

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030822\
 Data File : BM034164.D
 Acq On : 09 Mar 2022 00:43
 Operator : CG/JU
 Sample : N1733-24
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBRU2

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_M\Methods\SFAM-EPA-BM030822.M
 Title : SVOA CALIBRATION

Signal : TIC: BM034164.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.372	19	23	33	rVB	54079	72161	1.75%	0.159%
2	3.655	69	71	78	rVB	4695	6537	0.16%	0.014%
3	3.790	90	94	113	rBV	625549	885098	21.47%	1.952%
4	4.137	150	153	158	rVB2	9492	12044	0.29%	0.027%
5	4.366	189	192	197	rVB5	4689	7511	0.18%	0.017%
6	5.072	308	312	316	rVB4	9428	12160	0.29%	0.027%
7	5.149	321	325	327	rBV5	3555	4179	0.10%	0.009%
8	7.143	657	664	673	rBV	812148	1290497	31.30%	2.846%
9	7.313	687	693	701	rBV	630228	967015	23.46%	2.132%
10	7.384	701	705	710	rVB2	4517	6629	0.16%	0.015%
11	7.519	721	728	738	rBV	870954	1354971	32.87%	2.988%
12	7.990	801	808	816	rBV	518306	812624	19.71%	1.792%
13	8.054	816	819	824	rVB2	5322	7485	0.18%	0.017%
14	8.231	845	849	856	rVB6	3049	6633	0.16%	0.015%
15	8.695	920	928	939	rVV	1055907	1684448	40.86%	3.714%
16	8.784	939	943	949	rVB8	2607	4892	0.12%	0.011%
17	9.154	998	1006	1014	rBV	779766	1289238	31.27%	2.843%
18	9.242	1014	1021	1029	rVB2	11021	22855	0.55%	0.050%
19	9.331	1031	1036	1041	rVB8	4102	7325	0.18%	0.016%
20	9.607	1075	1083	1090	rBV	29065	46116	1.12%	0.102%
21	9.878	1121	1129	1137	rBV	693345	1119096	27.14%	2.468%
22	10.419	1212	1221	1232	rBV	1122496	1820484	44.16%	4.014%
23	10.583	1243	1249	1258	rBV2	20734	34027	0.83%	0.075%
24	10.807	1279	1287	1294	rBV	674750	1132330	27.47%	2.497%
25	10.931	1300	1308	1320	rBV	990054	1677457	40.69%	3.699%
26	11.748	1440	1447	1456	rBV8	4096	12010	0.29%	0.026%
27	12.236	1525	1530	1535	rBV2	12814	21287	0.52%	0.047%
28	12.389	1547	1556	1563	rVB2	19231	31335	0.76%	0.069%
29	12.901	1640	1643	1648	rBV4	3491	4386	0.11%	0.010%
30	13.048	1662	1668	1671	rVB7	5101	6480	0.16%	0.014%
31	13.124	1678	1681	1687	rVB7	3575	6046	0.15%	0.013%
32	13.177	1687	1690	1696	rVB4	4271	7220	0.18%	0.016%
33	13.466	1734	1739	1748	rVB2	32041	45991	1.12%	0.101%
34	13.595	1758	1761	1767	rVB6	3897	6307	0.15%	0.014%
35	14.030	1829	1835	1842	rBV	1822812	2532655	61.43%	5.585%
36	14.160	1853	1857	1862	rVB4	3932	7076	0.17%	0.016%

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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_M\Methods\SFAM-EPA-BM030822.M
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37	14.319	1873	1884	1896	rVV	2109145	3135581	76.06%	6.914%
38	14.530	1915	1920	1925	rVB2	10383	14951	0.36%	0.033%
39	14.630	1930	1937	1945	rVV	1030080	1528354	37.07%	3.370%
40	14.801	1960	1966	1979	rBV	701733	1001098	24.28%	2.208%
41	14.983	1993	1997	2001	rVV7	3473	5589	0.14%	0.012%
42	15.036	2001	2006	2010	rVB5	7669	12457	0.30%	0.027%
43	15.436	2068	2074	2078	rBV5	7474	12255	0.30%	0.027%
44	15.483	2078	2082	2086	rVB4	7710	10037	0.24%	0.022%
45	15.618	2098	2105	2117	rBV	2661170	3730324	90.48%	8.226%
46	15.724	2117	2123	2132	rBV	655341	903222	21.91%	1.992%
47	15.854	2142	2145	2155	rVB2	12910	22332	0.54%	0.049%
48	15.971	2162	2165	2170	rBV5	3569	4929	0.12%	0.011%
49	16.030	2170	2175	2180	rBV6	3812	8224	0.20%	0.018%
50	16.260	2210	2214	2216	rBV5	4381	6449	0.16%	0.014%
51	16.301	2216	2221	2224	rVV6	7196	12151	0.29%	0.027%
52	16.342	2225	2228	2230	rVV4	4889	6266	0.15%	0.014%
53	16.401	2234	2238	2243	rVB7	4687	9241	0.22%	0.020%
54	16.618	2271	2275	2279	rBV6	2572	4561	0.11%	0.010%
55	16.765	2294	2300	2307	rBV10	6486	13834	0.34%	0.031%
56	17.136	2358	2363	2364	rBV4	2762	4400	0.11%	0.010%
57	17.365	2396	2402	2408	rBV	1188344	1633967	39.63%	3.603%
58	17.465	2412	2419	2429	rVV	2667191	3712810	90.06%	8.187%
59	17.548	2430	2433	2439	rVB7	13718	21691	0.53%	0.048%
60	17.748	2463	2467	2472	rBV8	5230	8612	0.21%	0.019%
61	18.042	2515	2517	2522	rVB5	4006	5585	0.14%	0.012%
62	18.142	2530	2534	2538	rVB6	4278	6297	0.15%	0.014%
63	18.259	2549	2554	2561	rBV2	176139	243419	5.90%	0.537%
64	18.695	2625	2628	2632	rVB5	11509	13530	0.33%	0.030%
65	19.124	2697	2701	2702	rBV3	6350	6899	0.17%	0.015%
66	19.318	2729	2734	2739	rBV2	32152	47807	1.16%	0.105%
67	19.389	2742	2746	2749	rBV	33681	44998	1.09%	0.099%
68	19.530	2766	2770	2777	rBV7	34561	43640	1.06%	0.096%
69	19.753	2802	2808	2819	rBV	2994764	4036658	97.91%	8.901%
70	20.259	2891	2894	2898	rBV6	14681	21273	0.52%	0.047%
71	21.242	3058	3061	3067	rBV	283111	334632	8.12%	0.738%
72	21.447	3094	3096	3101	rBV	48042	59603	1.45%	0.131%
73	21.536	3106	3111	3117	rVV	1313782	1615351	39.18%	3.562%
74	21.865	3163	3167	3171	rBV	119862	158113	3.84%	0.349%
75	23.824	3493	3500	3508	rVB	2126169	4122743	100.00%	9.091%
76	23.983	3520	3527	3537	rVB	889343	1799662	43.65%	3.969%

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ClientSampleId :
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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_M\Methods\SFAM-EPA-BM030822.M
Title : SVOA CALIBRATION

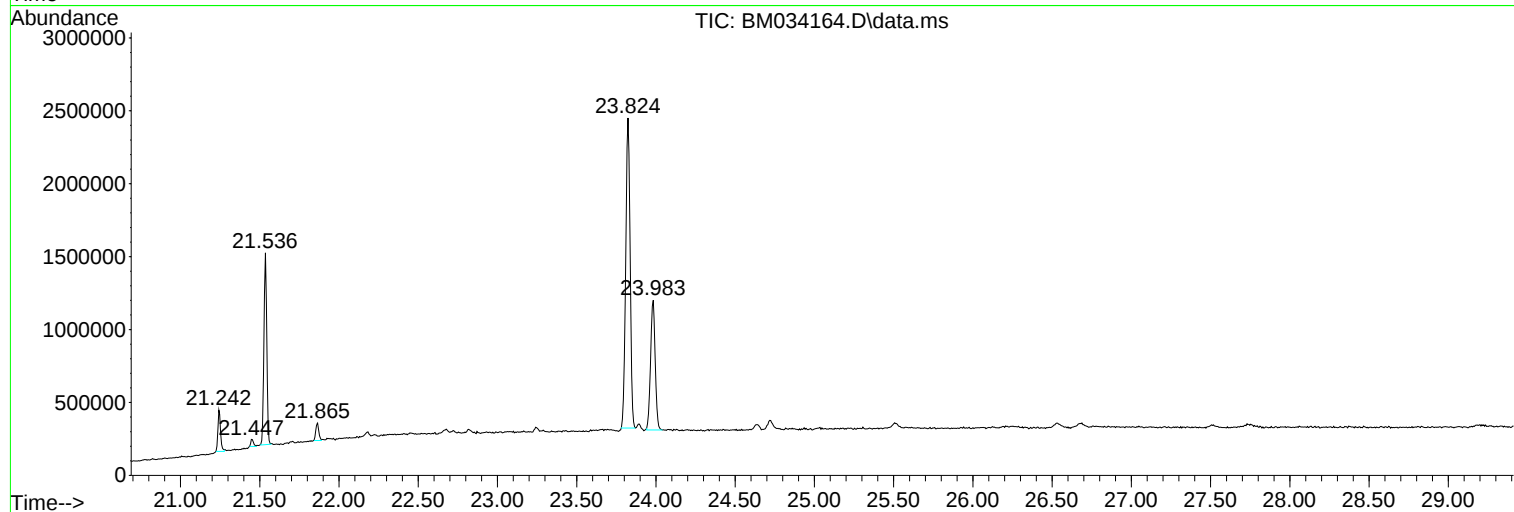
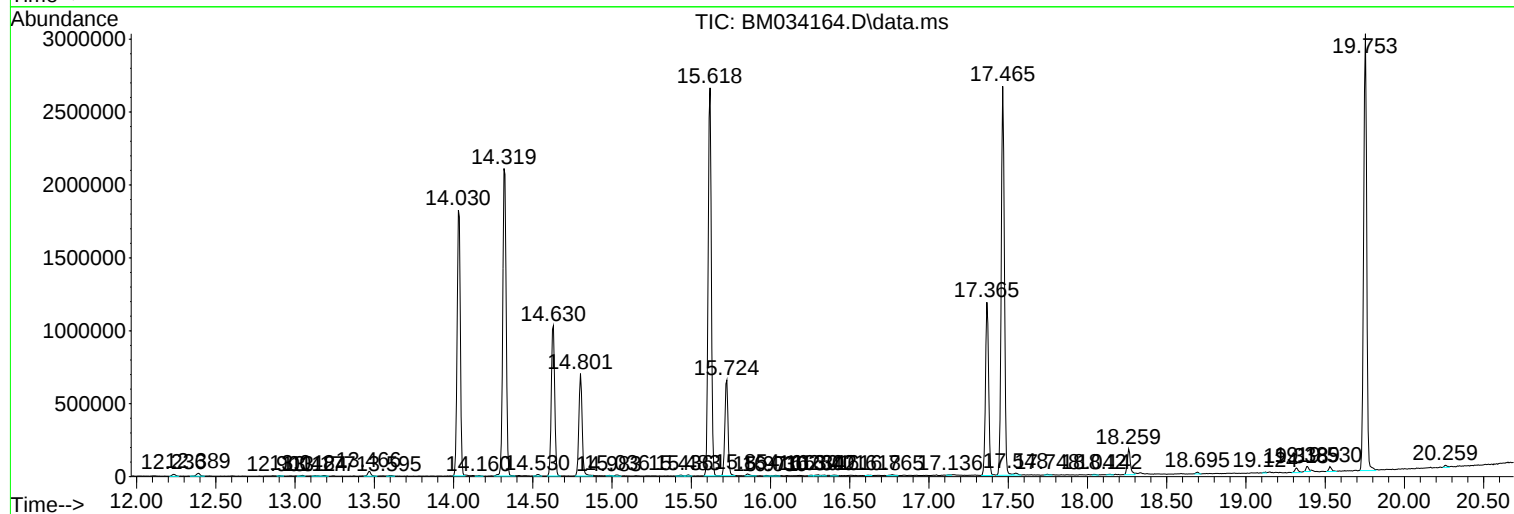
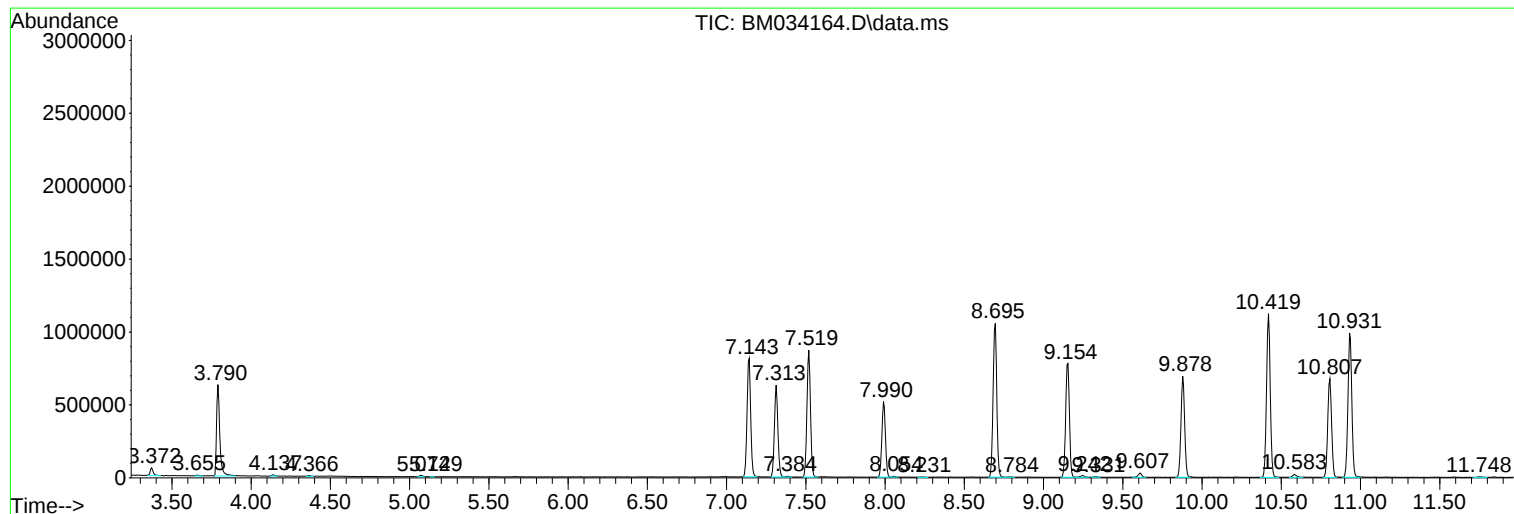
Sum of corrected areas: 45348150

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030822.M
Quant Title : SVOA CALIBRATION

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



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Acq On : 09 Mar 2022 00:43
Operator : CG/JU
Sample : N1733-24
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DBRU2

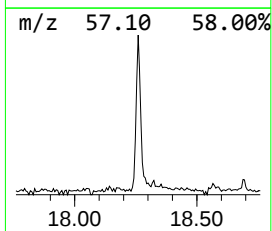
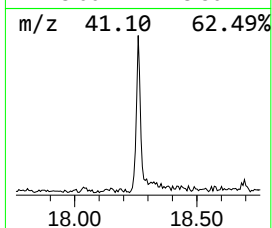
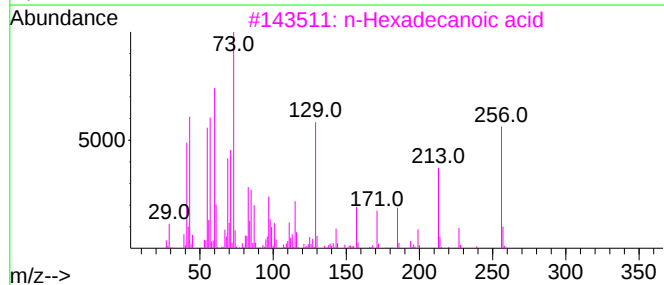
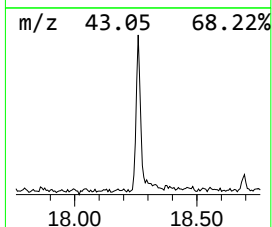
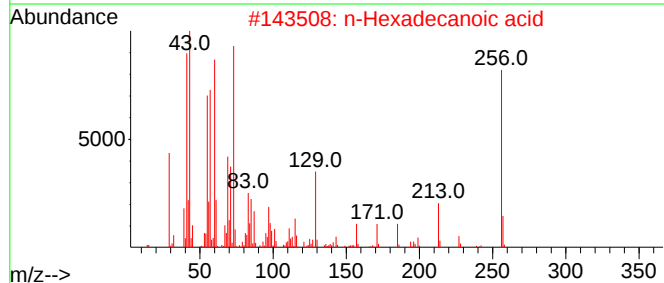
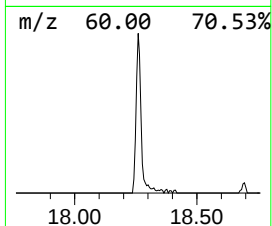
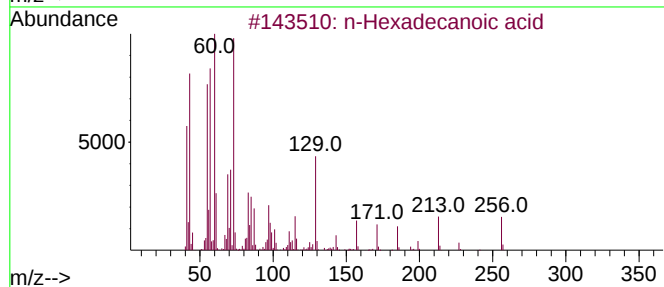
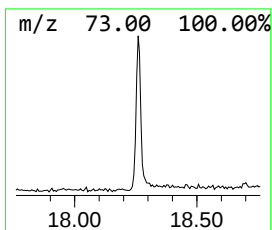
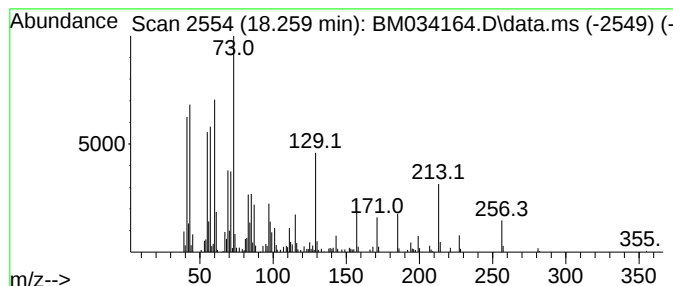
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030822.M
Quant Title : SVOA CALIBRATION

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.
18.259	2.98 ng/ul	243419	Phenanthrene-d10	17.365

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			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	89



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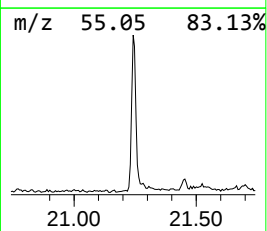
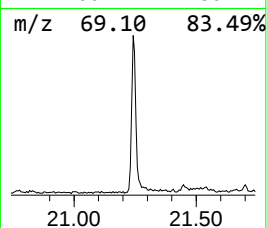
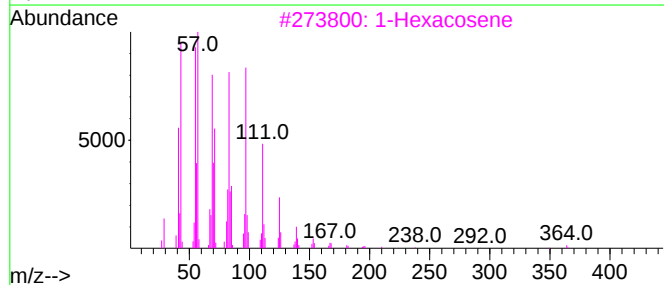
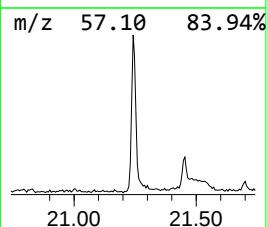
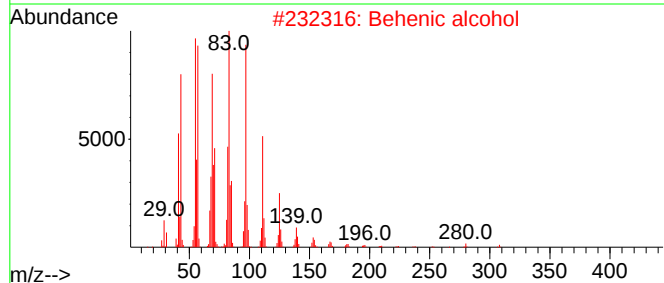
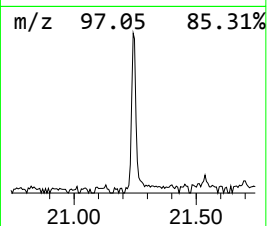
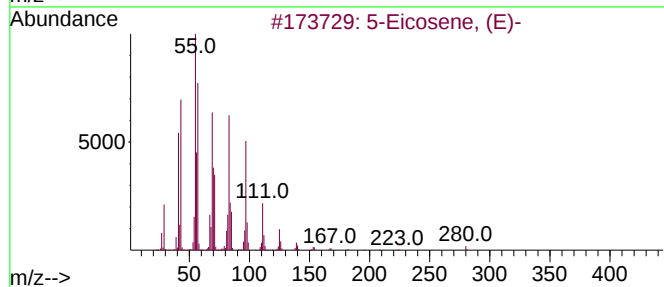
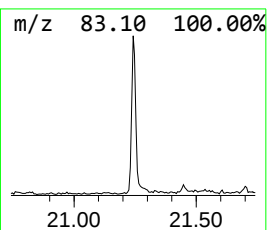
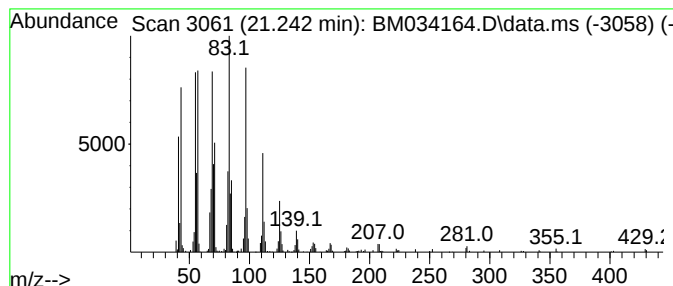
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030822.M
Quant Title : SVOA CALIBRATION

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.
21.242	4.14 ng/ul	334632	Chrysene-d12	21.536

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Eicosene, (E)-		280	C20H40	074685-30-6	99
2	Behenic alcohol		326	C22H46O	000661-19-8	95
3	1-Hexacosene		364	C26H52	018835-33-1	95
4	1-Heneicosanol		312	C21H44O	015594-90-8	95
5	1-Nonadecene		266	C19H38	018435-45-5	95



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Quant Title : SVOA CALIBRATION

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 ...	18.259	3.0	ng/u1	243419	4	17.365	1633970	20.0
(DEL) Alkane: S...	21.242	4.1	ng/u1	334632	5	21.536	1615350	20.0