

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071822\
 Data File : BP011018.D
 Acq On : 19 Jul 2022 06:09
 Operator : CG/JU
 Sample : N3710-14
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 H0B44

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-BP071422.M
 Title : SVOA CALIBRATION

Signal : TIC: BP011018.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.181	12	16	26	rVB	153841	237476	2.10%	0.212%
2	3.593	82	86	101	rVB	795272	1148122	10.16%	1.025%
3	3.811	119	123	128	rVB	59123	92792	0.82%	0.083%
4	3.899	135	138	144	rVB	40913	52176	0.46%	0.047%
5	4.616	258	260	267	rVB4	12727	18946	0.17%	0.017%
6	4.822	289	295	299	rBV3	19711	43538	0.39%	0.039%
7	5.046	331	333	338	rBV2	11503	12920	0.11%	0.012%
8	5.399	390	393	398	rVB2	17374	24398	0.22%	0.022%
9	6.122	511	516	519	rBV2	16461	22734	0.20%	0.020%
10	6.528	581	585	592	rBV6	11663	21710	0.19%	0.019%
11	6.687	607	612	617	rBV7	9538	20519	0.18%	0.018%
12	6.734	617	620	625	rVB7	6401	11874	0.11%	0.011%
13	6.857	633	641	651	rVB	441636	698385	6.18%	0.623%
14	7.022	662	669	678	rVV	1708652	2599845	23.00%	2.320%
15	7.099	679	682	687	rVB	17429	25687	0.23%	0.023%
16	7.210	693	701	710	rBV	1674818	2546006	22.52%	2.272%
17	7.299	713	716	727	rVB9	14745	27776	0.25%	0.025%
18	7.405	732	734	742	rVV9	8264	15187	0.13%	0.014%
19	7.675	774	780	788	rBV	1745312	2710444	23.98%	2.419%
20	7.746	788	792	797	rVB3	23044	37879	0.34%	0.034%
21	7.987	829	833	837	rBV7	15018	24004	0.21%	0.021%
22	8.110	845	854	860	rBV2	38054	83033	0.73%	0.074%
23	8.304	881	887	893	rVB3	53139	84321	0.75%	0.075%
24	8.393	893	902	911	rBV	1319150	2047962	18.12%	1.828%
25	8.504	917	921	926	rVB	22542	37863	0.33%	0.034%
26	8.787	964	969	971	rBV4	13893	23711	0.21%	0.021%
27	8.840	971	978	986	rVB	1865102	2961949	26.20%	2.643%
28	8.957	995	998	1003	rVB4	13526	22445	0.20%	0.020%
29	9.010	1003	1007	1013	rVB7	13431	21692	0.19%	0.019%
30	9.093	1014	1021	1029	rVB3	24878	56721	0.50%	0.051%
31	9.251	1046	1048	1054	rVB5	9419	14955	0.13%	0.013%
32	9.346	1054	1064	1067	rBV5	9620	27042	0.24%	0.024%
33	9.557	1091	1100	1108	rBV	1432733	2269288	20.07%	2.025%
34	9.722	1124	1128	1130	rBV5	9638	11759	0.10%	0.010%
35	9.793	1136	1140	1146	rVB6	13896	26734	0.24%	0.024%
36	9.981	1169	1172	1174	rBV4	12244	16003	0.14%	0.014%

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Integrator: RTE

Smoothing : OFF

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Sampling : 1

Min Area: 0.1 % of largest Peak

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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-BP071422.M
 Title : SVOA CALIBRATION

37	10.093	1184	1191	1199	rBV	2137173	3444796	30.47%	3.074%
38	10.169	1202	1204	1208	rVB5	11887	13933	0.12%	0.012%
39	10.263	1215	1220	1229	rBV5	21639	46530	0.41%	0.042%
40	10.340	1229	1233	1238	rVB8	9596	13610	0.12%	0.012%
41	10.469	1247	1255	1264	rBV	2323410	3775253	33.40%	3.369%
42	10.616	1271	1280	1295	rBV	2003775	3463589	30.64%	3.091%
43	10.775	1305	1307	1315	rVB8	10086	17658	0.16%	0.016%
44	10.887	1319	1326	1332	rVB10	13339	33578	0.30%	0.030%
45	10.957	1336	1338	1344	rVB7	8037	11907	0.11%	0.011%
46	11.128	1363	1367	1370	rBV3	10991	16492	0.15%	0.015%
47	11.281	1388	1393	1398	rVB9	8788	17099	0.15%	0.015%
48	11.451	1417	1422	1423	rBV4	10570	14375	0.13%	0.013%
49	11.716	1460	1467	1471	rBV5	14167	32632	0.29%	0.029%
50	11.757	1471	1474	1478	rVB5	10069	12732	0.11%	0.011%
51	11.845	1484	1489	1491	rBV6	7674	11605	0.10%	0.010%
52	12.069	1522	1527	1533	rVB	39412	60921	0.54%	0.054%
53	12.175	1540	1545	1550	rVB3	19736	32402	0.29%	0.029%
54	12.263	1550	1560	1569	rBV	114105	235171	2.08%	0.210%
55	12.622	1614	1621	1625	rBV9	7526	19421	0.17%	0.017%
56	12.692	1627	1633	1636	rBV4	19820	33592	0.30%	0.030%
57	12.781	1645	1648	1653	rBV4	16250	26336	0.23%	0.024%
58	12.951	1673	1677	1681	rVB7	13524	20988	0.19%	0.019%
59	13.063	1692	1696	1700	rBV7	11517	16631	0.15%	0.015%
60	13.304	1734	1737	1746	rBV9	10866	24605	0.22%	0.022%
61	13.463	1759	1764	1767	rBV7	10382	20301	0.18%	0.018%
62	13.522	1770	1774	1776	rBV5	11585	14221	0.13%	0.013%
63	13.663	1795	1798	1802	rVB6	12165	14611	0.13%	0.013%
64	13.757	1807	1814	1830	rBV	4052133	5665266	50.12%	5.056%
65	13.898	1836	1838	1844	rVB7	8454	14154	0.13%	0.013%
66	14.028	1850	1860	1872	rBV	4418481	6579885	58.21%	5.872%
67	14.145	1876	1880	1887	rVV	80256	137184	1.21%	0.122%
68	14.210	1887	1891	1895	rVV7	15028	36179	0.32%	0.032%
69	14.245	1895	1897	1903	rVV6	22142	34079	0.30%	0.030%
70	14.339	1906	1913	1922	rVV	3203263	4613982	40.82%	4.117%
71	14.422	1923	1927	1933	rVB3	36599	53407	0.47%	0.048%
72	14.557	1944	1950	1960	rBV	333129	561231	4.96%	0.501%
73	14.710	1972	1976	1982	rVV6	26213	37623	0.33%	0.034%
74	14.775	1983	1987	1992	rVB2	32341	44550	0.39%	0.040%
75	14.916	2008	2011	2016	rBV5	16697	23360	0.21%	0.021%
76	15.210	2058	2061	2065	rVB4	17537	23520	0.21%	0.021%

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 Sampling : 1 Min Area: 0.1 % of largest Peak
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 Stop Thrs : 0 Peak Location: TOP

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
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77	15.339	2077	2083	2090	rBV	5890890	8165057	72.23%	7.286%
78	15.463	2098	2104	2119	rBV	1776762	2428260	21.48%	2.167%
79	15.581	2120	2124	2128	rVB2	60494	88449	0.78%	0.079%
80	15.669	2136	2139	2144	rBV	36099	54481	0.48%	0.049%
81	15.904	2176	2179	2181	rBV4	11780	13205	0.12%	0.012%
82	16.033	2198	2201	2206	rBV3	24559	39828	0.35%	0.036%
83	16.081	2206	2209	2214	rVV6	20096	36872	0.33%	0.033%
84	16.133	2215	2218	2224	rVB6	31174	52338	0.46%	0.047%
85	16.881	2342	2345	2349	rBV4	45690	56783	0.50%	0.051%
86	17.110	2377	2384	2390	rBV	3536637	5149193	45.55%	4.595%
87	17.210	2394	2401	2413	rVB	6289403	8933729	79.03%	7.972%
88	17.433	2435	2439	2444	rVB4	28872	47995	0.42%	0.043%
89	17.492	2447	2449	2456	rVB7	22984	34424	0.30%	0.031%
90	17.616	2467	2470	2477	rVB2	72486	98765	0.87%	0.088%
91	17.798	2499	2501	2510	rVB7	27674	47437	0.42%	0.042%
92	18.022	2534	2539	2556	rBV	616271	1117951	9.89%	0.998%
93	19.039	2708	2712	2717	rVB2	95187	116986	1.03%	0.104%
94	19.104	2720	2723	2728	rBV	91575	116157	1.03%	0.104%
95	19.269	2747	2751	2761	rBV2	140863	262158	2.32%	0.234%
96	19.463	2778	2784	2791	rBV	8028493	10701404	94.67%	9.550%
97	20.951	3033	3037	3048	rBV	2846692	3654656	32.33%	3.261%
98	21.227	3079	3084	3089	rBV	4808542	5910076	52.28%	5.274%
99	23.433	3452	3459	3467	rVB2	5901223	11304363	100.00%	10.088%
100	23.586	3478	3485	3497	rVB2	3147923	6262081	55.40%	5.588%

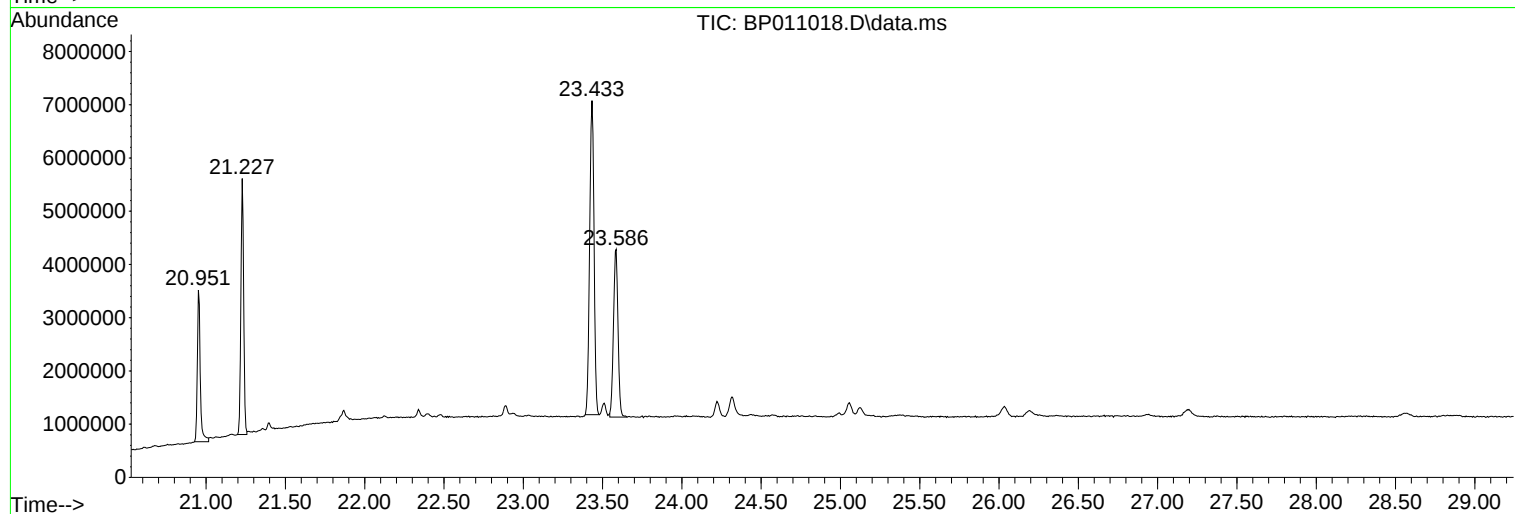
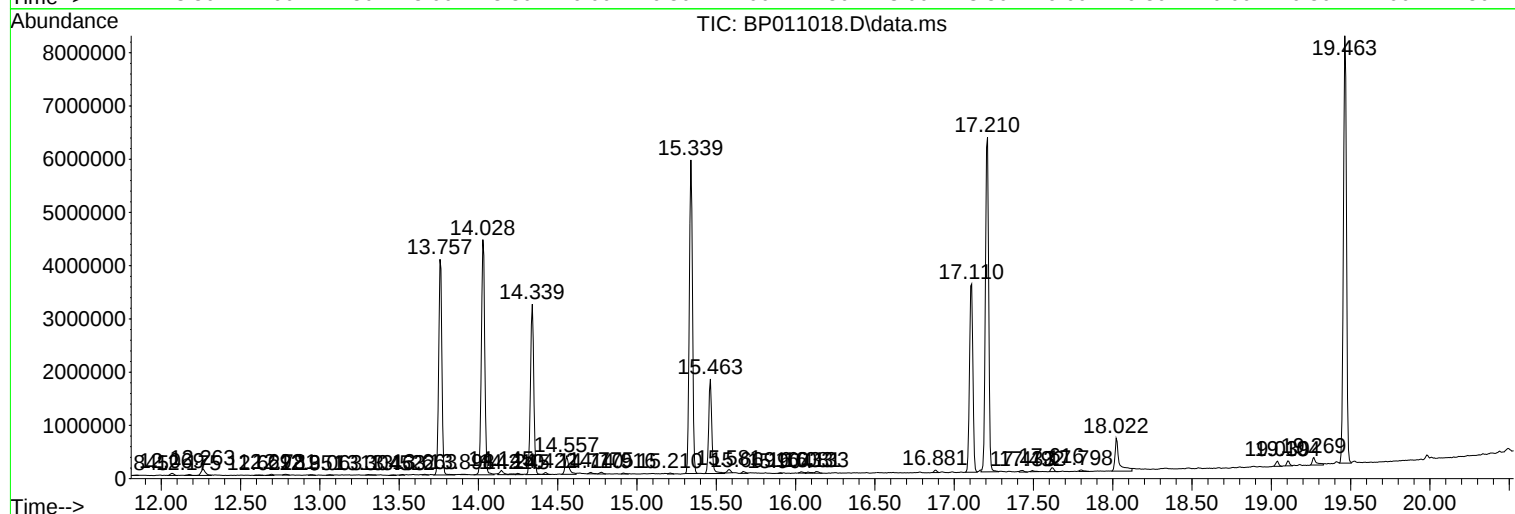
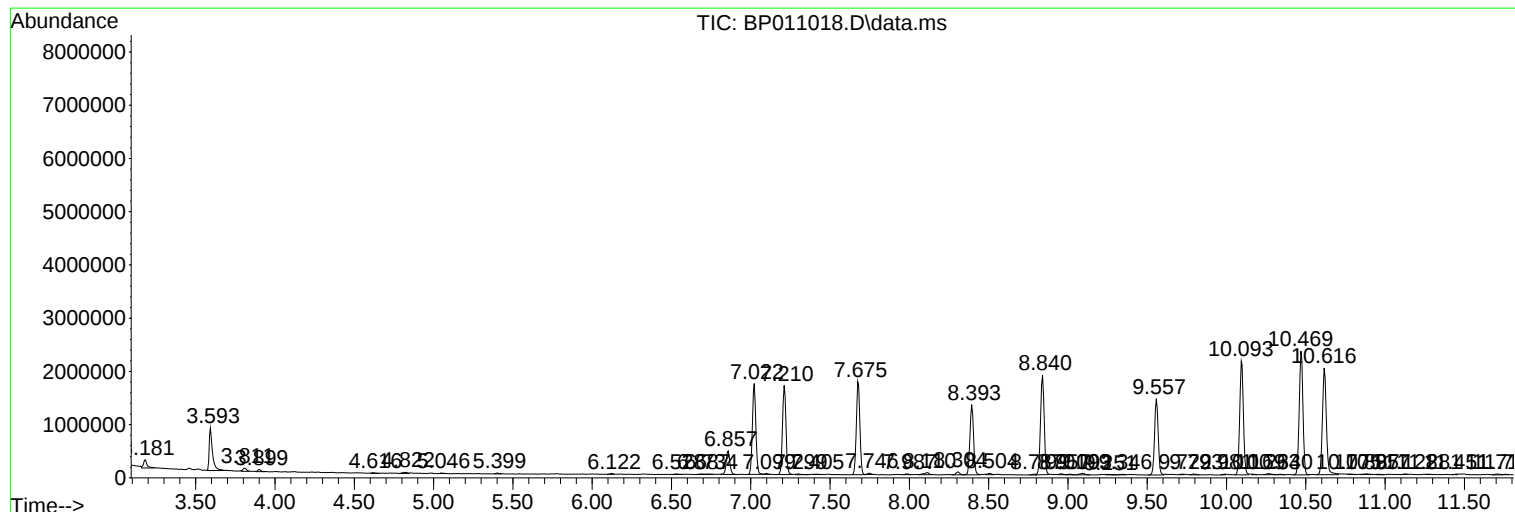
Sum of corrected areas: 112059953

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION

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



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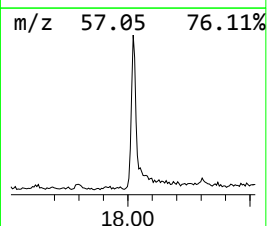
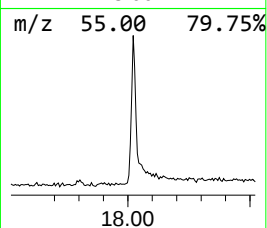
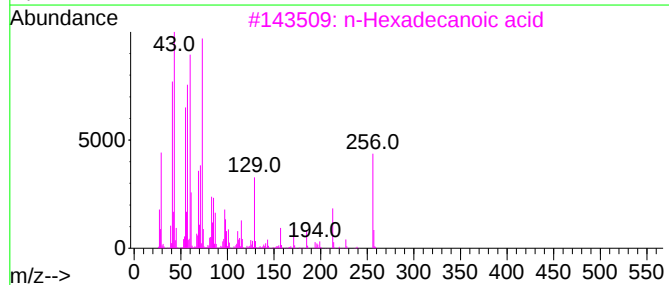
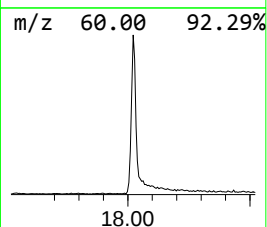
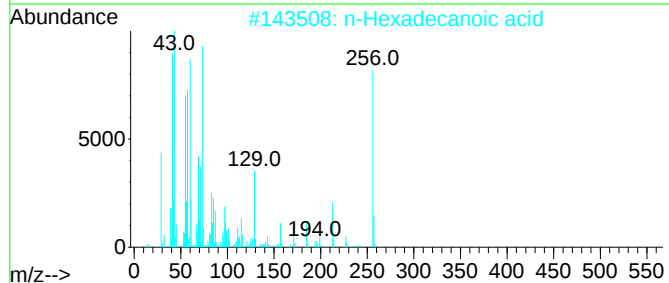
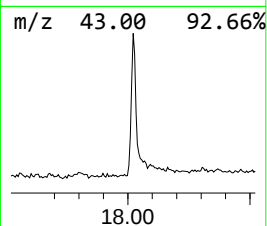
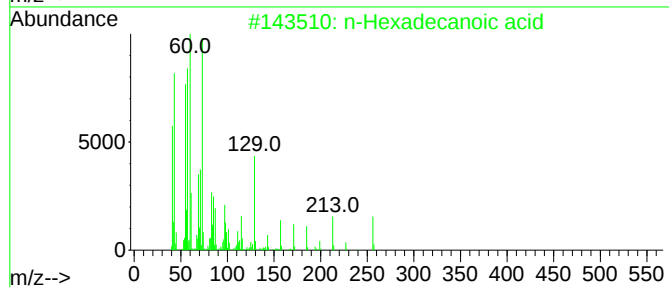
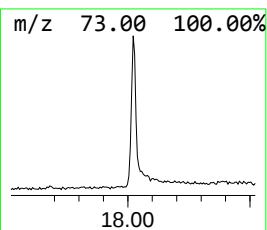
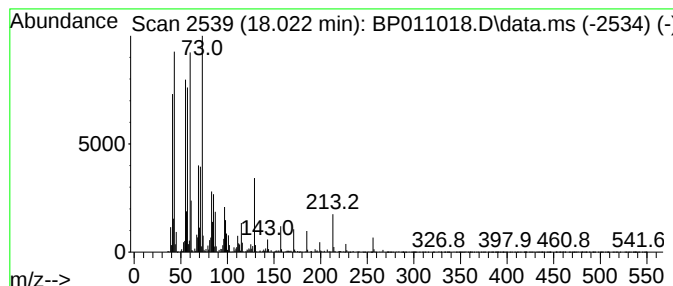
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.022	4.34 ng/ul	1117950	Phenanthrene-d10	17.110

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	99
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
5			Octadecanoic acid	284	C18H36O2	000057-11-4	90



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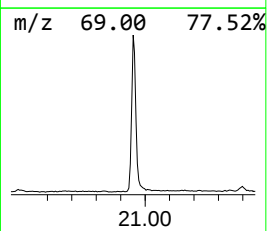
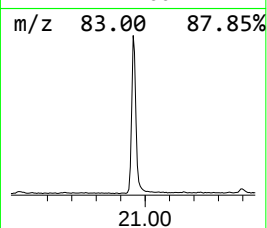
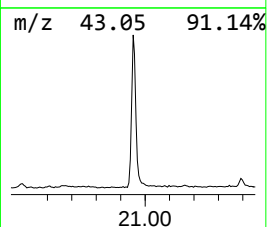
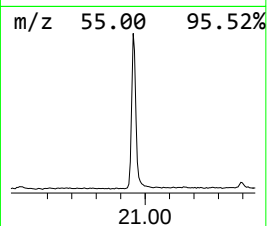
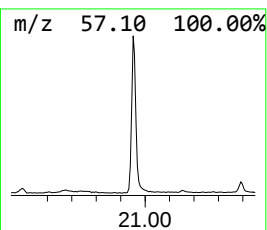
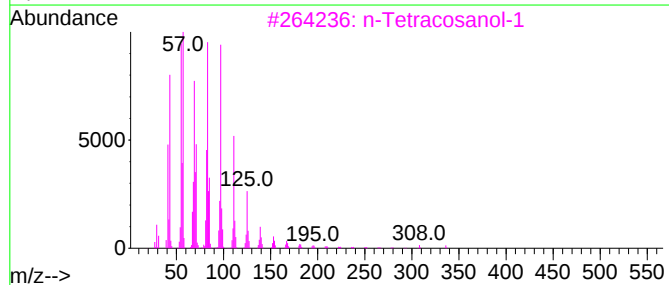
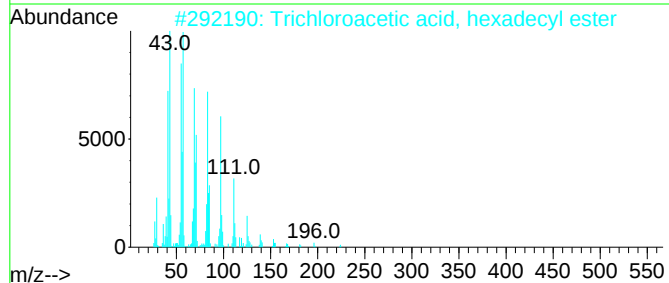
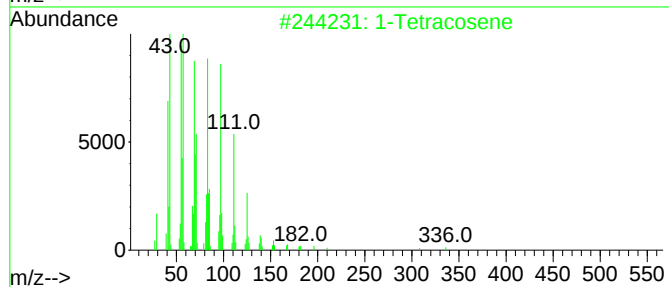
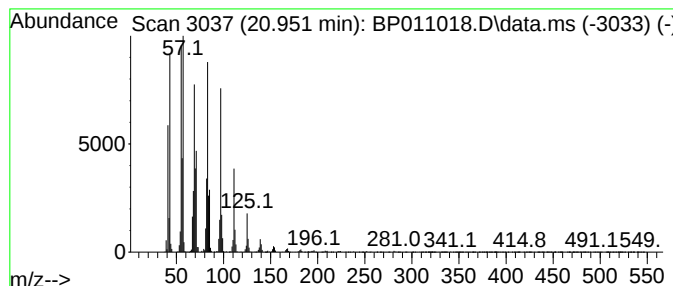
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.951	12.37 ng/ul	3654660	Chrysene-d12	21.227

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Tetracosene	336	C24H48	010192-32-2	94
2			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
3			n-Tetracosanol-1	354	C24H50O	000506-51-4	94
4			Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	94
5			1-Heneicosanol	312	C21H44O	015594-90-8	93



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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
n-Hexadecanoic ...	18.022	4.3	ng/ul	1117950	4	17.110	5149190	20.0
(DEL) Alkane: S...	20.951	12.4	ng/ul	3654660	5	21.227	5910080	20.0