



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 21, 2020 – 03:22 PM BST

PDB ID : 5GIY
Title : HSA-Palmitic acid-[RuCl5(ind)]2-
Authors : Yang, F.; Wang, T.
Deposited on : 2016-06-25
Resolution : 2.54 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.13.1
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) : 2.13.1

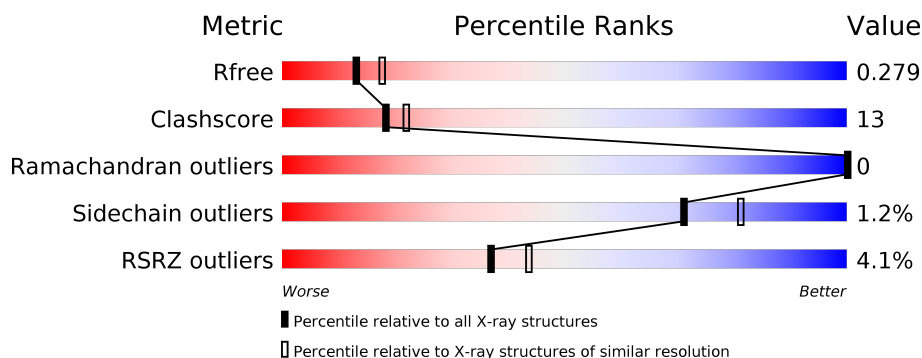
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.54 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	1284 (2.56-2.52)
Clashscore	141614	1332 (2.56-2.52)
Ramachandran outliers	138981	1315 (2.56-2.52)
Sidechain outliers	138945	1315 (2.56-2.52)
RSRZ outliers	127900	1272 (2.56-2.52)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. 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	581	<div> <div>4%</div> <div>72%</div> <div>27%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	PLM	A	1001	-	-	-	X
2	PLM	A	1004	-	-	X	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	PLM	A	1005	-	-	-	X
3	7GE	A	1007	-	-	X	-

2 Entry composition [i](#)

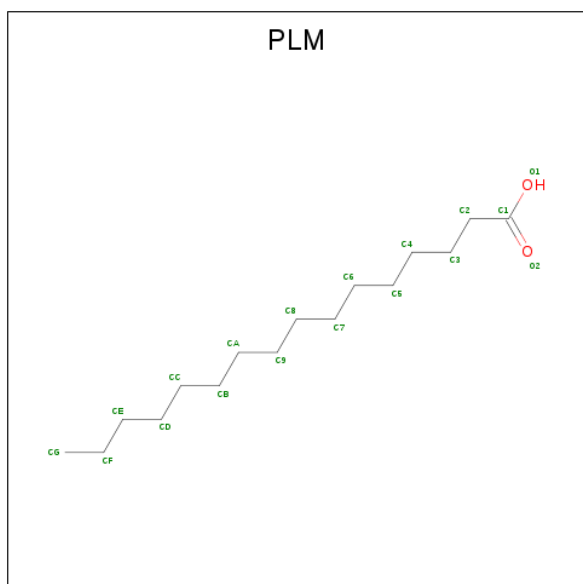
There are 4 unique types of molecules in this entry. The entry contains 4753 atoms, of which 10 are hydrogens and 0 are deuteriums.

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 Serum albumin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	581	Total	C	N	O	S	0	0	0
			4606	2909	776	880	41			

- Molecule 2 is PALMITIC ACID (three-letter code: PLM) (formula: $C_{16}H_{32}O_2$).



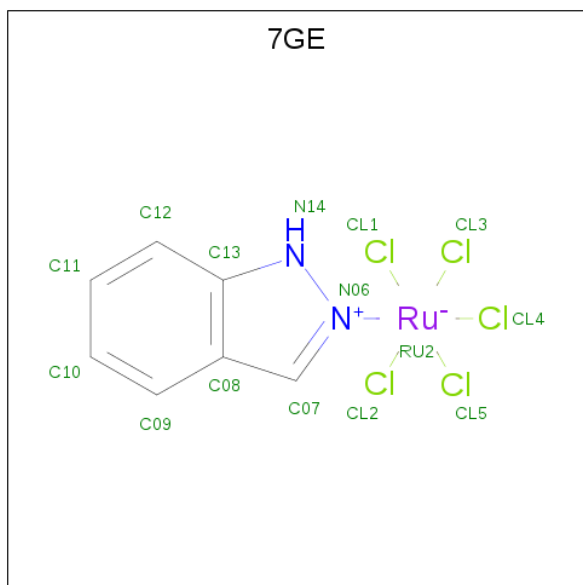
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			13	11	2		
2	A	1	Total	C	O	0	0
			18	16	2		
2	A	1	Total	C	O	0	0
			17	15	2		
2	A	1	Total	C	O	0	0
			18	16	2		
2	A	1	Total	C	O	0	0
			17	15	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			18	16	2		

- Molecule 3 is pentakis(chloranyl)-(1 {H}-indazol-2-ium-2-yl)ruthenium(1-) (three-letter code: 7GE) (formula: C₇H₆Cl₅N₂Ru).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	Cl	H	N	Ru	0	0
			18	7	4	4	2	1		
3	A	1	Total	C	Cl	H	N	Ru	0	0
			19	7	3	6	2	1		

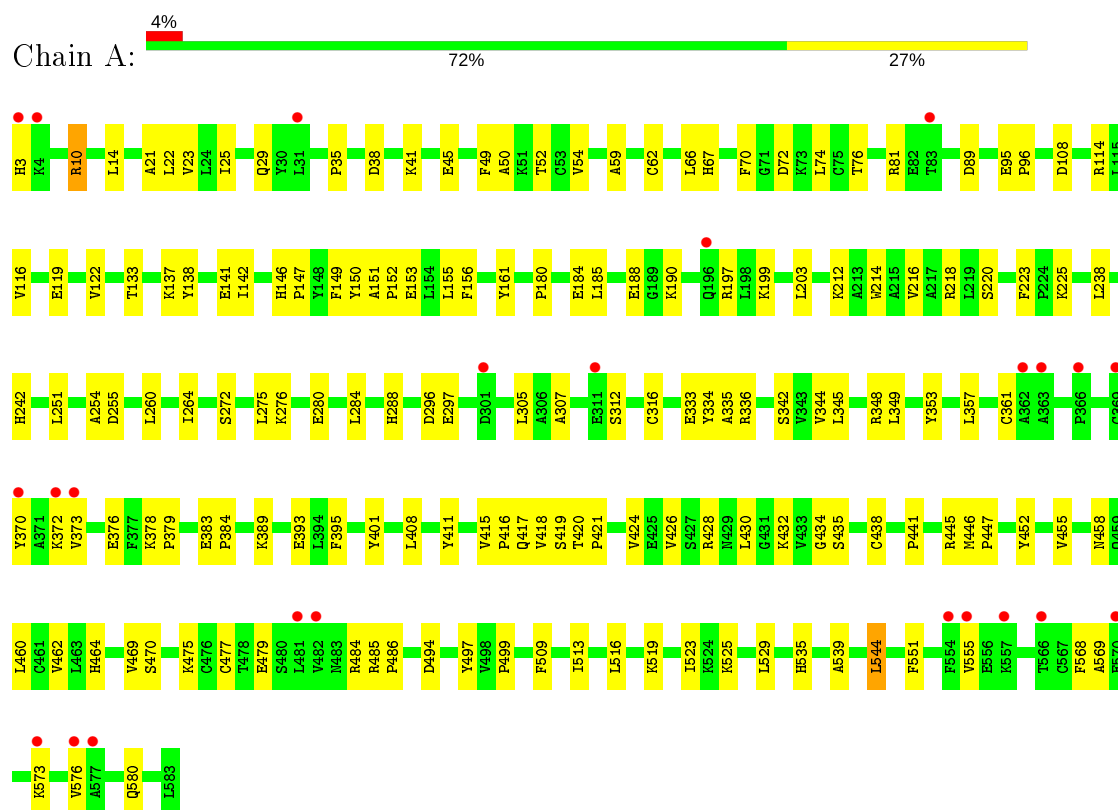
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	9	Total	O	0	0
			9	9		

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 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 ($\text{RSRZ} > 2$). 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: Serum albumin



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	185.84Å 38.71Å 96.07Å 90.00° 105.12° 90.00°	Depositor
Resolution (Å)	46.43 – 2.54 46.43 – 2.54	Depositor EDS
% Data completeness (in resolution range)	98.8 (46.43-2.54) 98.9 (46.43-2.54)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.49 (at 2.54Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.224 , 0.279 0.230 , 0.279	Depositor DCC
R_{free} test set	1129 reflections (5.12%)	wwPDB-VP
Wilson B-factor (Å ²)	67.4	Xtriage
Anisotropy	0.030	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 47.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4753	wwPDB-VP
Average B, all atoms (Å ²)	78.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.95% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 7GE, PLM

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	0.23	0/4696	0.36	0/6339

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4606	0	4495	120	0
2	A	101	0	163	13	0
3	A	27	10	0	12	0
4	A	9	0	0	1	0
All	All	4743	10	4658	125	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 13.

All (125) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:190:LYS:HD3	3:A:1007:7GE:CL3	1.35	1.63
1:A:146:HIS:HE1	3:A:1007:7GE:CL1	1.57	1.24
1:A:146:HIS:CD2	3:A:1007:7GE:CL4	2.33	1.18
1:A:190:LYS:CD	3:A:1007:7GE:CL3	2.31	1.13
1:A:146:HIS:CE1	3:A:1007:7GE:CL1	2.42	1.08
1:A:146:HIS:NE2	3:A:1007:7GE:CL4	2.29	1.03
1:A:460:LEU:HD21	2:A:1004:PLM:HC2	1.47	0.94
1:A:146:HIS:HD2	3:A:1007:7GE:CL4	1.98	0.84
2:A:1003:PLM:H81	2:A:1004:PLM:H22	1.61	0.83
1:A:59:ALA:HB3	1:A:62:CYS:SG	2.20	0.81
1:A:426:VAL:HG11	1:A:460:LEU:HD13	1.68	0.76
1:A:418:VAL:HG21	2:A:1004:PLM:HD2	1.67	0.75
1:A:242:HIS:NE2	3:A:1008:7GE:CL1	2.59	0.72
1:A:316:CYS:SG	1:A:370:TYR:OH	2.47	0.72
1:A:408:LEU:HD23	1:A:529:LEU:HD23	1.74	0.68
1:A:119:GLU:HB2	1:A:122:VAL:HG23	1.76	0.67
1:A:264:ILE:HG22	1:A:275:LEU:HD21	1.76	0.67
1:A:81:ARG:NH2	1:A:89:ASP:OD1	2.28	0.67
3:A:1008:7GE:CL1	4:A:1109:HOH:O	2.51	0.65
1:A:151:ALA:HB3	1:A:152:PRO:HD3	1.79	0.65
1:A:415:VAL:HG12	1:A:418:VAL:HG23	1.80	0.64
1:A:264:ILE:CG2	1:A:275:LEU:HD21	2.29	0.63
1:A:420:THR:HB	1:A:421:PRO:HD3	1.80	0.62
1:A:419:SER:OG	1:A:421:PRO:HD2	2.00	0.62
1:A:424:VAL:O	1:A:428:ARG:HG3	1.99	0.62
1:A:418:VAL:CG2	2:A:1004:PLM:HD2	2.30	0.60
1:A:155:LEU:HD23	1:A:284:LEU:HD21	1.84	0.60
1:A:225:LYS:HE3	1:A:297:GLU:O	2.01	0.60
1:A:485:ARG:HB3	1:A:486:PRO:HD3	1.84	0.60
1:A:416:PRO:HB2	1:A:497:TYR:CD1	2.37	0.60
1:A:417:GLN:NE2	1:A:494:ASP:OD2	2.30	0.59
1:A:254:ALA:HB1	2:A:1002:PLM:H51	1.84	0.58
1:A:95:GLU:HG3	1:A:96:PRO:HA	1.86	0.58
1:A:72:ASP:O	1:A:76:THR:HG23	2.04	0.57
1:A:464:HIS:HE1	1:A:470:SER:H	1.53	0.56
1:A:460:LEU:HD21	2:A:1004:PLM:CC	2.30	0.56
1:A:238:LEU:HD22	1:A:242:HIS:CE1	2.40	0.56
1:A:138:TYR:O	1:A:142:ILE:HG12	2.05	0.56
1:A:238:LEU:HD22	1:A:242:HIS:HE1	1.71	0.55
1:A:23:VAL:HG22	2:A:1002:PLM:HC1	1.87	0.55
1:A:446:MET:N	1:A:447:PRO:HD2	2.22	0.55
1:A:344:VAL:O	1:A:348:ARG:HG3	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:378:LYS:N	1:A:379:PRO:HD2	2.23	0.54
1:A:569:ALA:O	1:A:573:LYS:HG3	2.09	0.53
1:A:70:PHE:CE2	1:A:74:LEU:HD11	2.44	0.53
1:A:10:ARG:NH2	1:A:255:ASP:OD1	2.42	0.53
1:A:50:ALA:O	1:A:54:VAL:HG23	2.09	0.53
2:A:1003:PLM:HA1	2:A:1004:PLM:H21	1.91	0.52
1:A:276:LYS:O	1:A:280:GLU:HG3	2.09	0.52
1:A:242:HIS:CD2	3:A:1008:7GE:CL1	3.00	0.52
1:A:35:PRO:HG2	1:A:38:ASP:CG	2.31	0.51
1:A:108:ASP:OD2	1:A:197:ARG:NH1	2.43	0.51
1:A:519:LYS:O	1:A:523:ILE:HG12	2.10	0.51
1:A:137:LYS:O	1:A:141:GLU:HG2	2.10	0.51
1:A:460:LEU:CD1	2:A:1004:PLM:HE2	2.41	0.51
1:A:464:HIS:CE1	1:A:470:SER:H	2.29	0.51
1:A:161:TYR:CD2	1:A:185:LEU:HD23	2.46	0.50
1:A:161:TYR:CG	1:A:185:LEU:HD23	2.46	0.50
1:A:35:PRO:HG2	1:A:38:ASP:OD1	2.12	0.50
1:A:150:TYR:CZ	1:A:152:PRO:HG2	2.47	0.50
1:A:372:LYS:O	1:A:376:GLU:HG3	2.12	0.50
1:A:477:CYS:O	1:A:484:ARG:NH2	2.44	0.50
1:A:153:GLU:OE2	1:A:288:HIS:ND1	2.42	0.50
1:A:180:PRO:O	1:A:184:GLU:HG3	2.12	0.50
1:A:133:THR:O	1:A:137:LYS:HB2	2.11	0.49
1:A:14:LEU:HD13	1:A:22:LEU:HD12	1.95	0.49
1:A:296:ASP:OD1	1:A:297:GLU:N	2.45	0.49
1:A:307:ALA:O	1:A:312:SER:HB2	2.12	0.49
1:A:184:GLU:O	1:A:188:GLU:HG3	2.13	0.48
1:A:357:LEU:O	1:A:361:CYS:HB2	2.14	0.48
1:A:426:VAL:HG11	1:A:460:LEU:CD1	2.42	0.48
1:A:41:LYS:O	1:A:45:GLU:HG3	2.14	0.48
1:A:460:LEU:HD11	2:A:1004:PLM:HE2	1.96	0.48
1:A:114:ARG:CZ	1:A:116:VAL:HG12	2.43	0.48
1:A:576:VAL:O	1:A:580:GLN:HG3	2.14	0.48
3:A:1007:7GE:C07	3:A:1007:7GE:CL1	2.99	0.48
1:A:395:PHE:CZ	1:A:435:SER:HA	2.48	0.48
1:A:389:LYS:O	1:A:393:GLU:HG3	2.14	0.47
1:A:49:PHE:O	1:A:52:THR:HB	2.15	0.47
1:A:458:ASN:O	1:A:462:VAL:HG23	2.15	0.47
1:A:223:PHE:CD1	1:A:272:SER:HB2	2.50	0.46
1:A:551:PHE:O	1:A:555:VAL:HG23	2.15	0.46
1:A:415:VAL:CG1	1:A:418:VAL:HG23	2.44	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:333:GLU:HA	1:A:336:ARG:HD3	1.97	0.46
1:A:203:LEU:O	1:A:203:LEU:HD23	2.14	0.46
1:A:551:PHE:CD2	2:A:1005:PLM:H51	2.51	0.46
1:A:434:GLY:O	1:A:438:CYS:HB2	2.16	0.45
1:A:464:HIS:CE1	1:A:469:VAL:H	2.35	0.45
1:A:475:LYS:O	1:A:479:GLU:HG3	2.17	0.45
1:A:114:ARG:NH1	1:A:116:VAL:HG12	2.32	0.45
1:A:420:THR:O	1:A:424:VAL:HG23	2.17	0.44
1:A:539:ALA:CB	1:A:544:LEU:HD13	2.47	0.44
3:A:1007:7GE:CL4	3:A:1007:7GE:N14	2.87	0.44
1:A:411:TYR:HB3	2:A:1004:PLM:HA1	1.99	0.44
1:A:426:VAL:CG1	1:A:460:LEU:HD13	2.41	0.44
1:A:21:ALA:O	1:A:25:ILE:HG13	2.18	0.44
1:A:428:ARG:O	1:A:432:LYS:HG3	2.18	0.44
1:A:342:SER:HB3	1:A:446:MET:HG2	1.98	0.44
1:A:29:GLN:HG2	1:A:147:PRO:HA	2.00	0.44
1:A:305:LEU:HD22	1:A:333:GLU:HB3	2.00	0.44
1:A:150:TYR:CD2	1:A:152:PRO:HD2	2.54	0.43
1:A:383:GLU:HB3	1:A:384:PRO:HD3	2.00	0.43
1:A:383:GLU:HB3	1:A:384:PRO:CD	2.48	0.43
1:A:426:VAL:O	1:A:430:LEU:HG	2.19	0.43
1:A:214:TRP:CH2	1:A:218:ARG:HD2	2.54	0.43
1:A:513:ILE:HA	1:A:516:LEU:HD12	2.01	0.43
1:A:212:LYS:O	1:A:216:VAL:HG23	2.19	0.43
1:A:260:LEU:O	1:A:264:ILE:HG13	2.19	0.43
1:A:539:ALA:HB1	1:A:544:LEU:HD13	2.01	0.43
1:A:345:LEU:O	1:A:349:LEU:HG	2.18	0.42
1:A:441:PRO:O	1:A:445:ARG:HG3	2.19	0.42
1:A:383:GLU:N	1:A:384:PRO:HD2	2.35	0.42
1:A:216:VAL:HG21	2:A:1006:PLM:HE2	2.00	0.42
1:A:67:HIS:CD2	1:A:251:LEU:HD12	2.55	0.42
1:A:10:ARG:HG3	1:A:66:LEU:HD11	2.00	0.42
1:A:416:PRO:HB2	1:A:497:TYR:CE1	2.55	0.42
1:A:153:GLU:O	1:A:156:PHE:HB3	2.20	0.41
1:A:452:TYR:O	1:A:455:VAL:HG22	2.19	0.41
1:A:150:TYR:CE2	1:A:152:PRO:HG2	2.55	0.41
1:A:401:TYR:OH	1:A:525:LYS:HE2	2.20	0.41
1:A:499:PRO:HB3	1:A:535:HIS:O	2.21	0.41
1:A:353:TYR:HD1	1:A:373:VAL:HG11	1.85	0.41
1:A:220:SER:HB2	1:A:335:ALA:HB3	2.02	0.40
1:A:509:PHE:O	1:A:568:PHE:HB3	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:119:GLU:HB2	1:A:122:VAL:CG2	2.49	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	579/581 (100%)	563 (97%)	16 (3%)	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	503/509 (99%)	497 (99%)	6 (1%)	71 81

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	HIS
1	A	10	ARG
1	A	149	PHE
1	A	199	LYS
1	A	334	TYR

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Mol	Chain	Res	Type
1	A	544	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	67	HIS

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 ⓘ

8 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	PLM	A	1004	-	14,17,17	0.07	0	13,17,17	0.46	0
3	7GE	A	1007	1	8,15,16	6.81	6 (75%)	8,25,30	1.07	0
2	PLM	A	1003	1	13,16,17	0.08	0	12,16,17	0.48	0
2	PLM	A	1002	-	14,17,17	0.10	0	13,17,17	0.42	0
2	PLM	A	1006	1	14,17,17	0.09	0	13,17,17	0.42	0
2	PLM	A	1001	-	9,12,17	0.09	0	8,12,17	0.39	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	PLM	A	1005	-	13,16,17	0.07	0	12,16,17	0.42	0
3	7GE	A	1008	-	8,14,16	6.34	7 (87%)	8,21,30	0.95	0

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	PLM	A	1004	-	-	4/13/15/15	-
3	7GE	A	1007	1	-	-	0/2/2/2
2	PLM	A	1003	1	-	4/12/14/15	-
2	PLM	A	1002	-	-	2/13/15/15	-
2	PLM	A	1006	1	-	6/13/15/15	-
2	PLM	A	1001	-	-	4/8/10/15	-
2	PLM	A	1005	-	-	4/12/14/15	-
3	7GE	A	1008	-	-	-	0/2/2/2

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1007	7GE	C07-N06	-16.75	1.16	1.33
3	A	1008	7GE	C07-N06	-15.49	1.17	1.33
3	A	1007	7GE	C11-C10	-6.47	1.20	1.38
3	A	1008	7GE	C11-C10	-5.58	1.23	1.38
3	A	1007	7GE	C12-C13	4.73	1.50	1.41
3	A	1008	7GE	C12-C13	4.19	1.49	1.41
3	A	1008	7GE	C11-C12	-3.68	1.28	1.36
3	A	1007	7GE	C09-C08	3.24	1.49	1.41
3	A	1007	7GE	C11-C12	-2.82	1.30	1.36
3	A	1007	7GE	C10-C09	-2.78	1.30	1.36
3	A	1008	7GE	C10-C09	-2.72	1.30	1.36
3	A	1008	7GE	C09-C08	2.57	1.48	1.41
3	A	1008	7GE	C07-C08	-2.15	1.36	1.40

There are no bond angle outliers.

There are no chirality outliers.

All (24) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1004	PLM	C1-C2-C3-C4
2	A	1003	PLM	C1-C2-C3-C4
2	A	1001	PLM	C1-C2-C3-C4
2	A	1005	PLM	C1-C2-C3-C4
2	A	1005	PLM	CA-CB-CC-CD
2	A	1006	PLM	C5-C6-C7-C8
2	A	1004	PLM	C2-C3-C4-C5
2	A	1003	PLM	CA-CB-CC-CD
2	A	1006	PLM	C4-C5-C6-C7
2	A	1002	PLM	CA-CB-CC-CD
2	A	1006	PLM	C2-C3-C4-C5
2	A	1001	PLM	C6-C7-C8-C9
2	A	1003	PLM	CB-CC-CD-CE
2	A	1006	PLM	C6-C7-C8-C9
2	A	1001	PLM	C4-C5-C6-C7
2	A	1002	PLM	CC-CD-CE-CF
2	A	1003	PLM	CC-CD-CE-CF
2	A	1006	PLM	CC-CD-CE-CF
2	A	1005	PLM	CB-CC-CD-CE
2	A	1005	PLM	C9-CA-CB-CC
2	A	1001	PLM	C8-C9-CA-CB
2	A	1004	PLM	CB-CC-CD-CE
2	A	1004	PLM	C9-CA-CB-CC
2	A	1006	PLM	C3-C4-C5-C6

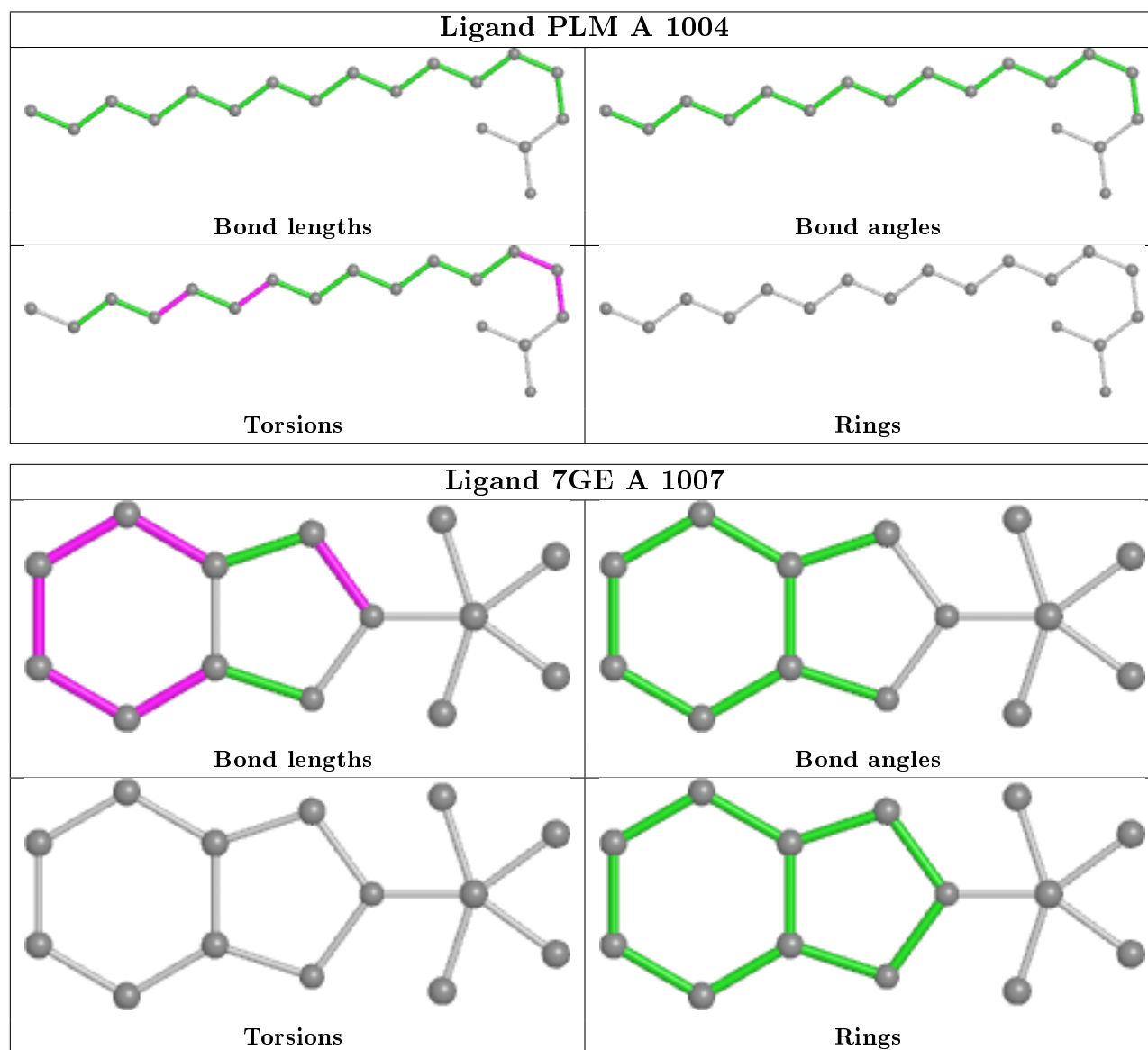
There are no ring outliers.

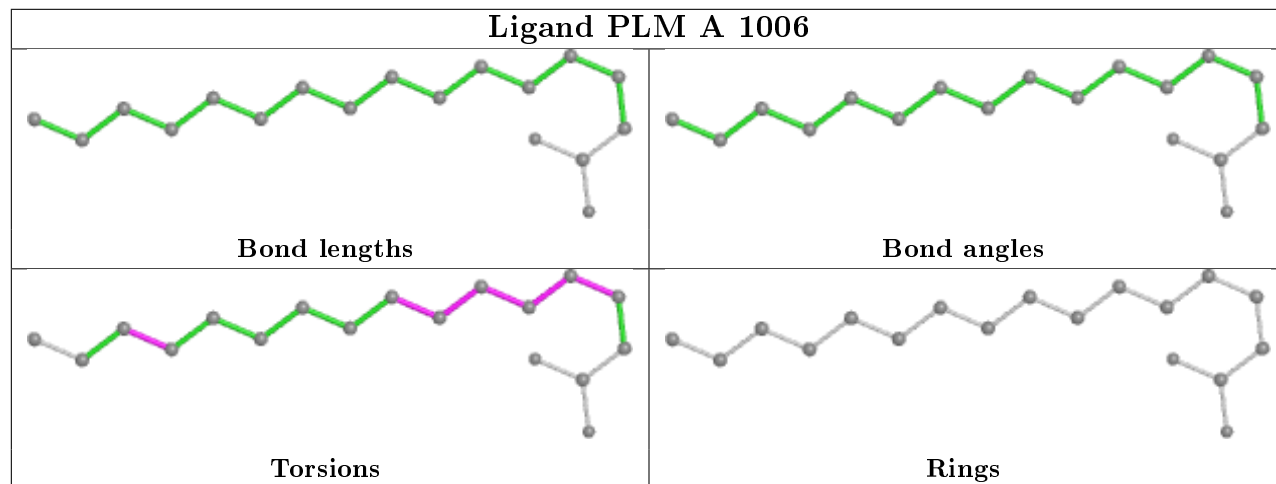
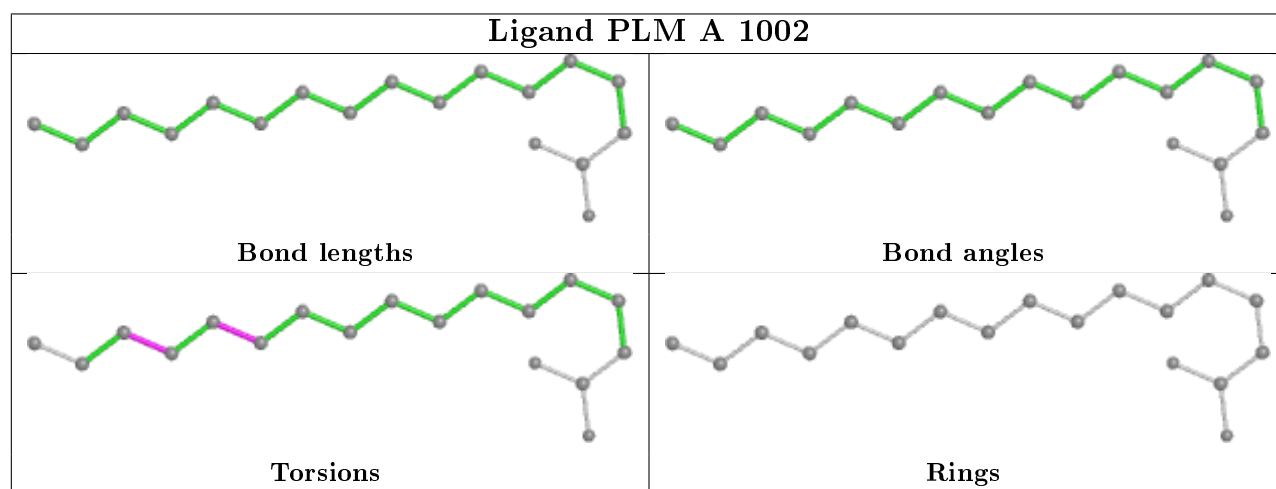
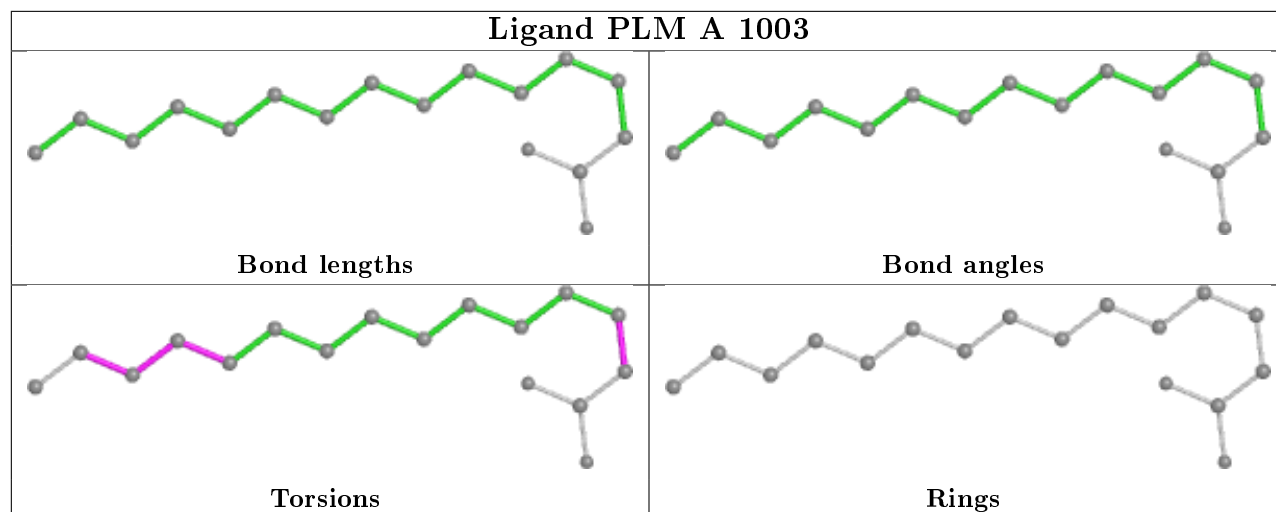
7 monomers are involved in 25 short contacts:

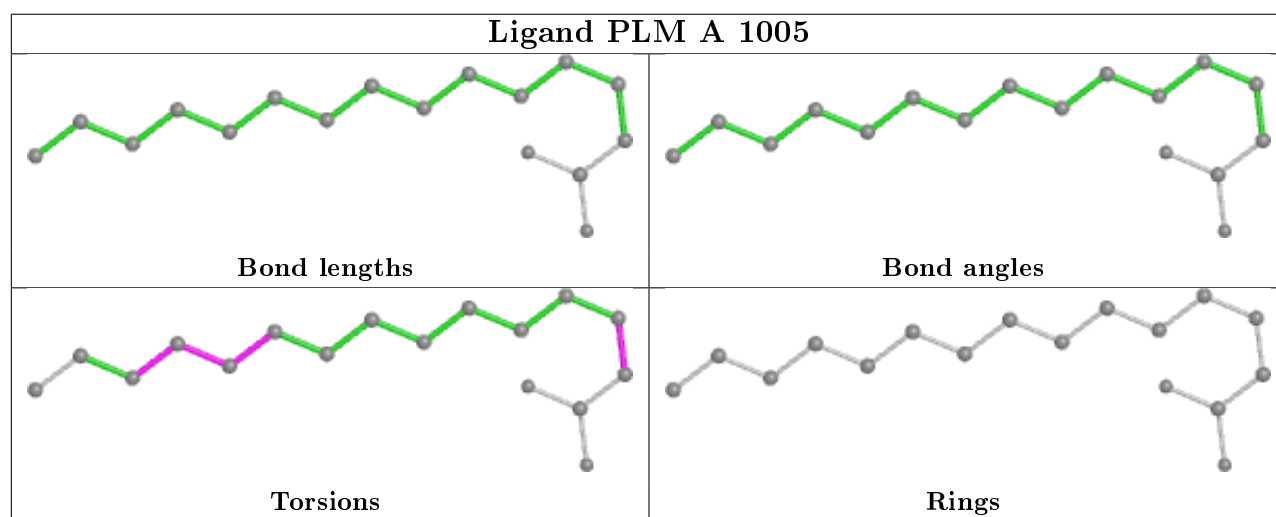
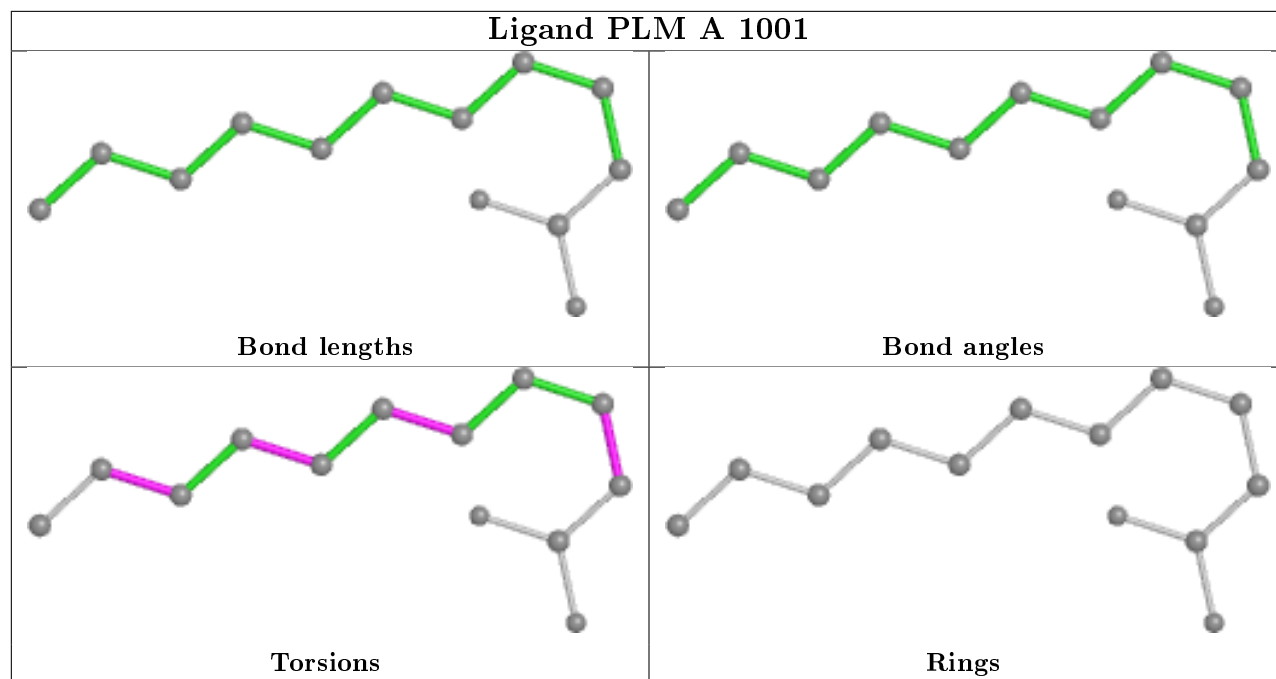
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1004	PLM	9	0
3	A	1007	7GE	9	0
2	A	1003	PLM	2	0
2	A	1002	PLM	2	0
2	A	1006	PLM	1	0
2	A	1005	PLM	1	0
3	A	1008	7GE	3	0

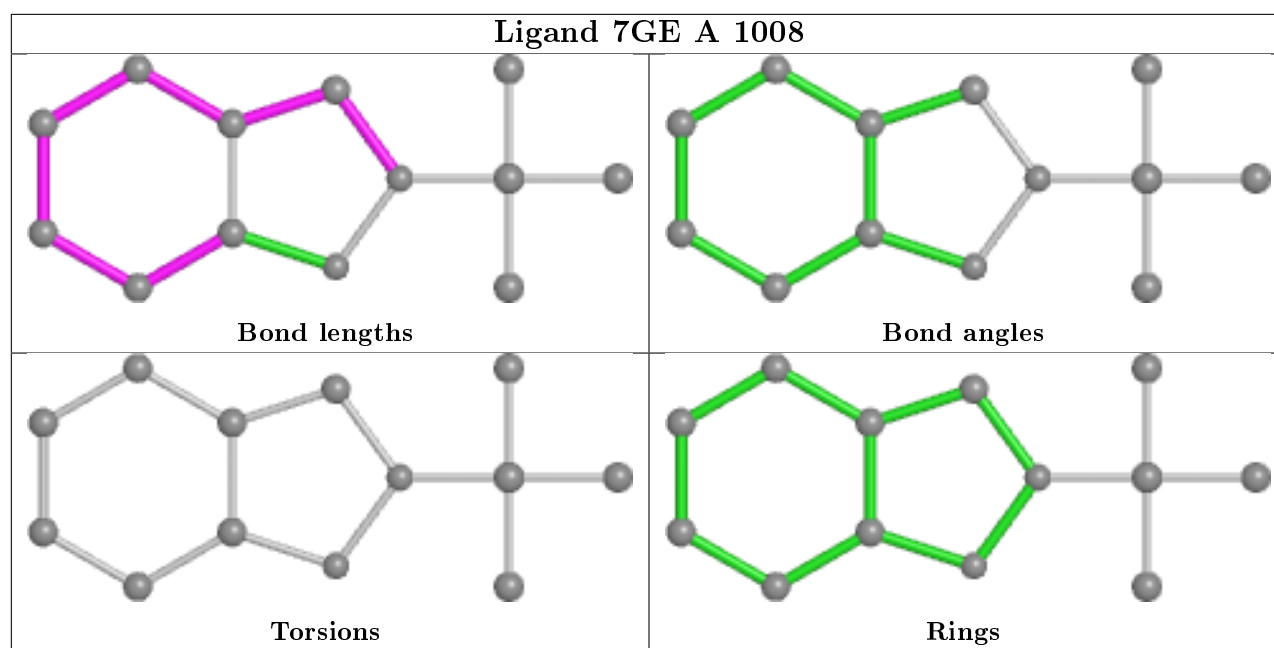
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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









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	581/581 (100%)	0.35	24 (4%)	37 44	46, 75, 112, 150	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	557	LYS	4.7
1	A	301	ASP	4.2
1	A	370	TYR	3.7
1	A	3	HIS	3.6
1	A	363	ALA	3.5
1	A	372	LYS	3.3
1	A	373	VAL	3.2
1	A	577	ALA	3.0
1	A	555	VAL	2.9
1	A	573	LYS	2.9
1	A	554	PHE	2.9
1	A	311	GLU	2.7
1	A	576	VAL	2.7
1	A	369	CYS	2.7
1	A	481	LEU	2.6
1	A	31	LEU	2.6
1	A	362	ALA	2.5
1	A	4	LYS	2.4
1	A	196	GLN	2.4
1	A	83	THR	2.3
1	A	566	THR	2.1
1	A	366	PRO	2.1
1	A	570	GLU	2.1
1	A	482	VAL	2.0

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

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

6.3 Carbohydrates ⓘ

There are no monosaccharides in this entry.

6.4 Ligands ⓘ

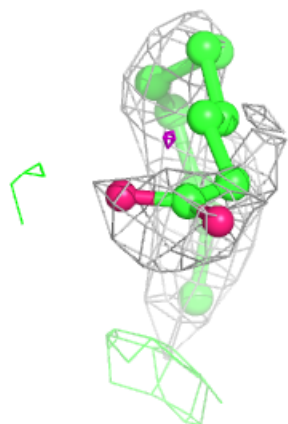
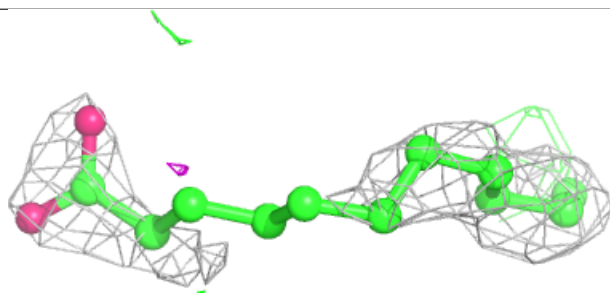
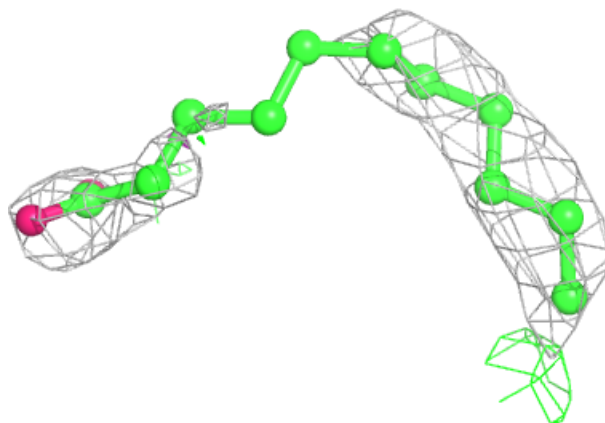
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	PLM	A	1001	13/18	0.79	0.46	55,79,101,104	0
2	PLM	A	1005	17/18	0.80	0.41	69,81,93,95	0
2	PLM	A	1006	18/18	0.81	0.47	52,70,114,119	0
2	PLM	A	1002	18/18	0.87	0.35	57,79,85,90	0
2	PLM	A	1003	17/18	0.91	0.31	55,71,81,86	0
2	PLM	A	1004	18/18	0.92	0.36	64,81,94,96	0
3	7GE	A	1008	13/15	0.93	0.18	131,137,165,168	0
3	7GE	A	1007	14/15	0.96	0.19	93,103,141,182	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.

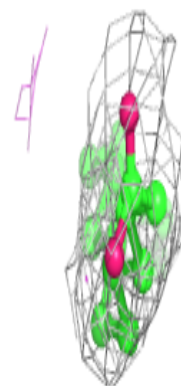
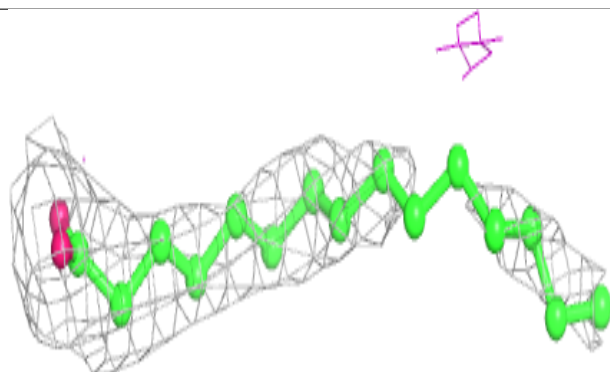
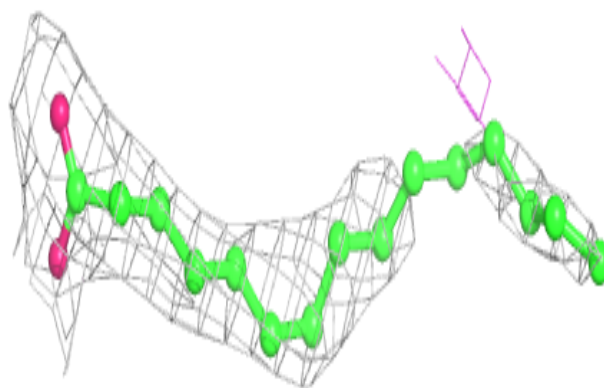
Electron density around PLM A 1001:

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

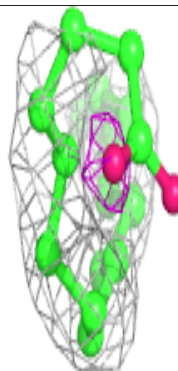
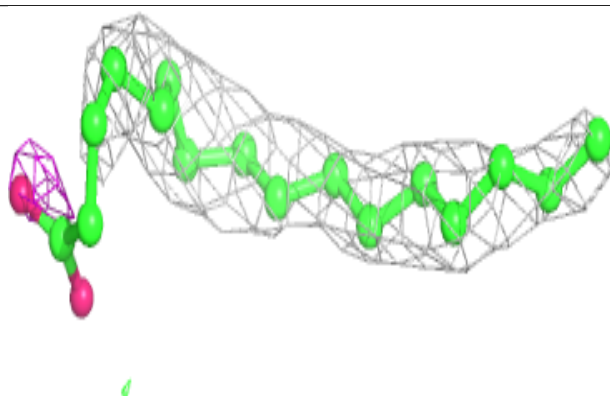
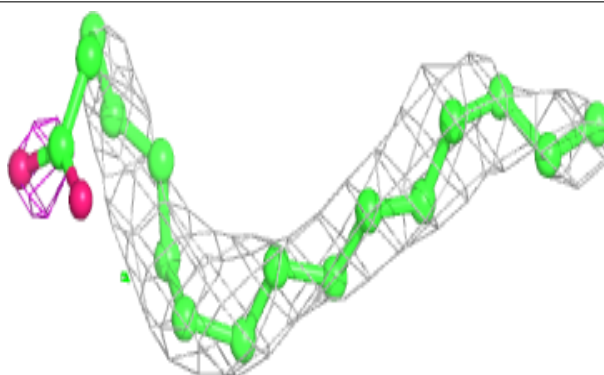


Electron density around PLM A 1005:

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

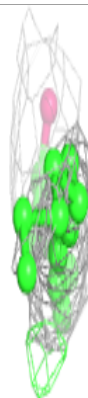
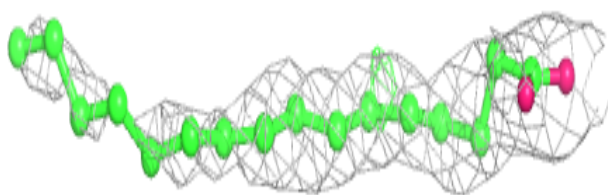
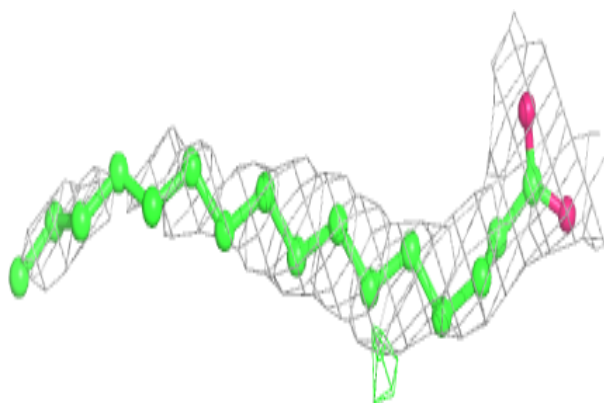
**Electron density around PLM A 1006:**

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

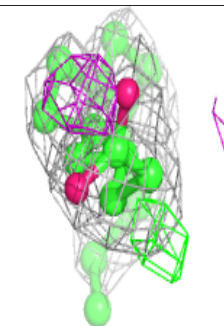
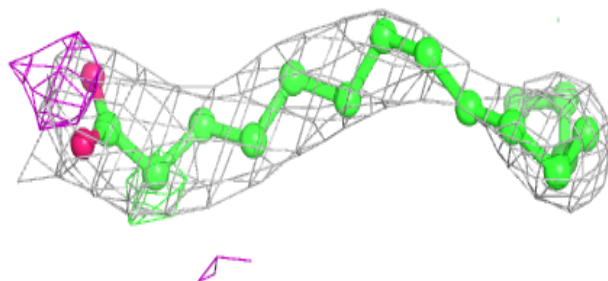
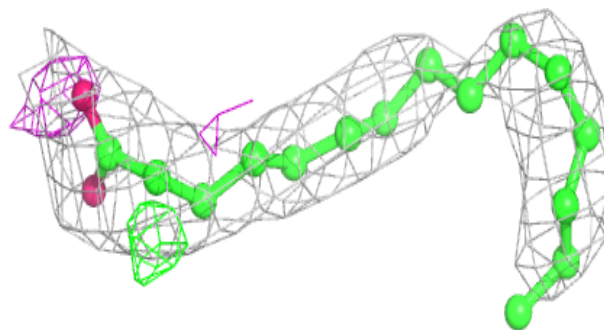


Electron density around PLM A 1002:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

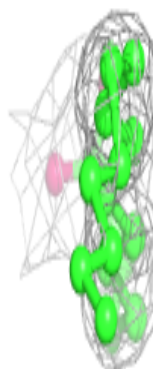
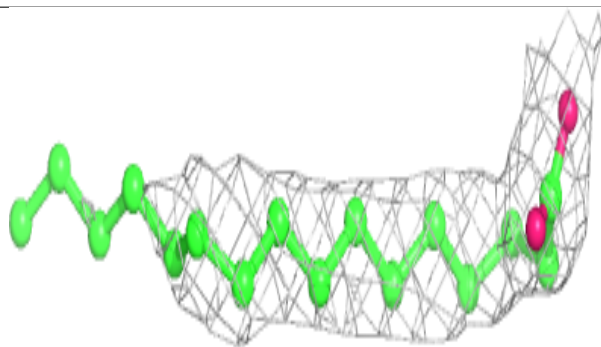
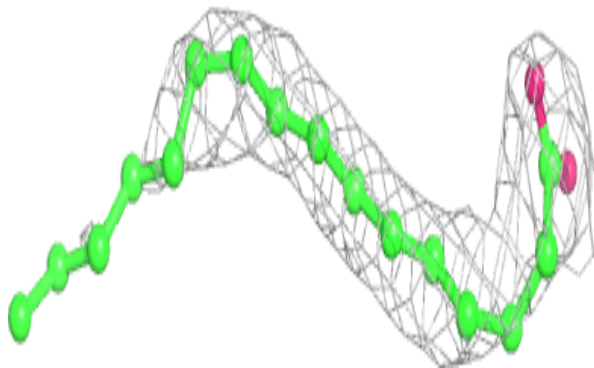
**Electron density around PLM A 1003:**

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



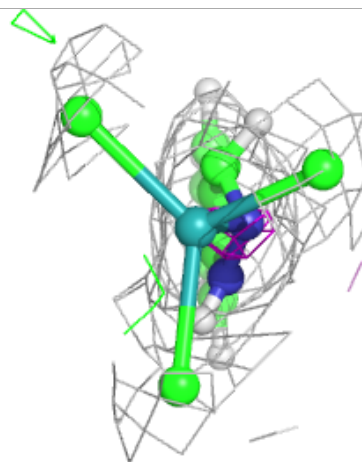
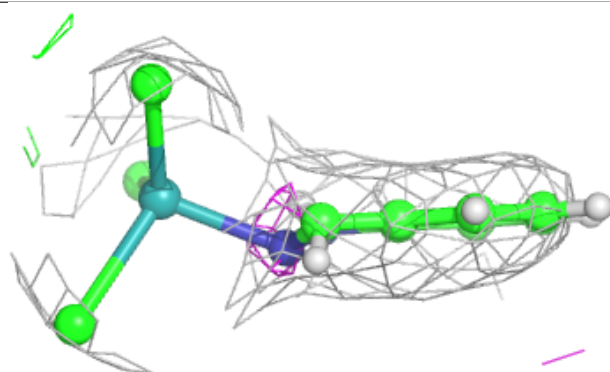
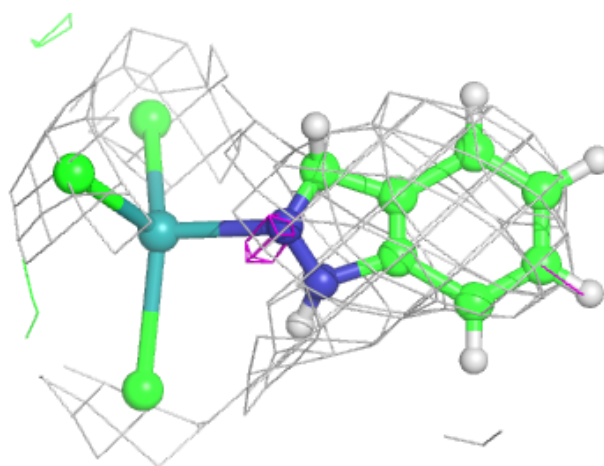
Electron density around PLM A 1004:

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



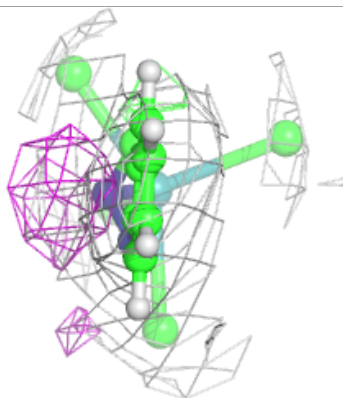
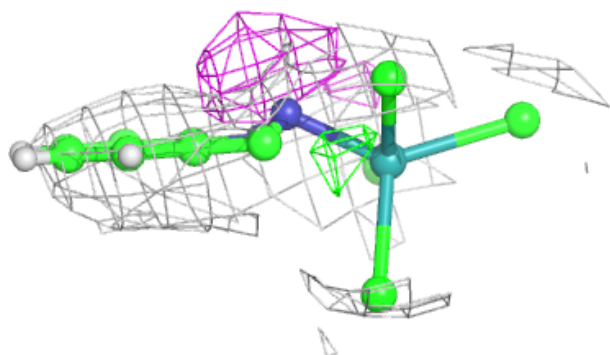
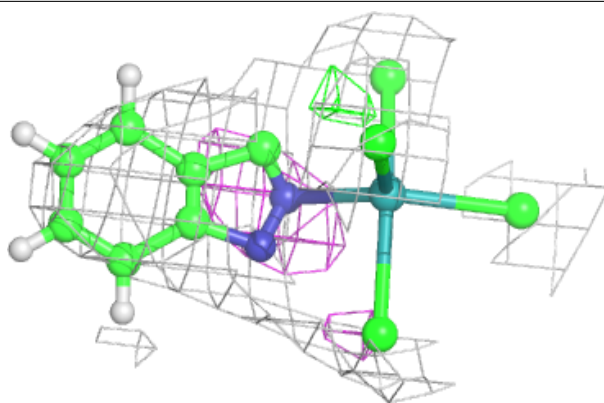
Electron density around 7GE A 1008:

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



Electron density around 7GE A 1007:

$2mF_o-DF_c$ (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.