

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071023\
 Data File : BP016169.D
 Acq On : 10 Jul 2023 22:39
 Operator : MA/JU
 Sample : 03404-10
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 YBW46

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP016169.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.340	23	26	30	rBV	34454	46588	0.90%	0.096%
2	3.799	103	104	132	rBV	222687	576868	11.12%	1.186%
3	3.993	135	137	140	rBV3	7705	8025	0.15%	0.017%
4	4.528	226	228	233	rVB4	10716	12091	0.23%	0.025%
5	6.893	623	630	633	rBV9	3494	6935	0.13%	0.014%
6	7.069	657	660	666	rVB8	3737	6285	0.12%	0.013%
7	7.234	680	688	703	rBV	257233	995891	19.20%	2.048%
8	7.352	703	708	738	rVB	289263	925351	17.84%	1.903%
9	7.563	738	744	763	rBV	381253	1022089	19.71%	2.102%
10	8.022	814	822	847	rBV	294093	1005555	19.39%	2.068%
11	8.793	943	953	987	rBV	295079	1447420	27.91%	2.977%
12	9.240	1021	1029	1056	rBV	316224	1106832	21.34%	2.276%
13	9.522	1075	1077	1081	rVB5	8692	9645	0.19%	0.020%
14	9.981	1147	1155	1178	rBV3	179320	915325	17.65%	1.883%
15	10.575	1249	1256	1273	rBV2	226944	1063352	20.50%	2.187%
16	10.852	1296	1303	1328	rVB	541163	1628254	31.40%	3.349%
17	11.075	1333	1341	1378	rBV4	176315	1008551	19.45%	2.074%
18	12.016	1498	1501	1507	rVB8	5011	8420	0.16%	0.017%
19	12.516	1576	1586	1596	rBV9	7036	24240	0.47%	0.050%
20	13.040	1670	1675	1677	rBV5	4127	6440	0.12%	0.013%
21	13.551	1755	1762	1766	rBV6	20724	42088	0.81%	0.087%
22	13.681	1777	1784	1786	rBV8	3627	6600	0.13%	0.014%
23	14.004	1836	1839	1843	rVB6	3871	5192	0.10%	0.011%
24	14.093	1847	1854	1878	rBV	1001412	2447343	47.19%	5.033%
25	14.369	1893	1901	1924	rBV	1508228	2937332	56.64%	6.041%
26	14.540	1928	1930	1937	rVB8	9327	17791	0.34%	0.037%
27	14.669	1945	1952	1972	rBV	1303957	2301856	44.39%	4.734%
28	15.092	2016	2024	2025	rBV4	60754	128649	2.48%	0.265%
29	15.498	2088	2093	2096	rVB7	8225	11743	0.23%	0.024%
30	15.681	2117	2124	2153	rBV	1563076	3814949	73.56%	7.846%
31	15.910	2156	2163	2164	rBV2	99232	173928	3.35%	0.358%
32	16.057	2183	2188	2208	rVB	389444	813760	15.69%	1.674%
33	16.310	2227	2231	2234	rVB6	9127	12553	0.24%	0.026%
34	16.610	2279	2282	2285	rBV4	11908	14757	0.28%	0.030%
35	16.916	2331	2334	2339	rBV6	7695	8730	0.17%	0.018%
36	17.092	2361	2364	2370	rBV8	7409	15223	0.29%	0.031%

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37	17.134	2370	2371	2376	rVB5	7601	7239	0.14%	0.015%
38	17.186	2376	2380	2383	rBV6	4209	6740	0.13%	0.014%
39	17.434	2415	2422	2434	rBV	1725407	2804734	54.08%	5.768%
40	17.545	2434	2441	2468	rVB2	1439458	4154431	80.11%	8.544%
41	18.286	2562	2567	2578	rBV2	272064	593618	11.45%	1.221%
42	19.351	2744	2748	2750	rBV2	35505	48100	0.93%	0.099%
43	19.516	2772	2776	2790	rBV3	105921	216230	4.17%	0.445%
44	19.769	2813	2819	2850	rBV2	2396634	5186073	100.00%	10.666%
45	20.710	2977	2979	2983	rBV	53664	53984	1.04%	0.111%
46	21.163	3054	3056	3059	rBV2	82745	78268	1.51%	0.161%
47	21.216	3059	3065	3078	rVV	348771	1041456	20.08%	2.142%
48	21.545	3116	3121	3130	rBV	1278467	2933769	56.57%	6.034%
49	23.898	3514	3521	3540	rBV2	1594266	4192632	80.84%	8.623%
50	24.074	3544	3551	3570	rVB	807883	2727666	52.60%	5.610%

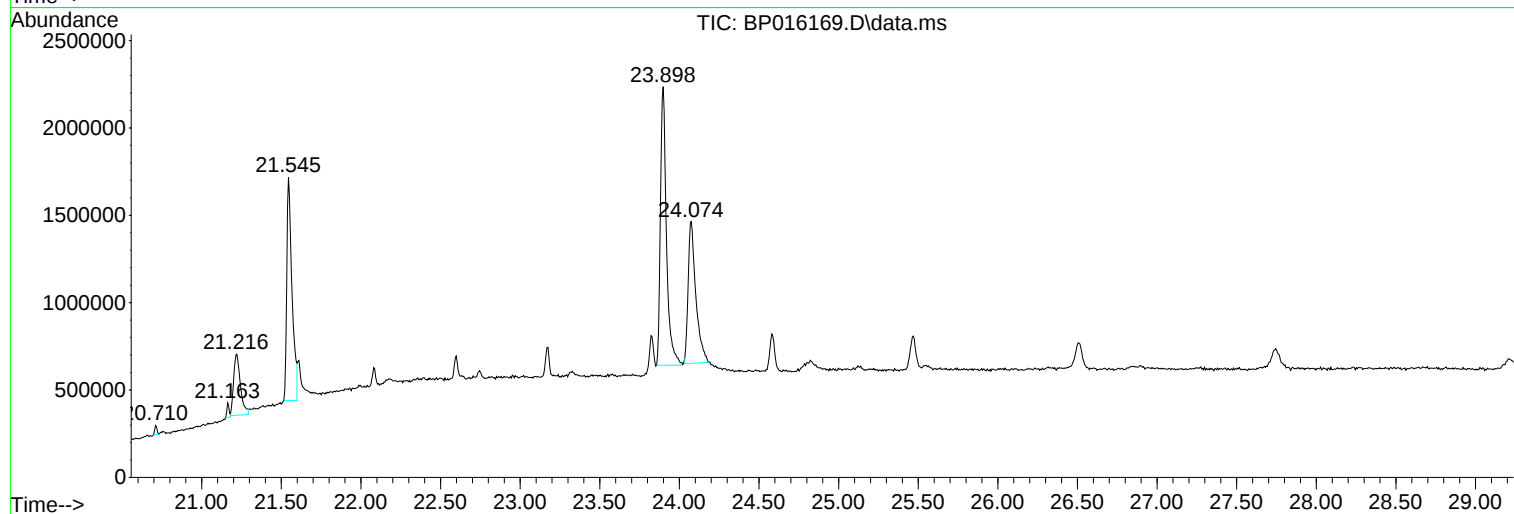
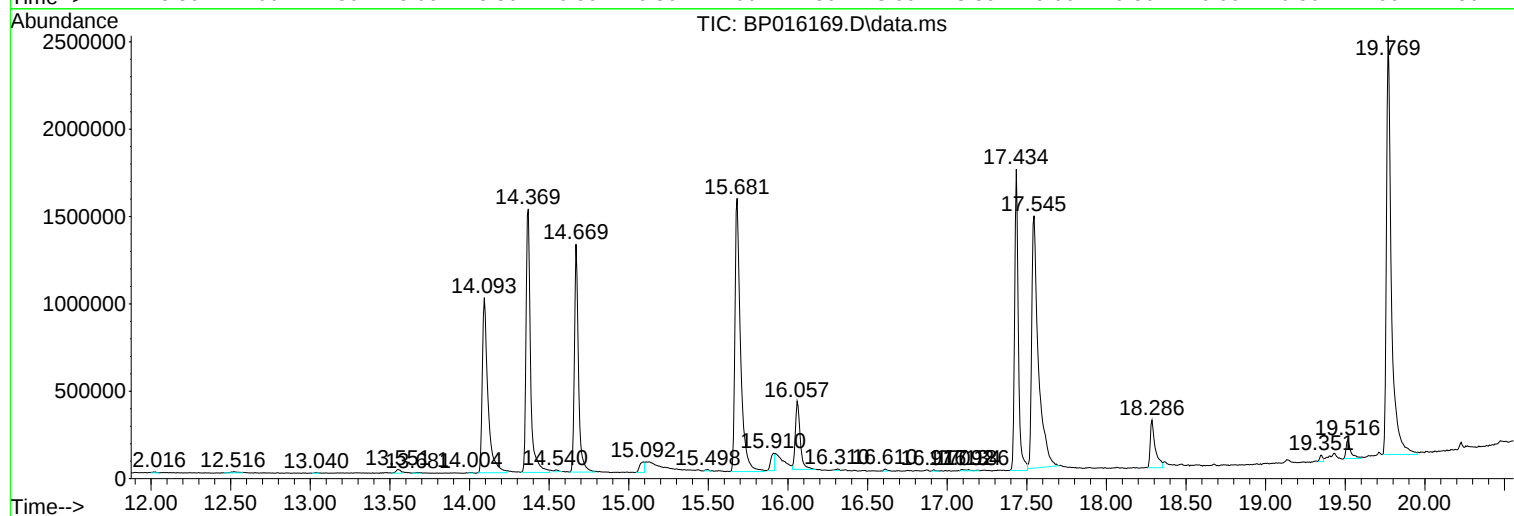
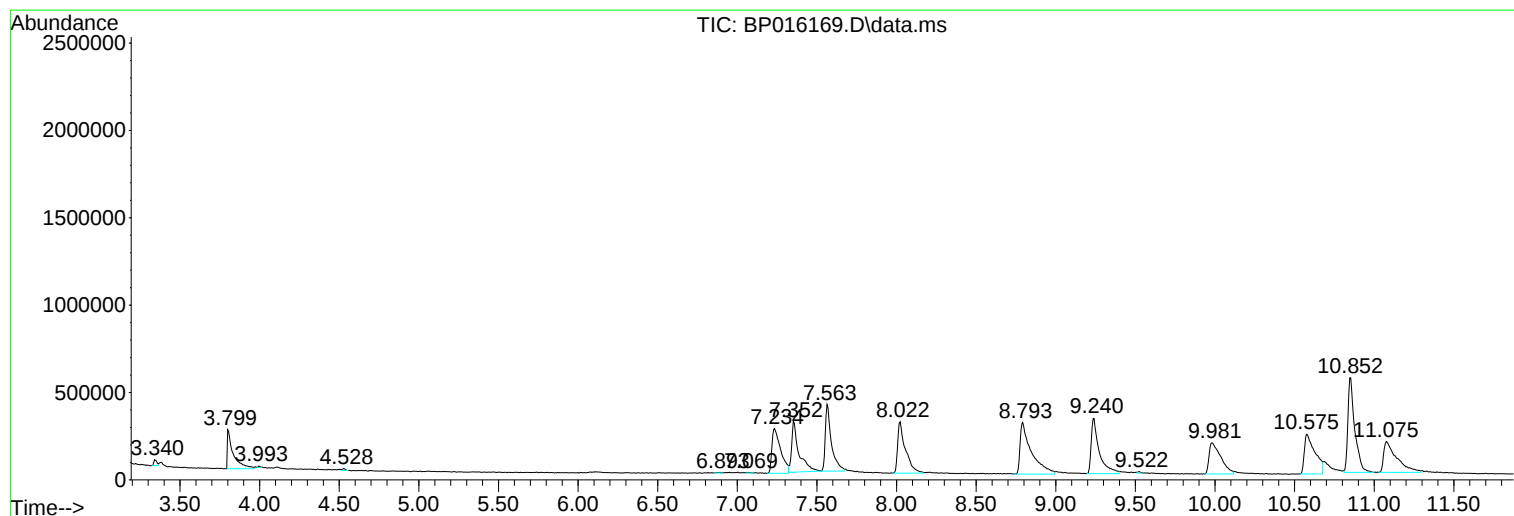
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P



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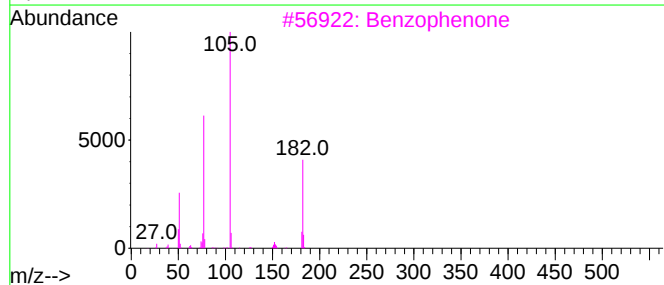
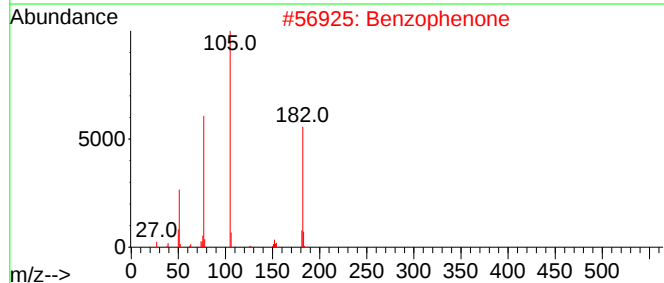
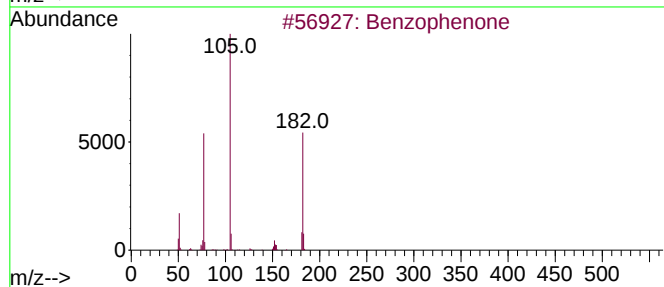
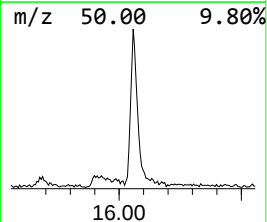
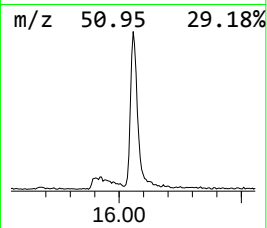
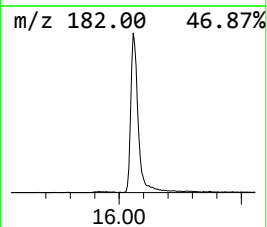
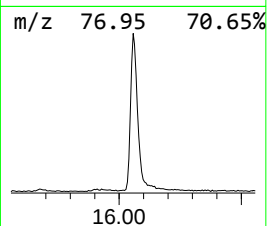
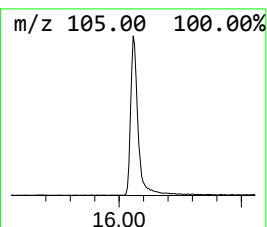
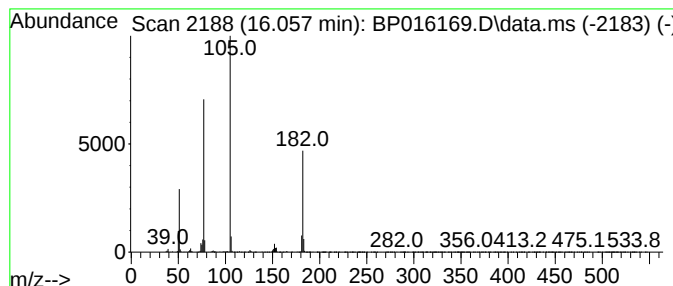
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Benzophenone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.057	5.80 ng/ul	813760	Phenanthrene-d10	17.433

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	97
2			Benzophenone	182	C13H10O	000119-61-9	96
3			Benzophenone	182	C13H10O	000119-61-9	96
4			Benzophenone	182	C13H10O	000119-61-9	91
5			Benzophenone	182	C13H10O	000119-61-9	90



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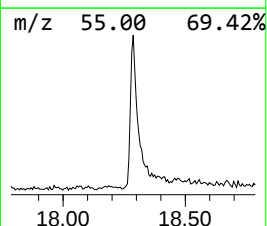
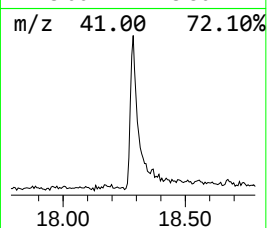
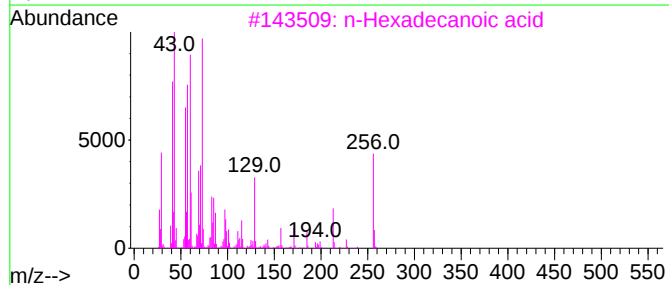
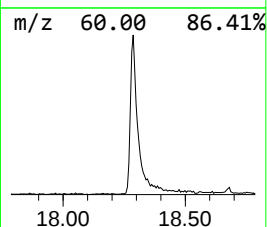
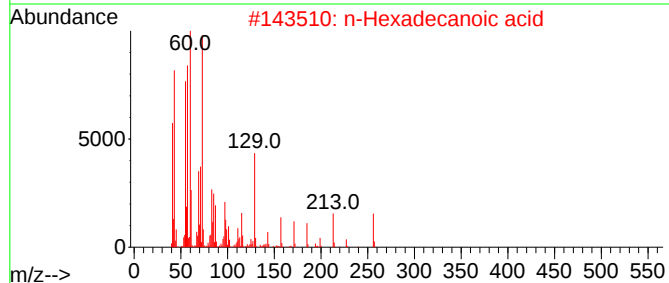
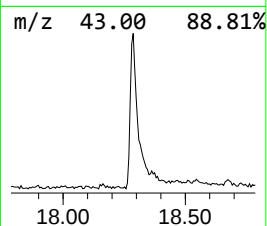
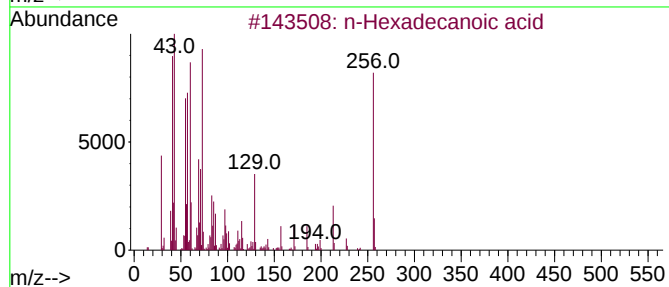
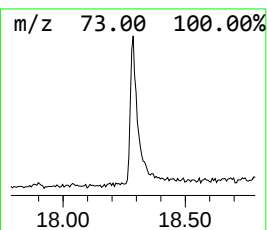
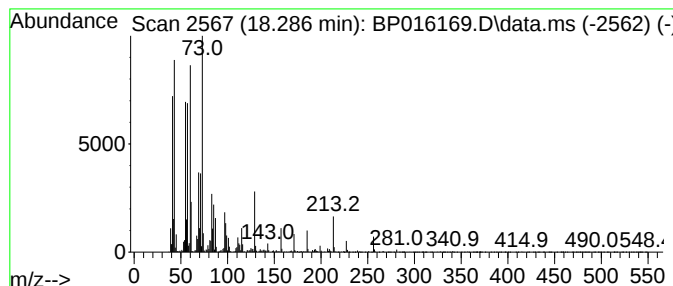
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.286	4.23 ng/ul	593618	Phenanthrene-d10	17.433

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	95
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91



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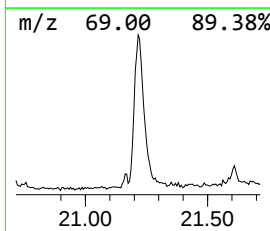
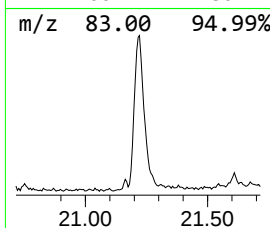
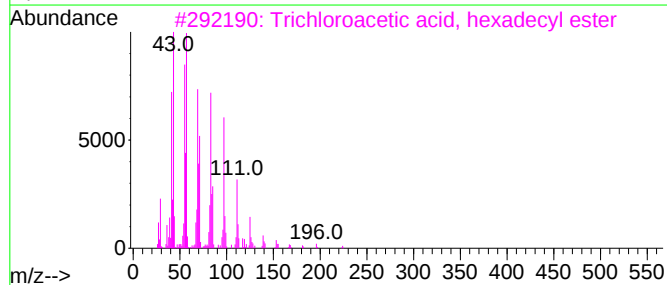
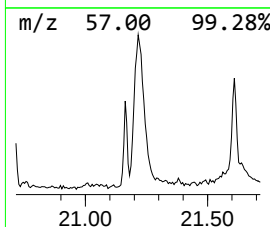
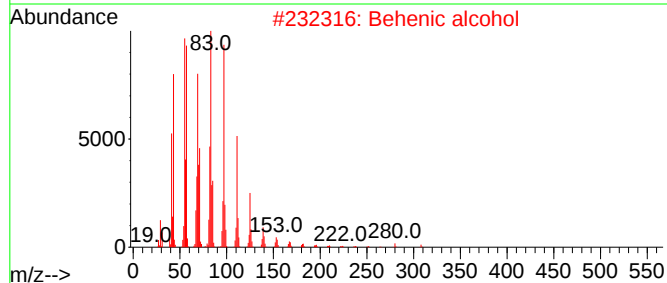
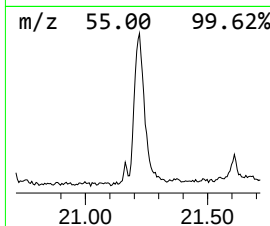
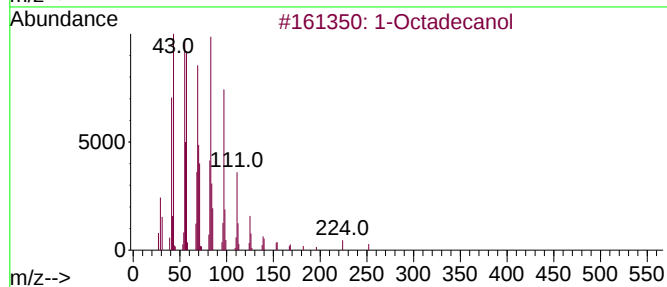
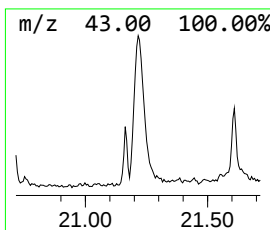
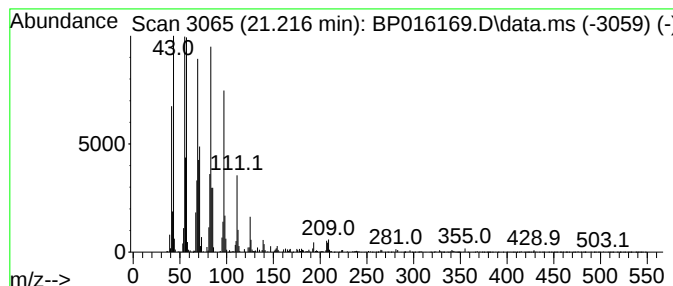
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Peak Number 3 1-Octadecanol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.216	7.10 ng/ul	1041460	Chrysene-d12	21.545

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Octadecanol	270	C18H38O	000112-92-5	94
2			Behenic alcohol	326	C22H46O	000661-19-8	94
3			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	93
4			1-Heneicosanol	312	C21H44O	015594-90-8	91
5			2-Chloropropionic acid, pentade...	318	C18H35ClO2	1000292-44-1	91



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzophenone	16.057	5.8	ng/ul	813760	4	17.433	2804730	20.0
n-Hexadecanoic ...	18.286	4.2	ng/ul	593618	4	17.433	2804730	20.0
1-Octadecanol	21.216	7.1	ng/ul	1041460	5	21.545	2933770	20.0