0.185 for h,-k,-l	Xtrriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	17249	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.53% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

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: SH3, 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	0.35	1/2102 (0.0%)	0.51	1/2844 (0.0%)
1	B	0.31	0/2096	0.50	0/2839
1	C	0.32	0/2067	0.49	0/2795
1	D	0.33	0/2089	0.49	0/2832
1	E	0.31	0/2020	0.48	0/2733
1	F	0.30	0/2051	0.48	0/2779
1	G	0.33	0/2087	0.51	0/2826
1	H	0.32	0/2088	0.51	0/2828
All	All	0.32	1/16600 (0.0%)	0.50	1/22476 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	266	GLU	CD-OE1	-5.27	1.19	1.25

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	20	ASN	CB-CA-C	-6.19	98.02	110.40

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	2067	0	2053	9	0
1	B	2061	0	2034	6	0
1	C	2034	0	2025	2	0
1	D	2053	0	2035	11	0
1	E	1988	0	1973	13	0
1	F	2016	0	1996	16	0
1	G	2052	0	2035	11	0
1	H	2053	0	2038	14	0
2	A	27	0	12	1	0
2	B	27	0	12	0	0
2	C	27	0	12	0	0
2	D	27	0	12	0	0
2	E	27	0	12	0	0
2	F	27	0	12	0	0
2	G	27	0	12	0	0
2	H	27	0	12	0	0
3	A	60	0	52	0	0
3	C	60	0	52	0	0
3	F	60	0	53	0	0
3	G	30	0	26	0	0
3	H	30	0	26	1	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
4	E	1	0	0	0	0
4	F	1	0	0	0	0
4	G	1	0	0	0	0
4	H	1	0	0	0	0
5	A	58	0	0	0	0
5	B	83	0	0	0	0
5	C	58	0	0	0	0
5	D	42	0	0	0	0
5	E	40	0	0	0	0
5	F	31	0	0	0	0
5	G	71	0	0	0	0
5	H	78	0	0	2	0
All	All	17249	0	16494	73	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 2.

All (73) 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:H:46:LYS:HG2	1:H:266:GLU:OE2	1.47	1.13
1:E:46:LYS:HE3	1:E:266:GLU:OE2	1.62	0.99
1:H:46:LYS:HG2	1:H:266:GLU:CD	1.83	0.97
1:E:46:LYS:CE	1:E:266:GLU:OE2	2.36	0.73
1:D:93:ILE:HD13	1:D:211:VAL:HG13	1.75	0.69
1:A:48:CYS:SG	1:A:266:GLU:HG3	2.35	0.67
1:G:205:THR:HG22	1:G:209:ILE:HD13	1.81	0.62
1:A:209:ILE:HD11	1:A:241:THR:HA	1.83	0.60
1:H:48:CYS:SG	1:H:266:GLU:HB2	2.42	0.59
1:G:157:ARG:NH2	1:G:170:ASP:OD1	2.37	0.58
1:E:117:ILE:HD11	1:F:203:VAL:HG22	1.88	0.56
1:D:114:VAL:CG2	1:D:211:VAL:HG21	2.35	0.55
1:H:46:LYS:CG	1:H:266:GLU:OE2	2.39	0.55
1:E:113:ARG:HD3	1:F:167:ILE:HD11	1.89	0.54
1:F:80:ILE:HD11	1:F:264:TYR:CD2	2.42	0.54
1:E:114:VAL:CG2	1:E:211:VAL:HG21	2.38	0.53
1:F:78:LEU:HD13	1:F:223:ILE:HG21	1.93	0.51
1:E:125:GLN:HB3	1:F:180:LEU:HD11	1.93	0.50
1:H:151:THR:HG23	1:H:152:ILE:HG13	1.94	0.50
1:E:126:GLY:HA3	1:F:176:PHE:O	2.12	0.49
1:D:93:ILE:HD12	1:D:212:ALA:HA	1.93	0.49
1:F:199:VAL:O	1:F:203:VAL:HG23	2.12	0.49
1:G:180:LEU:HD23	1:G:183:LEU:HD21	1.94	0.49
1:H:38:TRP:O	1:H:42:GLN:HG2	2.14	0.48
1:E:114:VAL:HG21	1:E:211:VAL:HG21	1.96	0.48
1:D:167:ILE:HG22	1:D:171:LEU:HD12	1.97	0.47
1:H:70:GLU:HG2	5:H:439:HOH:O	2.15	0.47
1:C:180:LEU:HD11	1:D:125:GLN:HB3	1.97	0.47
1:E:113:ARG:CD	1:F:167:ILE:HD11	2.44	0.47
1:G:73:ALA:O	1:G:261:GLY:HA3	2.15	0.47
1:G:199:VAL:O	1:G:203:VAL:HG23	2.14	0.47
1:A:209:ILE:HD13	1:A:220:ILE:CD1	2.44	0.46
1:G:16:GLN:HB2	1:G:23:THR:HB	1.97	0.46
1:B:209:ILE:HD13	1:B:244:ARG:HG3	1.98	0.46
1:H:46:LYS:HE2	1:H:266:GLU:OE1	2.16	0.46
1:D:78:LEU:HD11	1:D:221:VAL:HG11	1.98	0.46
1:G:176:PHE:O	1:H:126:GLY:HA3	2.16	0.46
1:G:237:VAL:CG2	1:G:250:TYR:OH	2.64	0.46
1:H:151:THR:HG21	5:H:435:HOH:O	2.15	0.46
1:G:209:ILE:HD11	1:G:241:THR:HA	1.99	0.45
1:E:180:LEU:HD11	1:F:125:GLN:HB3	1.98	0.45
1:A:18:GLN:O	1:A:19:ASP:HB2	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:234:ARG:NH1	1:D:250[B]:TYR:CE1	2.85	0.44
1:F:265:LEU:C	1:F:265:LEU:HD12	2.38	0.44
1:A:19:ASP:O	1:A:20:ASN:HB2	2.18	0.44
1:A:266:GLU:O	1:A:267:LYS:HG2	2.18	0.43
1:B:11:LEU:HD22	1:B:26:THR:CG2	2.47	0.43
1:F:225:SER:HA	1:F:228[B]:HIS:CE1	2.53	0.43
1:F:32:ILE:HD13	1:F:56:VAL:HG11	2.01	0.43
1:D:167:ILE:CG2	1:D:171:LEU:HD12	2.49	0.43
1:D:73:ALA:O	1:D:261:GLY:HA3	2.19	0.43
1:G:48:CYS:SG	1:G:266:GLU:HB2	2.59	0.43
1:D:144:ALA:HB1	1:D:197:ILE:HD12	2.01	0.43
1:D:114:VAL:HG21	1:D:211:VAL:HG21	1.99	0.42
1:F:234:ARG:HD3	1:F:250:TYR:CE2	2.53	0.42
1:E:188:THR:CG2	1:E:191:ASN:CG	2.88	0.42
1:C:103:LEU:HD11	1:C:207:MET:HB3	2.01	0.42
1:F:78:LEU:HD11	1:F:221:VAL:HG11	2.02	0.42
1:E:237:VAL:HG23	1:E:250:TYR:OH	2.20	0.42
1:B:33:ASP:O	1:B:37:GLU:HG2	2.20	0.41
1:F:144:ALA:HB1	1:F:197:ILE:HD12	2.02	0.41
1:E:57:ILE:O	1:E:61:ILE:HB	2.20	0.41
1:A:10:THR:HG21	2:A:301:ADP:H3'	2.02	0.41
1:A:157[A]:ARG:NH2	1:A:170:ASP:OD1	2.53	0.41
1:G:144:ALA:HB2	1:G:193:LEU:HB3	2.03	0.41
1:H:102:SER:HB2	3:H:302:SH3:OAM	2.20	0.41
1:B:265:LEU:HD23	1:B:265:LEU:HA	1.87	0.41
1:H:144:ALA:HB1	1:H:197:ILE:HD12	2.03	0.41
1:B:78:LEU:HD11	1:B:221:VAL:HG11	2.03	0.40
1:A:199:VAL:O	1:A:203:VAL:HG23	2.21	0.40
1:F:76:GLN:HG2	1:F:264:TYR:CD1	2.56	0.40
1:H:78:LEU:HD11	1:H:221:VAL:HG11	2.04	0.40
1:B:133:GLN:HB3	1:H:247:LYS:HE2	2.02	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	264/285 (93%)	257 (97%)	7 (3%)	0	100	100
1	B	264/285 (93%)	258 (98%)	6 (2%)	0	100	100
1	C	259/285 (91%)	254 (98%)	5 (2%)	0	100	100
1	D	265/285 (93%)	259 (98%)	6 (2%)	0	100	100
1	E	254/285 (89%)	248 (98%)	6 (2%)	0	100	100
1	F	259/285 (91%)	247 (95%)	12 (5%)	0	100	100
1	G	263/285 (92%)	256 (97%)	7 (3%)	0	100	100
1	H	266/285 (93%)	262 (98%)	4 (2%)	0	100	100
All	All	2094/2280 (92%)	2041 (98%)	53 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	219/233 (94%)	217 (99%)	2 (1%)	78	86
1	B	218/233 (94%)	217 (100%)	1 (0%)	88	92
1	C	215/233 (92%)	214 (100%)	1 (0%)	88	92
1	D	217/233 (93%)	214 (99%)	3 (1%)	67	76
1	E	210/233 (90%)	207 (99%)	3 (1%)	67	76
1	F	213/233 (91%)	212 (100%)	1 (0%)	88	92
1	G	217/233 (93%)	216 (100%)	1 (0%)	88	92
1	H	217/233 (93%)	215 (99%)	2 (1%)	78	86
All	All	1726/1864 (93%)	1712 (99%)	14 (1%)	81	88

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	178	HIS
1	A	267	LYS
1	B	178	HIS
1	C	178	HIS
1	D	43	GLN
1	D	178	HIS
1	D	244	ARG
1	E	112	ARG
1	E	157	ARG
1	E	178	HIS
1	F	178	HIS
1	G	178	HIS
1	H	88	ASP
1	H	178	HIS

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

Mol	Chain	Res	Type
1	A	125	GLN
1	C	158	HIS
1	D	43	GLN
1	D	76	GLN
1	D	139	GLN
1	E	125	GLN
1	F	42	GLN
1	F	109	GLN
1	F	150	ASN
1	G	76	GLN
1	G	158	HIS
1	H	158	HIS

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 24 ligands modelled in this entry, 8 are monoatomic - leaving 16 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
3	SH3	H	302	4	28,31,31	1.32	5 (17%)	42,44,44	1.74	6 (14%)
3	SH3	F	302	4	28,31,31	1.31	5 (17%)	42,44,44	1.09	4 (9%)
3	SH3	A	302	4	28,31,31	1.33	5 (17%)	42,44,44	1.01	4 (9%)
2	ADP	A	301	4	24,29,29	1.00	2 (8%)	29,45,45	1.28	3 (10%)
2	ADP	C	301	4	24,29,29	0.99	2 (8%)	29,45,45	1.34	3 (10%)
2	ADP	D	301	4	24,29,29	1.08	3 (12%)	29,45,45	1.26	2 (6%)
2	ADP	F	301	4	24,29,29	1.03	3 (12%)	29,45,45	1.24	4 (13%)
2	ADP	E	301	4	24,29,29	1.05	2 (8%)	29,45,45	1.23	1 (3%)
2	ADP	G	301	4	24,29,29	1.01	2 (8%)	29,45,45	1.36	4 (13%)
3	SH3	A	303	4	28,31,31	1.39	7 (25%)	42,44,44	1.34	6 (14%)
3	SH3	C	302	4	28,31,31	1.28	5 (17%)	42,44,44	1.29	5 (11%)
2	ADP	H	301	4	24,29,29	1.04	2 (8%)	29,45,45	1.28	4 (13%)
3	SH3	F	303	4	28,31,31	1.33	5 (17%)	42,44,44	1.16	4 (9%)
3	SH3	C	303	4	28,31,31	1.35	4 (14%)	42,44,44	1.12	2 (4%)
2	ADP	B	301	4	24,29,29	0.99	2 (8%)	29,45,45	1.24	1 (3%)
3	SH3	G	302	4	28,31,31	1.26	5 (17%)	42,44,44	1.21	5 (11%)

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
3	SH3	H	302	4	-	3/29/35/35	0/2/2/2
3	SH3	F	302	4	-	1/29/35/35	0/2/2/2
3	SH3	A	302	4	-	1/29/35/35	0/2/2/2
2	ADP	A	301	4	-	1/12/32/32	0/3/3/3
2	ADP	C	301	4	-	0/12/32/32	0/3/3/3
2	ADP	D	301	4	-	0/12/32/32	0/3/3/3
2	ADP	F	301	4	-	2/12/32/32	0/3/3/3
2	ADP	E	301	4	-	0/12/32/32	0/3/3/3
2	ADP	G	301	4	-	0/12/32/32	0/3/3/3
3	SH3	A	303	4	-	1/29/35/35	0/2/2/2
3	SH3	C	302	4	-	2/29/35/35	0/2/2/2
2	ADP	H	301	4	-	0/12/32/32	0/3/3/3
3	SH3	F	303	4	-	1/29/35/35	0/2/2/2
3	SH3	C	303	4	-	1/29/35/35	0/2/2/2
2	ADP	B	301	4	-	0/12/32/32	0/3/3/3
3	SH3	G	302	4	-	1/29/35/35	0/2/2/2

All (59) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	F	302	SH3	CAU-CAV	-3.56	1.41	1.51
3	A	303	SH3	CAU-CAV	-3.54	1.41	1.51
3	F	303	SH3	CAU-CAV	-3.45	1.41	1.51
3	A	302	SH3	CAU-CAV	-3.44	1.41	1.51
3	C	303	SH3	CAU-CAV	-3.38	1.41	1.51
3	C	302	SH3	CAU-CAV	-3.32	1.42	1.51
3	H	302	SH3	CAU-CAV	-3.21	1.42	1.51
3	G	302	SH3	CAU-CAV	-3.18	1.42	1.51
3	A	302	SH3	CAX-CAY	-2.78	1.33	1.39
3	G	302	SH3	CAX-CAY	-2.77	1.33	1.39
3	A	302	SH3	OAZ-CBA	2.71	1.48	1.43
3	F	303	SH3	CAX-CAY	-2.68	1.34	1.39
3	C	303	SH3	CAX-CAY	-2.66	1.34	1.39
3	C	302	SH3	CAX-CAY	-2.63	1.34	1.39
3	F	302	SH3	CAX-CAY	-2.61	1.34	1.39
2	D	301	ADP	C5-C4	2.60	1.47	1.40
2	H	301	ADP	C5-C4	2.59	1.47	1.40
3	C	303	SH3	CBD-CBC	-2.58	1.34	1.38
3	A	303	SH3	CBD-CBC	-2.56	1.34	1.38
2	C	301	ADP	C5-C4	2.53	1.47	1.40
2	E	301	ADP	C5-C4	2.53	1.47	1.40
3	H	302	SH3	CAX-CAY	-2.52	1.34	1.39
2	F	301	ADP	C5-C4	2.52	1.47	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	302	SH3	CBD-CBC	-2.50	1.34	1.38
2	D	301	ADP	O4'-C1'	2.49	1.44	1.41
2	A	301	ADP	C5-C4	2.45	1.47	1.40
2	G	301	ADP	C5-C4	2.45	1.47	1.40
2	E	301	ADP	C2-N3	2.44	1.36	1.32
3	A	303	SH3	CAY-CBC	-2.44	1.32	1.39
3	G	302	SH3	CBD-CBC	-2.42	1.34	1.38
3	A	303	SH3	CAX-CAY	-2.41	1.34	1.39
2	B	301	ADP	C5-C4	2.37	1.47	1.40
3	A	303	SH3	OAZ-CBA	2.33	1.47	1.43
3	H	302	SH3	CAP-CAQ	2.30	1.55	1.51
2	H	301	ADP	O4'-C1'	2.27	1.44	1.41
3	F	303	SH3	CAY-CBC	-2.26	1.33	1.39
2	F	301	ADP	C2-N3	2.26	1.35	1.32
3	C	303	SH3	CAY-CBC	-2.26	1.33	1.39
2	D	301	ADP	C2-N3	2.25	1.35	1.32
3	C	302	SH3	CAY-CBC	-2.25	1.33	1.39
3	H	302	SH3	CAY-CBC	-2.22	1.33	1.39
3	G	302	SH3	OAZ-CBA	2.21	1.47	1.43
3	F	303	SH3	CBD-CBC	-2.21	1.34	1.38
3	F	303	SH3	OAZ-CBA	2.21	1.47	1.43
2	B	301	ADP	C2-N3	2.20	1.35	1.32
2	G	301	ADP	O4'-C1'	2.20	1.44	1.41
3	H	302	SH3	CBD-CBC	-2.19	1.34	1.38
3	F	302	SH3	CAY-CBC	-2.18	1.33	1.39
3	G	302	SH3	CAY-CBC	-2.15	1.33	1.39
3	C	302	SH3	CBD-CBC	-2.13	1.34	1.38
3	A	303	SH3	PAB-OAD	2.12	1.63	1.54
3	F	302	SH3	CBD-CBC	-2.11	1.34	1.38
3	C	302	SH3	PAB-OAC	2.10	1.62	1.54
2	F	301	ADP	O4'-C1'	2.10	1.44	1.41
3	F	302	SH3	PAB-OAC	2.08	1.62	1.54
3	A	303	SH3	PAB-OAC	2.03	1.62	1.54
2	A	301	ADP	C2-N3	2.02	1.35	1.32
2	C	301	ADP	C2-N3	2.01	1.35	1.32
3	A	302	SH3	CAY-CBC	-2.00	1.34	1.39

All (58) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	H	302	SH3	CAO-NAN-CAL	6.29	133.81	122.59
3	A	303	SH3	OAZ-CBA-OBB	-4.75	100.49	108.08

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	H	302	SH3	CAJ-CAL-NAN	4.60	125.75	116.58
3	C	303	SH3	OAZ-CBA-OBB	-4.60	100.73	108.08
3	C	302	SH3	OAZ-CBA-OBB	-4.58	100.76	108.08
2	D	301	ADP	N3-C2-N1	-3.92	122.55	128.68
3	H	302	SH3	OAZ-CBA-OBB	-3.85	101.92	108.08
3	F	303	SH3	OAZ-CBA-OBB	-3.81	101.99	108.08
2	H	301	ADP	N3-C2-N1	-3.81	122.73	128.68
2	A	301	ADP	N3-C2-N1	-3.71	122.89	128.68
2	E	301	ADP	N3-C2-N1	-3.69	122.92	128.68
3	F	302	SH3	OAZ-CBA-OBB	-3.66	102.23	108.08
2	B	301	ADP	N3-C2-N1	-3.60	123.05	128.68
2	C	301	ADP	N3-C2-N1	-3.55	123.12	128.68
2	G	301	ADP	N3-C2-N1	-3.52	123.17	128.68
3	G	302	SH3	OAZ-CBA-OBB	-3.45	102.56	108.08
2	F	301	ADP	N3-C2-N1	-3.36	123.43	128.68
3	H	302	SH3	CAP-CAO-NAN	-3.33	105.18	111.90
3	C	302	SH3	OAE-PAB-OAA	3.04	115.01	106.47
3	G	302	SH3	CAO-CAP-CAQ	-3.02	107.33	112.36
3	H	302	SH3	OAM-CAL-NAN	-2.98	116.59	122.99
3	A	303	SH3	OAE-PAB-OAA	2.92	114.68	106.47
3	C	302	SH3	CAO-CAP-CAQ	-2.75	107.78	112.36
3	A	302	SH3	OAZ-CBA-OBB	-2.69	103.78	108.08
2	D	301	ADP	C4-C5-N7	-2.61	106.68	109.40
3	F	302	SH3	CAU-CAT-NAS	-2.57	104.22	111.99
3	F	303	SH3	CAU-CAT-NAS	-2.55	104.28	111.99
3	C	303	SH3	OAE-PAB-OAA	2.54	113.61	106.47
3	G	302	SH3	CAT-CAU-CAV	-2.53	107.01	112.87
2	C	301	ADP	PA-O3A-PB	-2.51	124.20	132.83
3	G	302	SH3	CAU-CAT-NAS	-2.48	104.51	111.99
3	A	302	SH3	CAT-CAU-CAV	-2.47	107.14	112.87
2	F	301	ADP	C4-C5-N7	-2.46	106.84	109.40
3	A	302	SH3	OAE-PAB-OAA	2.45	113.33	106.47
3	F	303	SH3	OAE-PAB-OAA	2.42	113.27	106.47
3	A	303	SH3	CAO-CAP-CAQ	-2.39	108.37	112.36
3	F	302	SH3	OAE-PAB-OAA	2.38	113.14	106.47
2	A	301	ADP	C4-C5-N7	-2.36	106.94	109.40
3	G	302	SH3	OAE-PAB-OAA	2.35	113.08	106.47
2	H	301	ADP	C4-C5-N7	-2.34	106.96	109.40
3	A	303	SH3	CAT-CAU-CAV	-2.28	107.57	112.87
2	G	301	ADP	O3B-PB-O2B	2.27	116.31	107.64
3	C	302	SH3	CAT-CAU-CAV	-2.26	107.62	112.87
3	F	302	SH3	CAO-CAP-CAQ	-2.26	108.60	112.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	H	302	SH3	CAO-CAP-CAQ	2.23	116.07	112.36
2	F	301	ADP	PA-O3A-PB	-2.16	125.41	132.83
2	G	301	ADP	N6-C6-N1	2.14	123.02	118.57
2	F	301	ADP	O3B-PB-O2B	2.13	115.77	107.64
3	C	302	SH3	CAU-CAT-NAS	-2.10	105.64	111.99
2	A	301	ADP	C1'-N9-C4	2.10	130.33	126.64
2	C	301	ADP	C4-C5-N7	-2.09	107.22	109.40
2	H	301	ADP	O3B-PB-O2B	2.08	115.59	107.64
2	H	301	ADP	C3'-C2'-C1'	2.07	104.09	100.98
3	F	303	SH3	CAO-CAP-CAQ	-2.06	108.93	112.36
3	A	303	SH3	OAZ-CAY-CAX	2.05	131.55	127.81
3	A	303	SH3	CAU-CAT-NAS	-2.04	105.83	111.99
3	A	302	SH3	CAU-CAT-NAS	-2.02	105.89	111.99
2	G	301	ADP	C4-C5-N7	-2.01	107.31	109.40

There are no chirality outliers.

All (14) torsion outliers are listed below:

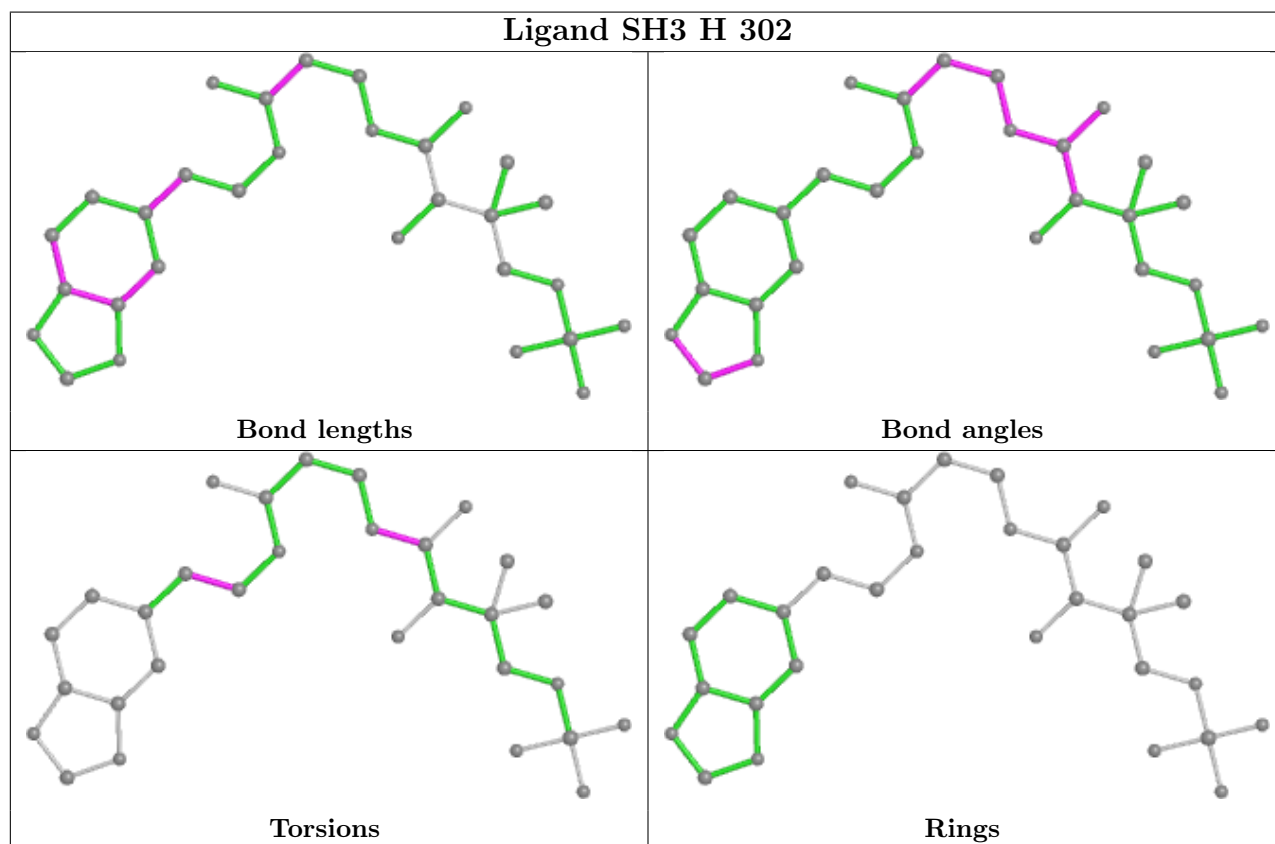
Mol	Chain	Res	Type	Atoms
3	A	302	SH3	NAS-CAT-CAU-CAV
3	A	303	SH3	NAS-CAT-CAU-CAV
3	C	302	SH3	NAS-CAT-CAU-CAV
3	C	303	SH3	NAS-CAT-CAU-CAV
3	F	302	SH3	NAS-CAT-CAU-CAV
3	F	303	SH3	NAS-CAT-CAU-CAV
3	G	302	SH3	NAS-CAT-CAU-CAV
3	H	302	SH3	CAJ-CAL-NAN-CAO
3	H	302	SH3	NAS-CAT-CAU-CAV
3	H	302	SH3	OAM-CAL-NAN-CAO
2	F	301	ADP	PB-O3A-PA-O2A
3	C	302	SH3	CAF-OAE-PAB-OAA
2	A	301	ADP	PB-O3A-PA-O1A
2	F	301	ADP	PB-O3A-PA-O1A

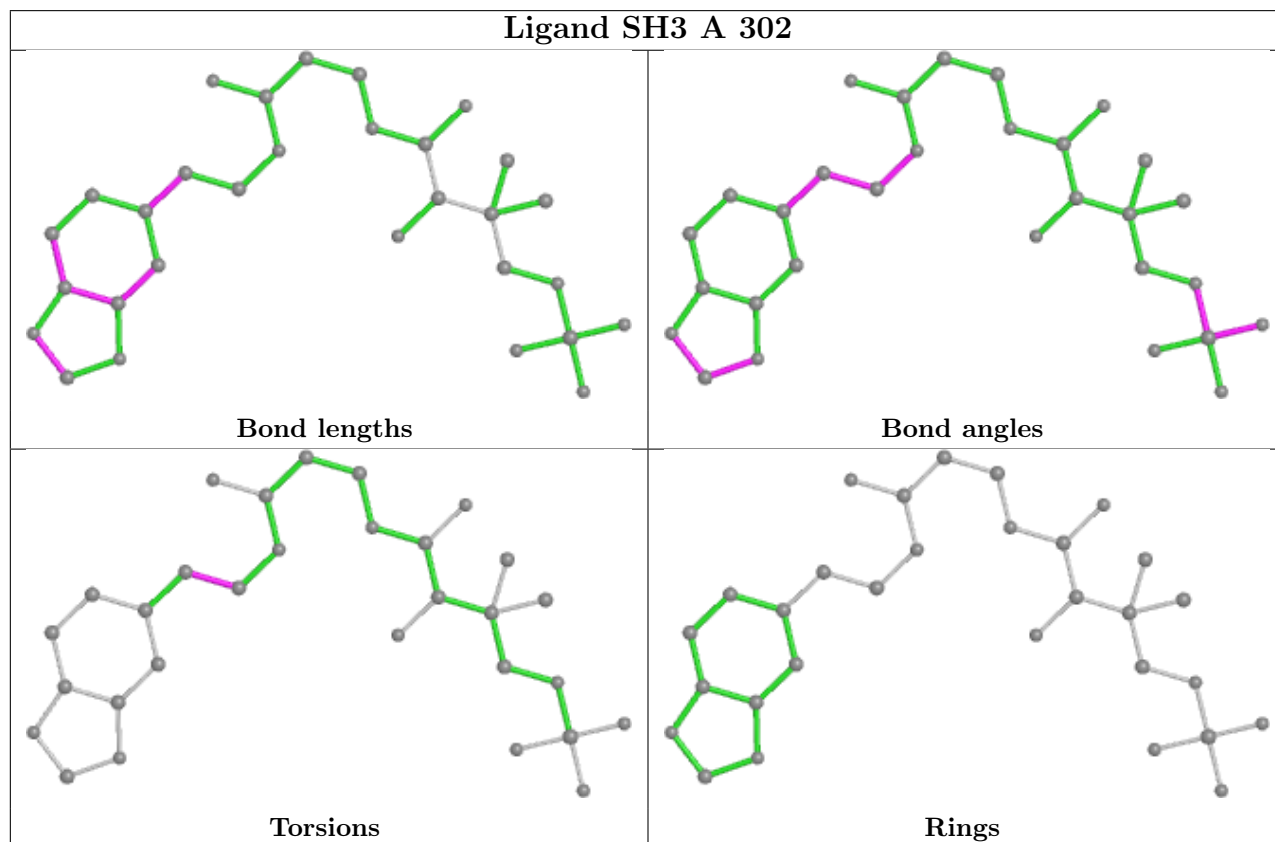
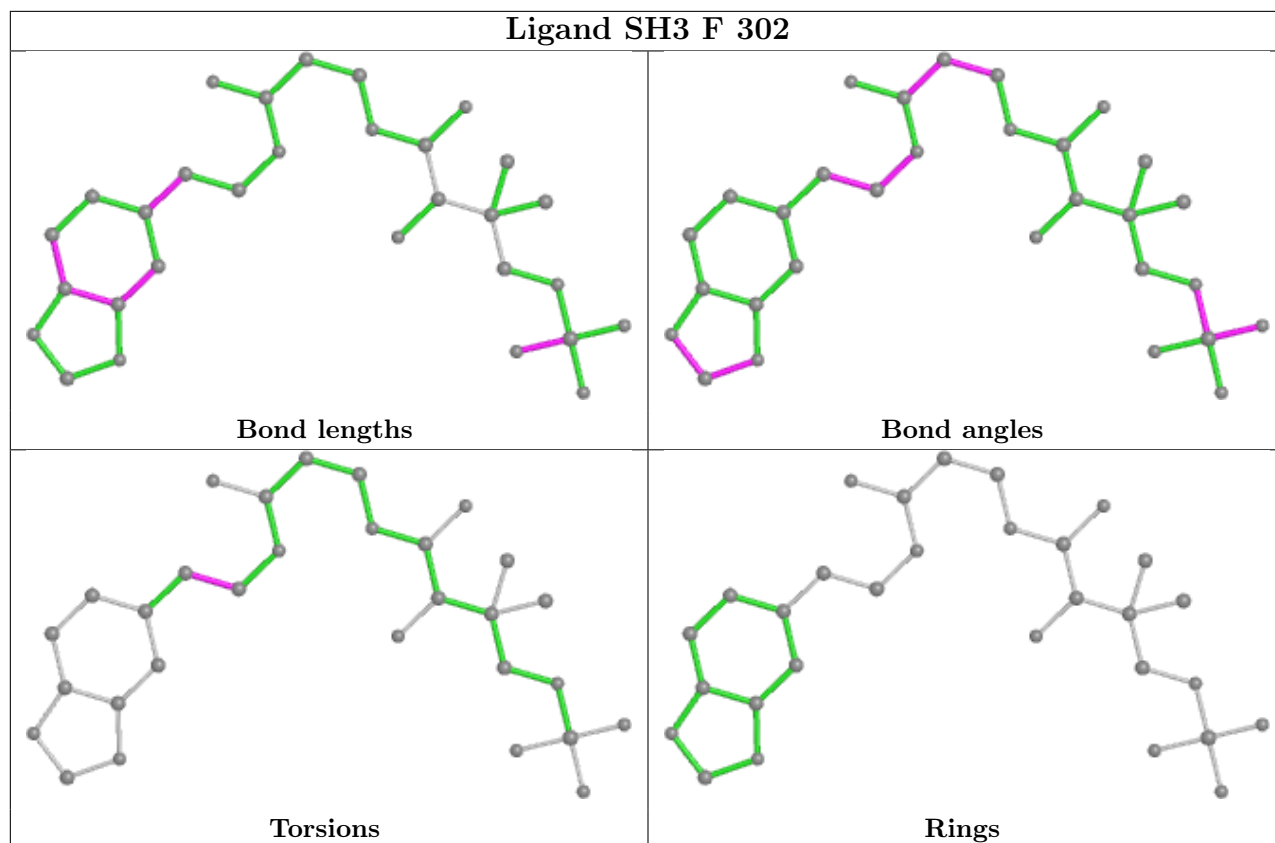
There are no ring outliers.

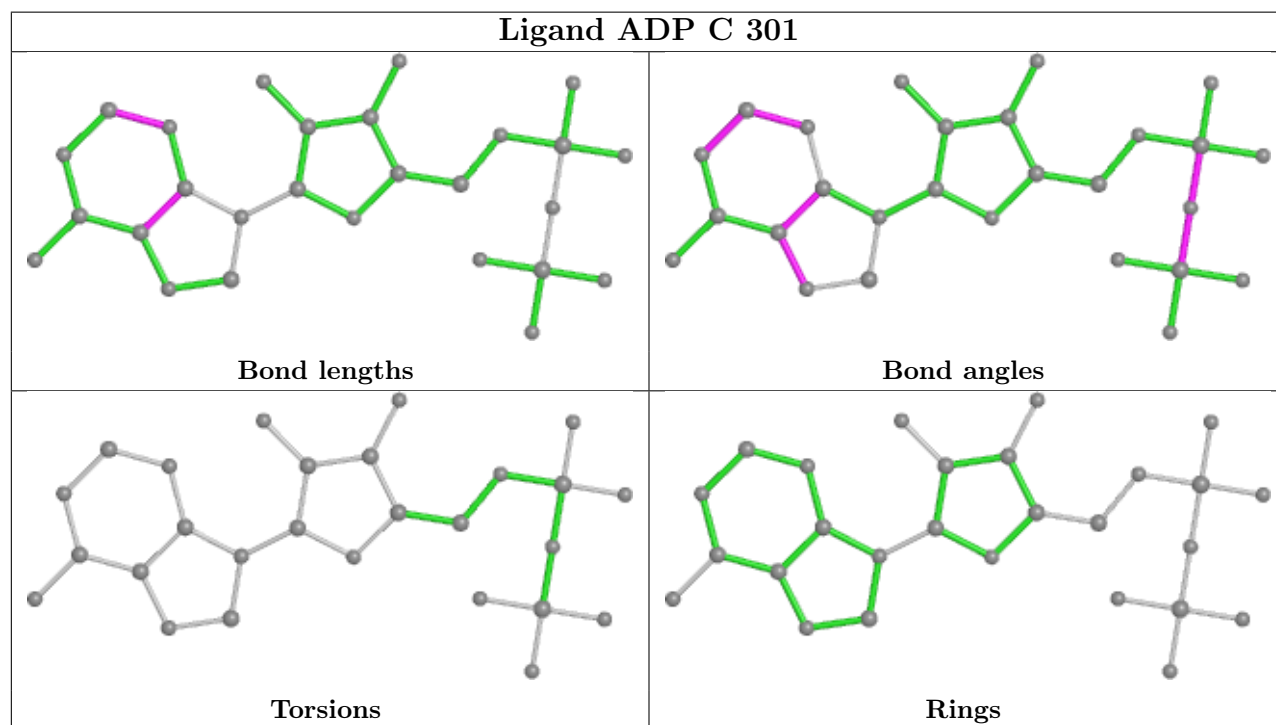
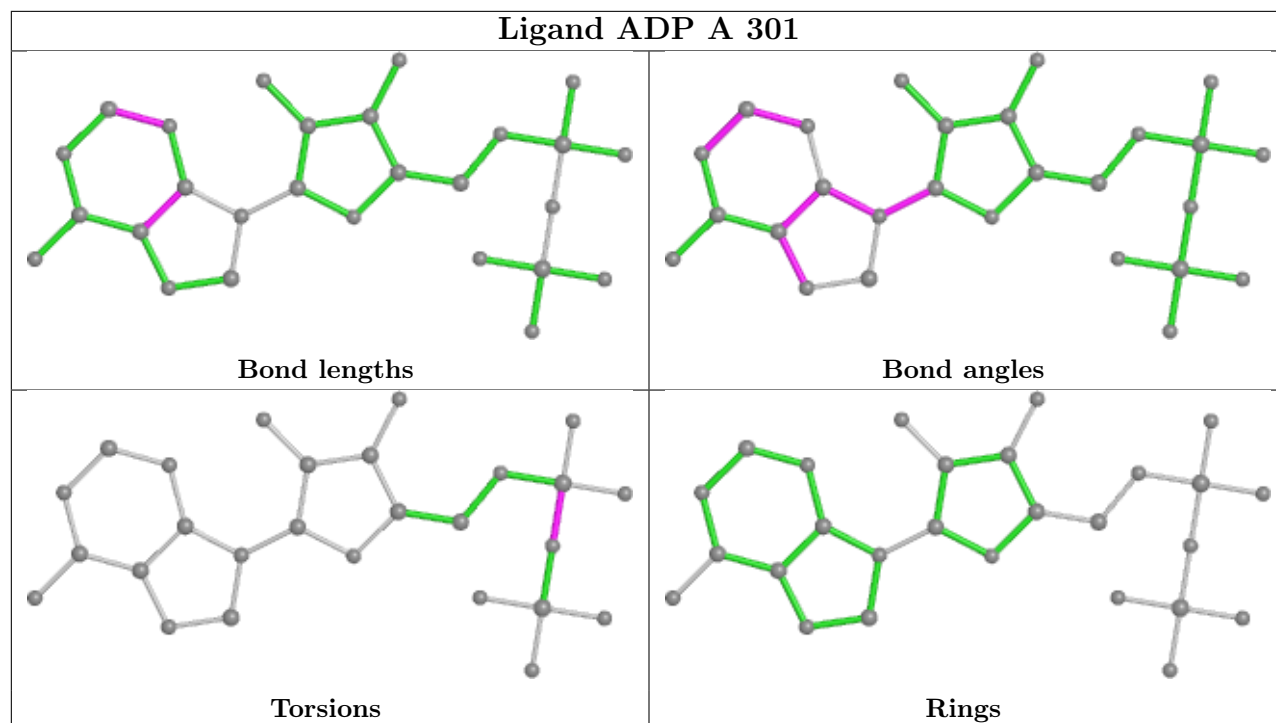
2 monomers are involved in 2 short contacts:

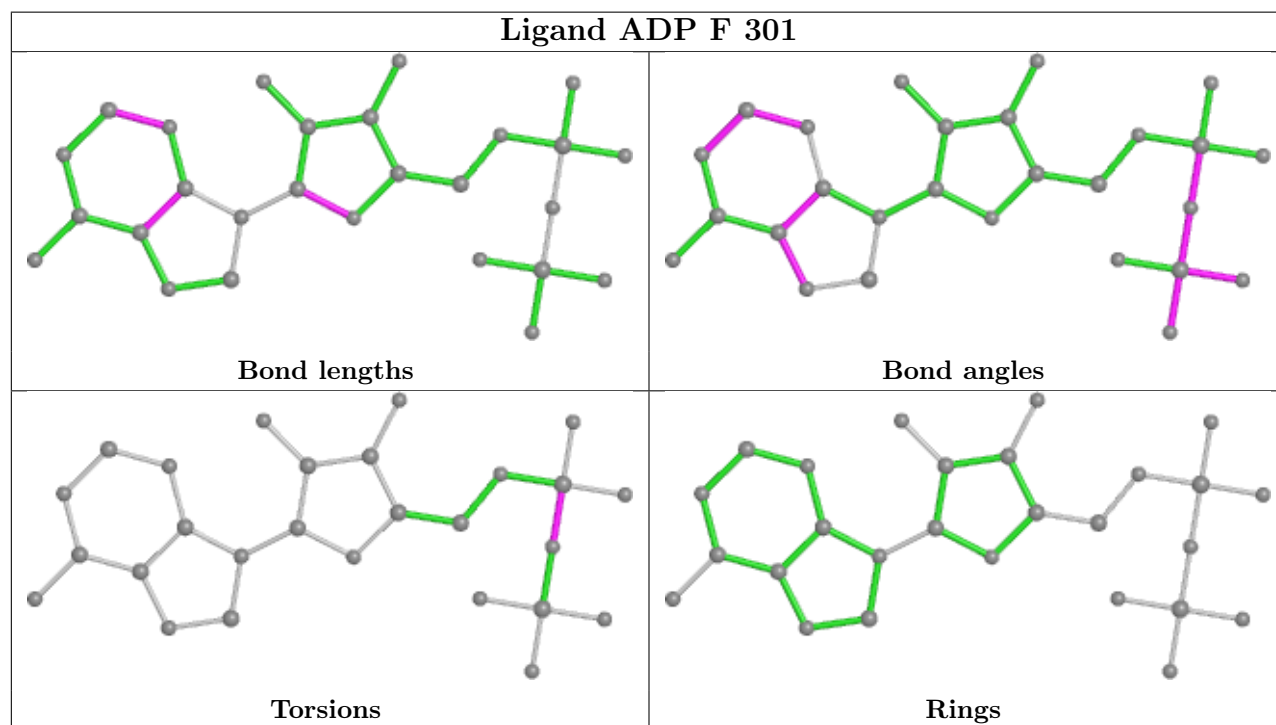
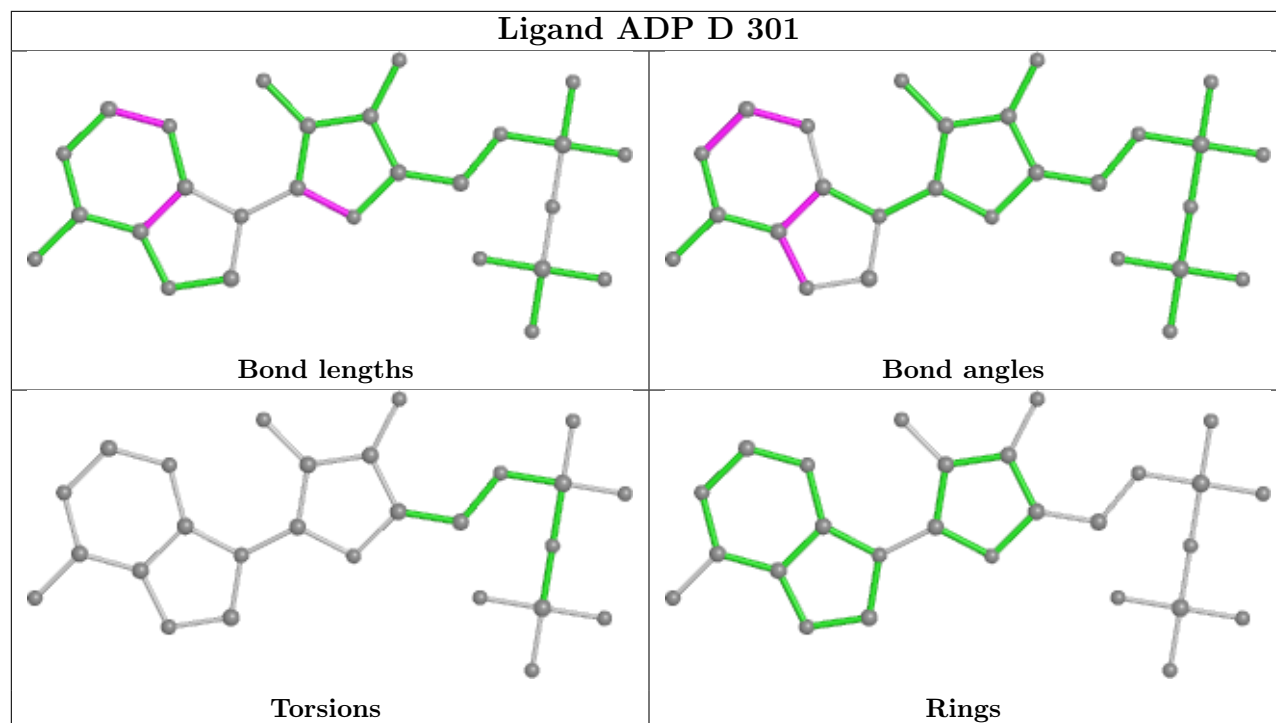
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	H	302	SH3	1	0
2	A	301	ADP	1	0

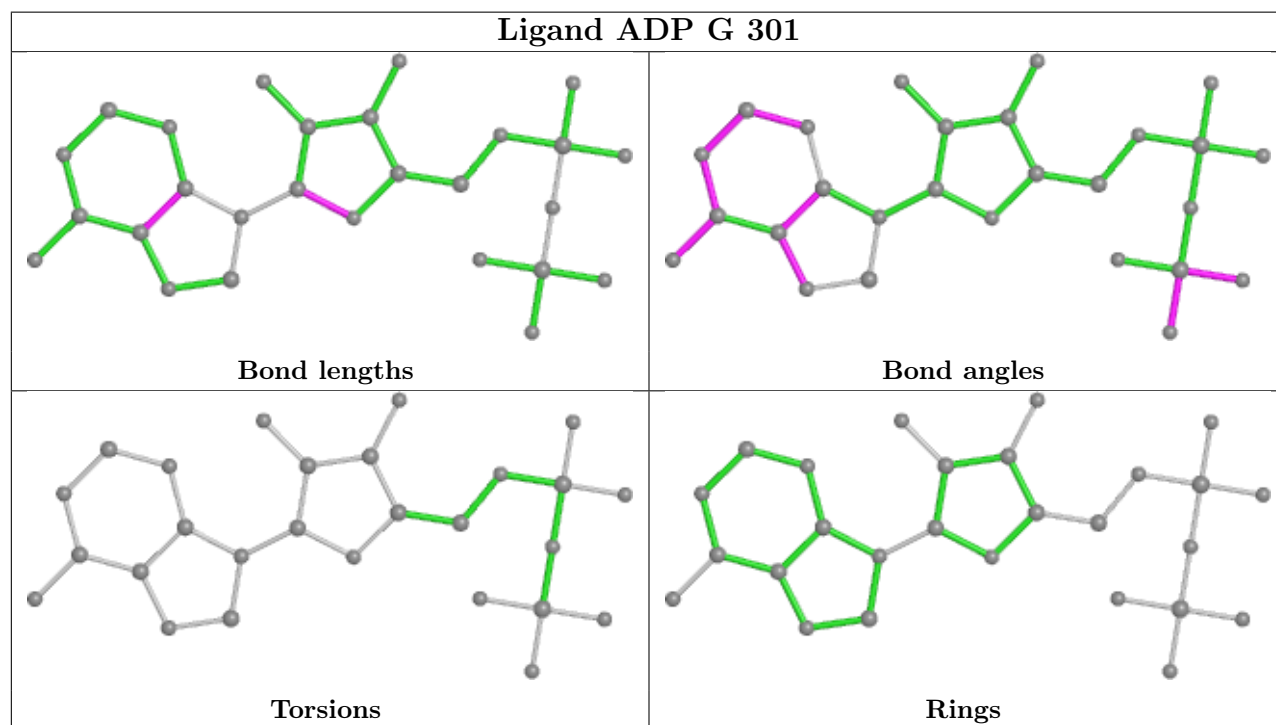
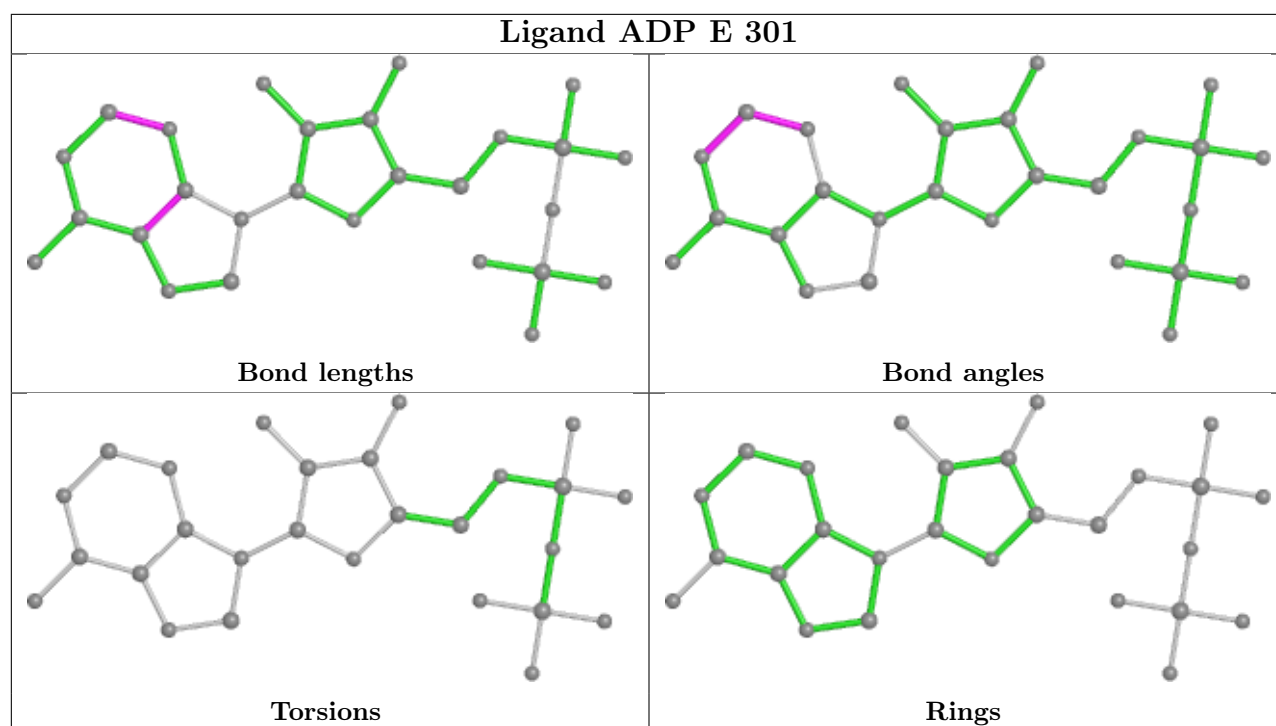
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

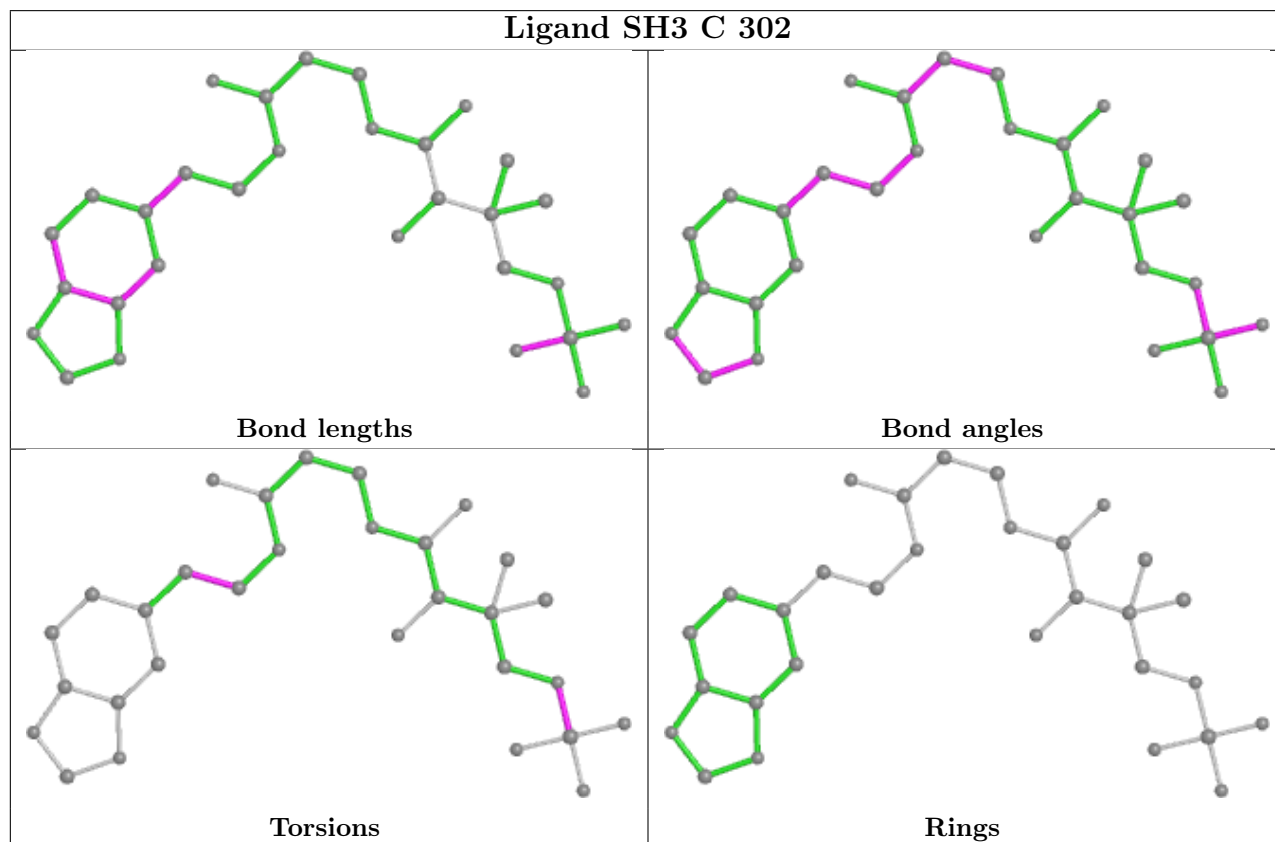
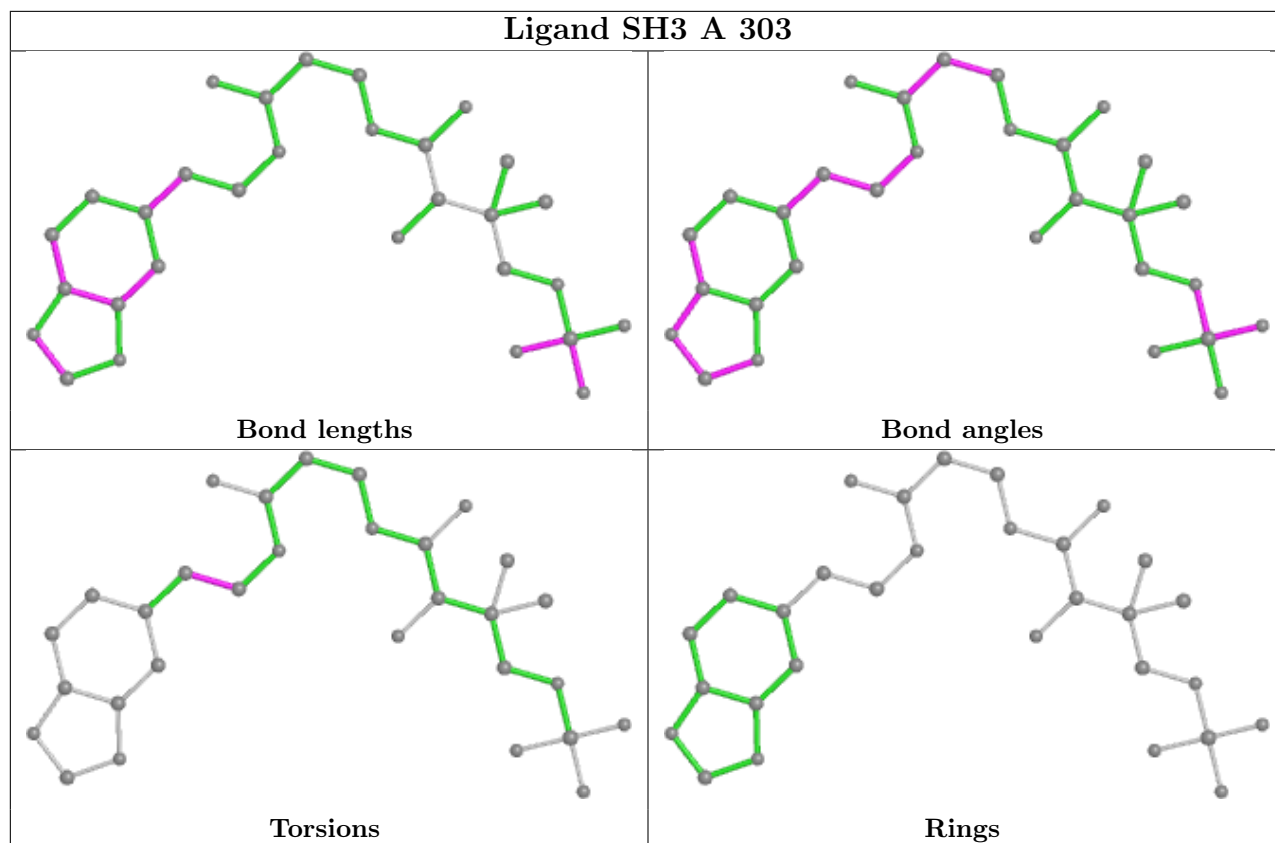


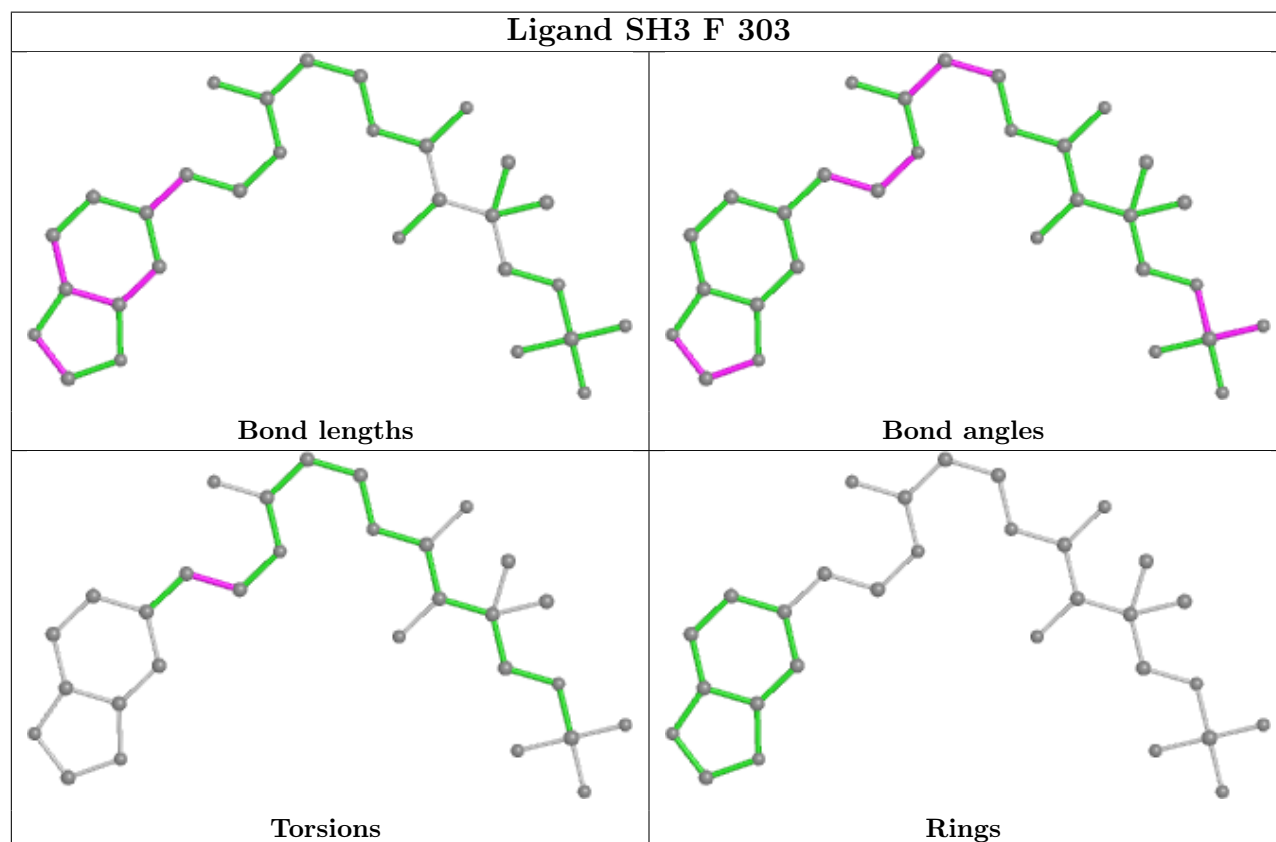
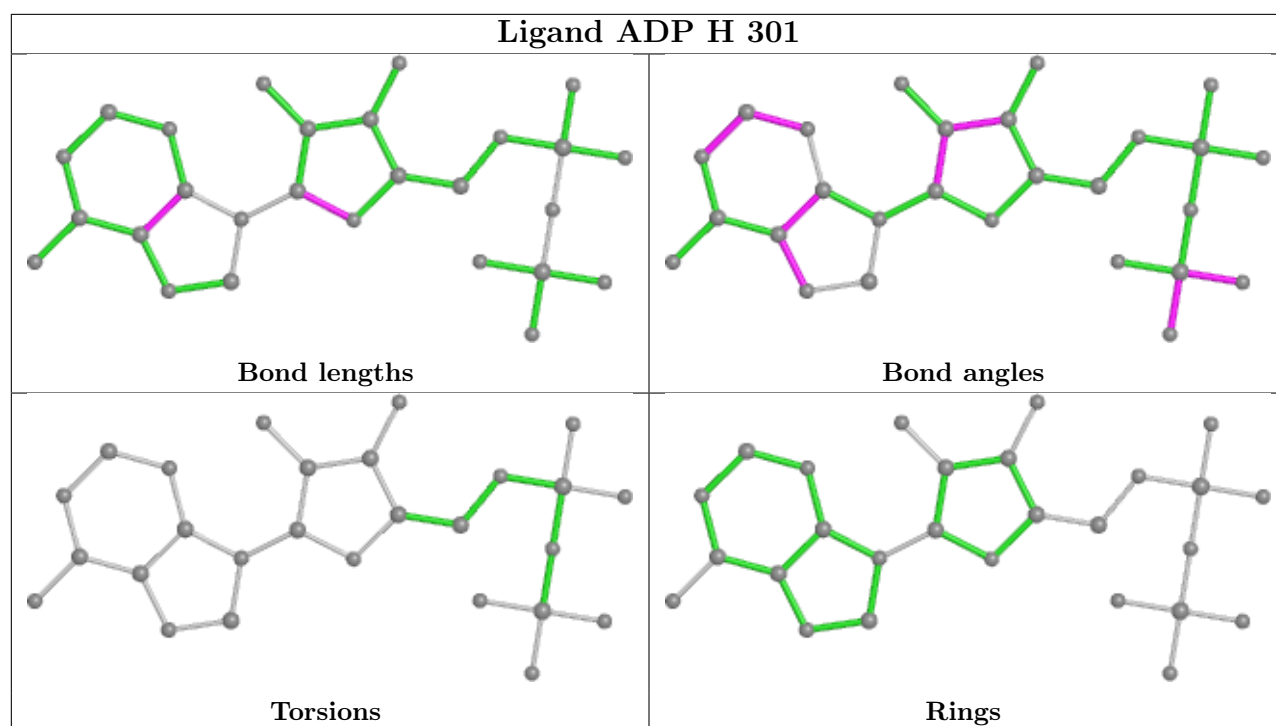


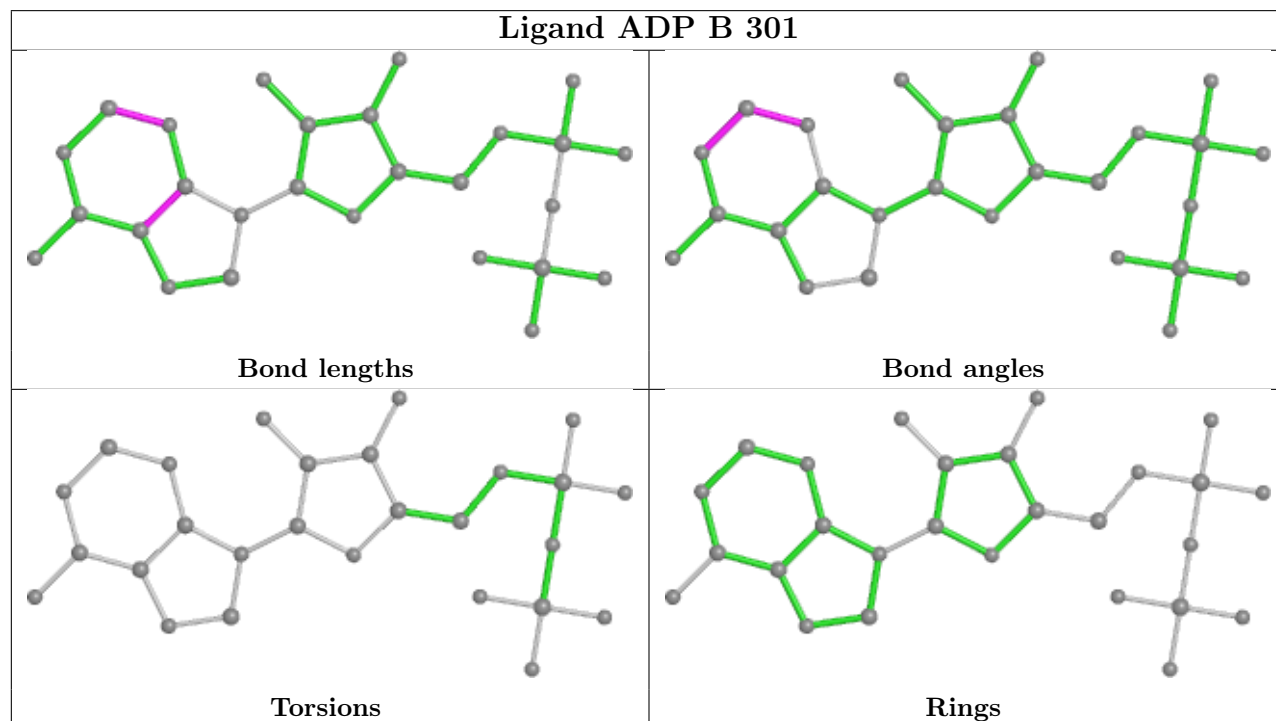
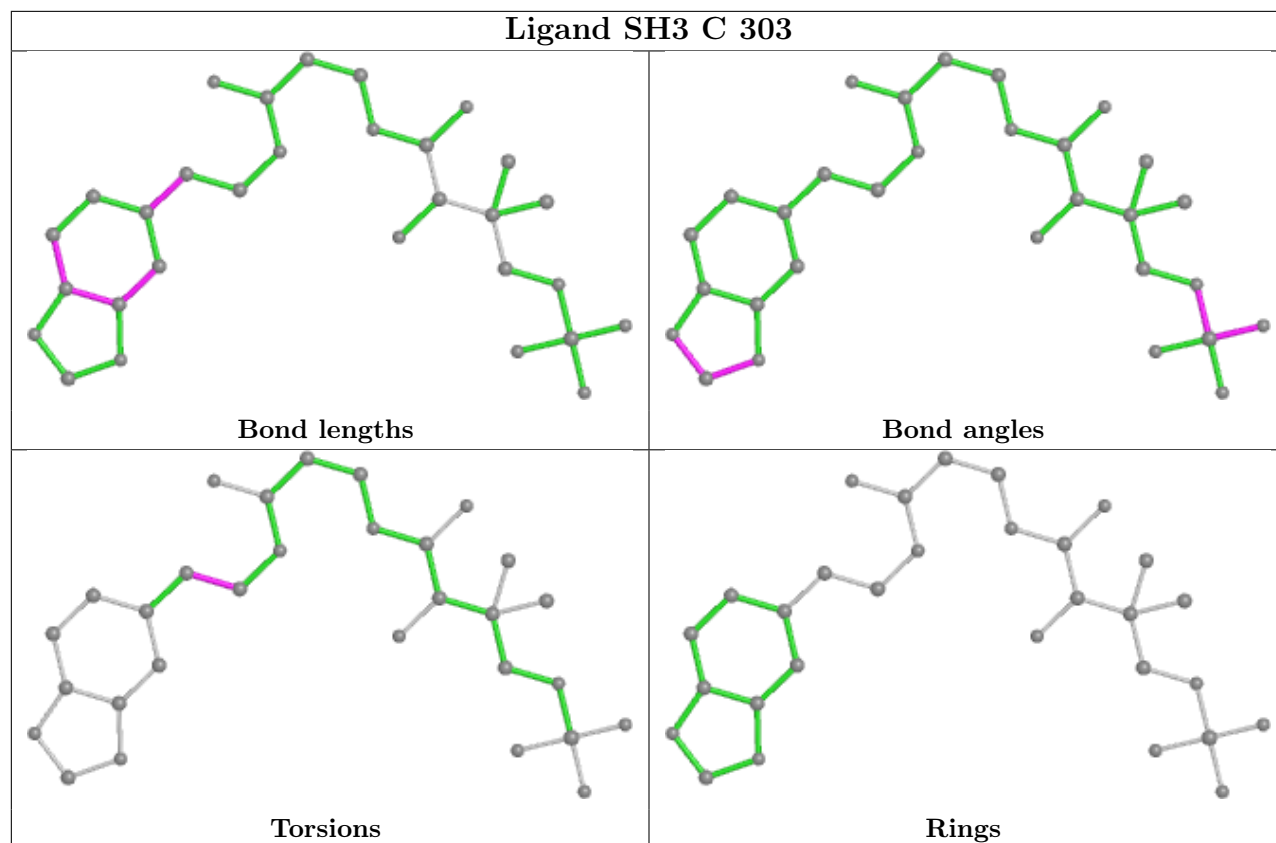


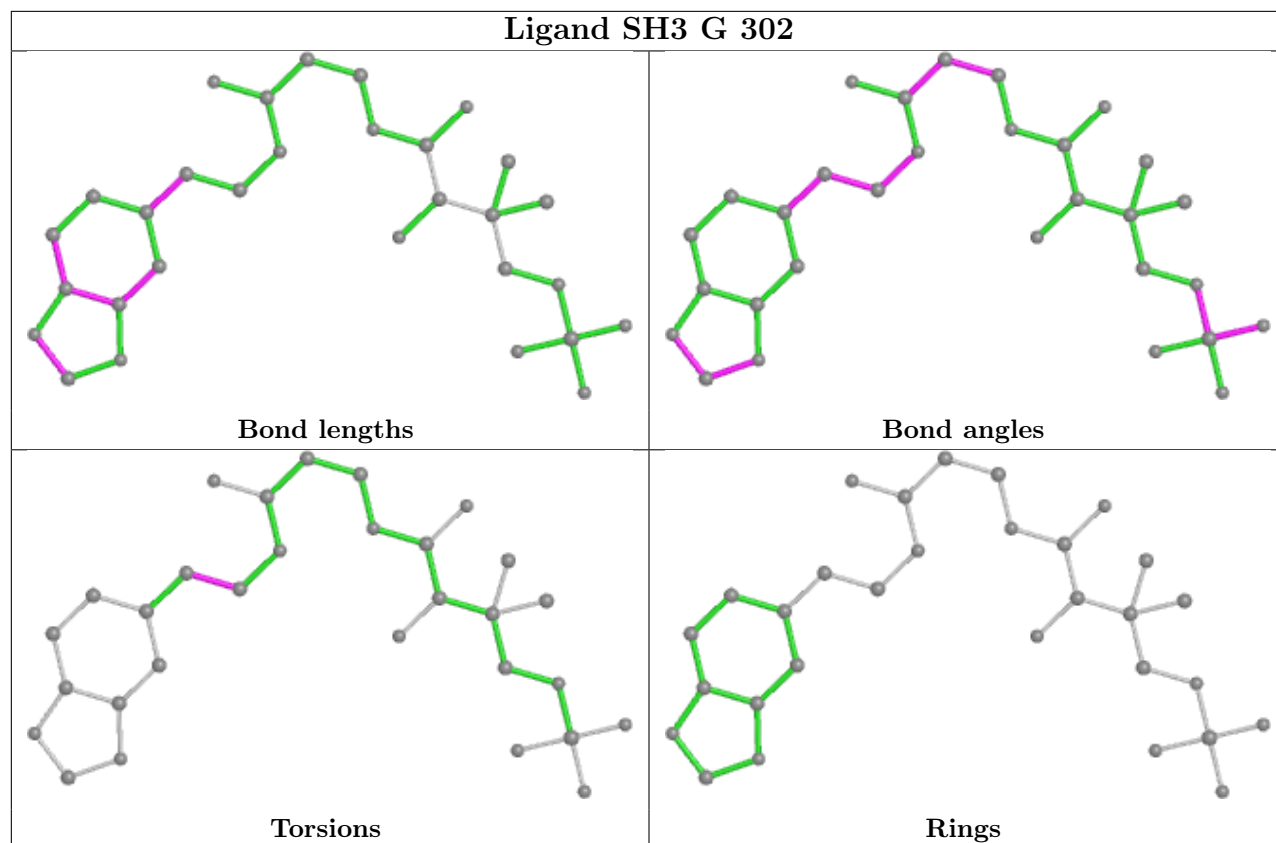












5.7 Other polymers [i](#)

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

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	265/285 (92%)	0.05	7 (2%) 56 59	8, 18, 33, 53	1 (0%)
1	B	265/285 (92%)	-0.10	1 (0%) 92 93	8, 15, 30, 38	0
1	C	265/285 (92%)	0.09	4 (1%) 73 75	11, 20, 34, 53	0
1	D	265/285 (92%)	0.31	11 (4%) 36 38	12, 26, 48, 65	0
1	E	259/285 (90%)	0.20	10 (3%) 39 42	15, 26, 41, 48	0
1	F	262/285 (91%)	0.48	17 (6%) 18 20	16, 31, 53, 56	0
1	G	265/285 (92%)	-0.06	2 (0%) 86 87	7, 16, 30, 44	0
1	H	267/285 (93%)	-0.03	4 (1%) 73 75	9, 16, 34, 43	0
All	All	2113/2280 (92%)	0.12	56 (2%) 54 57	7, 20, 41, 65	1 (0%)

All (56) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	264	TYR	6.1
1	F	186	ASP	4.8
1	E	157	ARG	4.6
1	D	250[A]	TYR	4.6
1	C	267	LYS	4.4
1	C	20	ASN	4.1
1	D	21	GLN	3.9
1	F	21	GLN	3.8
1	D	265	LEU	3.7
1	A	250	TYR	3.7
1	C	266	GLU	3.5
1	A	20	ASN	3.4
1	C	264	TYR	3.4
1	F	185	ALA	3.3
1	E	186	ASP	3.3
1	F	86	GLY	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	207	MET	3.2
1	F	33	ASP	3.2
1	E	266	GLU	3.1
1	G	163	THR	3.0
1	F	65	ALA	2.9
1	D	185	ALA	2.8
1	D	62	ASN	2.8
1	G	20	ASN	2.8
1	H	30	LYS	2.7
1	A	63	ILE	2.6
1	F	20	ASN	2.6
1	F	183	LEU	2.5
1	E	19	ASP	2.4
1	F	23	THR	2.4
1	F	250	TYR	2.4
1	F	88	ASP	2.4
1	A	184	ASP	2.4
1	D	88	ASP	2.4
1	E	170	ASP	2.4
1	D	69	VAL	2.3
1	F	1	MET	2.3
1	H	20	ASN	2.3
1	D	109	GLN	2.3
1	H	61	ILE	2.3
1	E	166	PRO	2.2
1	E	156	VAL	2.2
1	F	106	PHE	2.2
1	F	182	HIS	2.2
1	F	3	VAL	2.2
1	A	62	ASN	2.1
1	E	43	GLN	2.2
1	F	56	VAL	2.1
1	D	46	LYS	2.1
1	H	62	ASN	2.1
1	E	90	ALA	2.1
1	F	59	GLU	2.1
1	D	181	HIS	2.1
1	B	162	ASP	2.1
1	E	140	LEU	2.0
1	A	1	MET	2.0

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

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

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

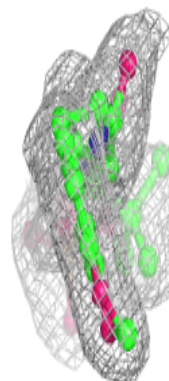
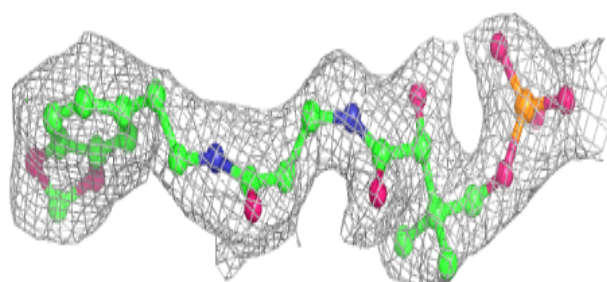
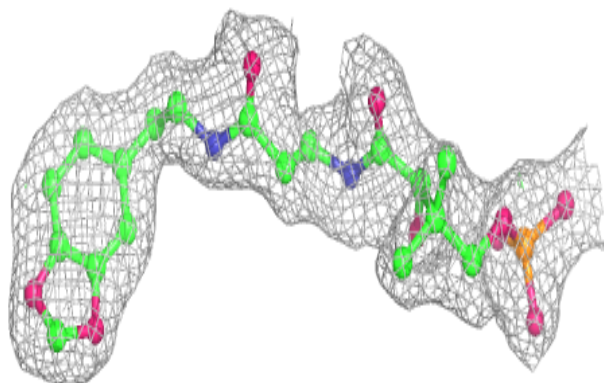
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	MG	D	302	1/1	0.82	0.08	19,19,19,19	0
4	MG	F	304	1/1	0.85	0.16	22,22,22,22	0
4	MG	E	302	1/1	0.88	0.07	25,25,25,25	0
4	MG	C	304	1/1	0.90	0.11	13,13,13,13	0
3	SH3	C	303	30/30	0.93	0.13	16,19,21,21	0
3	SH3	F	303	30/30	0.93	0.13	20,23,31,32	0
2	ADP	F	301	27/27	0.93	0.14	22,25,33,34	0
3	SH3	H	302	30/30	0.94	0.12	12,12,13,13	0
3	SH3	C	302	30/30	0.94	0.11	13,15,17,17	0
3	SH3	A	302	30/30	0.94	0.13	12,13,14,14	0
3	SH3	F	302	30/30	0.94	0.10	14,16,16,16	0
3	SH3	A	303	30/30	0.94	0.11	12,13,14,14	0
2	ADP	E	301	27/27	0.95	0.11	16,17,18,18	0
2	ADP	A	301	27/27	0.95	0.11	12,12,13,13	0
3	SH3	G	302	30/30	0.95	0.11	7,8,10,10	0
2	ADP	D	301	27/27	0.95	0.11	15,16,17,18	0
2	ADP	H	301	27/27	0.96	0.11	12,12,13,13	0
4	MG	B	302	1/1	0.96	0.09	11,11,11,11	0
2	ADP	C	301	27/27	0.96	0.11	10,10,11,11	0
4	MG	A	304	1/1	0.97	0.05	8,8,8,8	0
2	ADP	G	301	27/27	0.97	0.10	7,8,9,9	0
2	ADP	B	301	27/27	0.97	0.09	9,9,10,10	0
4	MG	G	303	1/1	0.97	0.11	2,2,2,2	0
4	MG	H	303	1/1	0.98	0.07	9,9,9,9	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different

orientation to approximate a three-dimensional view.

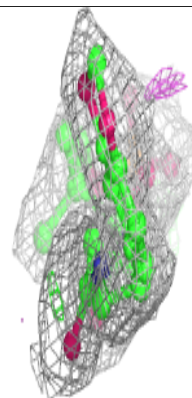
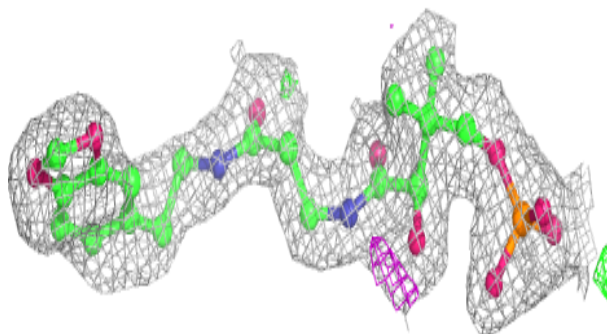
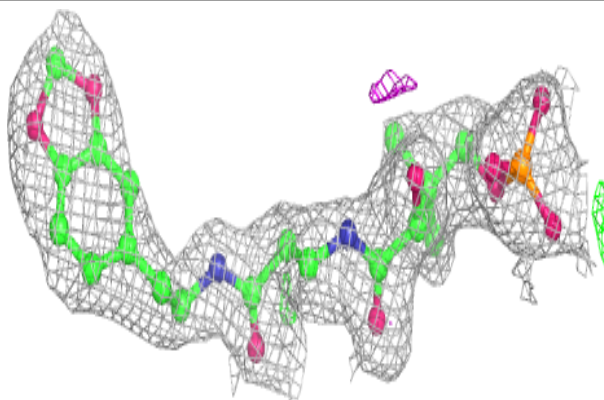
Electron density around SH3 C 303:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



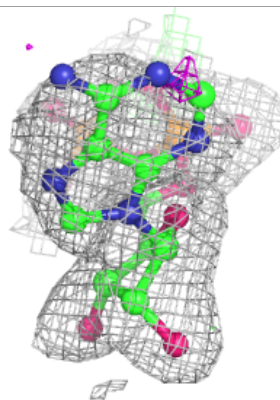
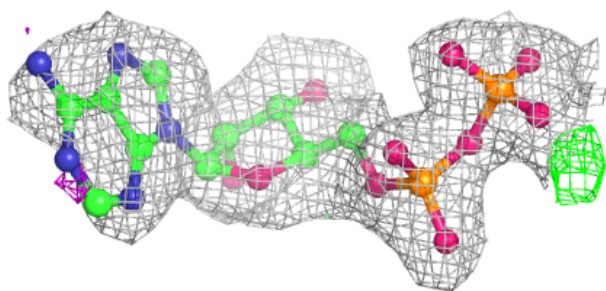
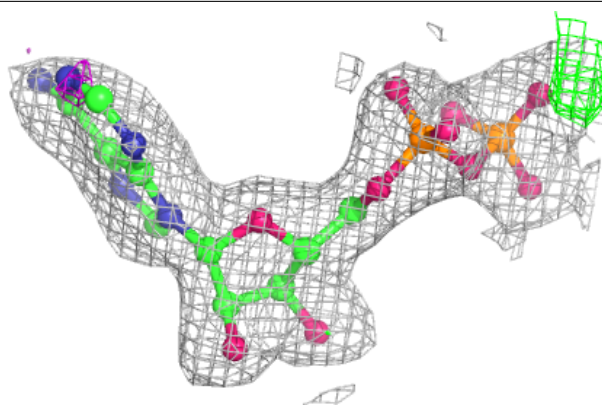
Electron density around SH3 F 303:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

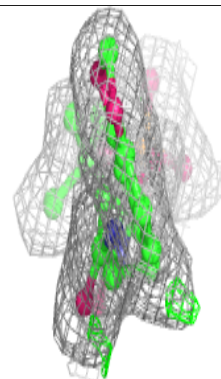
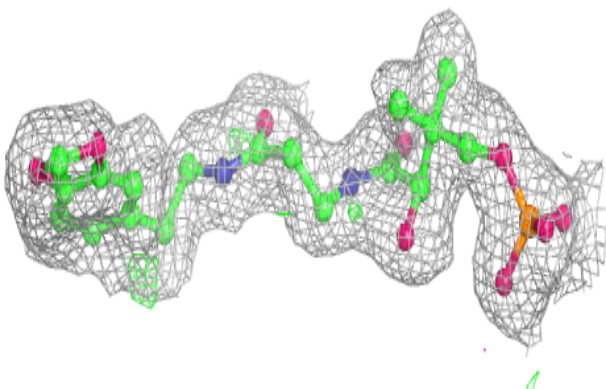
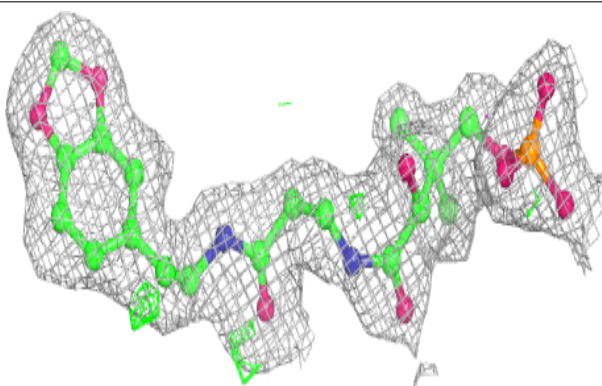


Electron density around ADP F 301:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

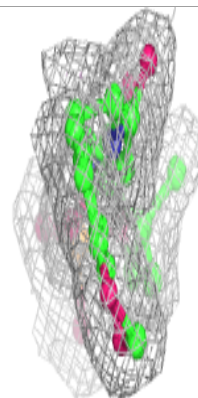
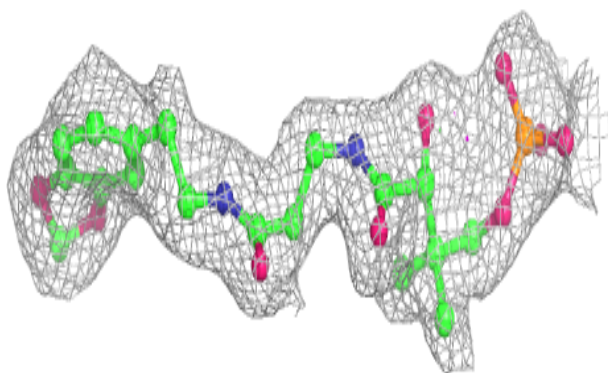
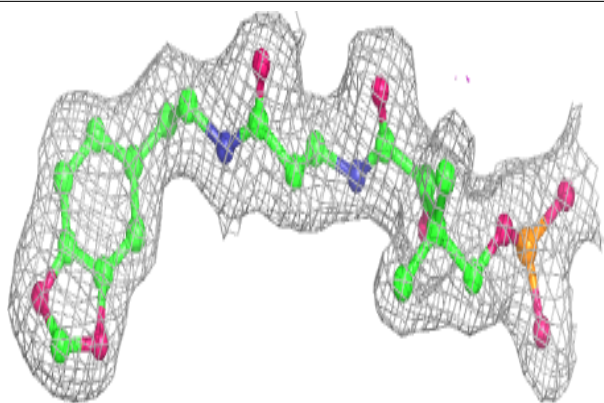
**Electron density around SH3 H 302:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

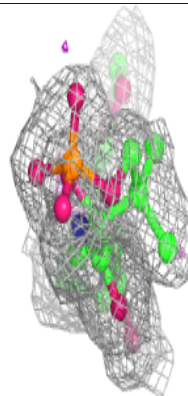
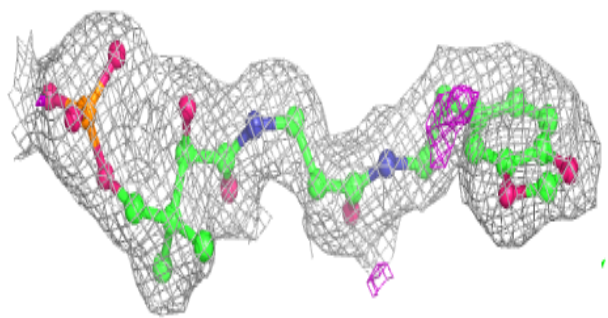
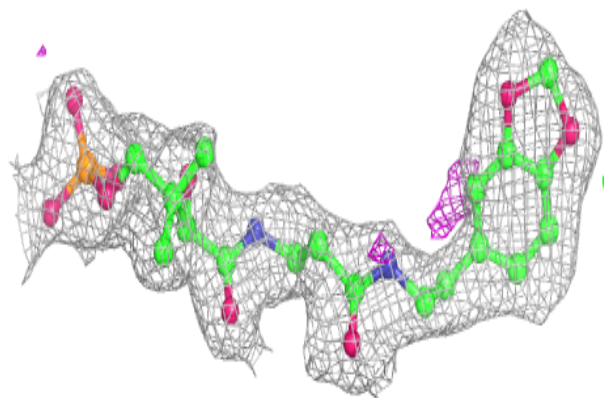


Electron density around SH3 C 302:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

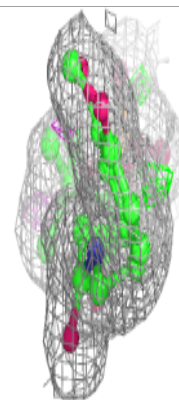
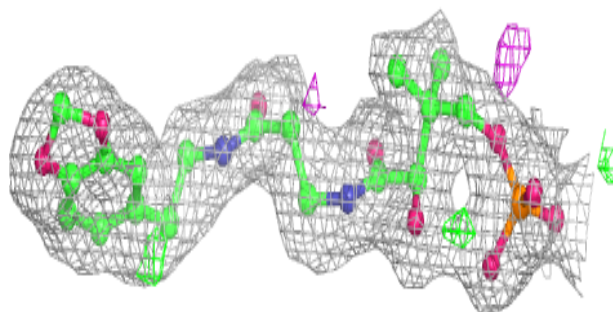
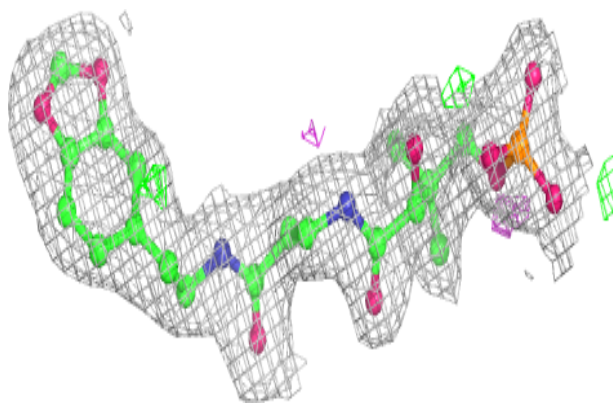
**Electron density around SH3 A 302:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

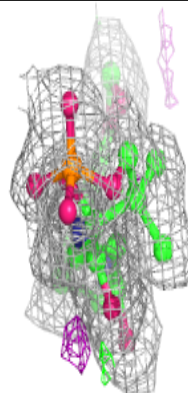
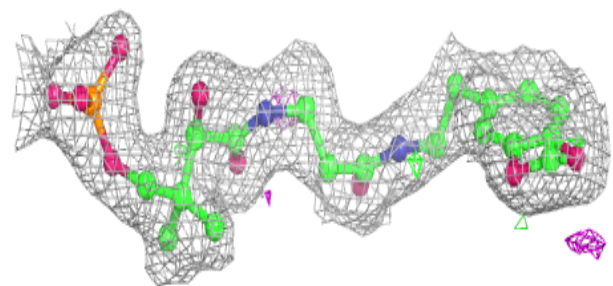
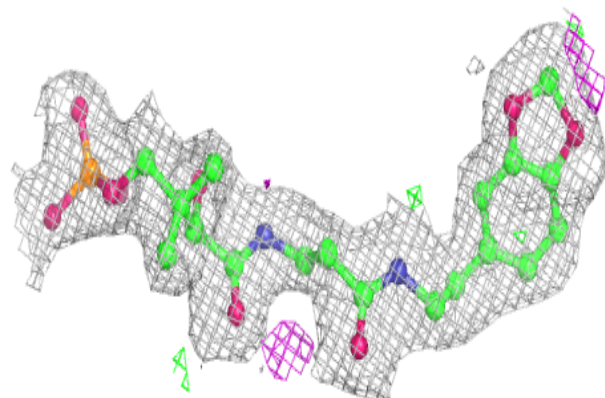


Electron density around SH3 F 302:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

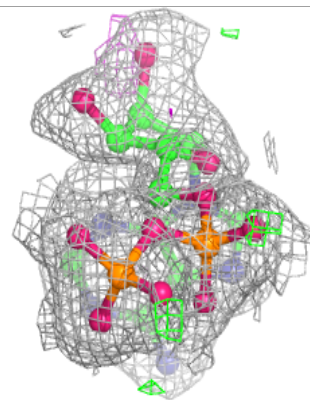
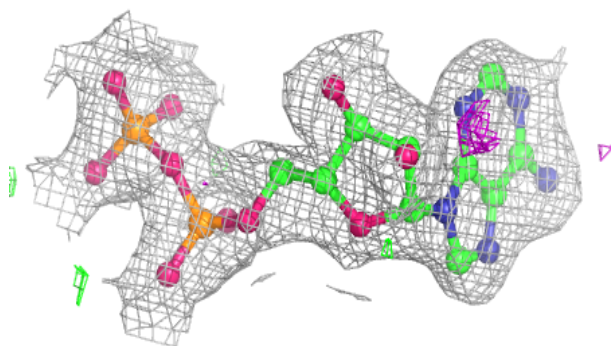
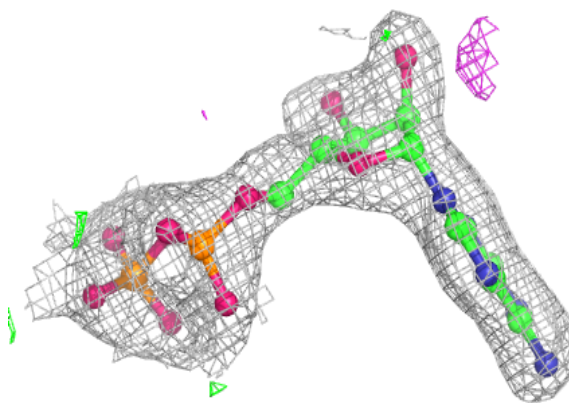
**Electron density around SH3 A 303:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

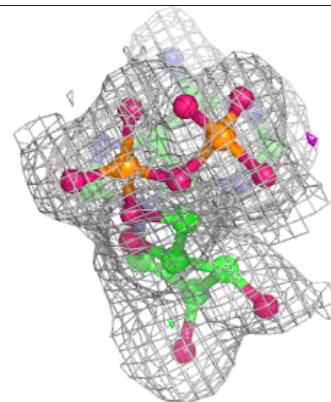
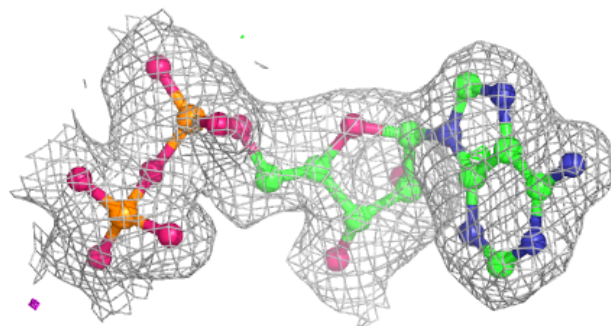
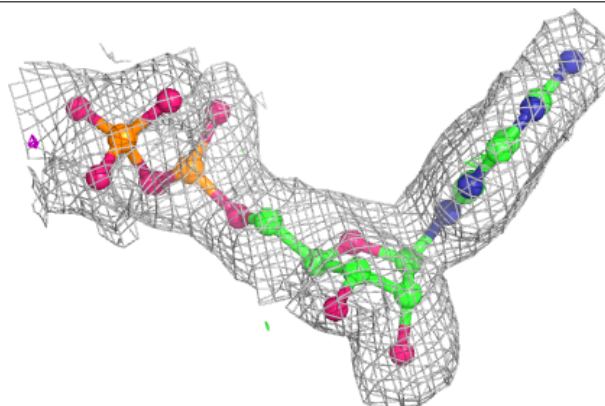


Electron density around ADP E 301:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

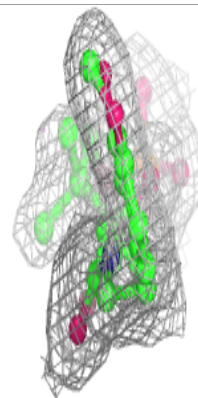
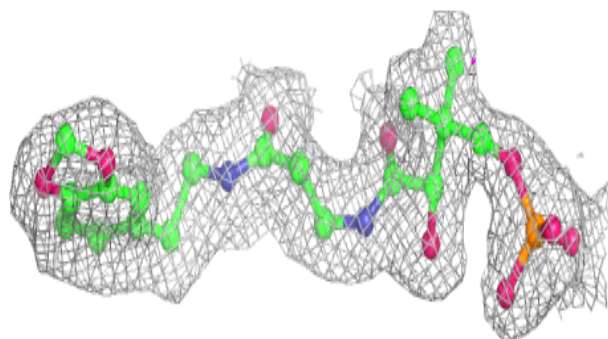
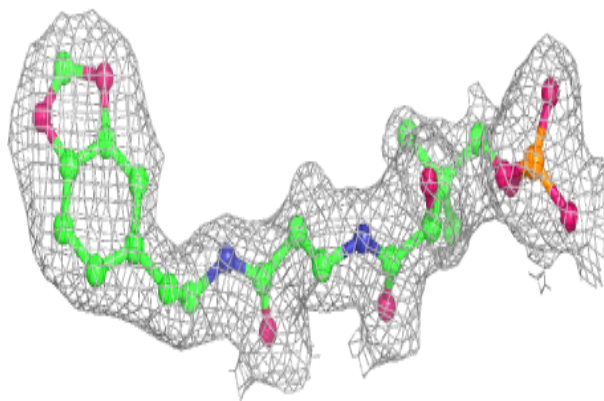
**Electron density around ADP A 301:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

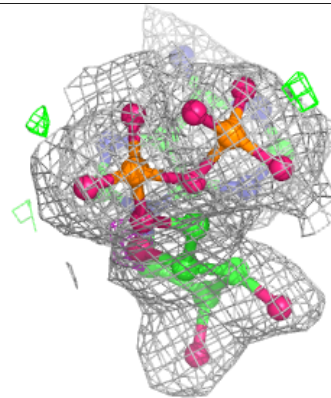
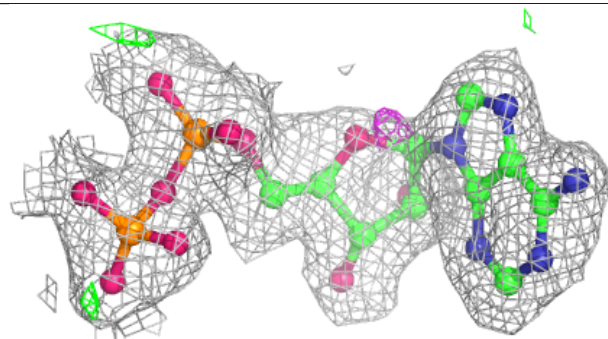
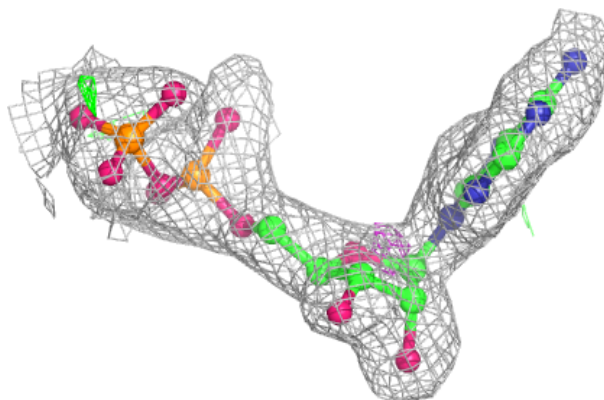


Electron density around SH3 G 302:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

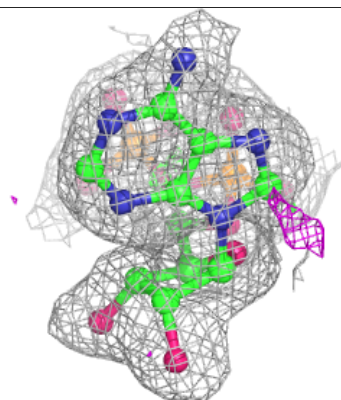
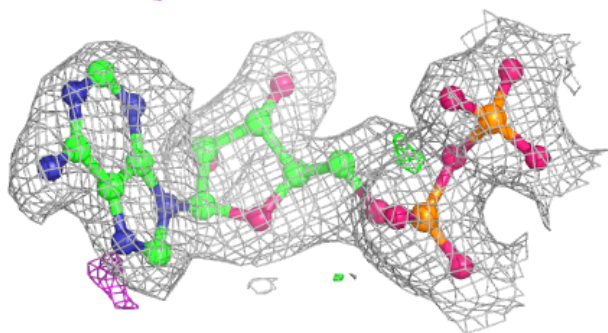
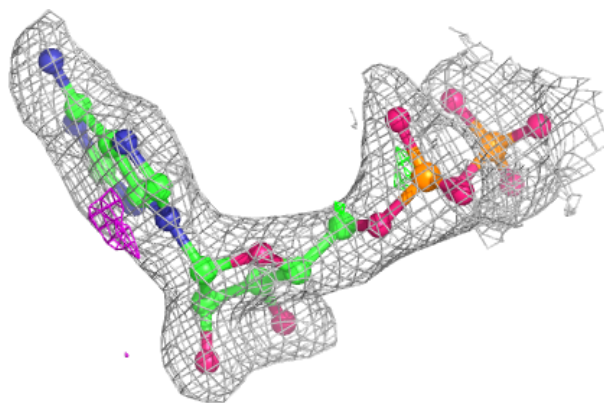
**Electron density around ADP D 301:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

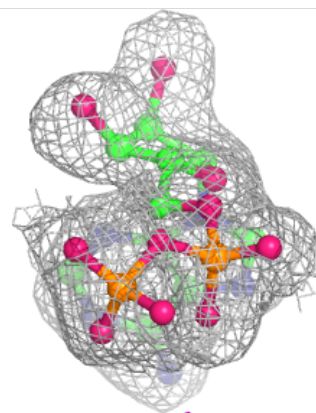
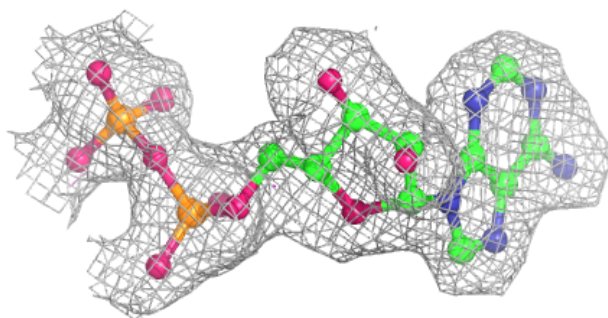
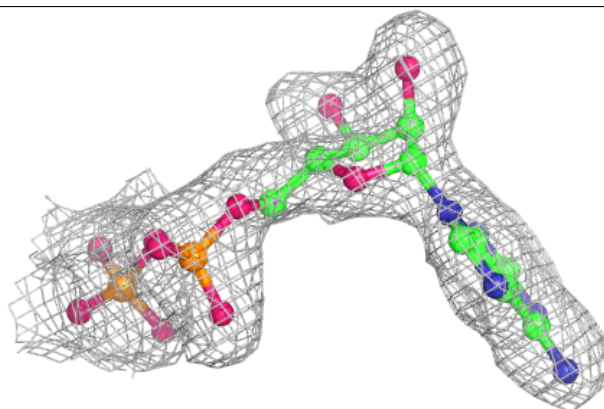


Electron density around ADP H 301:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

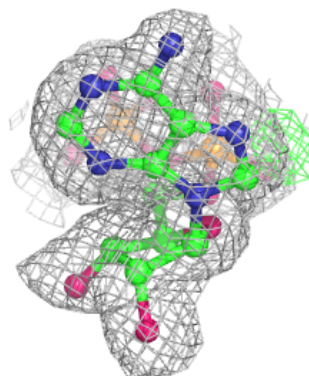
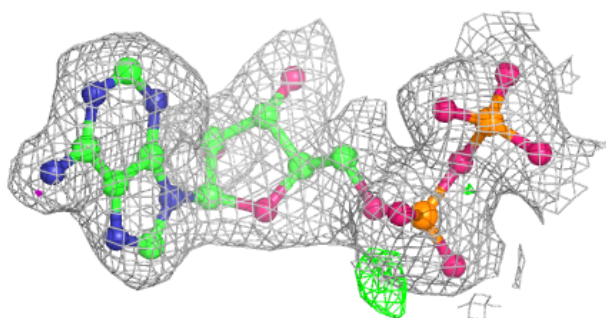
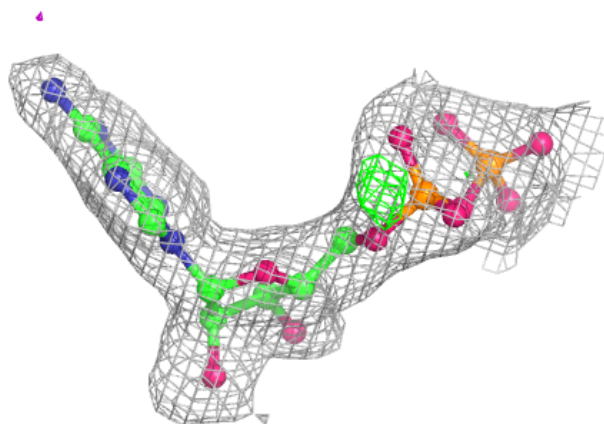
**Electron density around ADP C 301:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

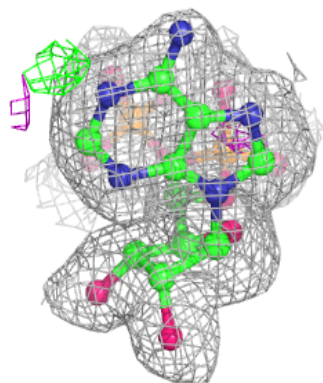
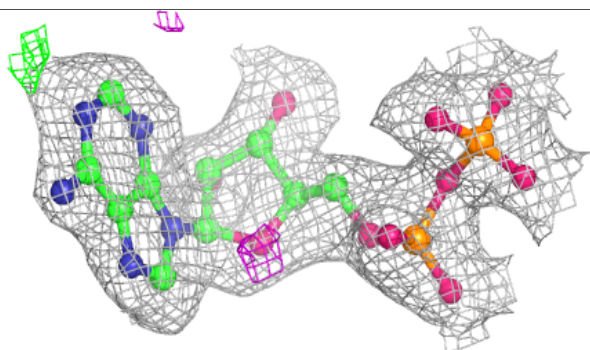
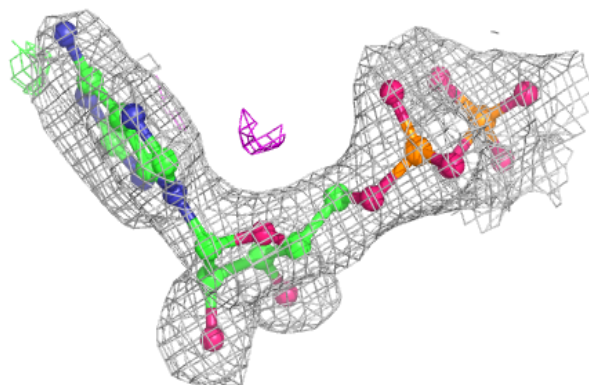


Electron density around ADP G 301:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around ADP B 301:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.