

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN070618\
 Data File : BN001984.D
 Acq On : 06 Jul 2018 11:41
 Operator : JU/SJ
 Sample : J3765-09
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F1H11

Integration Parameters: LSCINT.P

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

Filtering: 5
 Min Area: 0.1 % 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:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN061918.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	3.023	3	6	21	rVB	3052427	4571030	35.13%	2.522%
2	3.270	44	48	54	rVB	85309	117272	0.90%	0.065%
3	4.193	201	205	212	rVB	51820	72357	0.56%	0.040%
4	4.864	314	319	327	rVV3	32502	61470	0.47%	0.034%
5	6.881	655	662	676	rBV	1841344	3025623	23.25%	1.669%
6	7.046	684	690	704	rBV	1406954	2207954	16.97%	1.218%
7	7.246	717	724	735	rBV	1799941	2887634	22.19%	1.593%
8	7.711	795	803	810	rBV	1861947	3050318	23.44%	1.683%
9	8.411	912	922	933	rBV	2337599	3881141	29.83%	2.142%
10	8.852	990	997	1006	rBV	1752671	2923239	22.46%	1.613%
11	9.569	1111	1119	1133	rBV	1472041	2461376	18.92%	1.358%
12	10.105	1202	1210	1224	rBV	2495531	4167962	32.03%	2.300%
13	10.481	1266	1274	1281	rBV	3002237	5016820	38.55%	2.768%
14	10.616	1288	1297	1314	rBV	2164300	3722500	28.61%	2.054%
15	12.069	1538	1544	1551	rVB	50580	84404	0.65%	0.047%
16	12.699	1642	1651	1660	rBV5	89373	153278	1.18%	0.085%
17	12.946	1688	1693	1698	rBV	29718	46710	0.36%	0.026%
18	13.081	1710	1716	1717	rBV3	30570	48941	0.38%	0.027%
19	13.104	1717	1720	1725	rVV3	45024	63771	0.49%	0.035%
20	13.216	1735	1739	1750	rVV4	105360	215215	1.65%	0.119%
21	13.469	1778	1782	1787	rVV3	62237	97618	0.75%	0.054%
22	13.663	1809	1815	1822	rVV	1541277	2162064	16.61%	1.193%
23	13.751	1823	1830	1834	rVV	4687472	6428207	49.40%	3.547%
24	13.798	1834	1838	1849	rVB	1017139	1417980	10.90%	0.782%
25	14.028	1869	1877	1888	rBV	5089892	7924090	60.89%	4.372%
26	14.187	1898	1904	1909	rBV2	53516	76527	0.59%	0.042%
27	14.340	1915	1930	1937	rVV2	4305337	7164739	55.06%	3.953%
28	14.416	1937	1943	1950	rVB2	114654	203843	1.57%	0.112%
29	14.540	1957	1964	1971	rBV	2063340	3064688	23.55%	1.691%
30	14.598	1971	1974	1980	rVV2	155839	262403	2.02%	0.145%
31	14.663	1980	1985	1994	rVV	129312	277850	2.14%	0.153%
32	14.763	1994	2002	2009	rVB5	133687	258983	1.99%	0.143%
33	14.875	2014	2021	2029	rVB	31516	57501	0.44%	0.032%
34	15.069	2050	2054	2058	rVB5	30470	46757	0.36%	0.026%

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Integrator: RTE
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 Title : SVOA CALIBRATION

35	15.151	2063	2068	2074	rBV3	33943	67784	0.52%	0.037%
36	15.281	2085	2090	2094	rBV	649336	927312	7.13%	0.512%
37	15.334	2094	2099	2105	rVV	6691438	9648778	74.15%	5.324%
38	15.451	2112	2119	2131	rBV	2035384	3049185	23.43%	1.682%
39	15.569	2136	2139	2151	rVB	89356	156468	1.20%	0.086%
40	15.687	2153	2159	2162	rVV3	148694	265751	2.04%	0.147%
41	15.722	2162	2165	2170	rVV	225839	301049	2.31%	0.166%
42	15.822	2174	2182	2191	rVB2	291880	539126	4.14%	0.297%
43	15.916	2193	2198	2210	rVV6	140093	411270	3.16%	0.227%
44	16.063	2218	2223	2230	rVV3	114730	263272	2.02%	0.145%
45	16.128	2230	2234	2239	rVB	129129	187234	1.44%	0.103%
46	16.240	2249	2253	2258	rBV4	39114	56007	0.43%	0.031%
47	16.481	2290	2294	2300	rBV4	34602	67421	0.52%	0.037%
48	16.734	2330	2337	2345	rVB4	60570	123025	0.95%	0.068%
49	16.834	2349	2354	2357	rBV	116529	166122	1.28%	0.092%
50	16.863	2357	2359	2362	rVV3	52804	59902	0.46%	0.033%
51	16.922	2366	2369	2374	rVB4	67544	94553	0.73%	0.052%
52	16.998	2378	2382	2386	rBV3	36993	63311	0.49%	0.035%
53	17.081	2390	2396	2401	rVV2	5572203	8160475	62.71%	4.503%
54	17.122	2401	2403	2407	rVV	203632	244558	1.88%	0.135%
55	17.181	2407	2413	2417	rVV2	7042504	10443860	80.26%	5.763%
56	17.216	2417	2419	2426	rVV	999486	1129151	8.68%	0.623%
57	17.287	2426	2431	2441	rVV	55134	147196	1.13%	0.081%
58	17.416	2449	2453	2457	rVB2	62422	84888	0.65%	0.047%
59	17.486	2458	2465	2469	rBV	266629	396337	3.05%	0.219%
60	17.592	2480	2483	2488	rBV4	44997	68486	0.53%	0.038%
61	17.998	2546	2552	2560	rBV	2043190	3086318	23.72%	1.703%
62	18.298	2601	2603	2609	rVB5	45412	57157	0.44%	0.032%
63	18.498	2633	2637	2642	rVB	138357	186476	1.43%	0.103%
64	18.592	2648	2653	2659	rBV	178138	247914	1.91%	0.137%
65	18.881	2699	2702	2707	rVV2	74893	109300	0.84%	0.060%
66	18.933	2708	2711	2715	rVV5	47734	69670	0.54%	0.038%
67	19.016	2721	2725	2735	rVB	146454	250504	1.93%	0.138%
68	19.145	2737	2747	2757	rBV	1761124	3349505	25.74%	1.848%
69	19.281	2766	2770	2778	rBV2	985135	1483708	11.40%	0.819%
70	19.363	2782	2784	2792	rVB2	109389	191343	1.47%	0.106%
71	19.480	2798	2804	2808	rBV2	8522241	13012761	100.00%	7.180%

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72	19.692	2836	2840	2844	rBV	71821	110300	0.85%	0.061%
73	19.839	2859	2865	2871	rBV5	202538	537157	4.13%	0.296%
74	19.975	2883	2888	2892	rBV4	193746	476219	3.66%	0.263%
75	20.051	2897	2901	2905	rVB2	132558	180640	1.39%	0.100%
76	20.163	2914	2920	2925	rBV2	269933	458661	3.52%	0.253%
77	20.304	2940	2944	2949	rBV2	192475	326713	2.51%	0.180%
78	20.351	2949	2952	2957	rVB4	89392	165475	1.27%	0.091%
79	20.551	2983	2986	2989	rBV2	161929	181915	1.40%	0.100%
80	20.680	3005	3008	3011	rVB4	91429	101716	0.78%	0.056%
81	20.757	3017	3021	3027	rBV	322518	461262	3.54%	0.255%
82	20.839	3030	3035	3042	rVB	4768777	5484046	42.14%	3.026%
83	20.922	3043	3049	3053	rVV2	255568	476978	3.67%	0.263%
84	20.963	3053	3056	3059	rVV	247299	291144	2.24%	0.161%
85	21.016	3059	3065	3083	rVB5	1096599	3112123	23.92%	1.717%
86	21.280	3103	3110	3114	rBV2	6808654	10649250	81.84%	5.876%
87	21.316	3114	3116	3123	rVB	1261951	1492049	11.47%	0.823%
88	21.451	3136	3139	3143	rVB	387455	397216	3.05%	0.219%
89	21.539	3152	3154	3158	rVB	211227	233134	1.79%	0.129%
90	21.586	3159	3162	3168	rVV2	133232	237306	1.82%	0.131%
91	21.886	3209	3213	3217	rBV2	569610	679584	5.22%	0.375%
92	22.351	3288	3292	3299	rVB	571371	803226	6.17%	0.443%
93	22.851	3370	3377	3388	rBV2	1707571	4672689	35.91%	2.578%
94	23.316	3452	3456	3459	rBV	507350	844691	6.49%	0.466%
95	23.369	3459	3465	3470	rVV	4939753	9096598	69.91%	5.019%
96	23.421	3470	3474	3481	rVB2	1118073	2032369	15.62%	1.121%
97	23.510	3482	3489	3495	rBV	3706812	6860983	52.73%	3.786%
98	24.069	3578	3584	3593	rVB	822526	1548583	11.90%	0.854%
99	24.810	3705	3710	3721	rVB	571314	1180668	9.07%	0.651%
100	25.680	3852	3858	3862	rBV	335657	748445	5.75%	0.413%

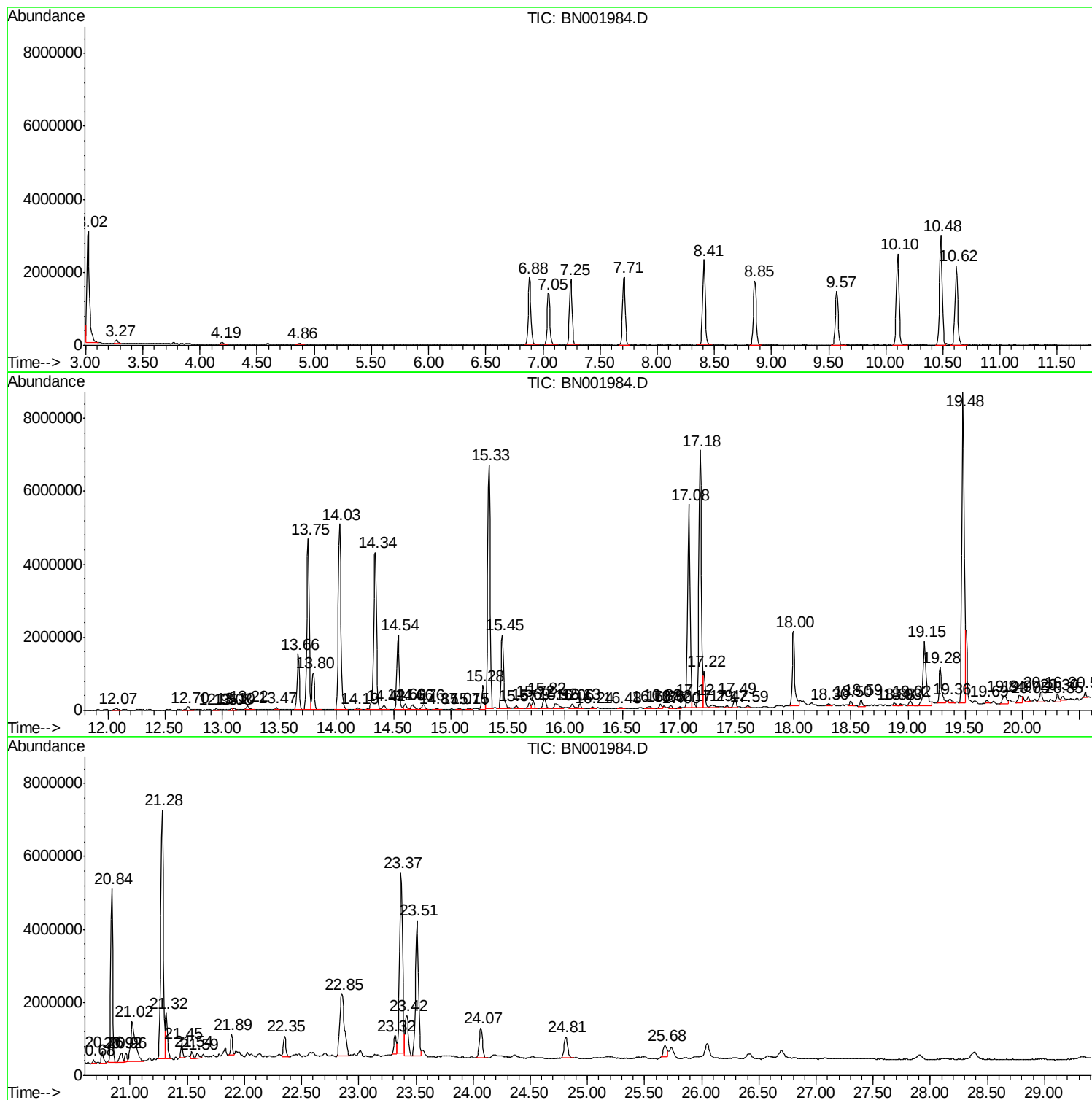
Sum of corrected areas: 181230582

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Instrument :
BNA_N
ClientSampled :
F1H11

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN061918.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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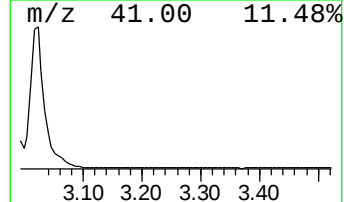
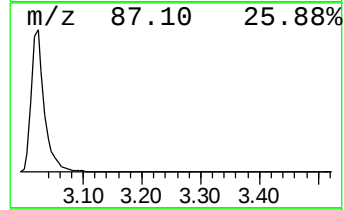
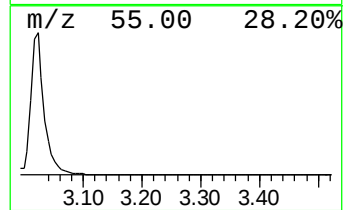
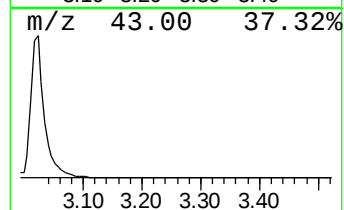
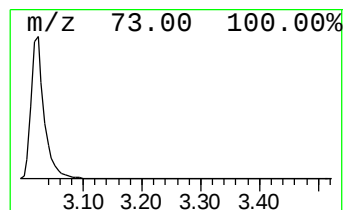
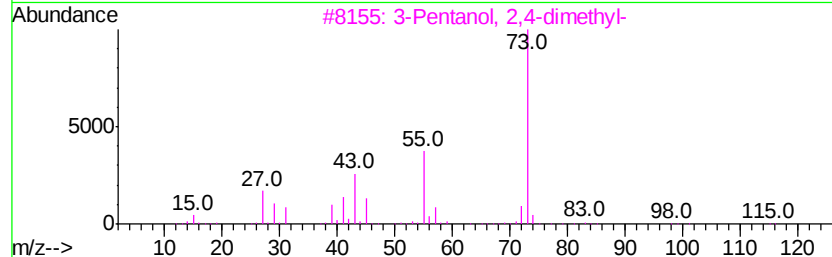
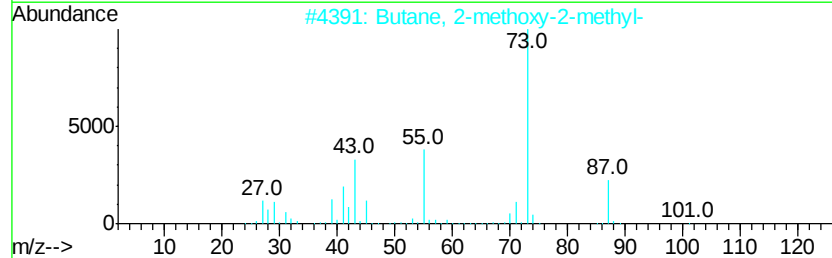
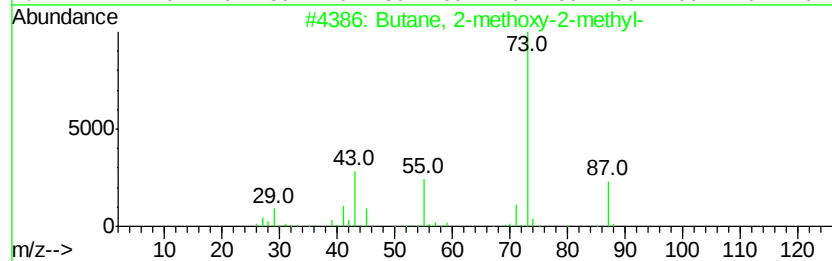
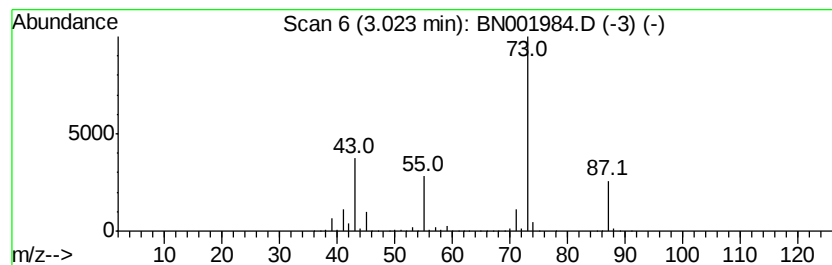
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN061918.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.02	29.97 ng/ul	4571030	1,4-Dichlorobenzene-d4	7.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	59
3		3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	37
4		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
5		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17



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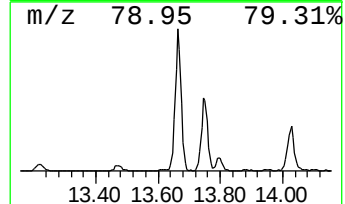
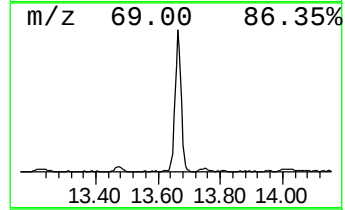
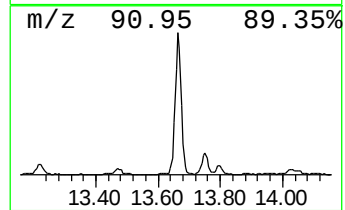
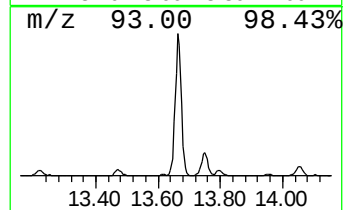
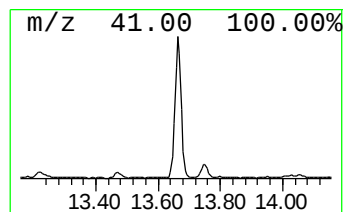
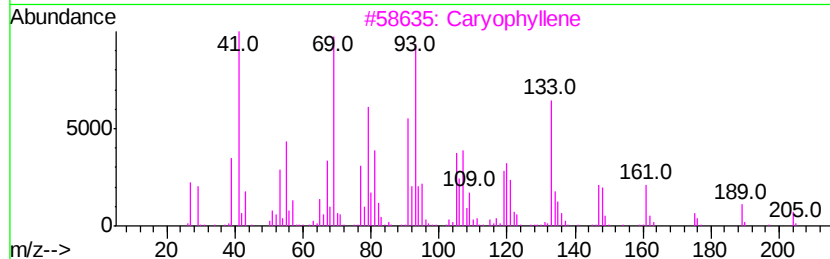
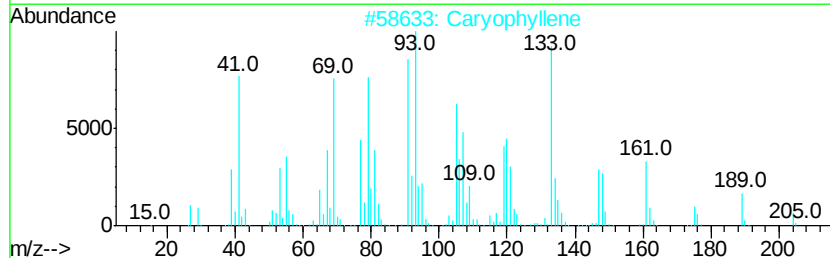
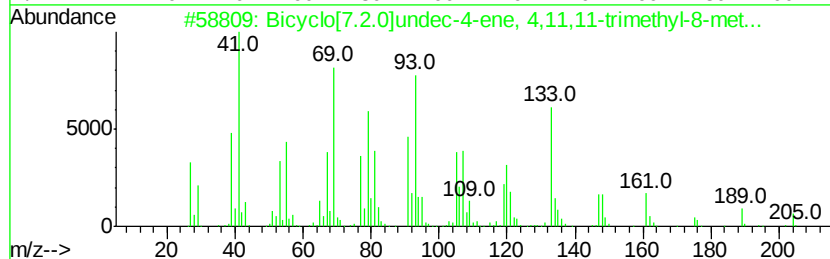
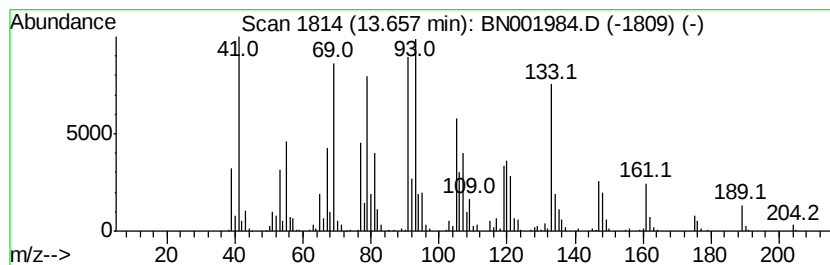
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Bicyclo[7.2.0]undec-4-ene, ... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.66	6.04 ng/ul	2162060	Acenaphthene-d10	14.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Bicyclo[7.2.0]undec-4-ene, 4,11,...	204	C15H24	000118-65-0	91
2		Carvophyllene	204	C15H24	000087-44-5	91
3		Carvophyllene	204	C15H24	000087-44-5	87
4		Bicyclo[7.2.0]undec-4-ene, 4,11,...	204	C15H24	013877-93-5	86
5		Bicyclo[7.2.0]undec-4-ene, 4,11,...	204	C15H24	013877-93-5	86



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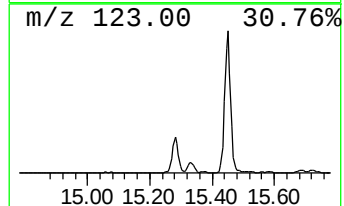
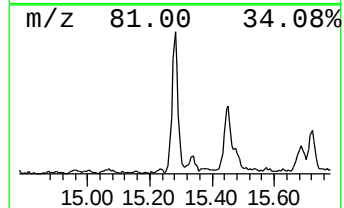
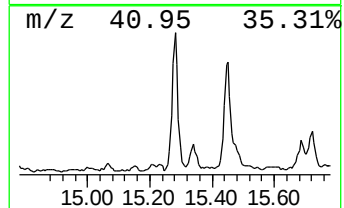
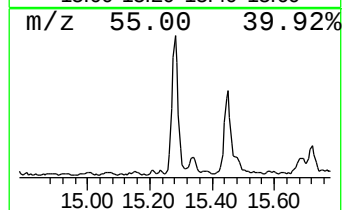
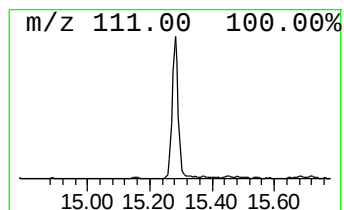
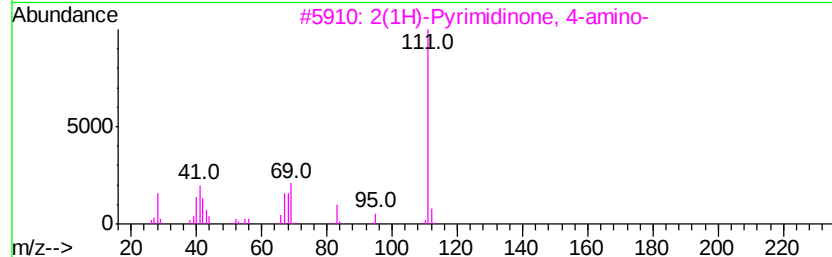
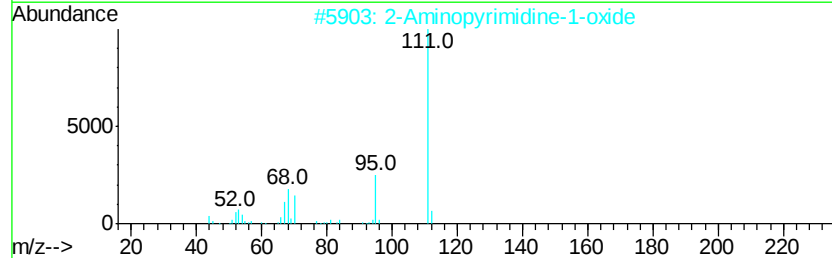
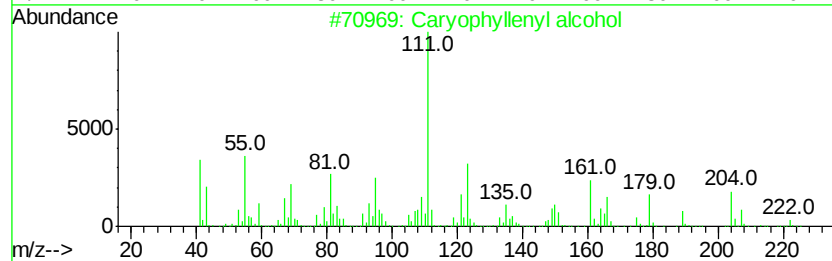
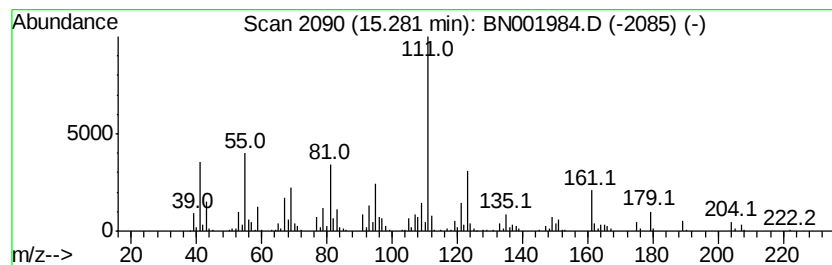
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Caryophyllenyl alcohol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.28	2.59 ng/ul	927312	Acenaphthene-d10	14.33

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Caryophyllenyl alcohol	222	C15H26O	1000131-90-4	83
2		2-Aminopyrimidine-1-oxide	111	C4H5N3O	035034-15-2	38
3		2(1H)-Pyrimidinone, 4-amino-	111	C4H5N3O	000071-30-7	38
4		Adipic acid, diphenyl ester	298	C18H18O4	003195-37-7	37
5		MONTANOL	352	C21H36O4	1000145-58-3	37



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Sample : J3765-09
Misc :
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ClientSampleId :
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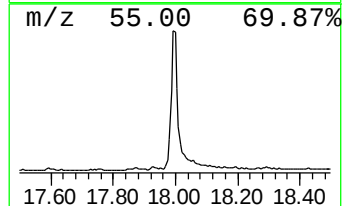
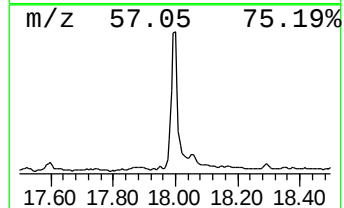
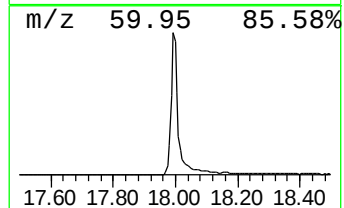
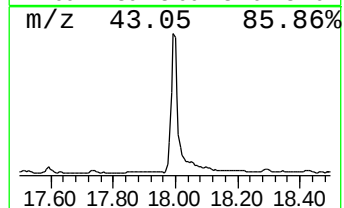
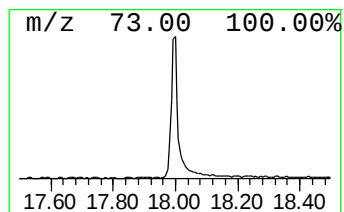
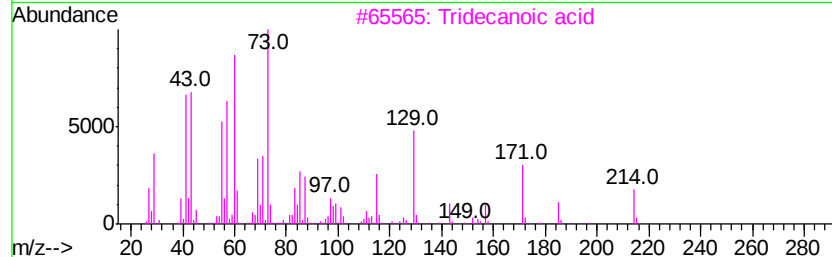
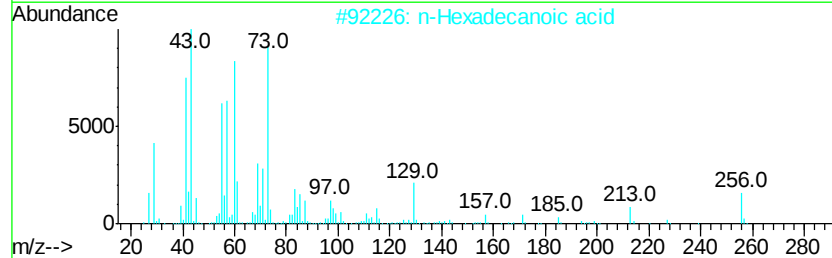
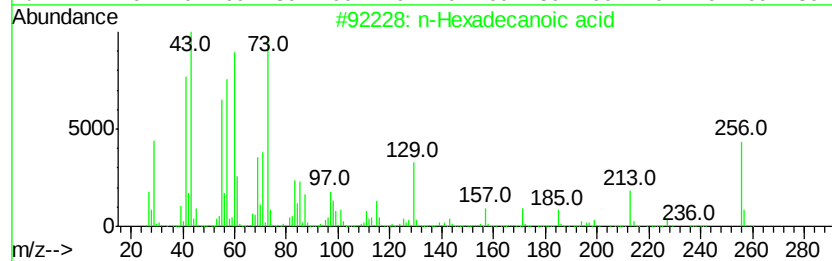
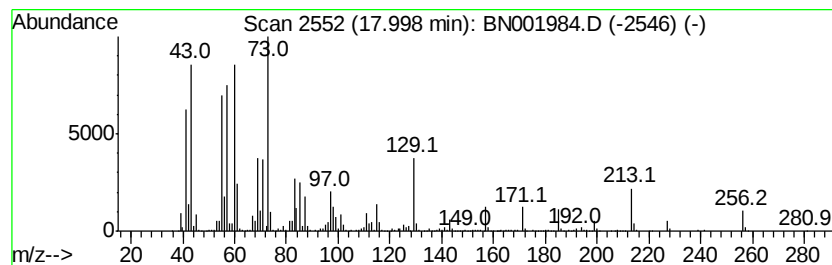
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.00	7.56 ng/ul	3086320	Phenanthrene-d10	17.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
3		Tridecanoic acid	214	C13H26O2	000638-53-9	92
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	89
5		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	89



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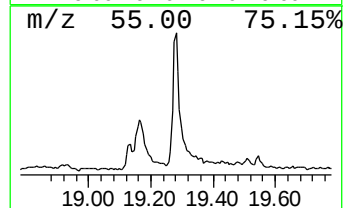
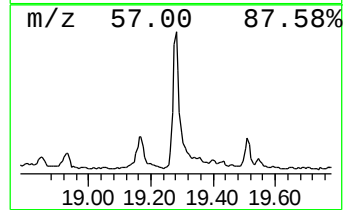
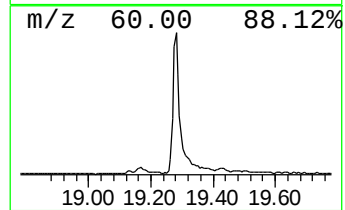
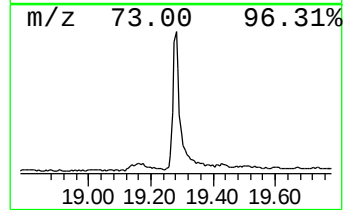
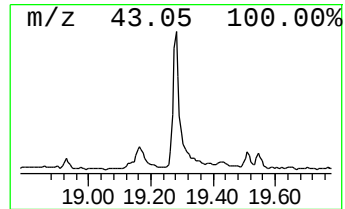
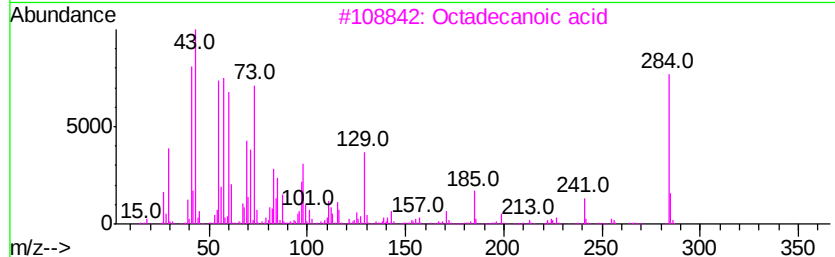
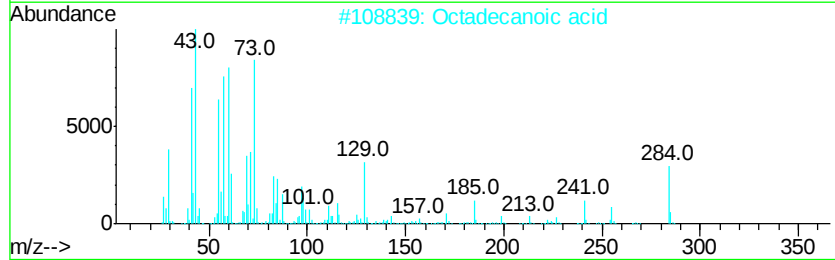
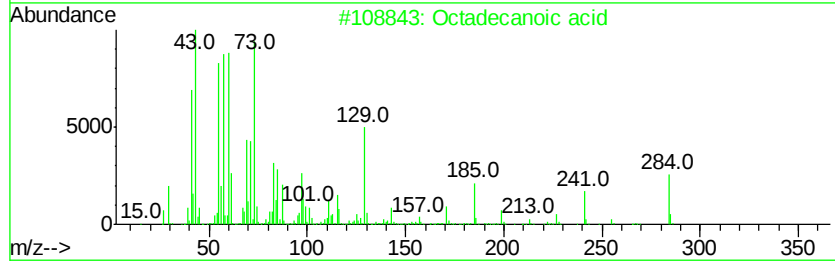
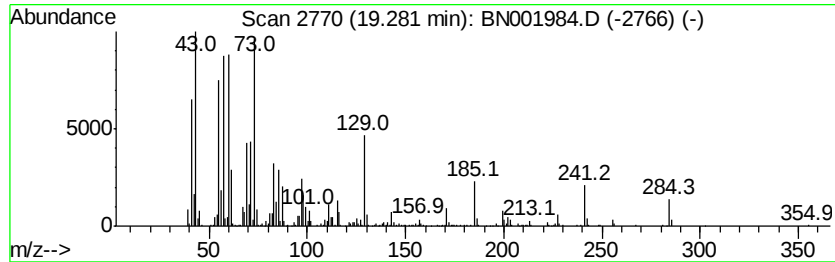
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Octadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.28	2.79 ng/ul	1483710	Chrysene-d12	21.28

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid		284	C18H36O2	000057-11-4	99
2	Octadecanoic acid		284	C18H36O2	000057-11-4	94
3	Octadecanoic acid		284	C18H36O2	000057-11-4	89
4	Pentadecanoic acid		242	C15H30O2	001002-84-2	89
5	Octadecanoic acid, 2-(2-hydroxye...		372	C22H44O4	000106-11-6	86



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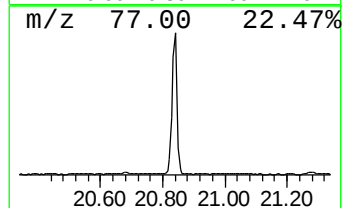
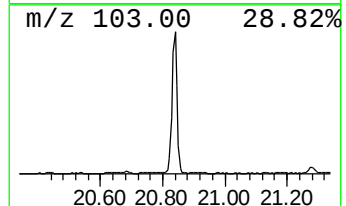
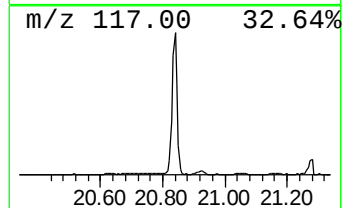
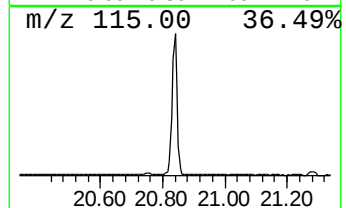
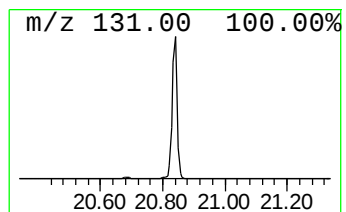
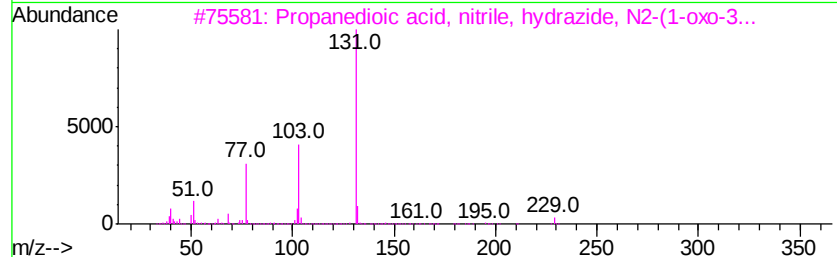
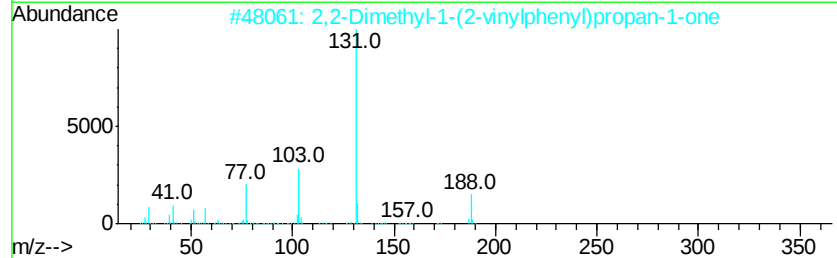
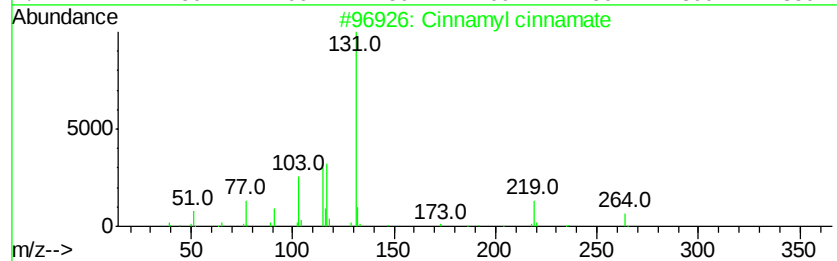
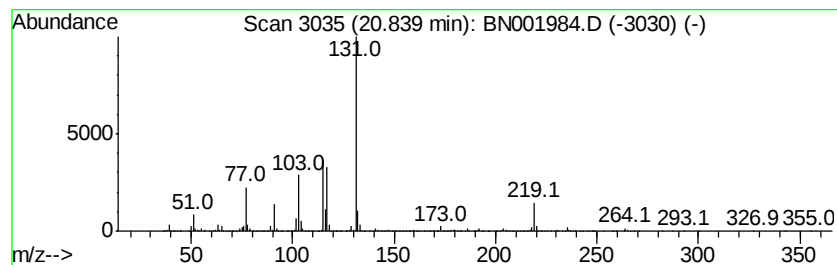
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Cinnamyl cinnamate Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.84	10.30 ng/ul	5484050	Chrysene-d12	21.28
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Cinnamyl cinnamate	264 C18H16O2	000122-69-0 87	
2	2,2-Dimethyl-1-(2-vinylphenyl)pr...	188 C13H16O	1000210-99-9 52	
3	Propanedioic acid, nitrile, hydr...	229 C12H11N3O2	294878-35-6 50	
4	Oxalic acid, monoamide monohydra...	323 C18H17N3O3	1000294-01-4 50	
5	1-Penten-3-one, 4,4-dimethyl-1-p...	188 C13H16O	000538-44-3 50	



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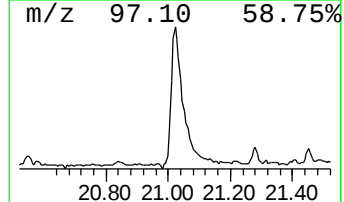
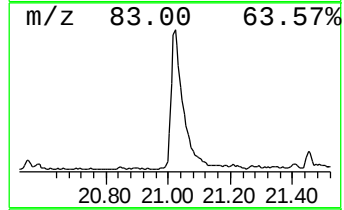
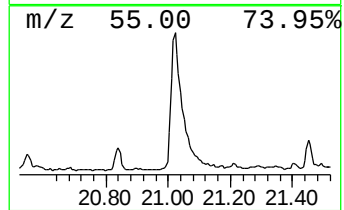
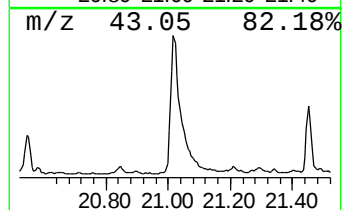
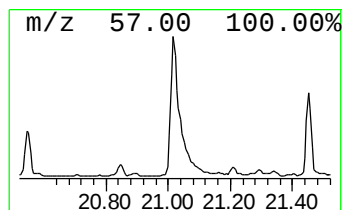
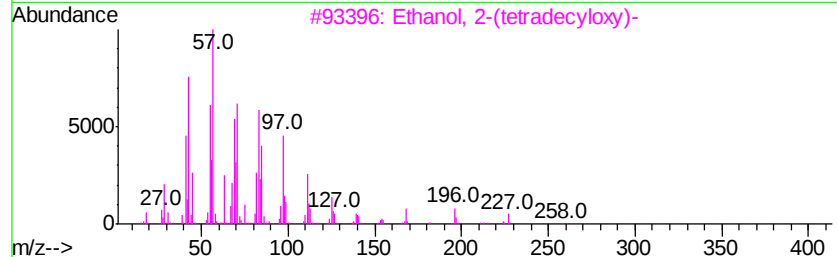
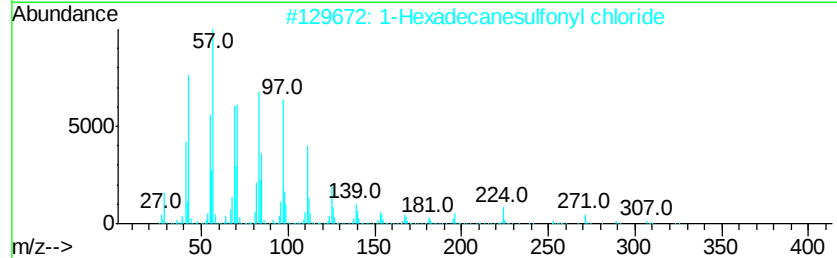
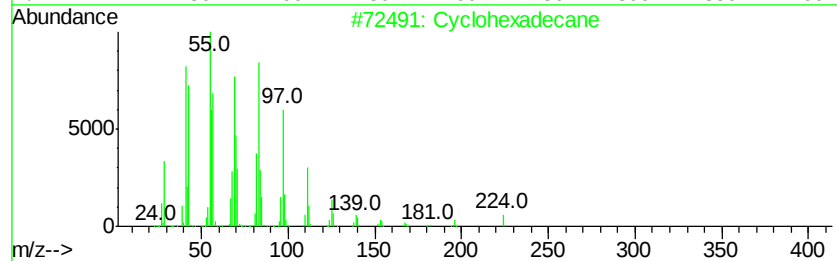
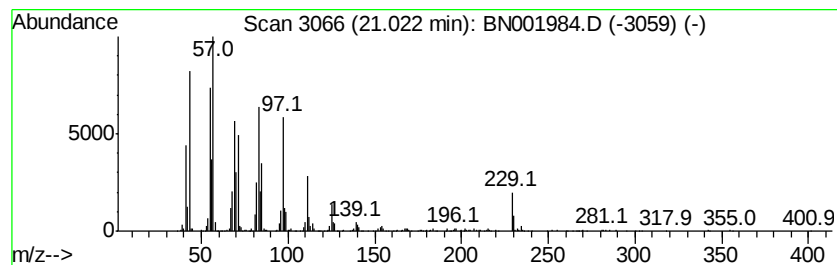
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 (DEL) Alkane: Branched21.02 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.02	5.84 ng/ul	3112120	Chrysene-d12	21.28

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexadecane	224	C16H32	000295-65-8	90
2		1-Hexadecanesulfonyl chloride	324	C16H33ClO2S	038775-38-1	83
3		Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	76
4		Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	64
5		Cyclotetracosane	336	C24H48	000297-03-0	64



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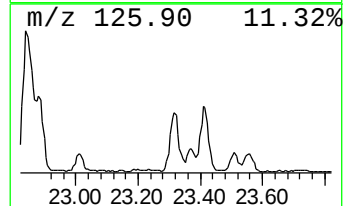
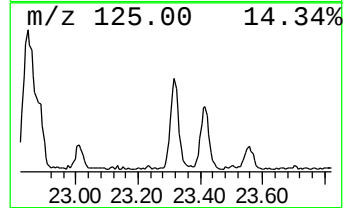
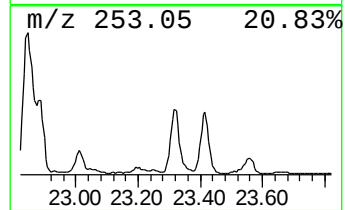
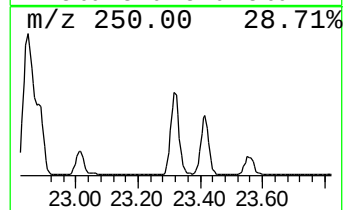
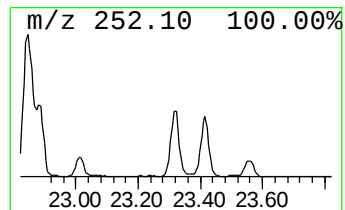
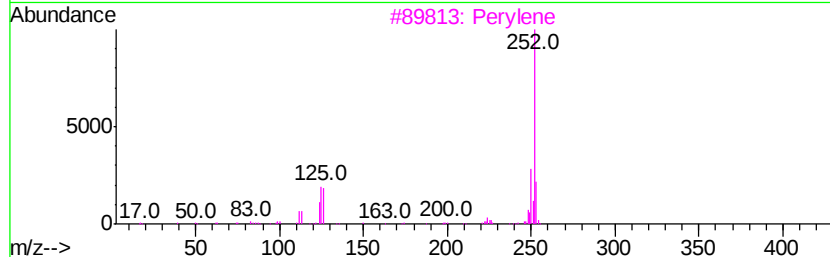
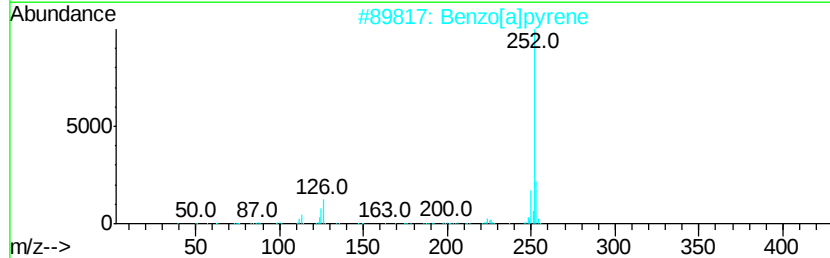
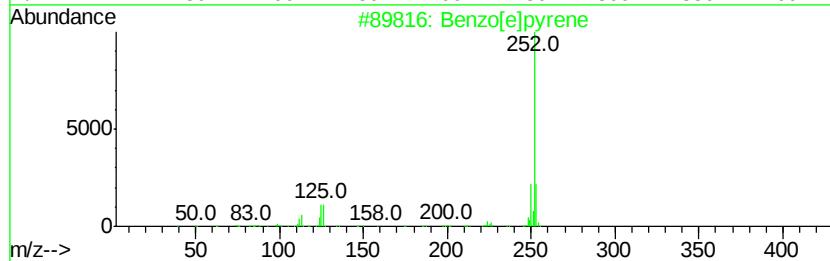
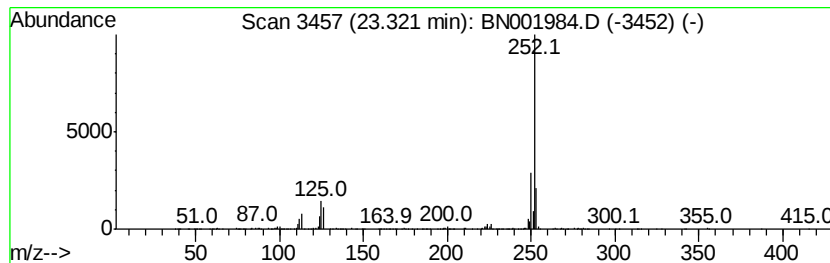
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Benzo[e]pyrene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.32	2.46 ng/ul	844691	Perylene-d12	23.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[e]pyrene	252	C20H12	000192-97-2	99
2		Benzo[a]pyrene	252	C20H12	000050-32-8	95
3		Perylene	252	C20H12	000198-55-0	95
4		Benzo[k]fluoranthene	252	C20H12	000207-08-9	93
5		Benz[e]acephenanthrylene	252	C20H12	000205-99-2	90



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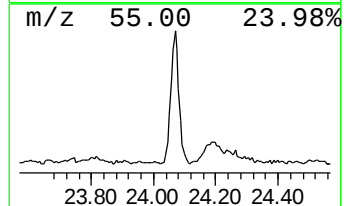
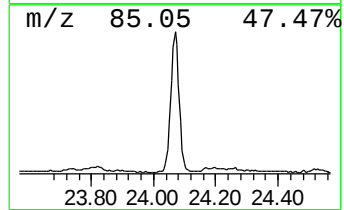
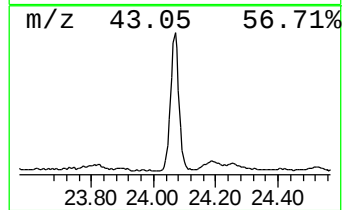
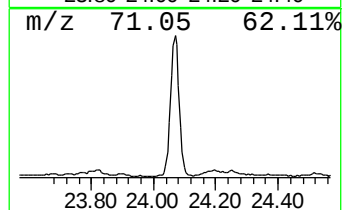
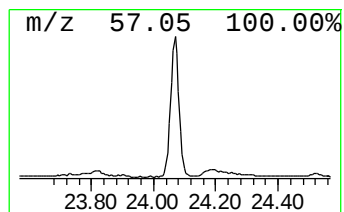
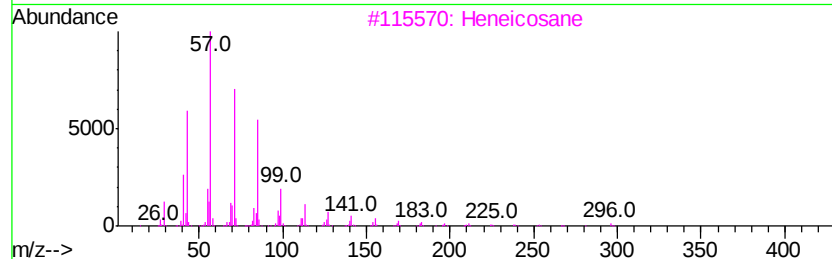
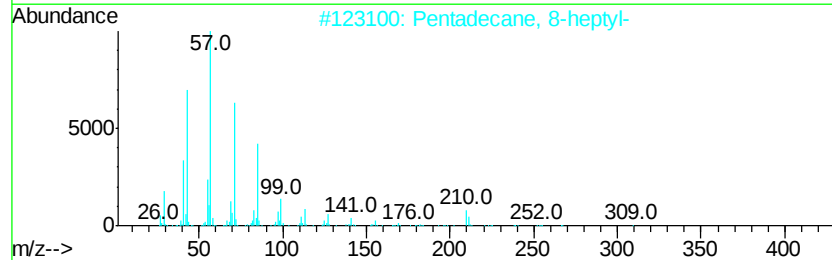
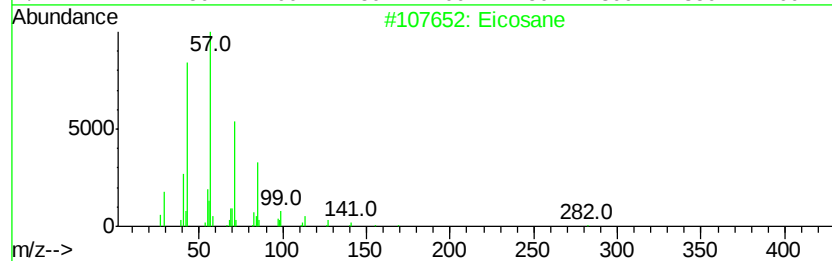
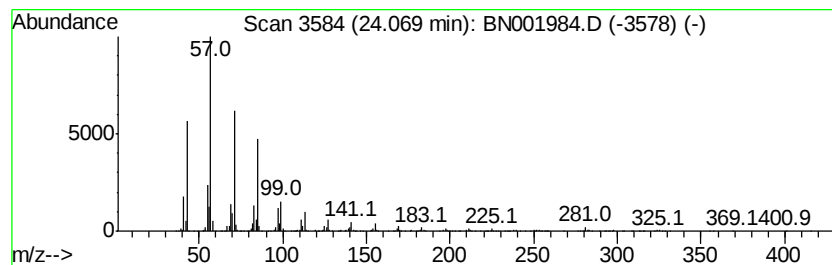
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TIC Integration Parameters: LSCINT.P

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.07	4.51 ng/ul	1548580	Perylene-d12	23.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	94
2		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	91
3		Heneicosane	296	C21H44	000629-94-7	91
4		Docosane, 11-butyl-	366	C26H54	013475-76-8	91
5		Tetracosane	338	C24H50	000646-31-1	91



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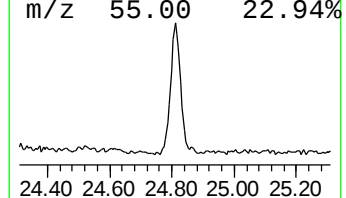
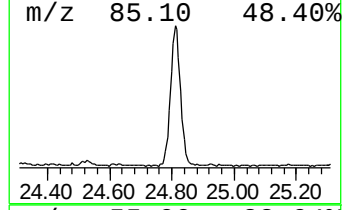
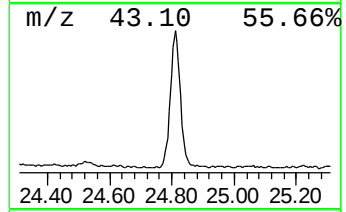
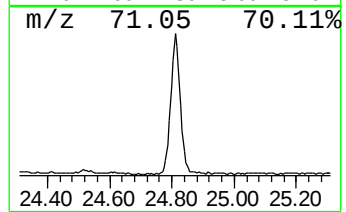
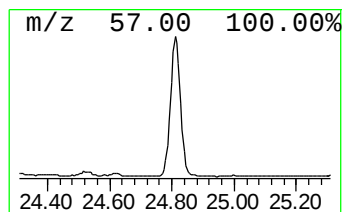
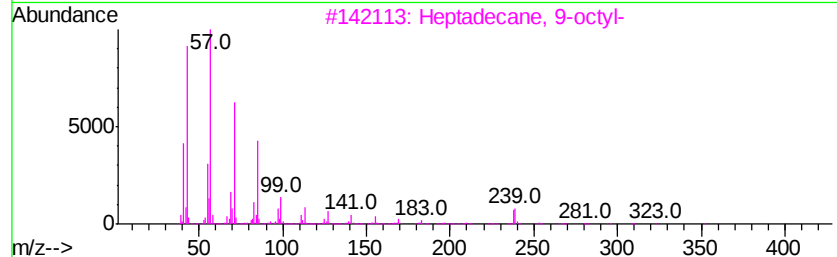
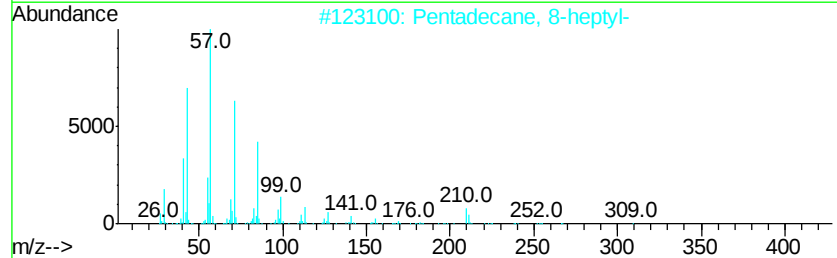
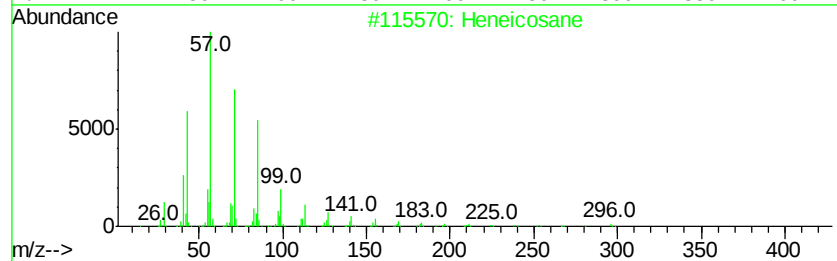
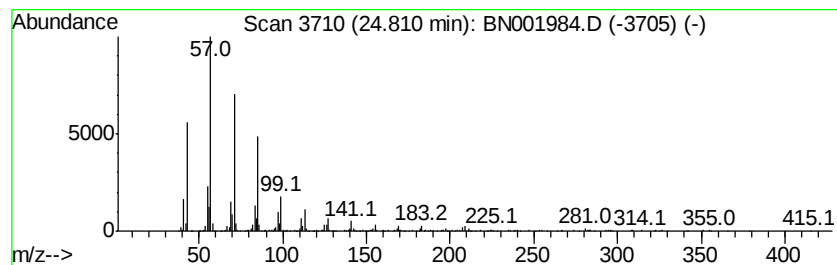
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.81	3.44 ng/ul	1180670	Perylene-d12	23.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	91
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4		Tetracosane	338	C24H50	000646-31-1	91
5		Tetratetracontane	619	C44H90	007098-22-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN070618\
Data File : BN001984.D
Acq On : 06 Jul 2018 11:41
Operator : JU/SJ
Sample : J3765-09
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
F1H11

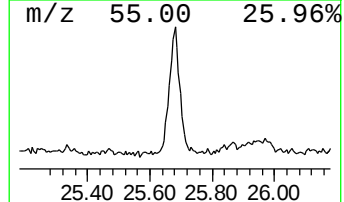
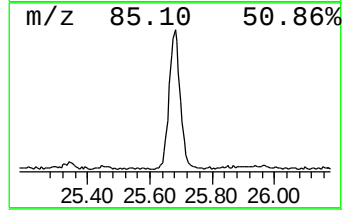
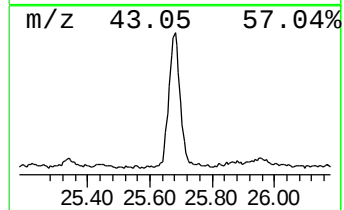
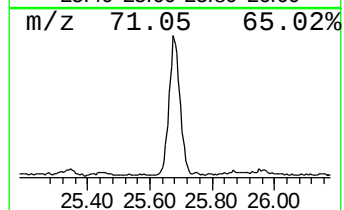
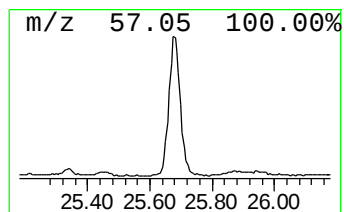
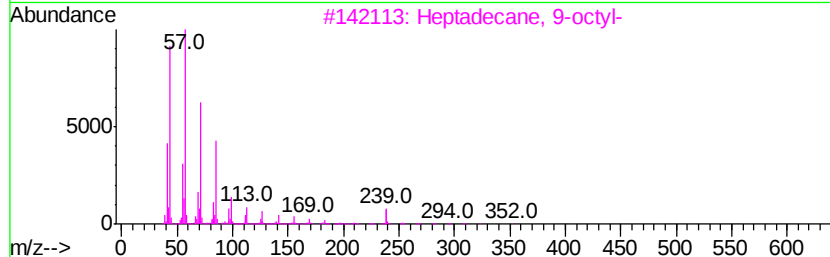
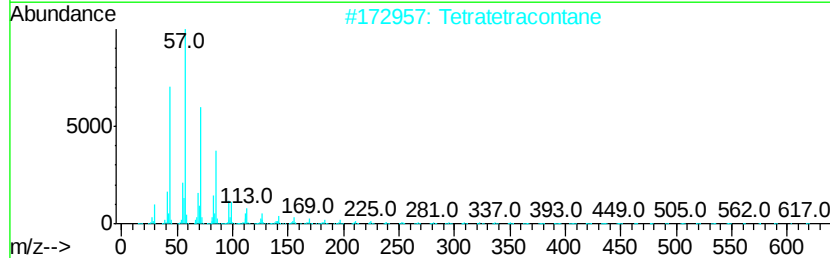
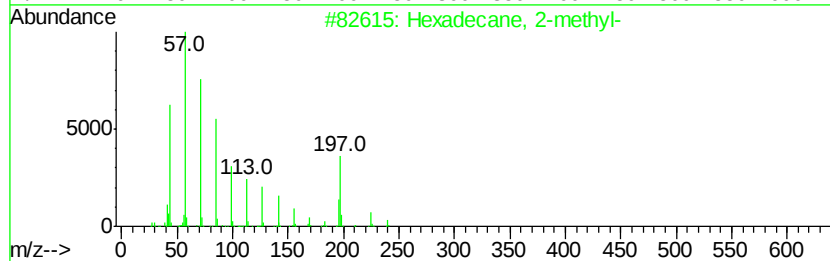
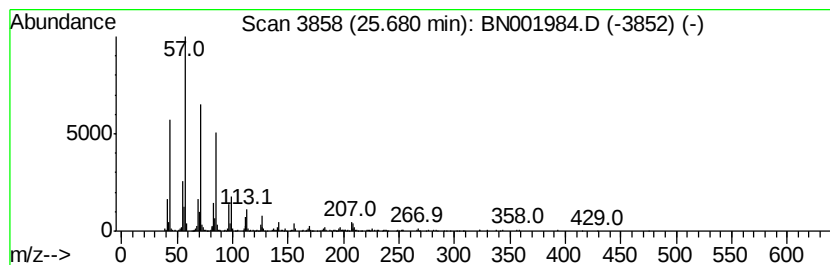
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.68	2.18 ng/ul	748445	Perylene-d12	23.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane, 2-methyl-	240	C17H36	001560-92-5	90
2		Tetratetracontane	619	C44H90	007098-22-8	90
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	90
4		Nonacosane	408	C29H60	000630-03-5	90
5		Octacosane	394	C28H58	000630-02-4	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN070618\
 Data File : BN001984.D
 Acq On : 06 Jul 2018 11:41
 Operator : JU/SJ
 Sample : J3765-09
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F1H11

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN061918.M
 Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	3.02	30.0	ng/ul	4571030	1	7.71	3050320	20.0
Bicyclo[7.2.0]und...	13.66	6.0	ng/ul	2162060	3	14.33	7164740	20.0
Caryophyllenyl al...	15.28	2.6	ng/ul	927312	3	14.33	7164740	20.0
n-Hexadecanoic acid	18.00	7.6	ng/ul	3086320	4	17.08	8160480	20.0
Octadecanoic acid	19.28	2.8	ng/ul	1483710	5	21.28	10649300	20.0
Cinnamyl cinnamate	20.84	10.3	ng/ul	5484050	5	21.28	10649300	20.0
(DEL) Alkane: Bra...	21.02	5.8	ng/ul	3112120	5	21.28	10649300	20.0
Benzo[e]pyrene	23.32	2.5	ng/ul	844691	6	23.51	6860980	20.0
(DEL) Alkane: Str...	24.07	4.5	ng/ul	1548580	6	23.51	6860980	20.0
(DEL) Alkane: Str...	24.81	3.4	ng/ul	1180670	6	23.51	6860980	20.0
(DEL) Alkane: Str...	25.68	2.2	ng/ul	748445	6	23.51	6860980	20.0