

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM060319\
 Data File : BM020697.D
 Acq On : 04 Jun 2019 09:30
 Operator : HP/JU
 Sample : K3139-07
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 TP-1A-D

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM053119.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.057	8	12	25	rVB	2313602	2836311	18.27%	2.848%
2	5.410	405	412	425	rBV	5879588	8359277	53.85%	8.395%
3	6.992	672	681	692	rBV	5805568	9273400	59.74%	9.312%
4	7.357	735	743	753	rBV	5848634	9417406	60.67%	9.457%
5	7.828	815	823	830	rBV	1143427	1833299	11.81%	1.841%
6	8.145	869	877	885	rBV	4392850	7068840	45.54%	7.099%
7	8.986	1010	1020	1030	rBV	3460470	5716762	36.83%	5.741%
8	10.616	1287	1297	1305	rBV	1470425	2419674	15.59%	2.430%
9	13.080	1709	1716	1730	rBV	8605213	12573557	81.00%	12.627%
10	13.916	1851	1858	1868	rBV	472284	682840	4.40%	0.686%
11	14.463	1944	1951	1961	rBV2	1939043	2974495	19.16%	2.987%
12	15.951	2197	2204	2218	rBV	6023449	8677603	55.90%	8.714%
13	17.204	2410	2417	2430	rBV	2377131	3344895	21.55%	3.359%
14	18.092	2563	2568	2578	rBV	244371	386781	2.49%	0.388%
15	19.827	2857	2863	2869	rBV	12424135	15522585	100.00%	15.588%
16	21.092	3074	3078	3088	rBV	634993	792545	5.11%	0.796%
17	21.380	3122	3127	3133	rBV	2792770	3445921	22.20%	3.460%
18	22.450	3306	3309	3314	rVB	137077	173356	1.12%	0.174%
19	22.974	3395	3398	3404	rVB2	142499	185887	1.20%	0.187%
20	23.562	3494	3498	3505	rVB	160387	265531	1.71%	0.267%
21	23.668	3509	3516	3523	rVB	1942308	3629262	23.38%	3.645%

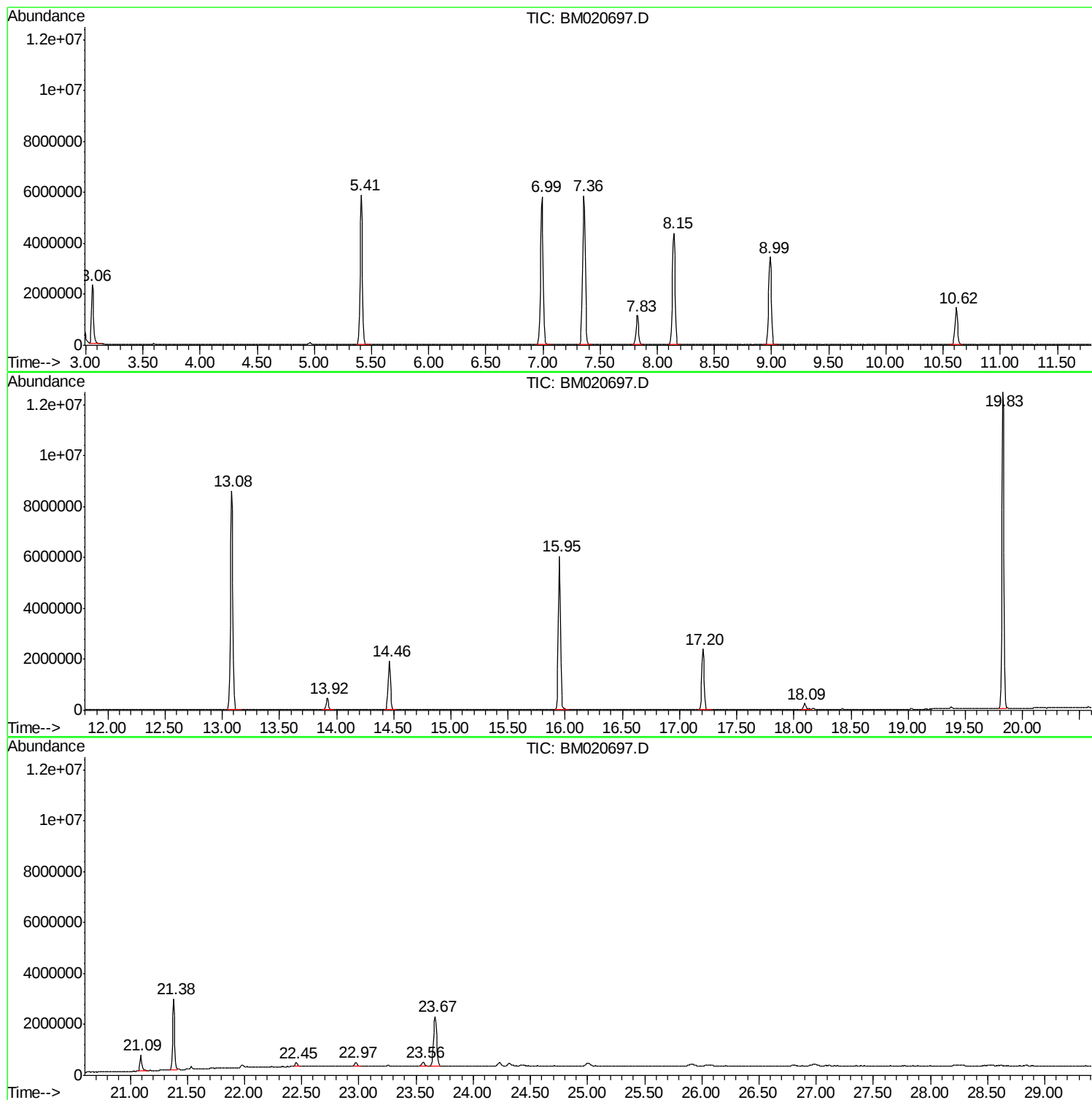
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM053119.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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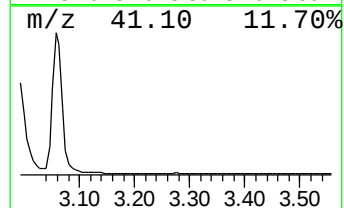
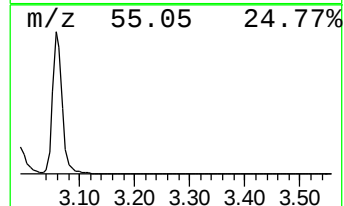
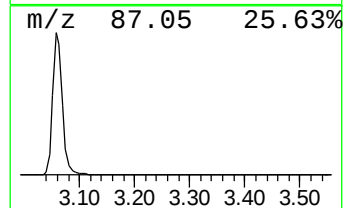
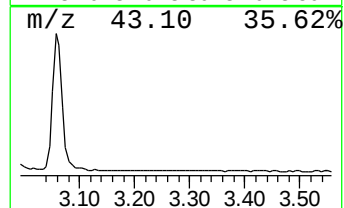
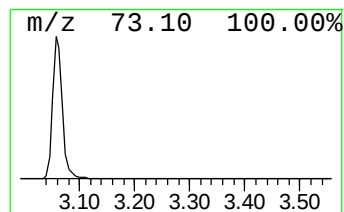
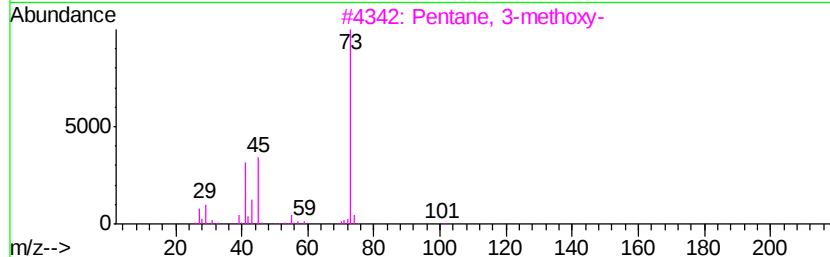
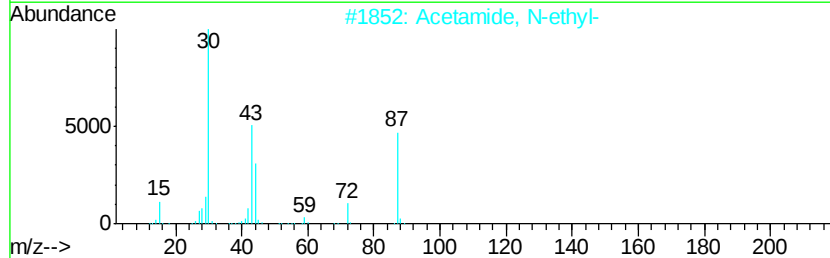
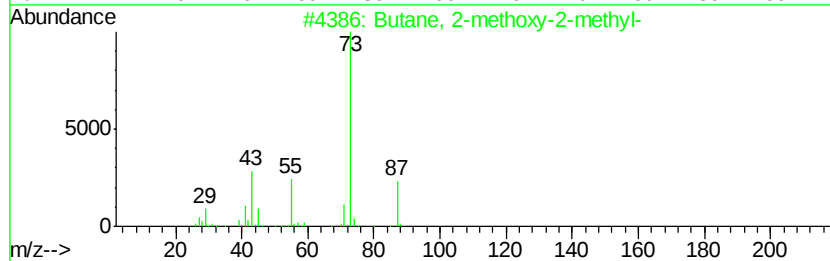
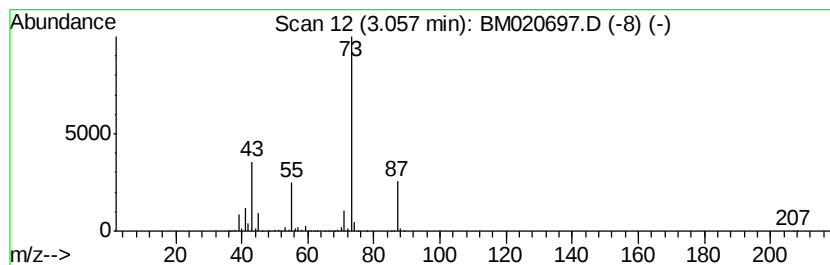
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM053119.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.06	30.94 ng	2836310	1,4-Dichlorobenzene-d4	7.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	22
3		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
4		Silane, tetramethyl-	88	C4H12Si	000075-76-3	12
5		N-Ethylformamide	73	C3H7NO	000627-45-2	9



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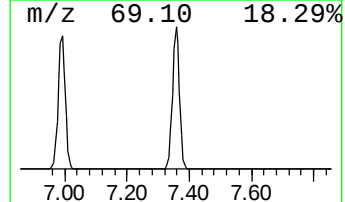
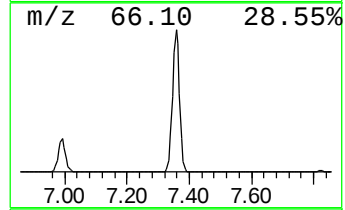
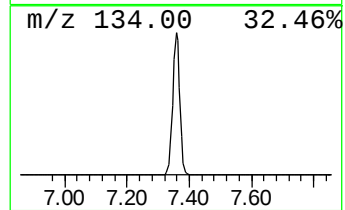
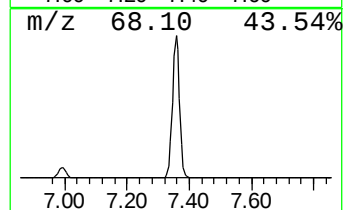
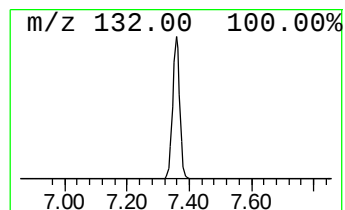
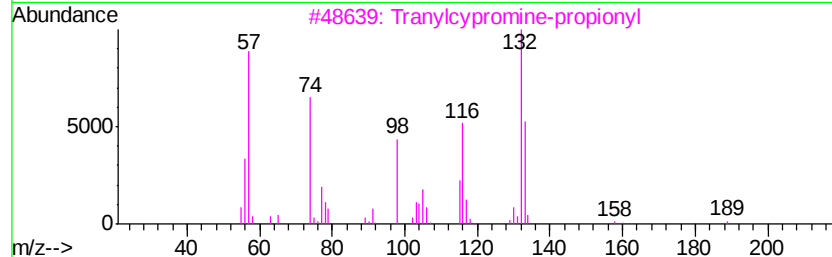
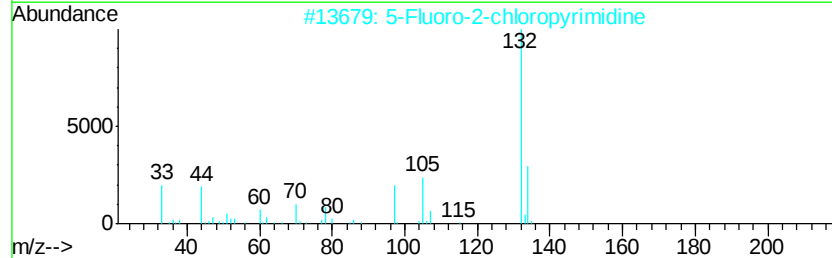
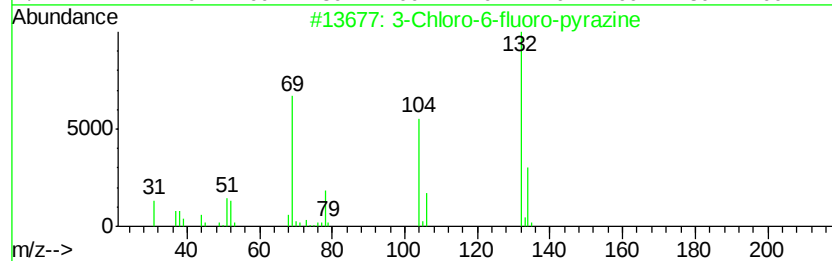
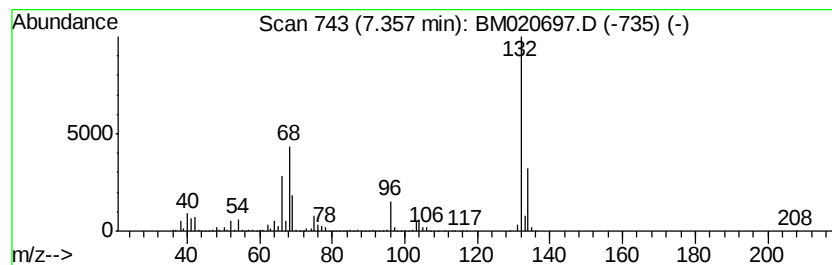
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Peak Number 2 unknown7.36 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.36	102.74 ng	9417410	1,4-Dichlorobenzene-d4	7.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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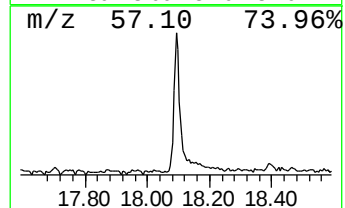
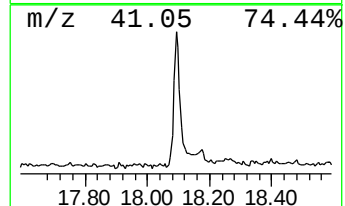
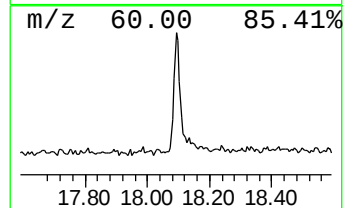
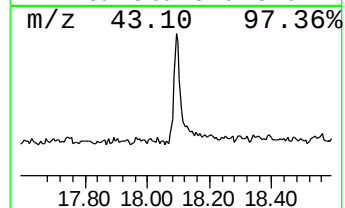
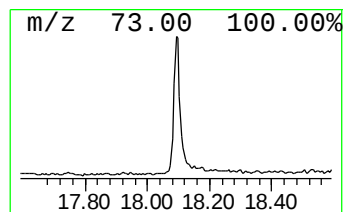
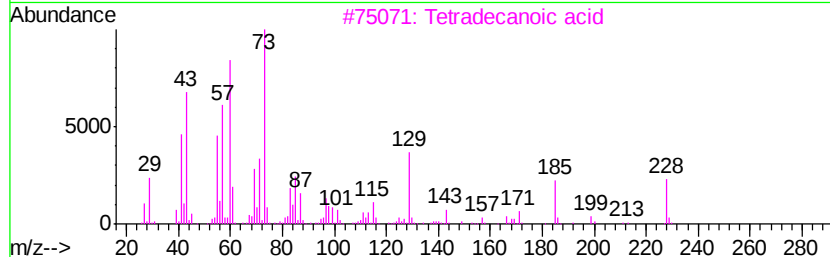
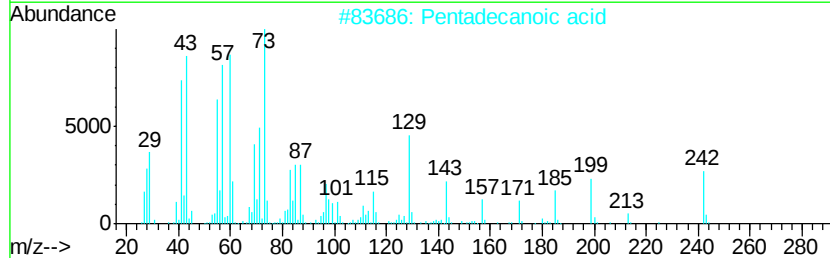
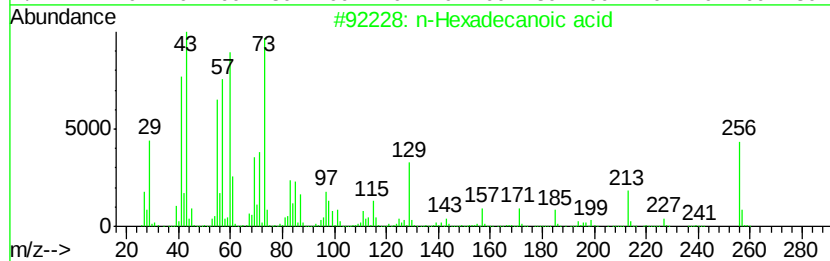
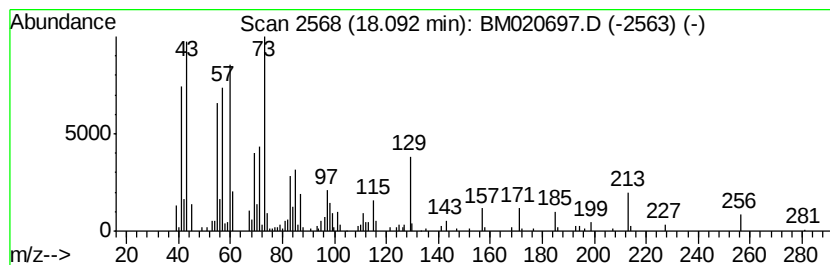
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Peak Number 4 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.09	2.31 ng	386781	Phenanthrene-d10	17.20

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	83
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	80
4		Tridecanoic acid	214	C13H26O2	000638-53-9	76
5		n-Decanoic acid	172	C10H20O2	000334-48-5	76



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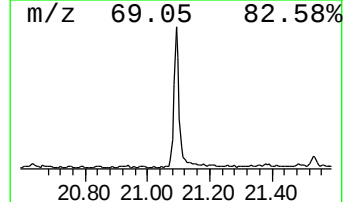
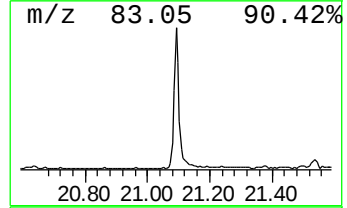
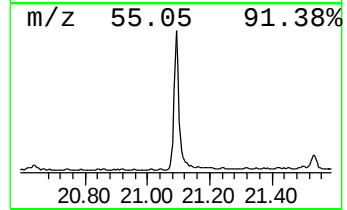
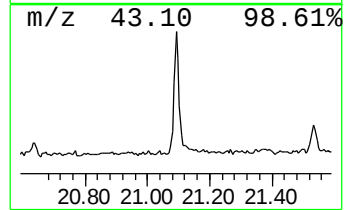
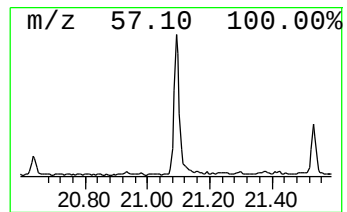
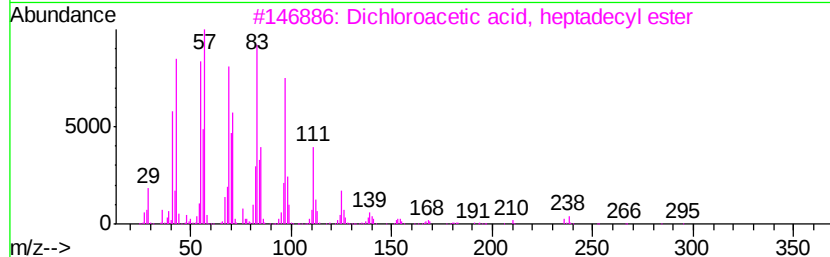
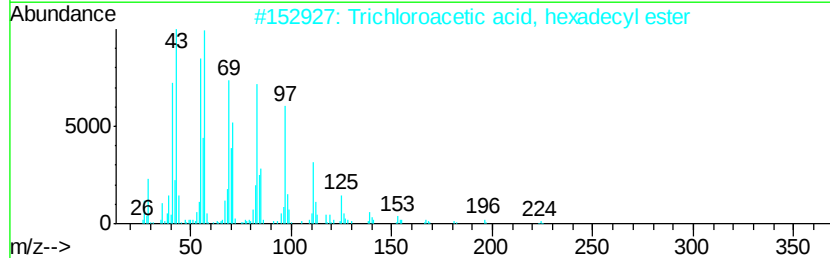
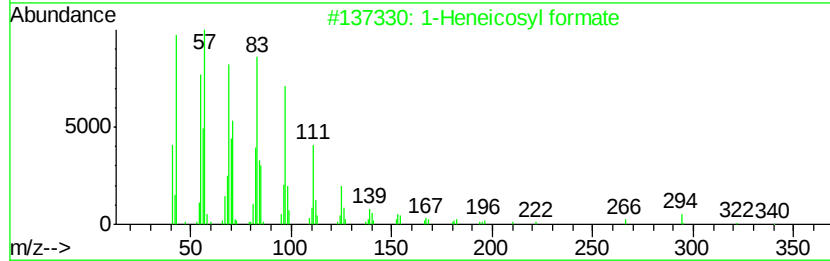
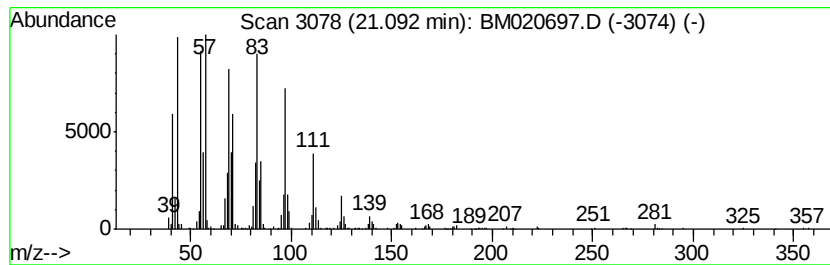
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Peak Number 5 1-Heneicosyl formate Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.09	4.60 ng	792545	Chrysene-d12	21.38

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosyl formate		340	C22H44O2	077899-03-7	94
2	Trichloroacetic acid, hexadecyl ...		386	C18H33Cl3O2	074339-54-1	94
3	Dichloroacetic acid, heptadecyl ...		366	C19H36Cl2O2	1000282-98-2	94
4	Trifluoroacetic acid, n-octadecyl ...		366	C20H37F3O2	079392-43-1	91
5	17-Pentatriacontene		491	C35H70	006971-40-0	91



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	3.06	30.9	ng	2836310	1	7.83	1833300	20.0
unknown7.36	7.36	102.7	ng	9417410	1	7.83	1833300	20.0
n-Hexadecanoic acid	18.09	2.3	ng	386781	4	17.20	3344900	20.0
1-Heneicosyl formate	21.09	4.6	ng	792545	5	21.38	3445920	20.0