

Data Path : U:\HPCHEM1\BNA M\DATA\BM020218\
 Data File : BM014141.D
 Acq On : 03 Feb 2018 07:41
 Operator : SJ/JU
 Sample : J1339-04
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 F6C94

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 1 % 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:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM013118.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	6.739	633	638	654	rBV	356672	693093	23.00%	2.199%
2	6.898	660	665	673	rBV	295738	470634	15.62%	1.493%
3	7.087	691	697	716	rVB2	447460	763700	25.35%	2.423%
4	7.551	769	776	789	rBV	413684	712324	23.64%	2.260%
5	8.269	892	898	914	rBV	458854	873337	28.99%	2.770%
6	8.710	966	973	985	rBV	426184	758289	25.17%	2.405%
7	9.422	1088	1094	1104	rBV	361513	695212	23.08%	2.205%
8	9.963	1179	1186	1196	rBV2	477216	930300	30.88%	2.951%
9	10.322	1240	1247	1255	rBV	606170	973046	32.30%	3.087%
10	10.480	1268	1274	1286	rBV	374510	814013	27.02%	2.582%
11	13.627	1803	1809	1814	rBV	1052844	1532490	50.87%	4.861%
12	13.674	1814	1817	1825	rVB	142610	233518	7.75%	0.741%
13	13.898	1848	1855	1869	rBV	1173189	1801844	59.81%	5.716%
14	14.110	1886	1891	1896	rBV3	35772	50663	1.68%	0.161%
15	14.204	1901	1907	1918	rVV3	867620	1370009	45.47%	4.346%
16	14.457	1945	1950	1964	rBV2	180851	537272	17.83%	1.704%
17	15.068	2049	2054	2059	rVB3	44340	58568	1.94%	0.186%
18	15.210	2071	2078	2089	rBV	1498856	2221678	73.74%	7.047%
19	15.357	2098	2103	2113	rBV2	338950	598562	19.87%	1.899%
20	15.457	2116	2120	2129	rVB8	23587	57109	1.90%	0.181%
21	15.933	2197	2201	2209	rBV2	41055	72427	2.40%	0.230%
22	16.727	2332	2336	2339	rBV2	24519	35688	1.18%	0.113%
23	16.962	2370	2376	2385	rBV	1208677	1699167	56.40%	5.390%
24	17.062	2385	2393	2408	rVV	1652169	2478273	82.26%	7.861%
25	17.651	2487	2493	2496	rBV6	22602	33963	1.13%	0.108%
26	17.868	2526	2530	2540	rBV2	195072	333733	11.08%	1.059%
27	19.009	2719	2724	2728	rBV4	28308	48049	1.59%	0.152%
28	19.162	2746	2750	2761	rBV3	128683	244637	8.12%	0.776%
29	19.262	2762	2767	2770	rVB4	21838	37659	1.25%	0.119%
30	19.368	2779	2785	2798	rBV	2030423	2881607	95.65%	9.141%
31	20.886	3039	3043	3052	rBV	329653	470315	15.61%	1.492%
32	21.168	3086	3091	3099	rBV2	1586767	2104018	69.84%	6.674%
33	23.209	3431	3438	3449	rVB2	1641698	3012796	100.00%	9.557%
34	23.344	3455	3461	3470	rVB2	1007559	1927080	63.96%	6.113%

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Stop Thrs : 0 Peak Location: TOP

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Peak separation: 5

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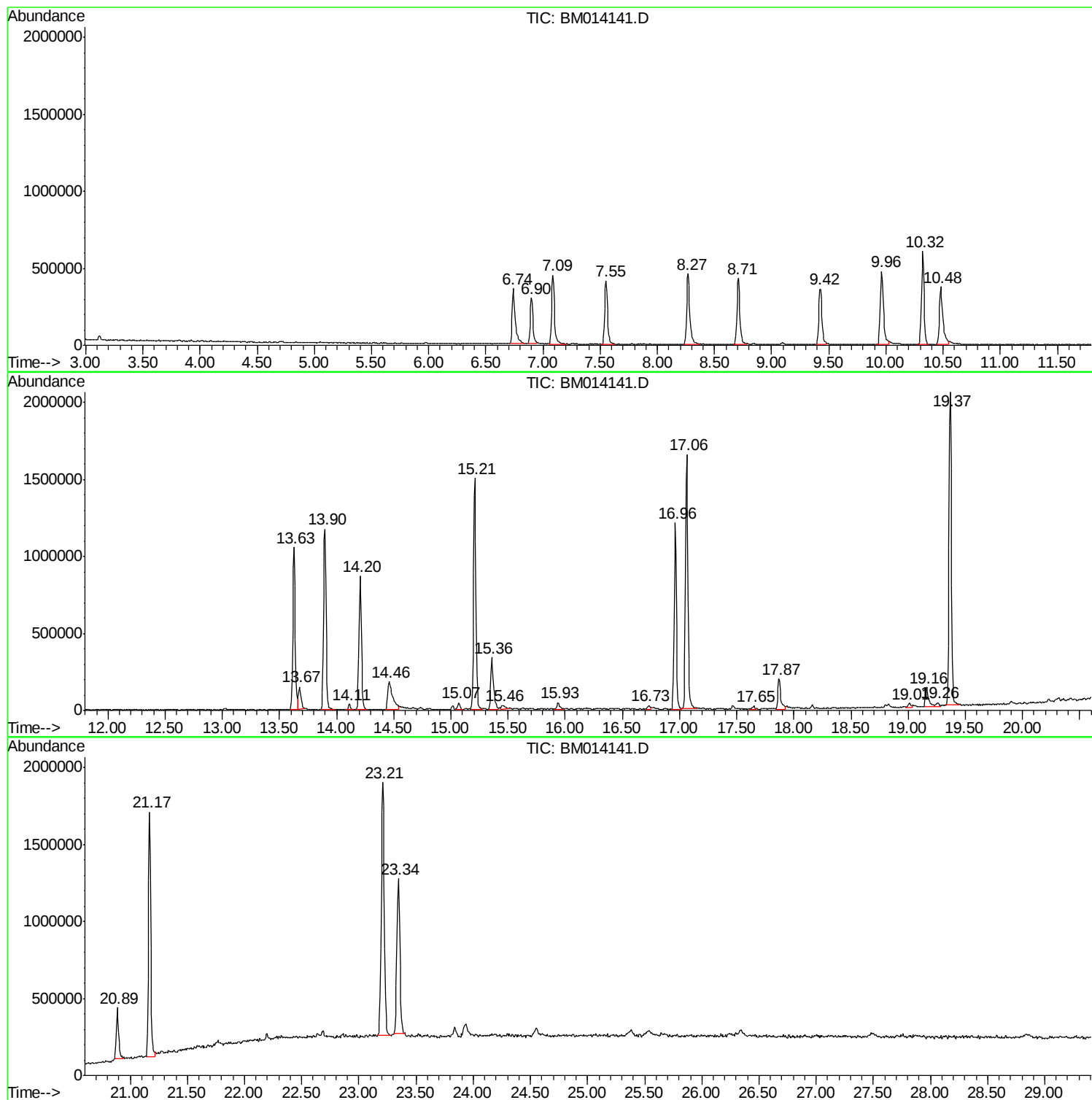
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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM013118.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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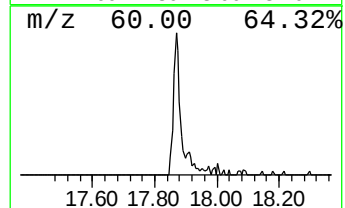
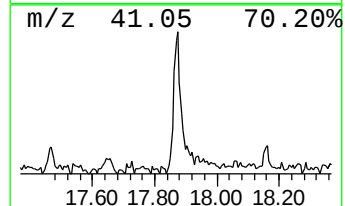
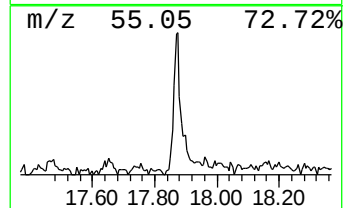
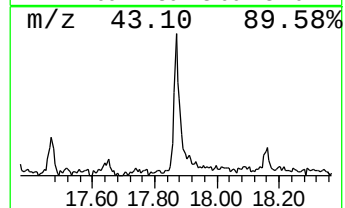
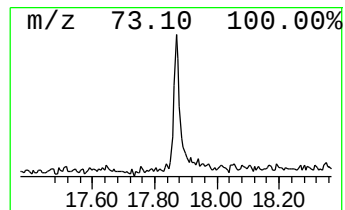
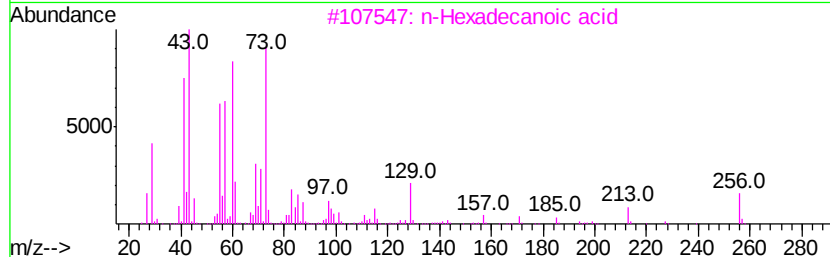
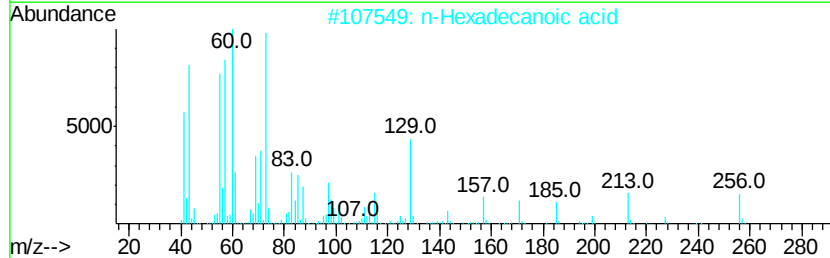
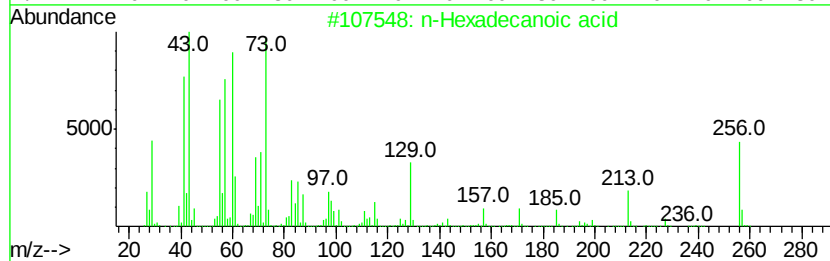
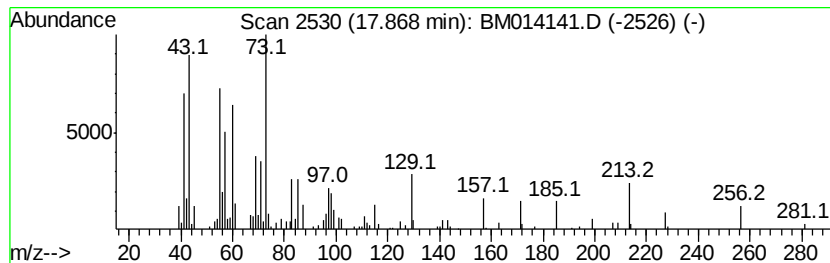
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.87	3.93 ng/ul	333733	Phenanthrene-d10	16.96

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	68	
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	60	
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	43	
4	Tridecanoic acid	214	C13H26O2	000638-53-9	41	
5	n-Decanoic acid	172	C10H20O2	000334-48-5	38	



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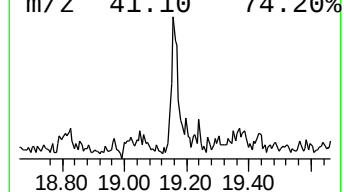
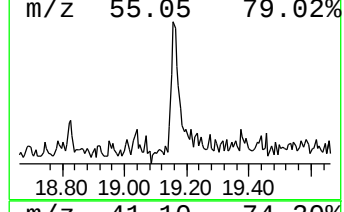
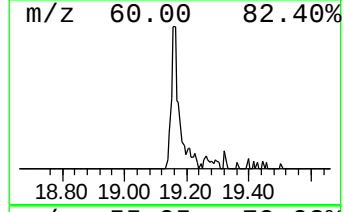
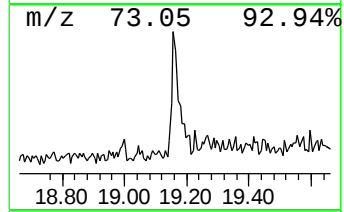
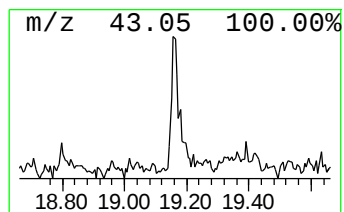
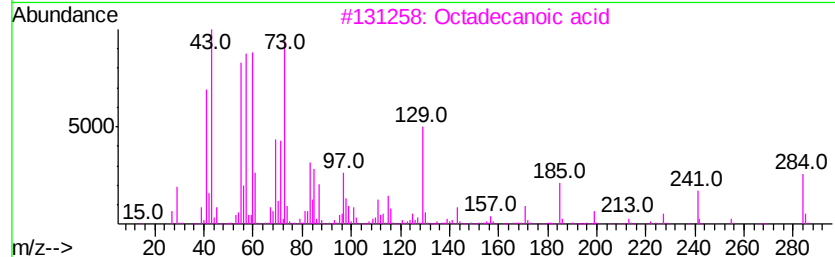
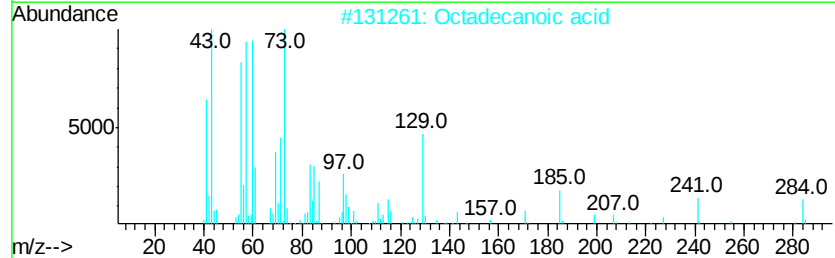
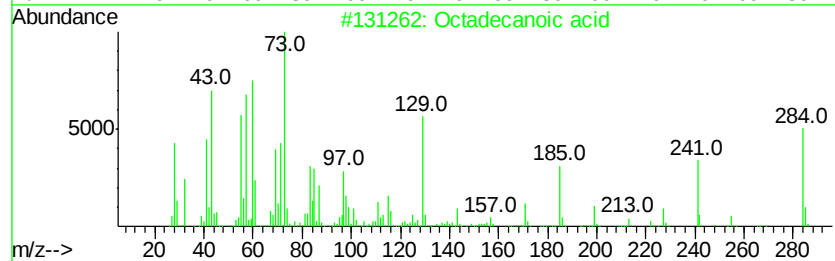
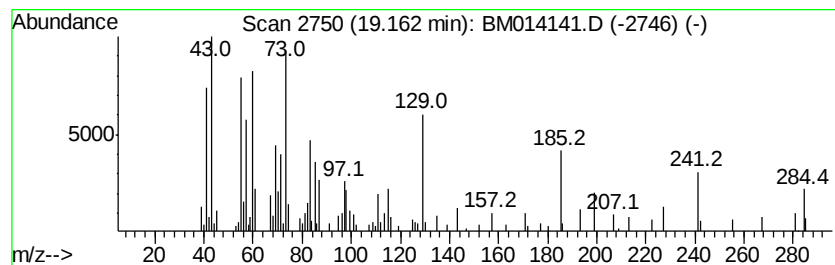
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Peak Number 2 Octadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.16	2.33 ng/ul	244637	Chrysene-d12	21.17

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	98
2			Octadecanoic acid	284	C18H36O2	000057-11-4	93
3			Octadecanoic acid	284	C18H36O2	000057-11-4	83
4			Octadecanoic acid	284	C18H36O2	000057-11-4	62
5			Octadecanoic acid	284	C18H36O2	000057-11-4	50



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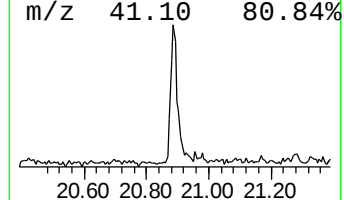
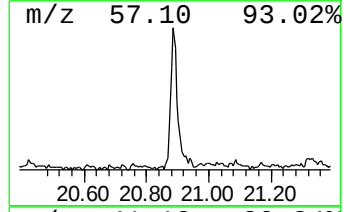
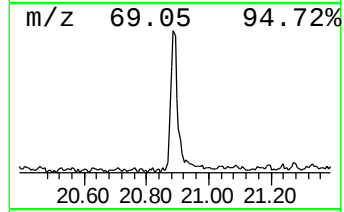
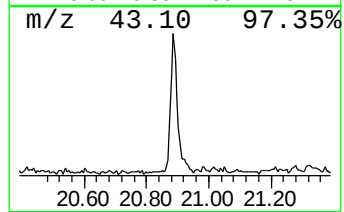
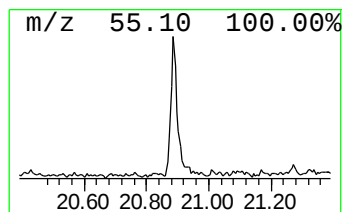
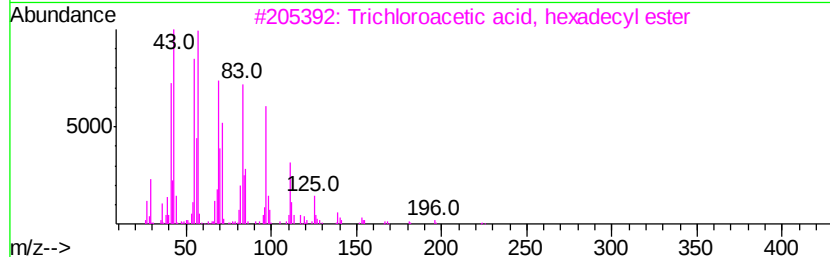
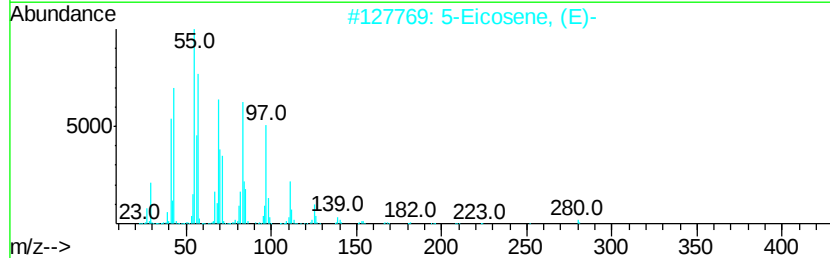
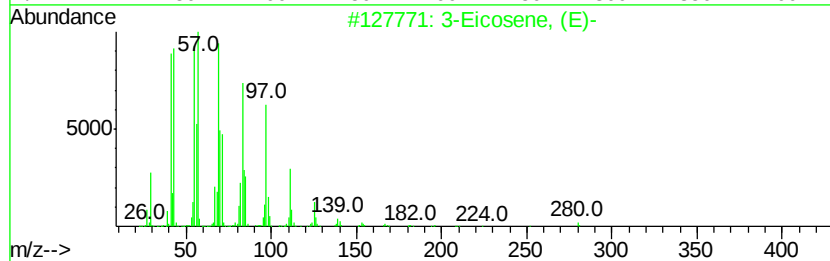
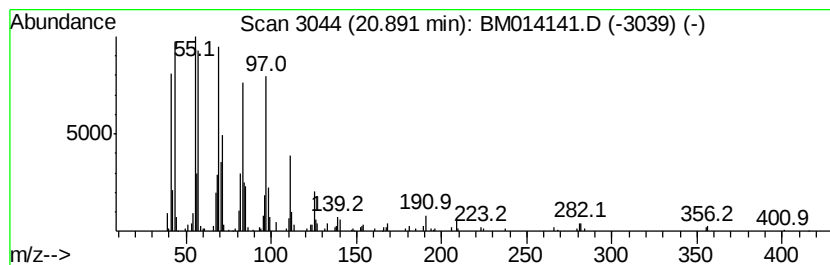
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Peak Number 3 (DEL) Alkane: Cyclic20.89 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.89	4.47 ng/ul	470315	Chrysene-d12	21.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Eicosene, (E)-	280	C20H40	074685-33-9	98
2		5-Eicosene, (E)-	280	C20H40	074685-30-6	96
3		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
4		9-Eicosene, (E)-	280	C20H40	074685-29-3	93
5		Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	91



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
n-Hexadecanoic acid	17.87	3.9	ng/ul	333733	4	16.96	1699170	20.0
Octadecanoic acid	19.16	2.3	ng/ul	244637	5	21.17	2104020	20.0
(DEL) Alkane: Cyc...	20.89	4.5	ng/ul	470315	5	21.17	2104020	20.0