

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN040120\
 Data File : BN010366.D
 Acq On : 01 Apr 2020 18:49
 Operator : CG/JU
 Sample : L2093-01
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBEW7

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-BN040120MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BN010367.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.893	14	18	26	rBV	98486	186771	0.55%	0.036%
2	2.964	26	30	47	rVB	247838	450355	1.33%	0.087%
3	3.211	68	72	73	rBV	66118	73639	0.22%	0.014%
4	3.240	73	77	96	rVB	324791	568289	1.67%	0.110%
5	3.917	189	192	197	rVB2	30489	40823	0.12%	0.008%
6	4.393	268	273	280	rVB4	21582	52750	0.16%	0.010%
7	4.793	336	341	355	rBV	689224	1053906	3.10%	0.204%
8	6.764	669	676	679	rBV	2760518	4811156	14.16%	0.931%
9	6.793	679	681	683	rVV	2043870	2761427	8.13%	0.534%
10	6.822	683	686	695	rVB	2790586	3966195	11.68%	0.767%
11	6.958	701	709	724	rVB	1517202	2391616	7.04%	0.463%
12	7.152	734	742	745	rBV	2110896	3718463	10.95%	0.719%
13	7.181	745	747	758	rVB	1937348	2492594	7.34%	0.482%
14	7.611	811	820	828	rBV	1739402	2665947	7.85%	0.516%
15	7.693	828	834	841	rVB2	38161	71370	0.21%	0.014%
16	8.311	931	939	949	rBV	2610096	4204868	12.38%	0.814%
17	8.411	949	956	965	rVB	1837753	2698423	7.94%	0.522%
18	8.746	1005	1013	1017	rBV	1888534	3156058	9.29%	0.611%
19	8.787	1017	1020	1029	rVB	1535544	2301356	6.77%	0.445%
20	9.234	1091	1096	1104	rVB	95617	145767	0.43%	0.028%
21	9.463	1127	1135	1137	rBV	1557490	2459039	7.24%	0.476%
22	9.493	1137	1140	1156	rVB	1421777	2311817	6.81%	0.447%
23	9.999	1218	1226	1228	rBV	3015558	5136406	15.12%	0.994%
24	10.369	1282	1289	1293	rBV	2572186	4242859	12.49%	0.821%
25	10.422	1293	1298	1305	rVV	2674445	4331890	12.75%	0.838%
26	10.505	1305	1312	1324	rVV	1566665	2714903	7.99%	0.525%
27	10.722	1338	1349	1357	rVB	3271917	5036517	14.83%	0.974%
28	11.657	1498	1508	1516	rBV	4946855	7655412	22.54%	1.481%
29	11.910	1543	1551	1556	rVV	207781	329016	0.97%	0.064%
30	11.963	1556	1560	1566	rVB	62402	105520	0.31%	0.020%
31	12.257	1600	1610	1619	rBV	4465030	7078726	20.84%	1.370%
32	12.416	1629	1637	1644	rBV	3359967	5000751	14.72%	0.967%
33	12.657	1671	1678	1685	rBV	5149151	7520168	22.14%	1.455%
34	13.063	1739	1747	1751	rVV	6117795	9812098	28.88%	1.898%

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35	13.099	1751	1753	1758	rVV	3283111	4243610	12.49%	0.821%
36	13.157	1758	1763	1776	rVB	1814449	2780684	8.19%	0.538%
37	13.657	1841	1848	1852	rBV	5248879	7680335	22.61%	1.486%
38	13.704	1852	1856	1864	rVB	5124513	7245551	21.33%	1.402%
39	13.810	1868	1874	1882	rVB	3068381	3955845	11.64%	0.765%
40	13.928	1886	1894	1903	rBV	6061894	9209853	27.11%	1.782%
41	14.175	1932	1936	1940	rBV3	33250	40660	0.12%	0.008%
42	14.240	1940	1947	1952	rBV	4059966	6154253	18.12%	1.191%
43	14.299	1952	1957	1967	rVB	5548175	8127777	23.93%	1.572%
44	14.446	1970	1982	2001	rBV2	3367895	6874656	20.24%	1.330%
45	14.640	2008	2015	2021	rBV	4206221	5986117	17.62%	1.158%
46	14.869	2047	2054	2068	rBV	3646204	5217099	15.36%	1.009%
47	15.075	2084	2089	2094	rVB3	43302	64047	0.19%	0.012%
48	15.134	2094	2099	2103	rBV	77977	103355	0.30%	0.020%
49	15.234	2109	2116	2121	rBV	8389987	11885346	34.99%	2.299%
50	15.293	2121	2126	2131	rVV	3881202	5000616	14.72%	0.967%
51	15.363	2131	2138	2151	rVV2	4731940	8330428	24.52%	1.612%
52	15.469	2152	2156	2163	rVB2	159704	258700	0.76%	0.050%
53	15.898	2223	2229	2234	rVV	262239	330388	0.97%	0.064%
54	15.957	2235	2239	2242	rVV6	25462	44623	0.13%	0.009%
55	16.022	2243	2250	2258	rVB6	43501	111126	0.33%	0.021%
56	16.175	2269	2276	2279	rBV	112531	155449	0.46%	0.030%
57	16.298	2291	2297	2303	rVV	4346958	5946798	17.51%	1.151%
58	16.640	2345	2355	2367	rBV	6598290	8827557	25.99%	1.708%
59	16.934	2400	2405	2408	rBV	183175	245000	0.72%	0.047%
60	16.981	2408	2413	2417	rVV	5817398	8464796	24.92%	1.638%
61	17.028	2417	2421	2425	rVV	6020156	8467326	24.93%	1.638%
62	17.081	2425	2430	2433	rVV2	9375874	13325206	39.23%	2.578%
63	17.116	2433	2436	2445	rVV	7395724	9594878	28.24%	1.856%
64	17.192	2445	2449	2455	rVB	48793	73574	0.22%	0.014%
65	17.387	2474	2482	2488	rVV	6193471	8897819	26.19%	1.721%
66	17.463	2489	2495	2503	rVB	5830761	7679038	22.60%	1.486%
67	17.634	2521	2524	2527	rVV3	36630	41301	0.12%	0.008%
68	17.675	2527	2531	2539	rVB9	32835	68452	0.20%	0.013%
69	17.910	2566	2571	2576	rBV	592791	894051	2.63%	0.173%
70	17.969	2576	2581	2588	rVB	5346145	6590736	19.40%	1.275%
71	18.216	2619	2623	2631	rVB7	81986	142905	0.42%	0.028%

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72	18.392	2648	2653	2658	rVB	89776	137317	0.40%	0.027%
73	18.775	2714	2718	2724	rBV5	79291	132108	0.39%	0.026%
74	18.851	2727	2731	2737	rVB	139759	202308	0.60%	0.039%
75	19.051	2754	2765	2771	rBV	16362635	22394245	65.92%	4.333%
76	19.198	2786	2790	2797	rBV	347515	552679	1.63%	0.107%
77	19.269	2798	2802	2803	rBV	100972	126450	0.37%	0.024%
78	19.381	2815	2821	2824	rBV2	11710330	17822505	52.46%	3.448%
79	19.410	2824	2826	2835	rVB	11931510	14951827	44.01%	2.893%
80	19.539	2845	2848	2852	rBV6	56883	70123	0.21%	0.014%
81	19.933	2912	2915	2919	rBV5	49081	56745	0.17%	0.011%
82	19.975	2919	2922	2925	rBV	96128	102176	0.30%	0.020%
83	20.333	2979	2983	2991	rBV	6143576	7043706	20.73%	1.363%
84	20.475	3004	3007	3016	rVB4	173804	215140	0.63%	0.042%
85	20.928	3079	3084	3099	rBV3	2471146	3590437	10.57%	0.695%
86	21.128	3114	3118	3121	rBV	9169239	10133340	29.83%	1.961%
87	21.175	3121	3126	3131	rVV2	18242469	32038593	94.31%	6.199%
88	21.222	3131	3134	3146	rVB	13332432	15732806	46.31%	3.044%
89	21.375	3156	3160	3167	rBV	1725681	1874786	5.52%	0.363%
90	21.804	3229	3233	3241	rBV	3343572	3743316	11.02%	0.724%
91	22.257	3305	3310	3323	rVB	4409024	5465715	16.09%	1.057%
92	22.716	3381	3388	3392	rBV	10336727	18318662	53.92%	3.544%
93	22.751	3392	3394	3400	rVB	4524472	4904083	14.44%	0.949%
94	23.227	3467	3475	3478	rBV	9994308	18700170	55.05%	3.618%
95	23.269	3478	3482	3486	rVV	5811848	8930547	26.29%	1.728%
96	23.357	3492	3497	3514	rVB2	6012317	11108469	32.70%	2.149%
97	23.922	3587	3593	3600	rBV	3095430	5499569	16.19%	1.064%
98	23.992	3601	3605	3614	rVB2	593572	1033288	3.04%	0.200%
99	24.639	3708	3715	3724	rVB	1558465	3415577	10.05%	0.661%
100	25.521	3851	3865	3875	rBV2	11704668	33971083	100.00%	6.572%

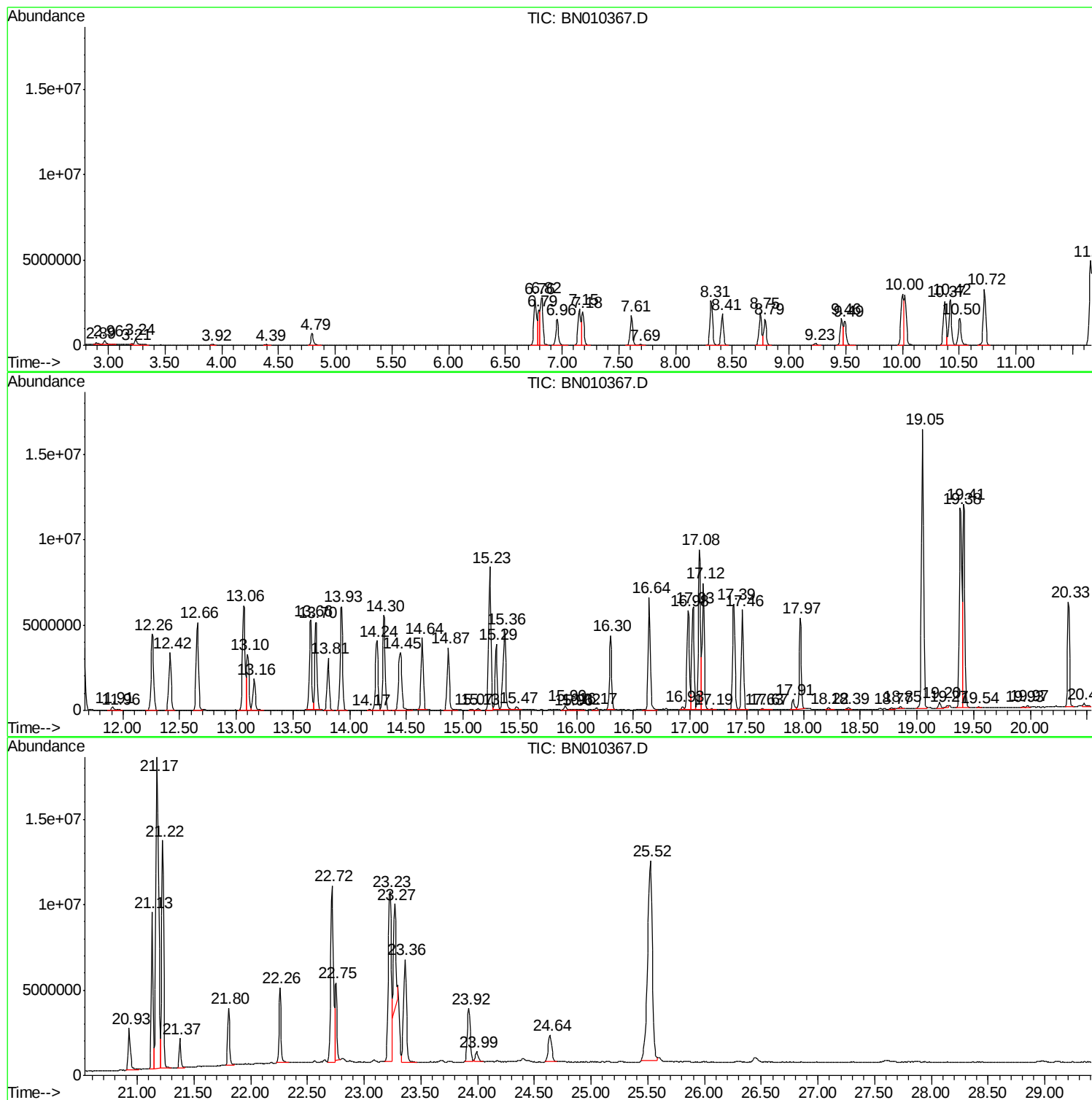
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.M
Quant Title : SVOA CALIBRATION

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



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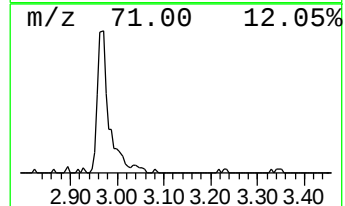
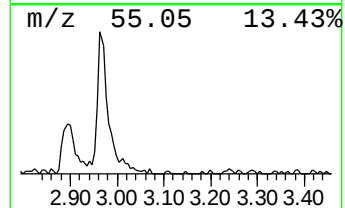
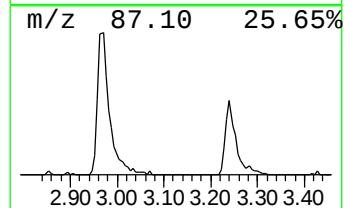
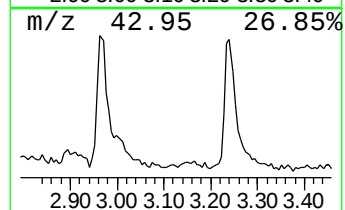
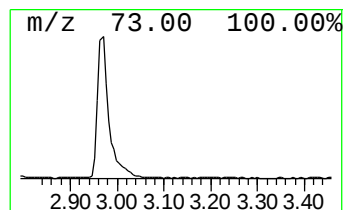
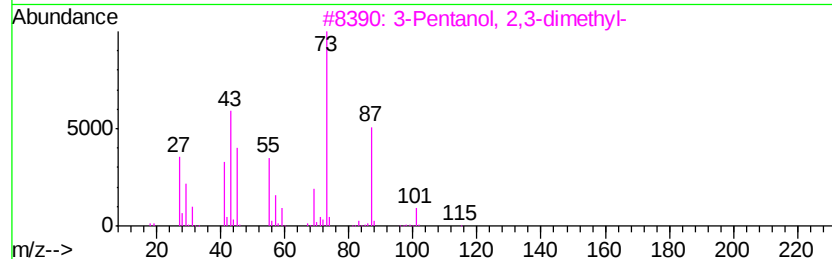
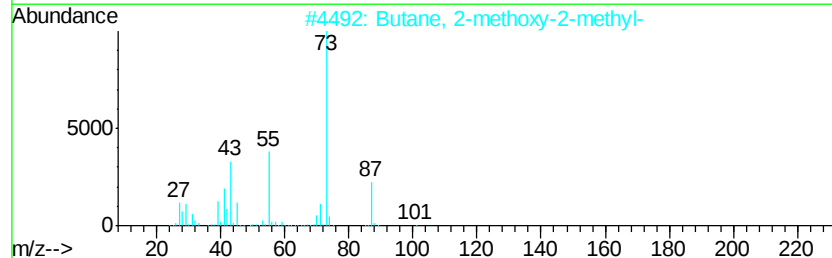
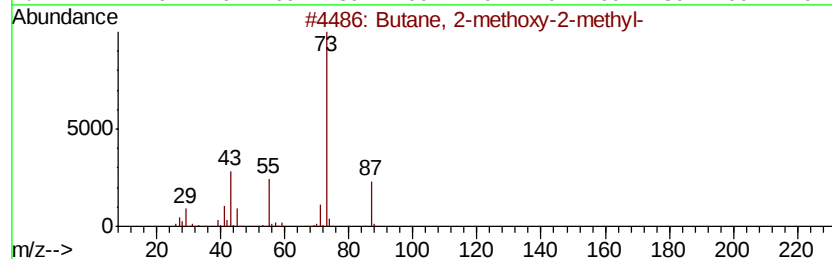
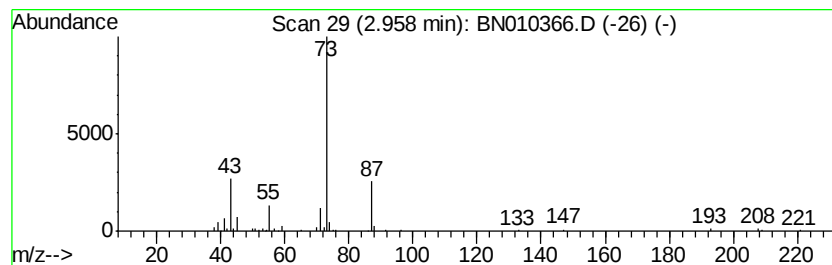
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.96	3.38 ng/ul	450355	1,4-Dichlorobenzene-d4	7.61

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	56
2		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	56
3		3-Pentanol, 2,3-dimethyl-	116	C7H16O	000595-41-5	28
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
5		N-Ethylformamide	73	C3H7NO	000627-45-2	9



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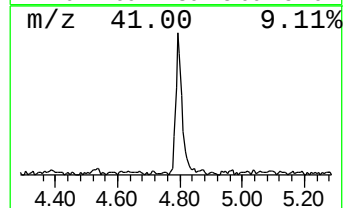
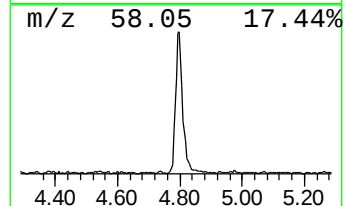
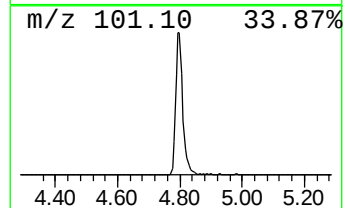
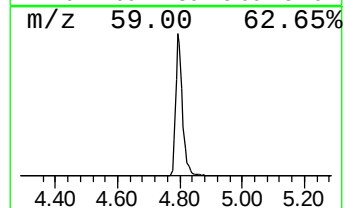
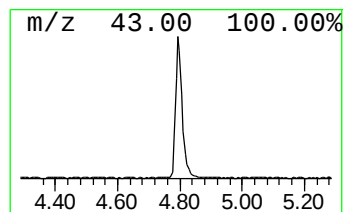
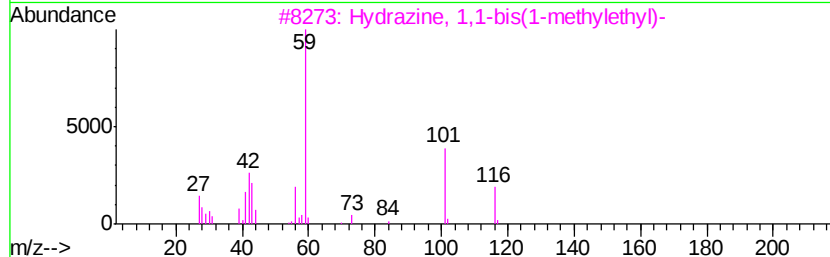
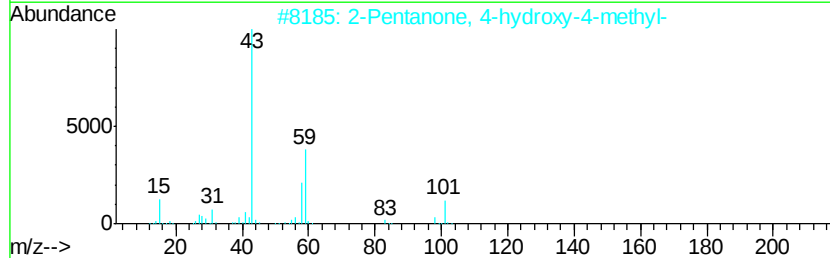
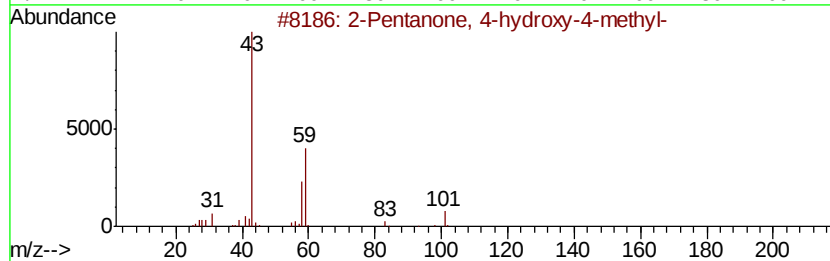
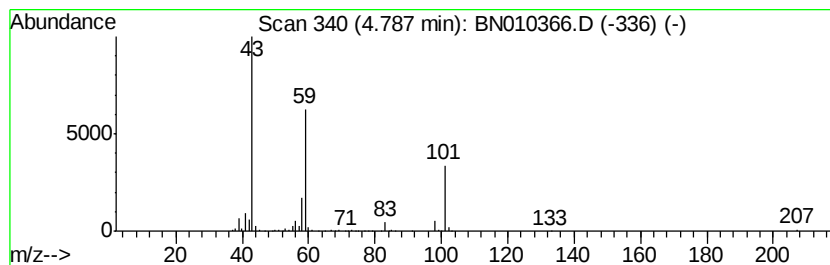
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.79	7.91 ng/ul	1053910	1,4-Dichlorobenzene-d4	7.61

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
3		Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	47
4		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38
5		Dimethadione	129	C5H7NO3	000695-53-4	33



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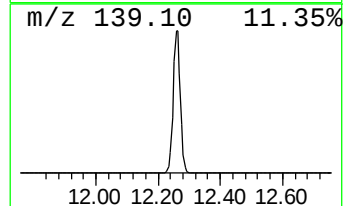
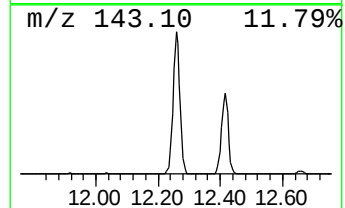
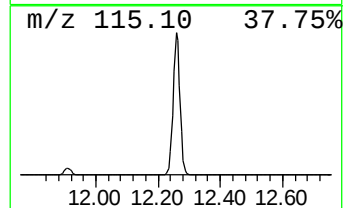
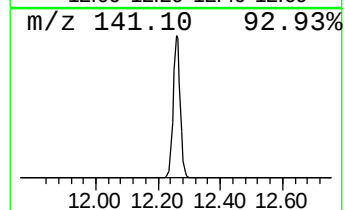
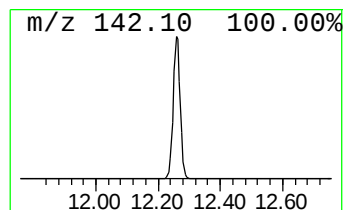
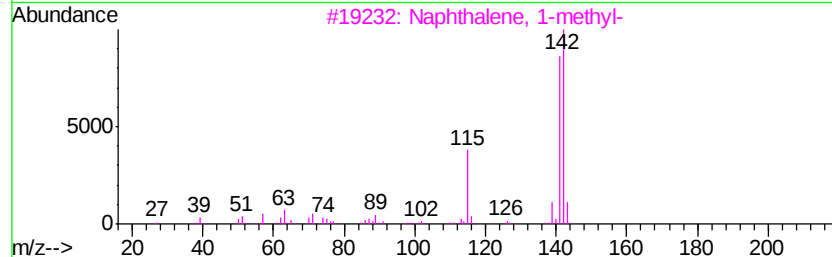
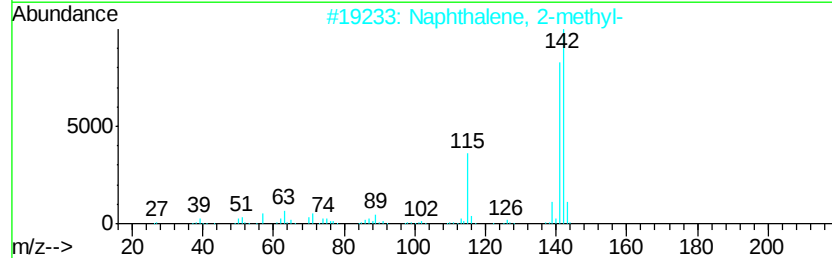
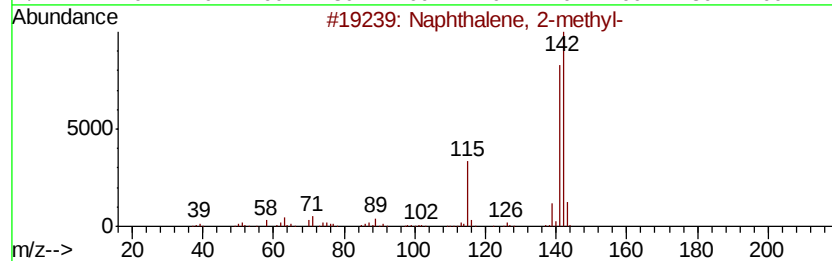
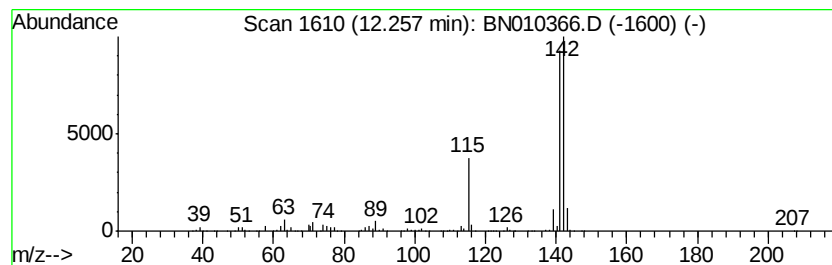
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Naphthalene, 1-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.26	33.37 ng/ul	7078730	Naphthalene-d8	10.37

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
4			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
5			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91



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Sample : L2093-01
Misc :
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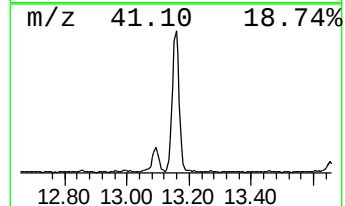
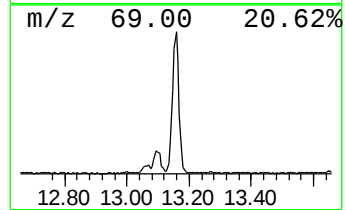
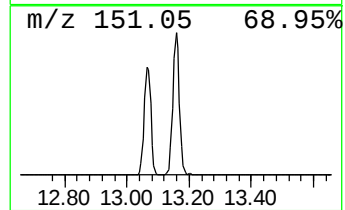
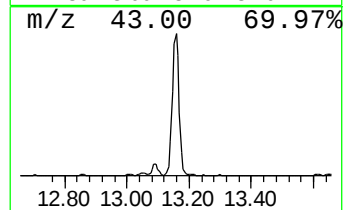
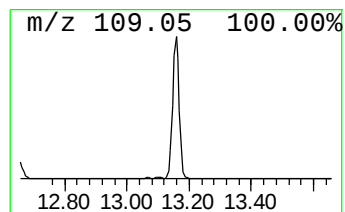
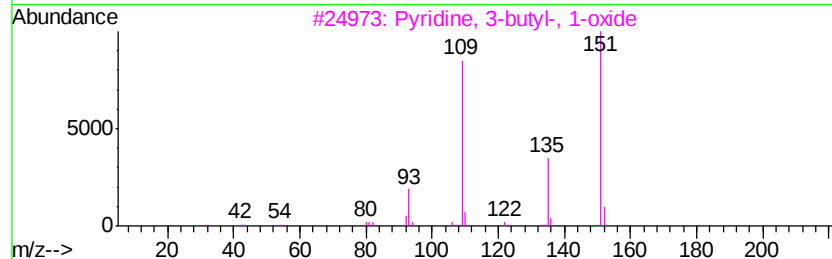
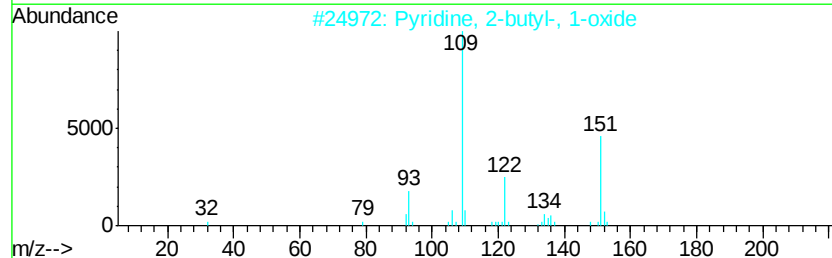
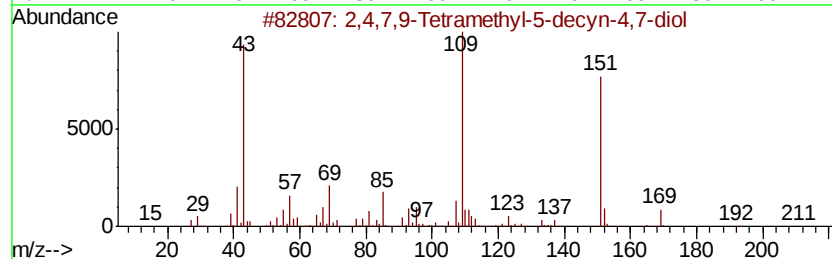
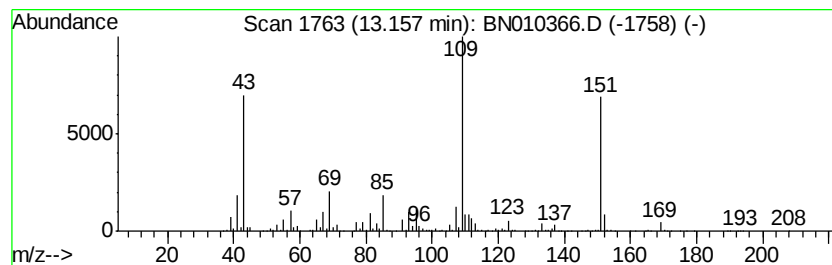
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.16	9.04 ng/ul	2780680	Acenaphthene-d10	14.24

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2		000126-86-3	64
2	Pyridine, 2-butyl-, 1-oxide	151	C9H13NO		031396-32-4	59
3	Pyridine, 3-butyl-, 1-oxide	151	C9H13NO		031396-33-5	53
4	m-Isopropoxyvaniline	151	C9H13NO		041406-00-2	53
5	2-Propenal, 3-(2,2,6-trimethyl-7...	194	C12H18O2		055759-91-6	50



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN040120\
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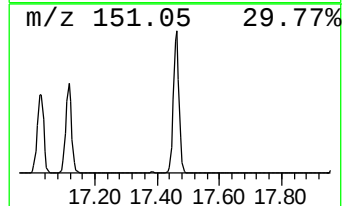
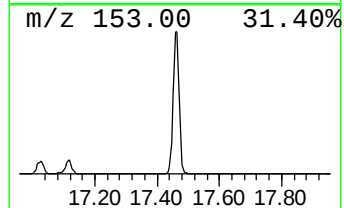
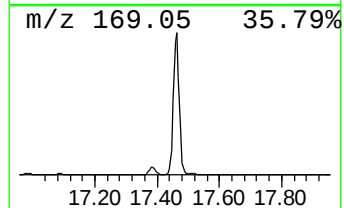
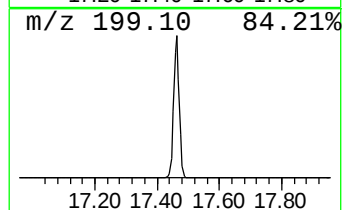
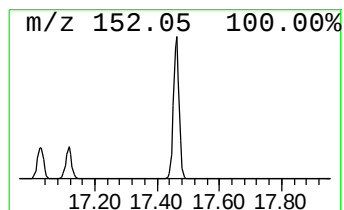
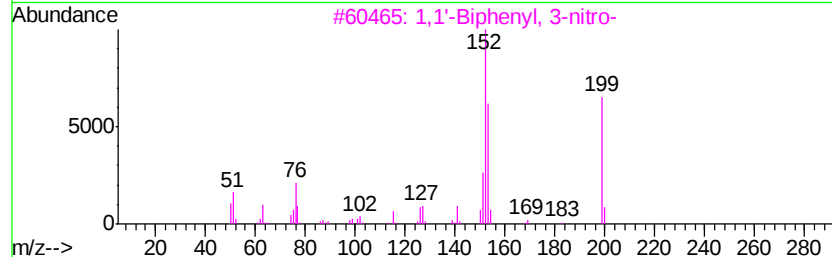
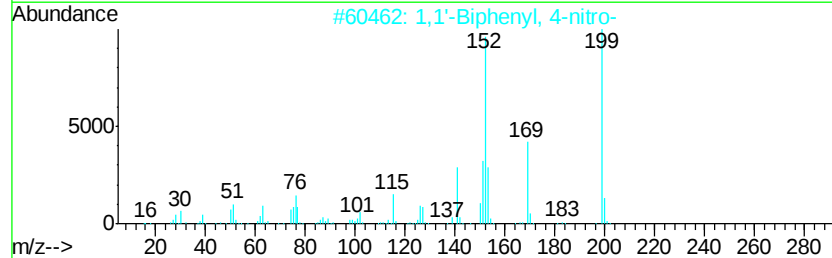
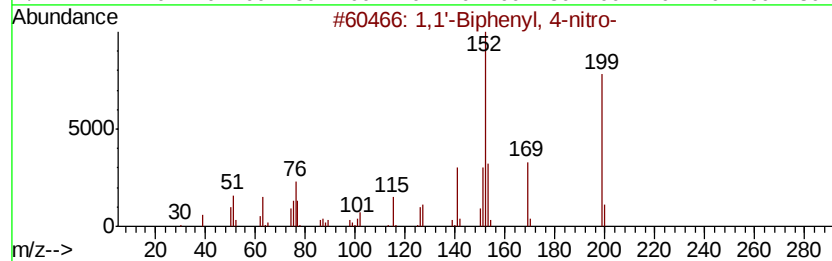
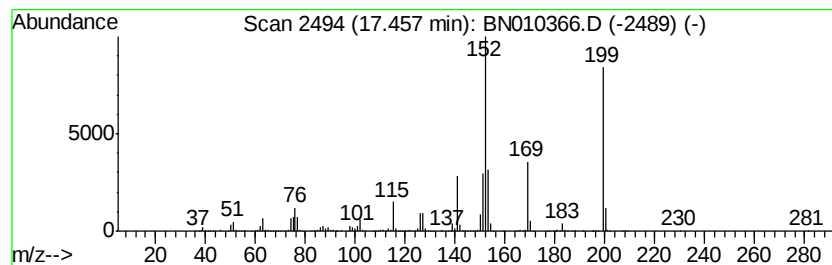
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 1,1'-Biphenyl, 4-nitro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.46	18.14 ng/ul	7679040	Phenanthrene-d10	16.98

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 4-nitro-		199	C12H9NO2	000092-93-3	98
2	1,1'-Biphenyl, 4-nitro-		199	C12H9NO2	000092-93-3	96
3	1,1'-Biphenyl, 3-nitro-		199	C12H9NO2	002113-58-8	76
4	1,1'-Biphenyl, 3-nitro-		199	C12H9NO2	002113-58-8	72
5	1,1'-Biphenyl, 3-nitro-		199	C12H9NO2	002113-58-8	68



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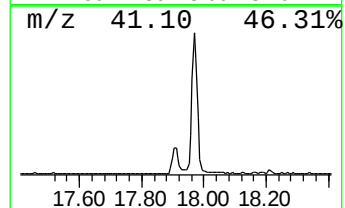
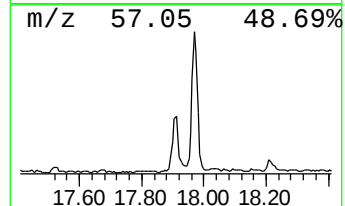
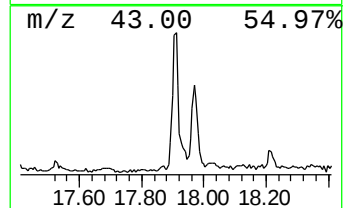
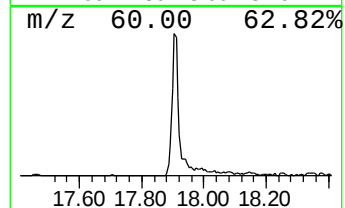
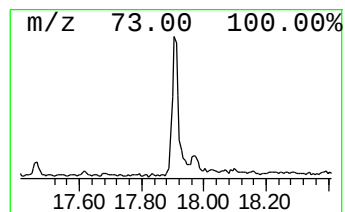
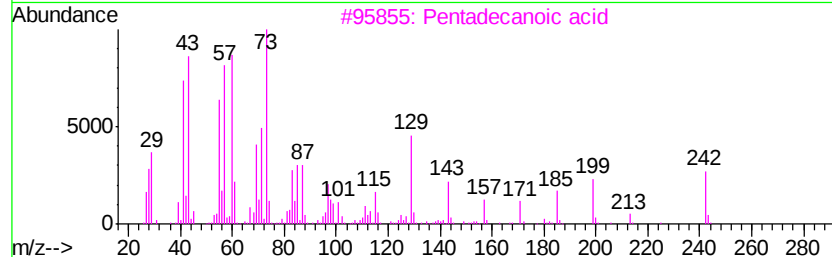
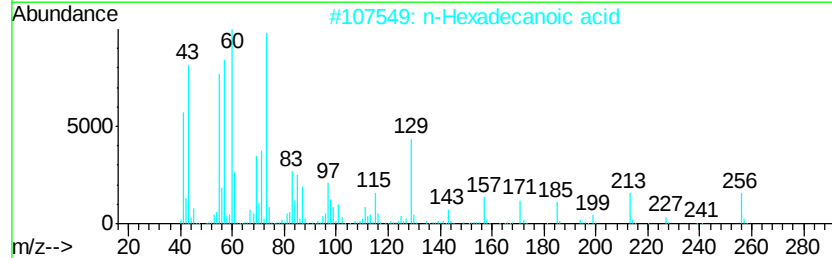
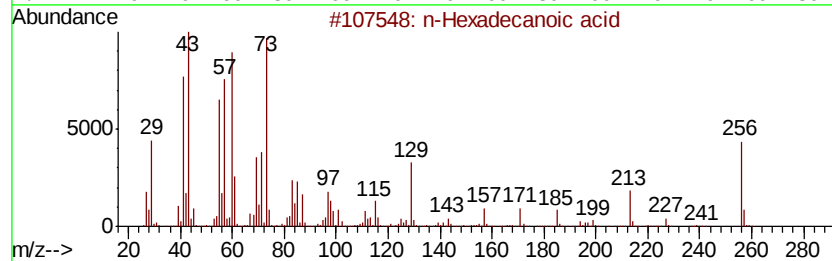
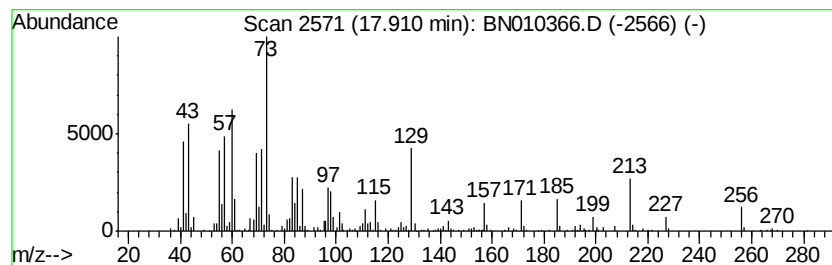
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 n-Hexadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.
17.91	2.11 ng/ul	894051	Phenanthrene-d10		16.98
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	68
4	n-Decanoic acid	172	C10H20O2	000334-48-5	64
5	Tridecanoic acid	214	C13H26O2	000638-53-9	62



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN040120\
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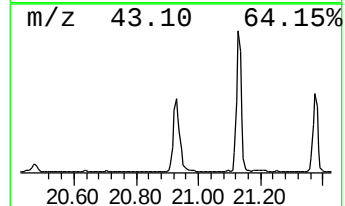
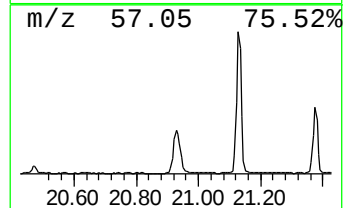
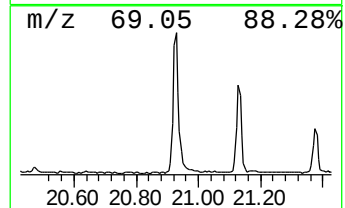
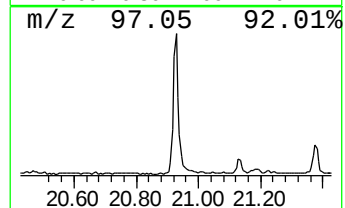
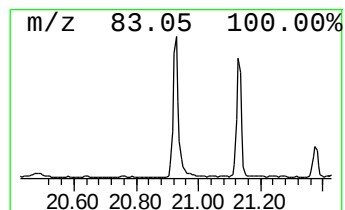
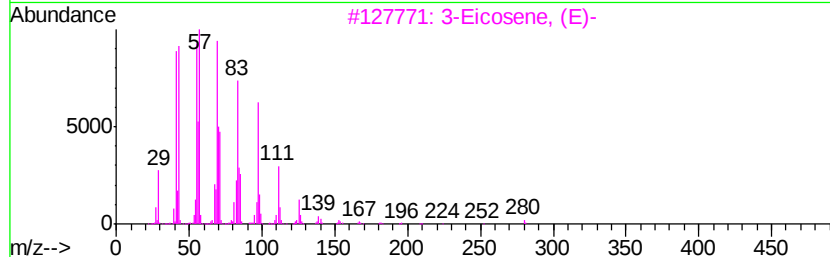
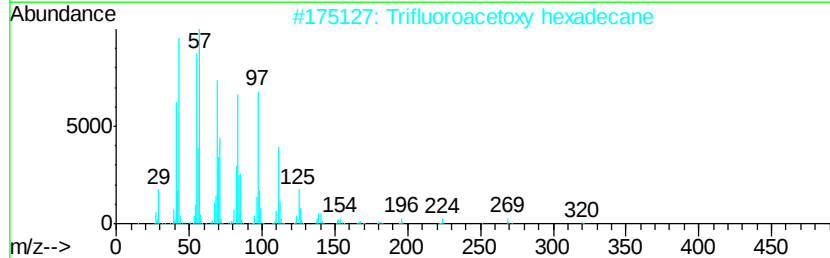
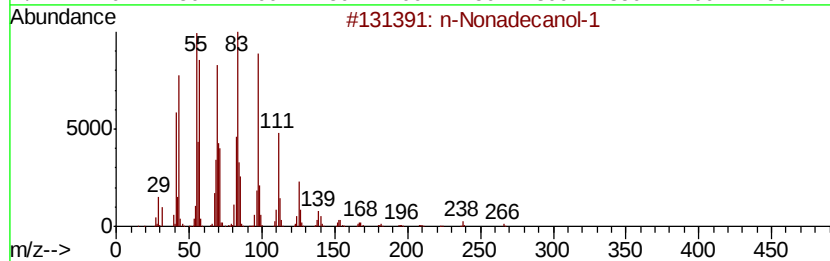
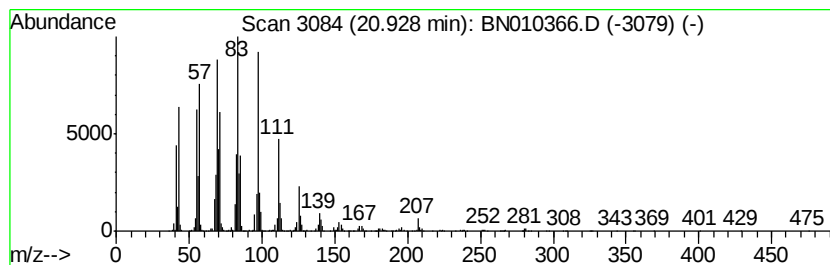
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 n-Nonadecanol-1 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.93	2.24 ng/ul	3590440	Chrysene-d12	21.19

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Nonadecanol-1	284	C19H40O	001454-84-8	91
2		Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	91
3		3-Eicosene, (E)-	280	C20H40	074685-33-9	91
4		n-Tetracosanol-1	354	C24H50O	000506-51-4	91
5		Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	91



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Data File : BN010366.D
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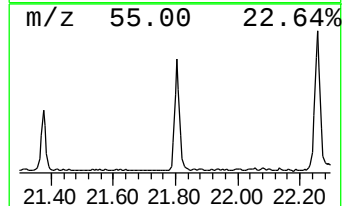
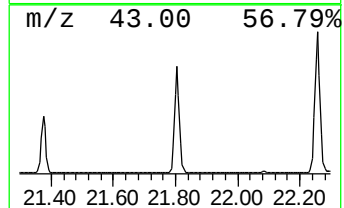
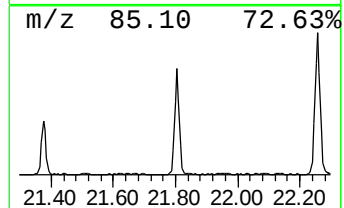
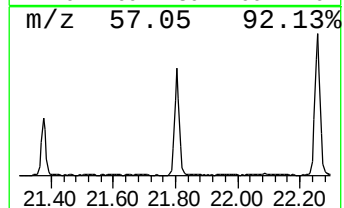
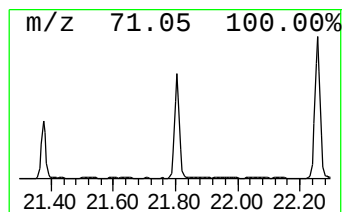
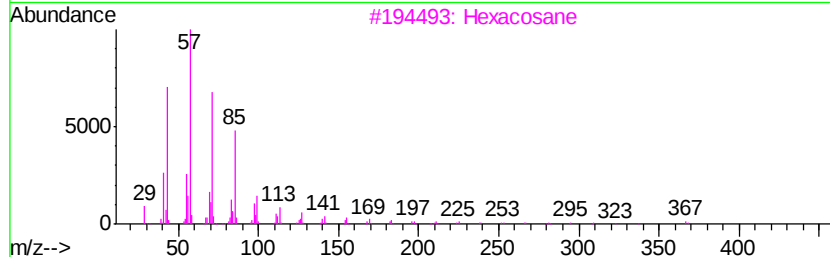
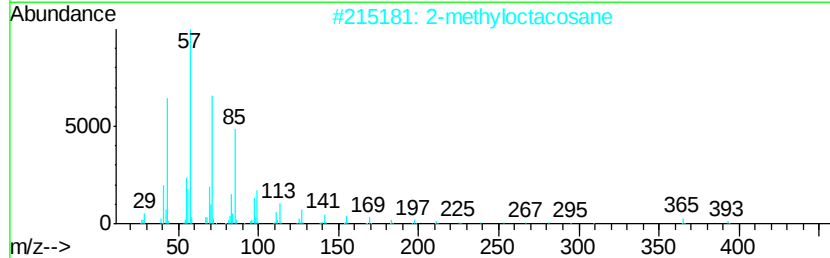
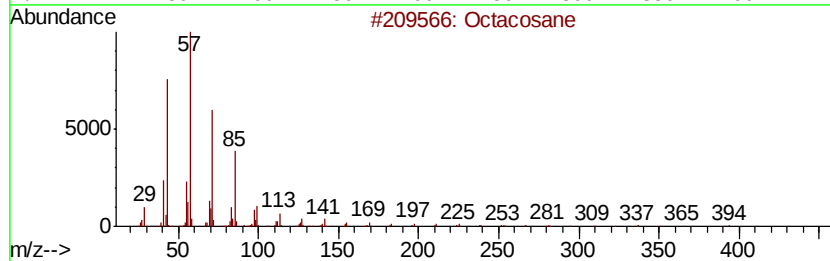
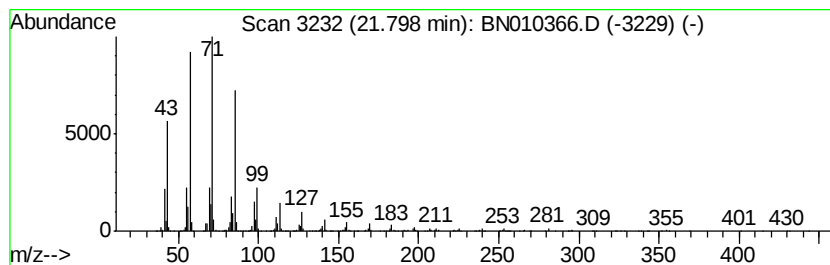
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.80	2.34 ng/ul	3743320	Chrysene-d12	21.19

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosane	394	C28H58	000630-02-4	90
2			2-methyloctacosane	408	C29H60	1000376-72-8	90
3			Hexacosane	366	C26H54	000630-01-3	90
4			Tetracosane, 11-decyl-	479	C34H70	055429-84-0	87
5			Heneicosane	296	C21H44	000629-94-7	87



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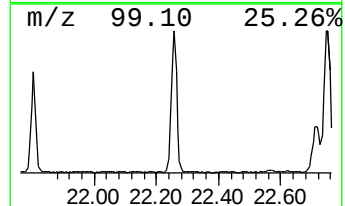
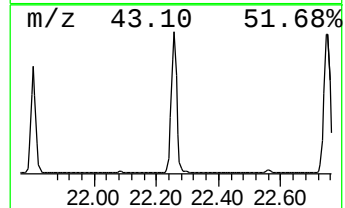
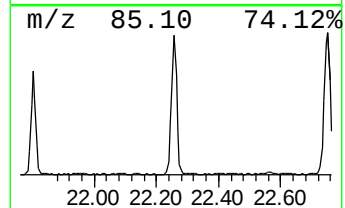
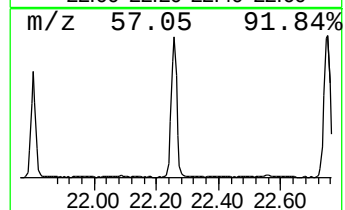
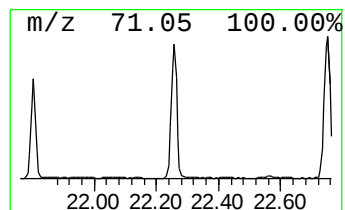
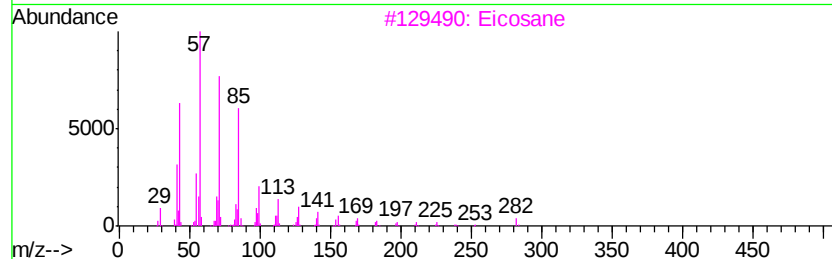
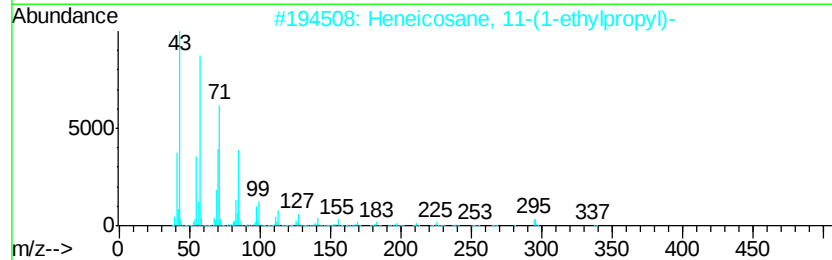
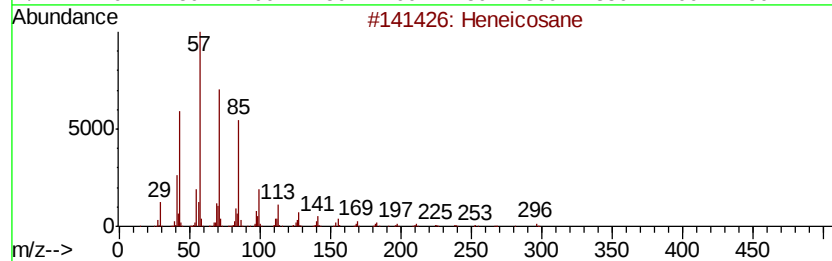
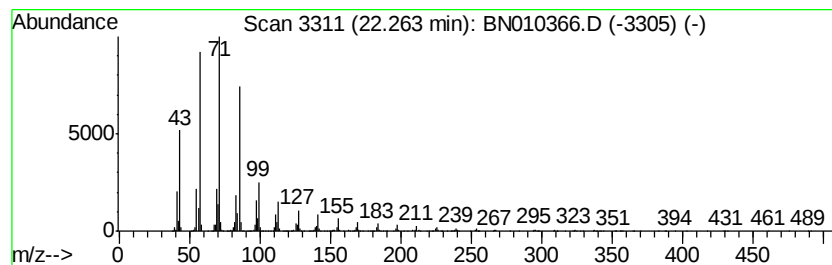
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.26	3.41 ng/ul	5465720	Chrysene-d12	21.19

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91
3		Eicosane	282	C20H42	000112-95-8	91
4		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	91
5		2-methyloctacosane	408	C29H60	1000376-72-8	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN040120\
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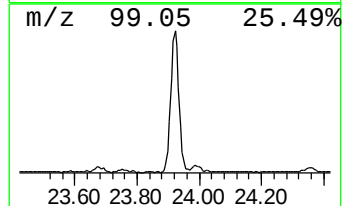
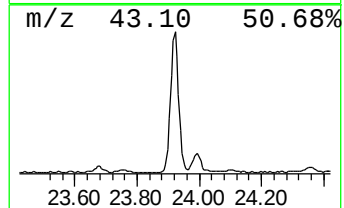
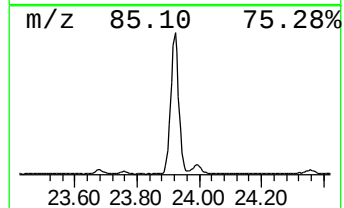
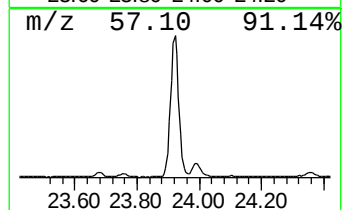
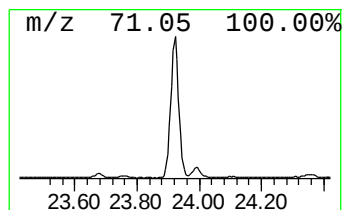
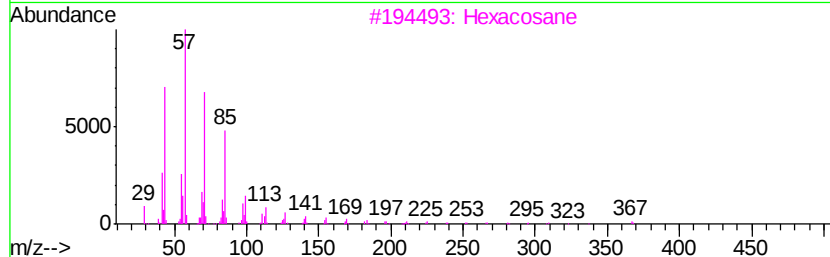
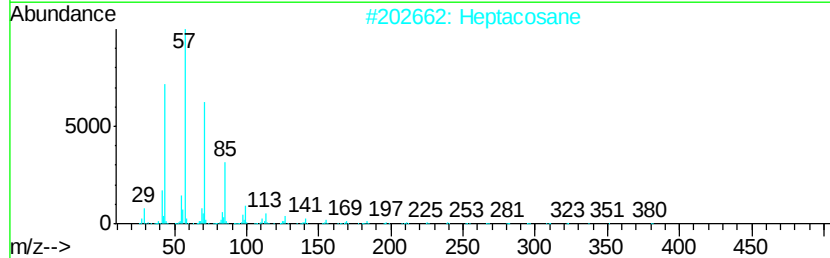
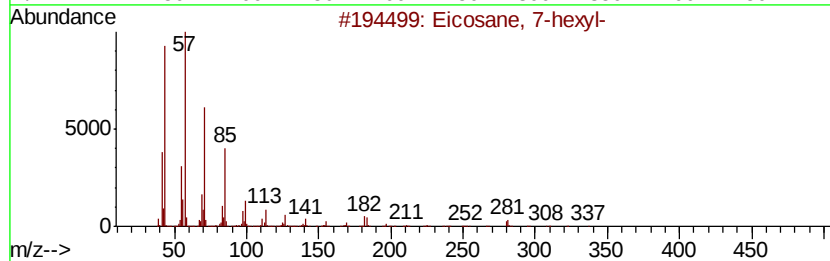
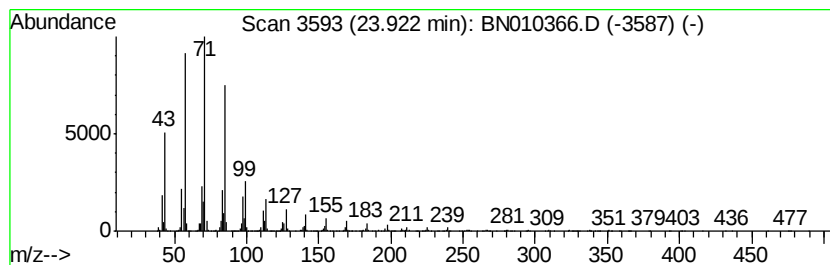
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.92	9.90 ng/ul	5499570	Perylene-d12	23.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane, 7-hexyl-	366	C26H54	055333-99-8	91
2		Heptacosane	380	C27H56	000593-49-7	91
3		Hexacosane	366	C26H54	000630-01-3	90
4		Hentriacontane	437	C31H64	000630-04-6	90
5		Octacosane	394	C28H58	000630-02-4	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN040120\
Data File : BN010366.D
Acq On : 01 Apr 2020 18:49
Operator : CG/JU
Sample : L2093-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBEW7

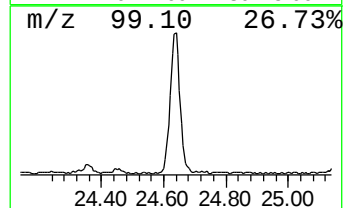
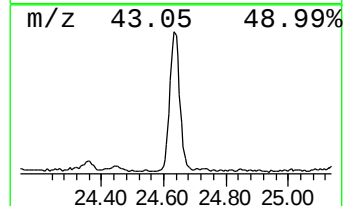
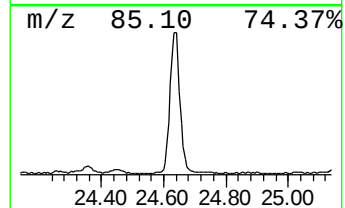
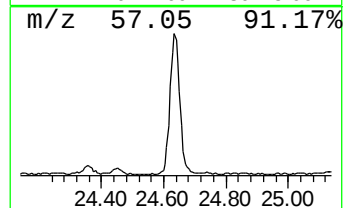
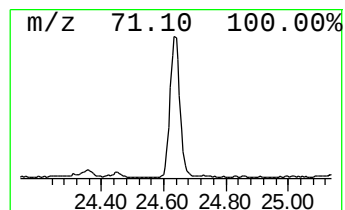
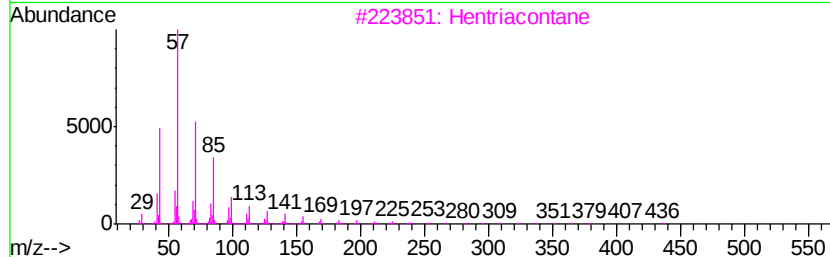
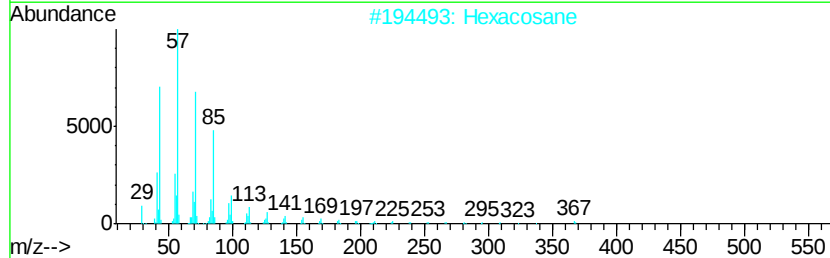
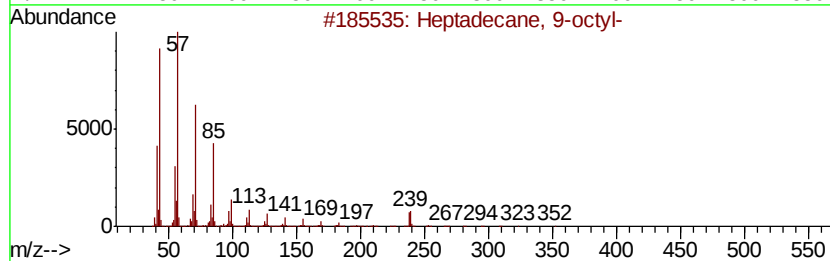
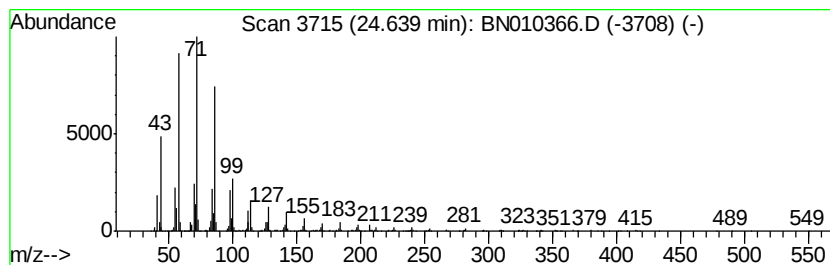
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.64	6.15 ng/ul	3415580	Perylene-d12	23.36

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	94
2			Hexacosane	366	C26H54	000630-01-3	91
3			Hentriacontane	437	C31H64	000630-04-6	91
4			Triacontane, 1-bromo-	500	C30H61Br	004209-22-7	91
5			Tetratetracontane	619	C44H90	007098-22-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN040120\
 Data File : BN010366.D
 Acq On : 01 Apr 2020 18:49
 Operator : CG/JU
 Sample : L2093-01
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBEW7

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.M
 Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	2.96	3.4	ng/ul	450355	1	7.61	2665950	20.0
2-Pentanone, 4-hy...	4.79	7.9	ng/ul	1053910	1	7.61	2665950	20.0
Naphthalene, 1-me...	12.26	33.4	ng/ul	7078730	2	10.37	4242860	20.0
2,4,7,9-Tetrameth...	13.16	9.0	ng/ul	2780680	3	14.24	6154250	20.0
1,1'-Biphenyl, 4-...	17.46	18.1	ng/ul	7679040	4	16.98	8464800	20.0
n-Hexadecanoic acid	17.91	2.1	ng/ul	894051	4	16.98	8464800	20.0
n-Nonadecanol-1	20.93	2.2	ng/ul	3590440	5	21.19	32038600	20.0
(DEL) Alkane: Str...	21.80	2.3	ng/ul	3743320	5	21.19	32038600	20.0
(DEL) Alkane: Str...	22.26	3.4	ng/ul	5465720	5	21.19	32038600	20.0
(DEL) Alkane: Str...	23.92	9.9	ng/ul	5499570	6	23.36	11108500	20.0
(DEL) Alkane: Str...	24.64	6.2	ng/ul	3415580	6	23.36	11108500	20.0