

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\
 Data File : BM021278.D
 Acq On : 30 Jun 2019 00:54
 Operator : HP/JU
 Sample : K3590-01
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C5BA0

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_M\METHODS\SOM-EPA-BM061419MA.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	3.016	3	5	11	rVB	2819106	3117238	4.25%	0.632%
2	4.404	236	241	250	rVB	376931	500264	0.68%	0.101%
3	6.275	554	559	571	rVB	233811	368440	0.50%	0.075%
4	6.434	579	586	596	rVB	1473860	2292210	3.13%	0.465%
5	6.816	641	651	654	rVV2	218233	564510	0.77%	0.114%
6	6.857	654	658	664	rVV	267576	386384	0.53%	0.078%
7	6.934	664	671	681	rVB	1057699	1823590	2.49%	0.370%
8	7.057	681	692	695	rBV2	274968	455339	0.62%	0.092%
9	7.098	695	699	708	rVV	801774	1211113	1.65%	0.246%
10	7.187	708	714	725	rVB	692374	1190897	1.63%	0.241%
11	7.292	725	732	743	rBV	1104126	1735539	2.37%	0.352%
12	7.757	804	811	818	rBV	1210101	1863819	2.54%	0.378%
13	8.463	925	931	947	rVV	1329415	2279702	3.11%	0.462%
14	8.922	1002	1009	1019	rVV	1074162	1868750	2.55%	0.379%
15	9.281	1064	1070	1074	rBV	202671	318881	0.44%	0.065%
16	9.633	1120	1130	1145	rBV	929135	1579099	2.16%	0.320%
17	9.857	1162	1168	1180	rVB4	307959	746266	1.02%	0.151%
18	10.169	1213	1221	1239	rVB	1572639	2738326	3.74%	0.555%
19	10.545	1278	1285	1291	rBV	1662849	2735614	3.73%	0.555%
20	12.739	1651	1658	1670	rVV4	137760	291808	0.40%	0.059%
21	13.263	1741	1747	1753	rBV2	242751	412280	0.56%	0.084%
22	13.710	1818	1823	1831	rVB	369175	582729	0.80%	0.118%
23	13.816	1831	1841	1845	rBV	2617935	3978153	5.43%	0.807%
24	13.863	1845	1849	1854	rVV	726846	1004709	1.37%	0.204%
25	14.092	1882	1888	1894	rVB	3107212	4568934	6.24%	0.926%
26	14.404	1935	1941	1946	rVV2	2499855	3842379	5.24%	0.779%
27	14.463	1946	1951	1963	rVB3	362062	810262	1.11%	0.164%
28	14.633	1970	1980	1990	rBV4	1149895	2831980	3.87%	0.574%
29	14.821	2007	2012	2016	rBV2	483024	645037	0.88%	0.131%
30	15.310	2086	2095	2099	rBV2	380389	699596	0.95%	0.142%
31	15.398	2103	2110	2116	rVV	4282678	6257162	8.54%	1.269%
32	15.527	2126	2132	2139	rBV	695283	948292	1.29%	0.192%
33	15.963	2199	2206	2213	rVB2	656503	1056272	1.44%	0.214%
34	16.110	2226	2231	2239	rVB6	124292	322319	0.44%	0.065%

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 Title : SVOA CALIBRATION

35	16.257	2251	2256	2261	rBV3	202608	358920	0.49%	0.073%
36	16.474	2289	2293	2301	rVB	204286	337835	0.46%	0.069%
37	16.557	2301	2307	2311	rBV	1013885	1397923	1.91%	0.283%
38	16.651	2319	2323	2337	rVB	460949	700014	0.96%	0.142%
39	16.904	2359	2366	2373	rBV8	245385	740025	1.01%	0.150%
40	16.974	2373	2378	2383	rVB2	312809	615160	0.84%	0.125%
41	17.051	2386	2391	2397	rBV	1283759	1658736	2.26%	0.336%
42	17.115	2397	2402	2404	rVV	578534	758142	1.03%	0.154%
43	17.151	2404	2408	2412	rVV2	2998126	4316668	5.89%	0.875%
44	17.192	2412	2415	2419	rVV	1244369	1787702	2.44%	0.362%
45	17.251	2419	2425	2431	rVV	3720326	6121183	8.35%	1.241%
46	17.315	2431	2436	2446	rVB2	2314375	3560605	4.86%	0.722%
47	17.533	2468	2473	2476	rBV	336085	510592	0.70%	0.104%
48	17.668	2492	2496	2502	rVB3	235878	367191	0.50%	0.074%
49	17.786	2511	2516	2518	rBV	281307	424827	0.58%	0.086%
50	17.939	2534	2542	2547	rBV	3798256	6363753	8.69%	1.290%
51	17.998	2547	2552	2558	rVV	4706610	7366068	10.05%	1.494%
52	18.080	2558	2566	2576	rVB	13492157	24686756	33.70%	5.006%
53	18.315	2601	2606	2610	rBV	1902666	2540470	3.47%	0.515%
54	18.351	2610	2612	2617	rVB3	293846	428024	0.58%	0.087%
55	18.480	2631	2634	2639	rVB2	393844	501103	0.68%	0.102%
56	18.533	2639	2643	2647	rBV3	282165	379845	0.52%	0.077%
57	18.662	2660	2665	2670	rVB	1103402	1584023	2.16%	0.321%
58	18.715	2670	2674	2679	rBV3	431223	677842	0.93%	0.137%
59	18.768	2680	2683	2687	rVB4	247331	356762	0.49%	0.072%
60	18.986	2718	2720	2727	rVB2	411192	785704	1.07%	0.159%
61	19.062	2727	2733	2739	rBV	3403394	4675433	6.38%	0.948%
62	19.233	2749	2762	2772	rBV4	11207511	35900541	49.00%	7.280%
63	19.339	2775	2780	2790	rVB	4134574	5742536	7.84%	1.164%
64	19.545	2810	2815	2818	rBV	4254271	6168524	8.42%	1.251%
65	19.574	2818	2820	2825	rVB	2204804	2539747	3.47%	0.515%
66	19.774	2851	2854	2859	rVB4	662987	841610	1.15%	0.171%
67	20.021	2892	2896	2900	rBV	3777034	4961990	6.77%	1.006%
68	20.227	2926	2931	2941	rVB5	1871916	2928321	4.00%	0.594%
69	20.427	2961	2965	2968	rBV3	889494	1065196	1.45%	0.216%
70	20.521	2977	2981	2982	rBV	772767	973938	1.33%	0.197%
71	20.592	2982	2993	3008	rVV	29239089	73265223	100.00%	14.856%

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72	20.921	3045	3049	3055	rBV	2390846	3015641	4.12%	0.611%
73	21.039	3064	3069	3079	rVB3	5168650	6907819	9.43%	1.401%
74	21.203	3095	3097	3102	rBV2	405681	493119	0.67%	0.100%
75	21.339	3114	3120	3124	rBV2	3635517	6604780	9.01%	1.339%
76	21.374	3124	3126	3132	rVB	1641114	1832903	2.50%	0.372%
77	21.656	3172	3174	3177	rVB	528443	530208	0.72%	0.108%
78	21.692	3178	3180	3185	rVB3	382369	446166	0.61%	0.090%
79	21.915	3215	3218	3219	rBV	916232	1002205	1.37%	0.203%
80	21.939	3219	3222	3232	rVB3	2524014	3640574	4.97%	0.738%
81	22.380	3294	3297	3301	rBV2	523326	786627	1.07%	0.160%
82	22.427	3303	3305	3311	rVB	852293	1129405	1.54%	0.229%
83	22.621	3332	3338	3344	rBV	10388702	14570329	19.89%	2.954%
84	22.897	3380	3385	3388	rBV	2823455	4590723	6.27%	0.931%
85	22.956	3388	3395	3404	rVB2	16075899	29770723	40.63%	6.037%
86	23.415	3469	3473	3476	rBV	731909	1162996	1.59%	0.236%
87	23.468	3477	3482	3487	rVV	3110680	5544220	7.57%	1.124%
88	23.515	3487	3490	3499	rVB	1068509	1980228	2.70%	0.402%
89	23.609	3501	3506	3512	rBV	1948472	3557802	4.86%	0.721%
90	23.791	3532	3537	3547	rVB	3585839	6259002	8.54%	1.269%
91	24.127	3588	3594	3601	rVB	4117799	8383001	11.44%	1.700%
92	24.227	3602	3611	3624	rVB	16346733	35846400	48.93%	7.269%
93	24.885	3717	3723	3737	rVB5	2016393	4833398	6.60%	0.980%
94	25.344	3794	3801	3814	rVB	3280471	7991041	10.91%	1.620%
95	25.768	3866	3873	3878	rBV	1857997	4884744	6.67%	0.990%
96	26.062	3915	3923	3942	rVB5	2424477	9323081	12.73%	1.890%
97	26.703	4022	4032	4051	rVB2	8228942	26135054	35.67%	5.299%
98	27.474	4155	4163	4174	rVB3	2381385	7526530	10.27%	1.526%
99	27.603	4177	4185	4209	rVB	2637313	9104027	12.43%	1.846%
100	28.297	4291	4303	4320	rVB2	5546701	20898234	28.52%	4.238%

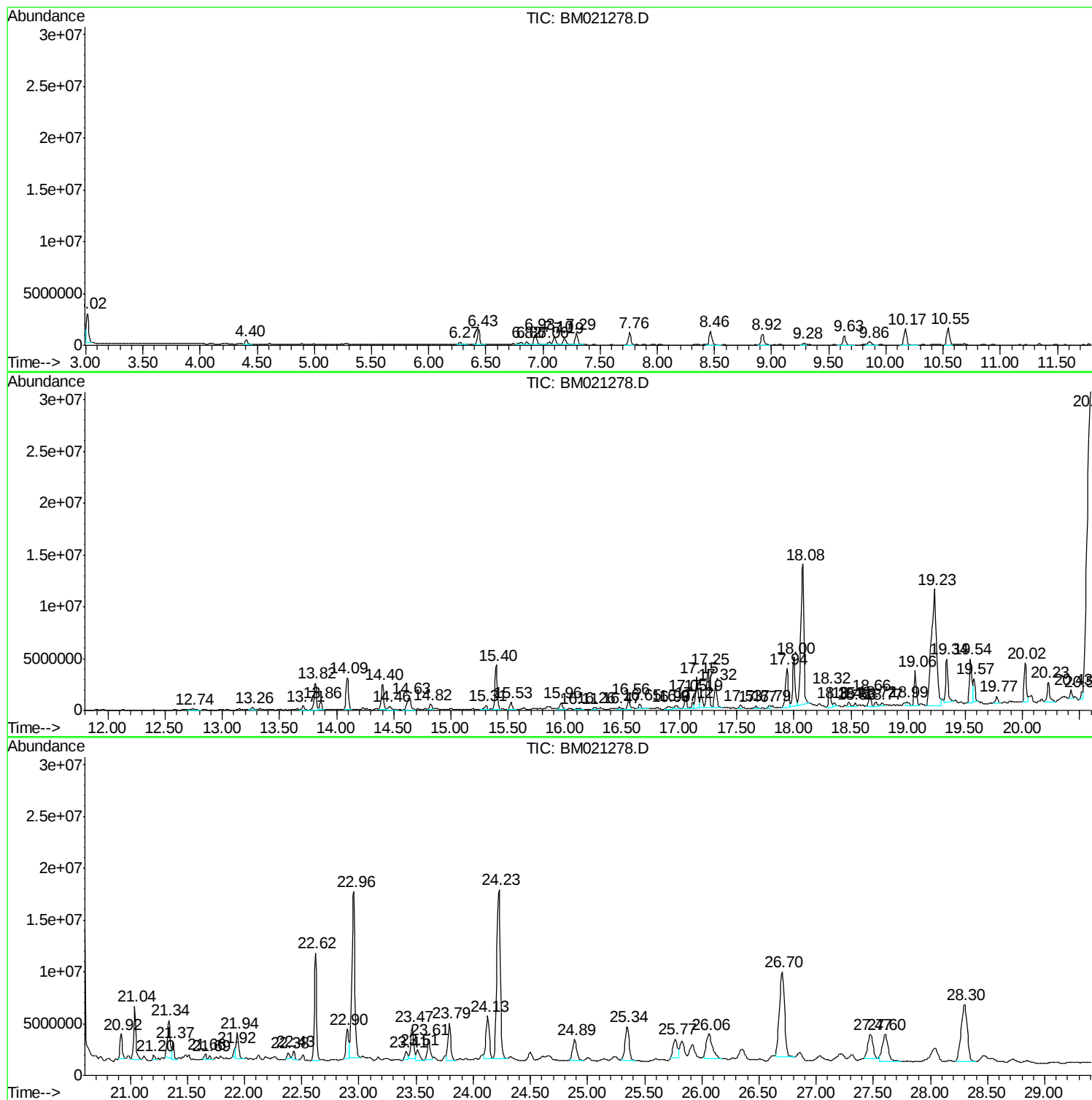
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Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M
Quant Title : SVOA CALIBRATION

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



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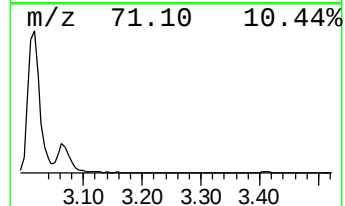
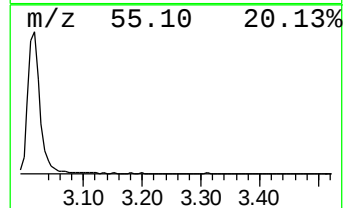
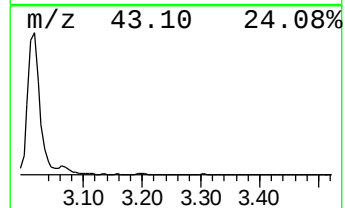
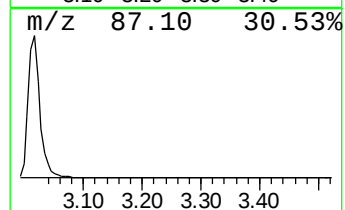
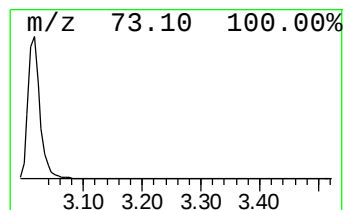
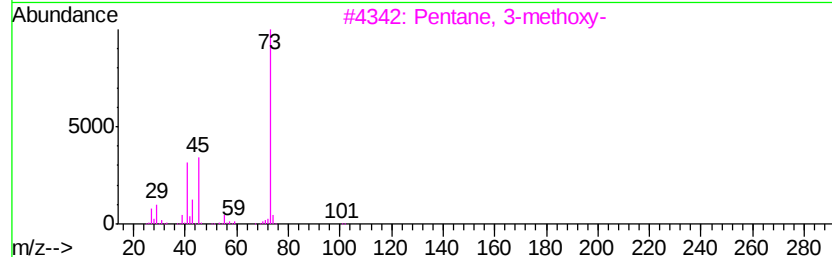
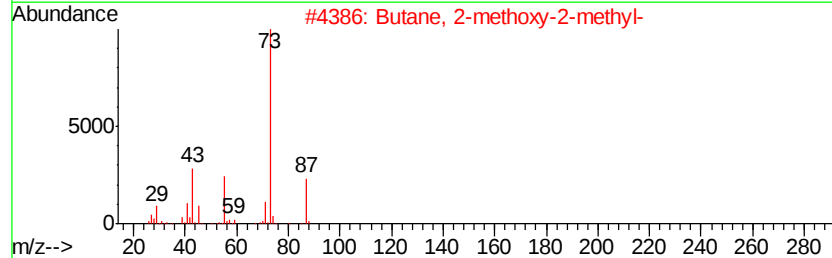
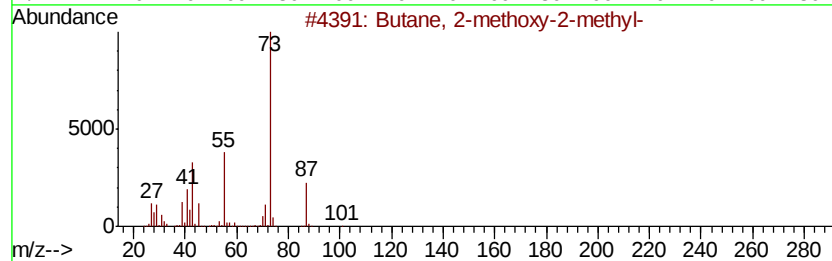
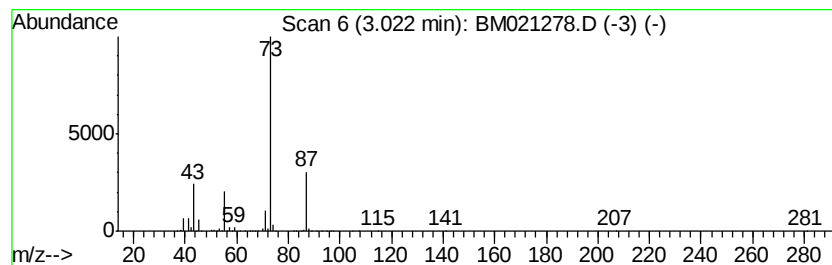
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.02	33.45 ng/ul	3117240	1,4-Dichlorobenzene-d4	7.76

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
2		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
3		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
4		N-Ethylformamide	73	C3H7NO	000627-45-2	9
5		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9



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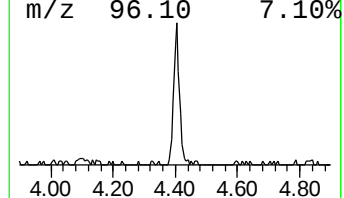
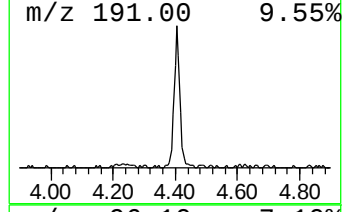
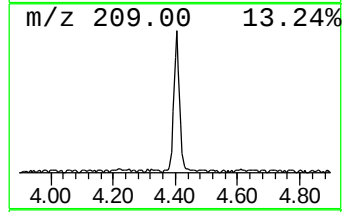
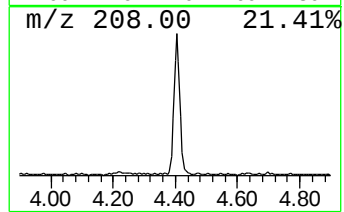
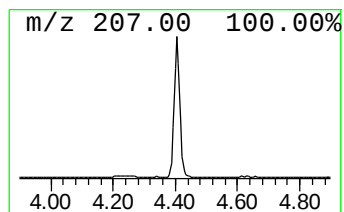
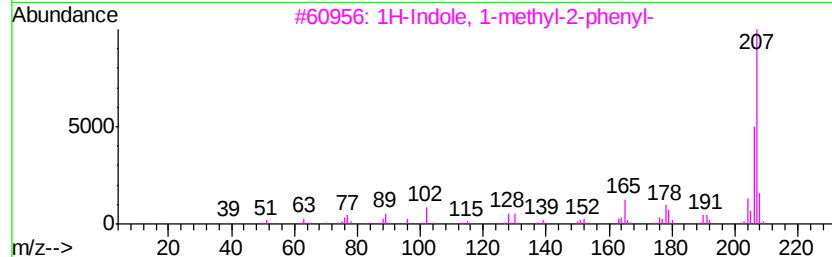
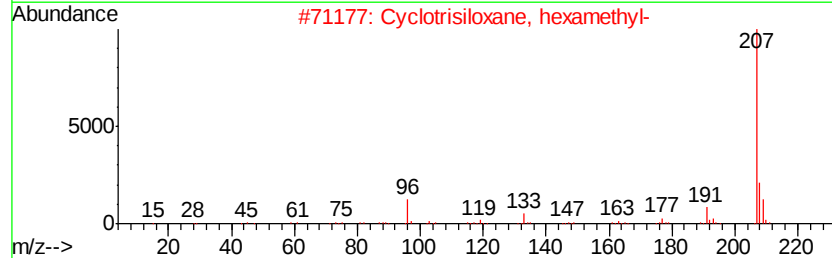
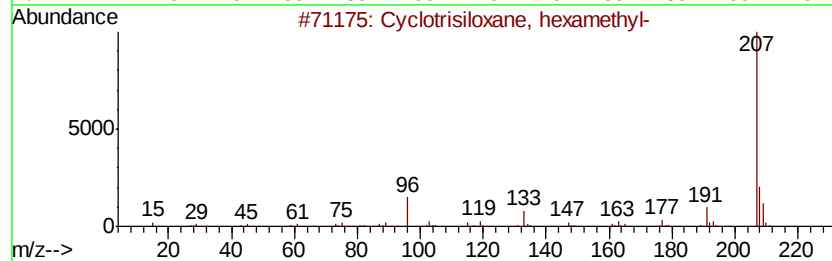
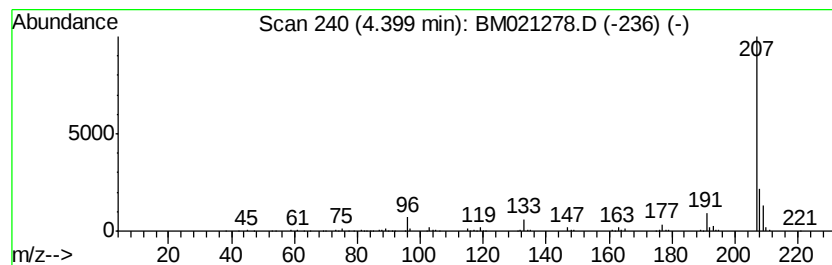
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 Cyclotrisiloxane, hexamethyl- Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.40	5.37 ng/ul	500264	1,4-Dichlorobenzene-d4	7.76

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclotrisiloxane, hexamethyl-	222	C6H18O3Si3	000541-05-9	90
2			Cyclotrisiloxane, hexamethyl-	222	C6H18O3Si3	000541-05-9	87
3			1H-Indole, 1-methyl-2-phenyl-	207	C15H13N	003558-24-5	39
4			2-Ethylacridine	207	C15H13N	055751-83-2	38
5			Cyclotrisiloxane, hexamethyl-	222	C6H18O3Si3	000541-05-9	35



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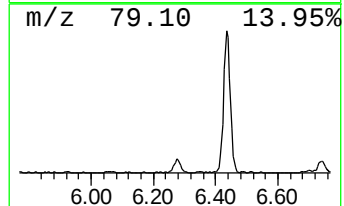
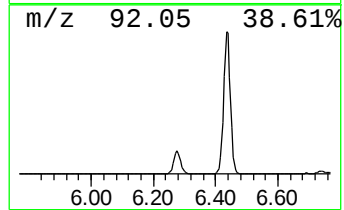
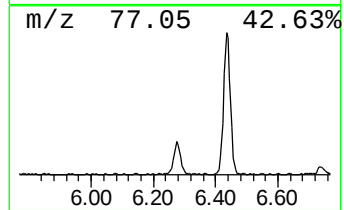
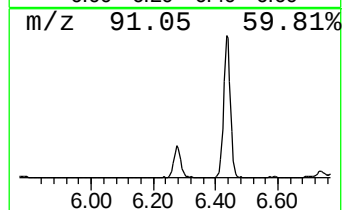
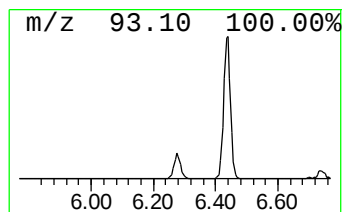
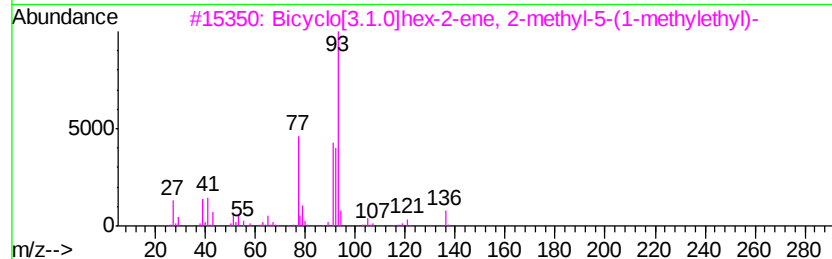
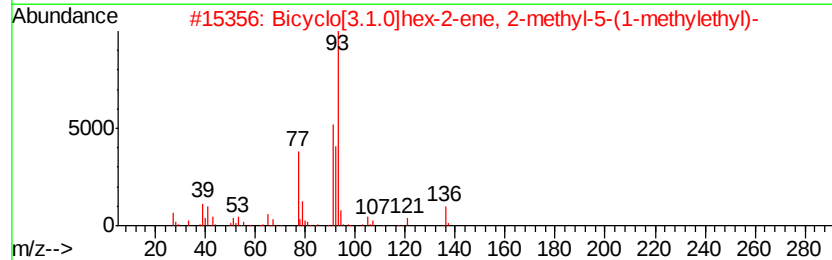
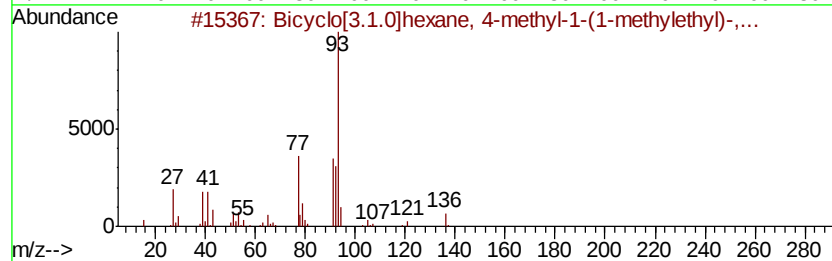
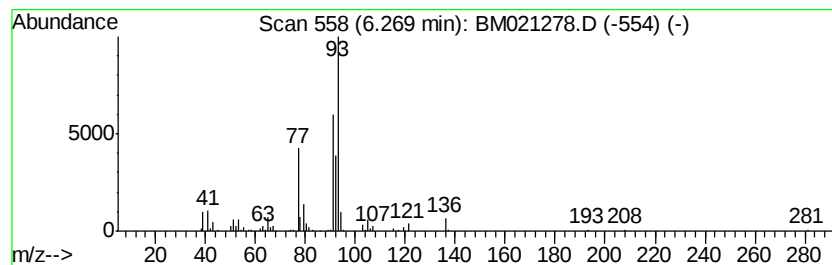
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Bicyclo[3.1.0]hexane, 4-met... Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.27	3.95 ng/ul	368440	1,4-Dichlorobenzene-d4	7.76

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Bicyclo[3.1.0]hexane, 4-methyl-1...	136	C10H16	058037-87-9	94
2		Bicyclo[3.1.0]hex-2-ene, 2-methy...	136	C10H16	002867-05-2	94
3		Bicyclo[3.1.0]hex-2-ene, 2-methv...	136	C10H16	002867-05-2	91
4		Bicyclo[3.1.0]hex-2-ene, 2-methv...	136	C10H16	002867-05-2	90
5		.alpha.-Phellandrene	136	C10H16	000099-83-2	78



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\
Data File : BM021278.D
Acq On : 30 Jun 2019 00:54
Operator : HP/JU
Sample : K3590-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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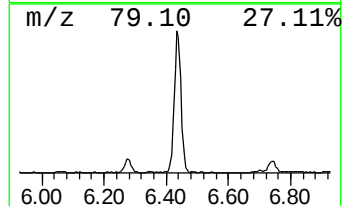
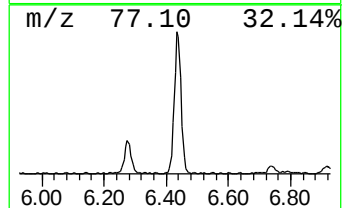
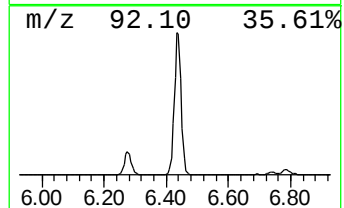
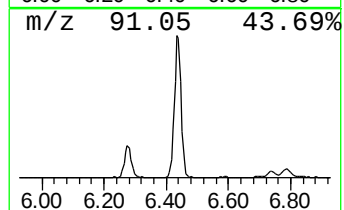
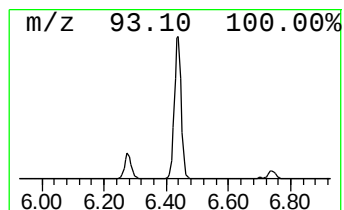
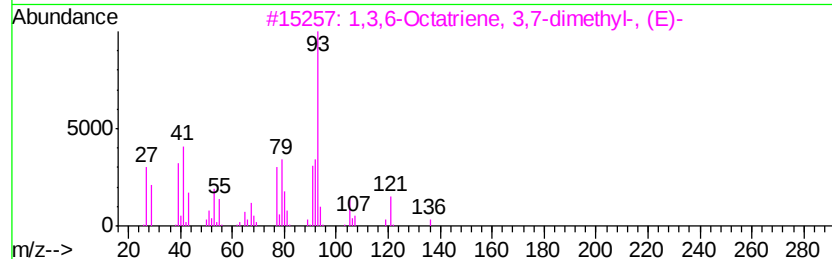
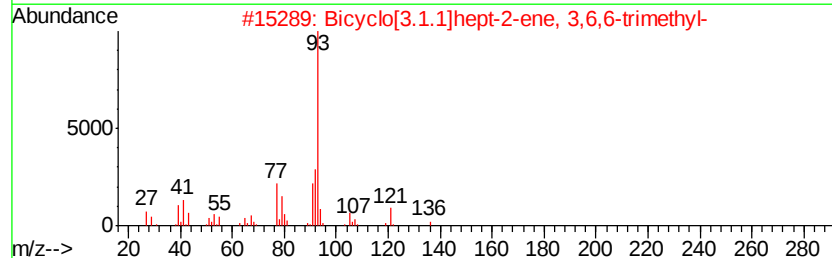
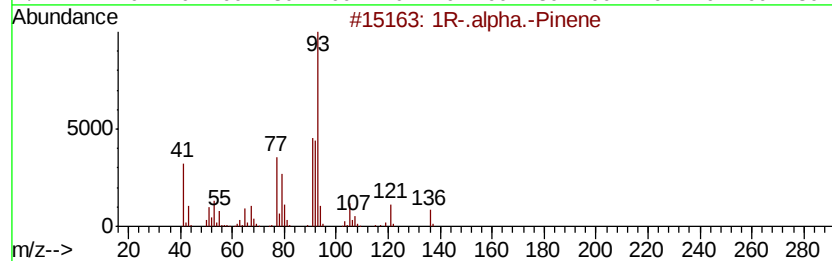
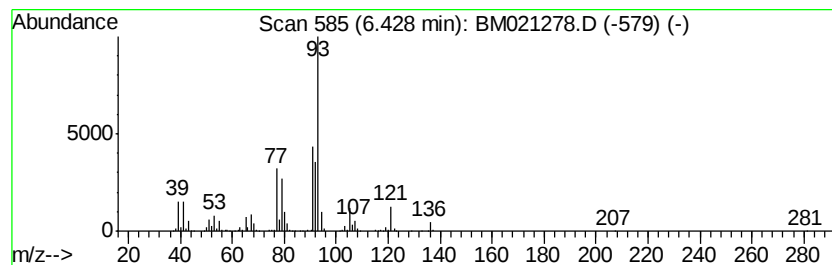
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 1R-.alpha.-Pinene Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.43	24.60 ng/ul	2292210	1,4-Dichlorobenzene-d4	7.76

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1R-.alpha.-Pinene	136	C10H16	007785-70-8	95
2		Bicyclo[3.1.1]hept-2-ene, 3,6,6-...	136	C10H16	004889-83-2	94
3		1,3,6-Octatriene, 3,7-dimethyl-,...	136	C10H16	003779-61-1	94
4		.alpha.-Pinene	136	C10H16	000080-56-8	94
5		1R-.alpha.-Pinene	136	C10H16	007785-70-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\
Data File : BM021278.D
Acq On : 30 Jun 2019 00:54
Operator : HP/JU
Sample : K3590-01
Misc :
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Instrument :
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ClientSampleId :
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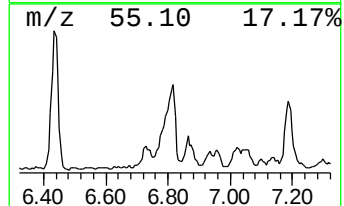
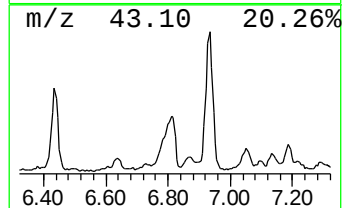
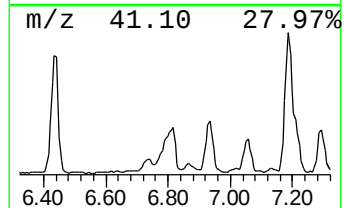
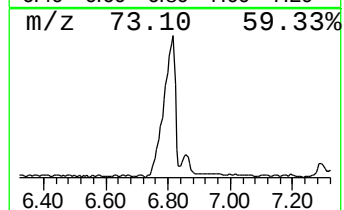
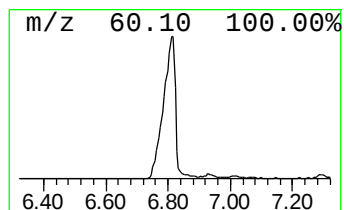
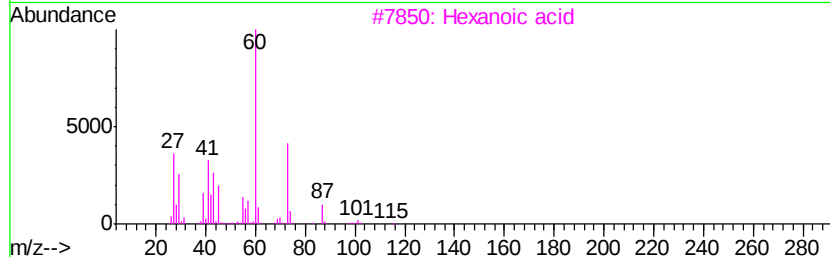
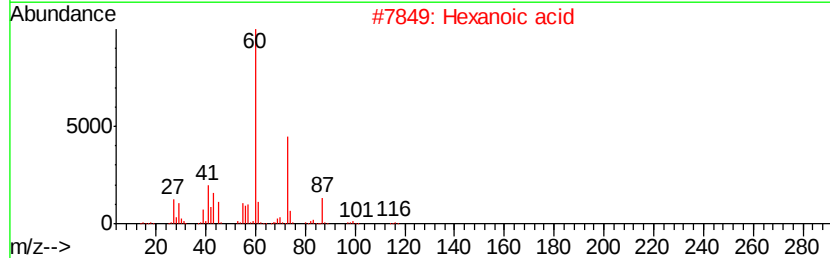
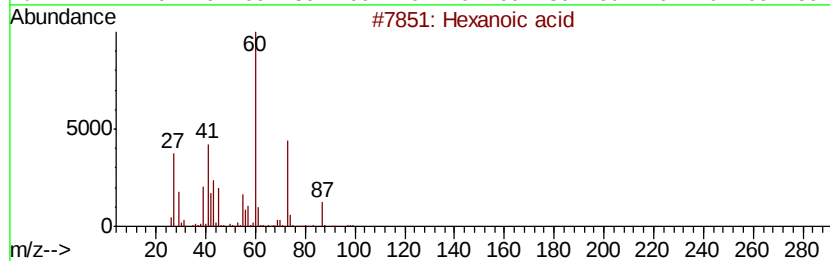
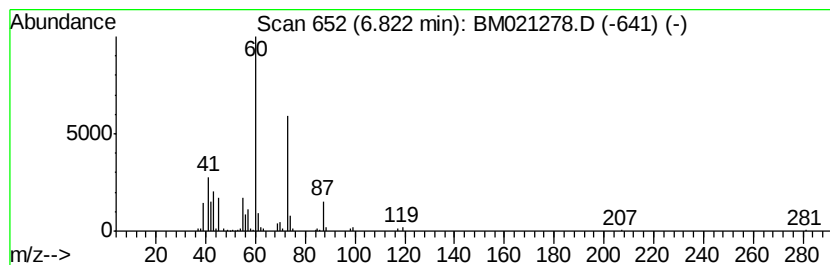
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Hexanoic acid Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.82	6.06 ng/ul	564510	1,4-Dichlorobenzene-d4	7.76

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexanoic acid	116	C6H12O2	000142-62-1	90
2		Hexanoic acid	116	C6H12O2	000142-62-1	83
3		Hexanoic acid	116	C6H12O2	000142-62-1	78
4		Propanedioic acid, propyl-	146	C6H10O4	000616-62-6	78
5		Octanoic Acid	144	C8H16O2	000124-07-2	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\
Data File : BM021278.D
Acq On : 30 Jun 2019 00:54
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Sample : K3590-01
Misc :
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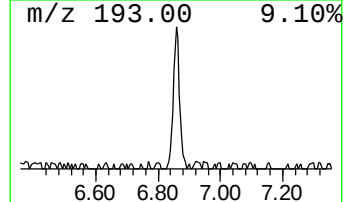
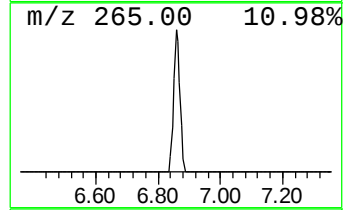
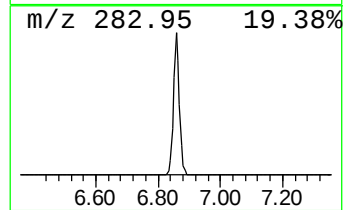
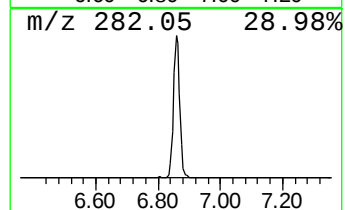
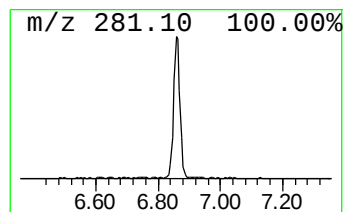
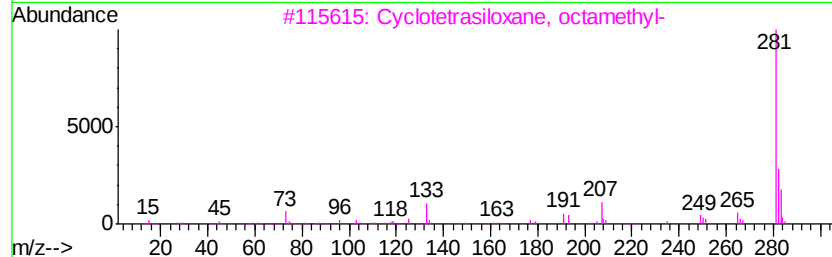
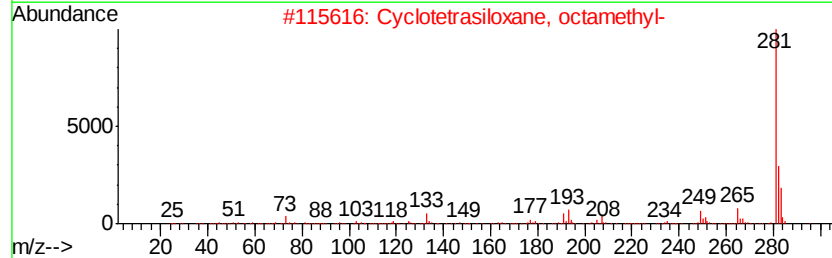
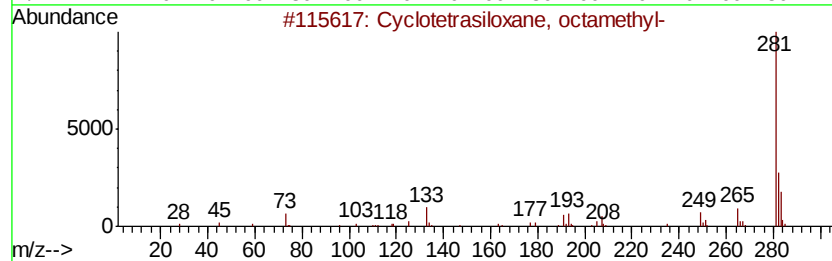
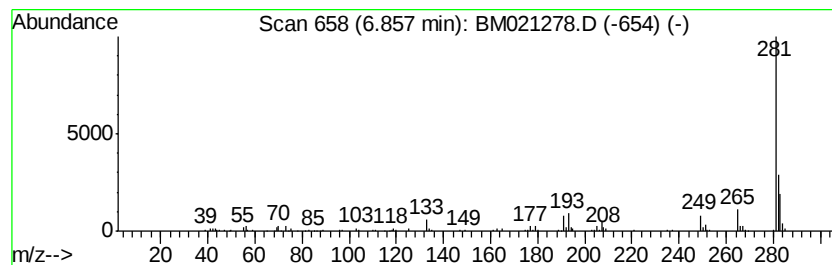
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Cyclotetrasiloxane, octamet... Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.86	4.15 ng/ul	386384	1,4-Dichlorobenzene-d4	7.76

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclotetrasiloxane, octamethyl-	296	C8H24O4Si4	000556-67-2	91
2		Cyclotetrasiloxane, octamethyl-	296	C8H24O4Si4	000556-67-2	91
3		Cyclotetrasiloxane, octamethyl-	296	C8H24O4Si4	000556-67-2	90
4		4-Nitro-4'-chlorodiphenylsulfoxide	281	C12H8ClNO3S	024535-53-3	53
5		Benzene, 1-phenyl-4-(2-cyano-2-p...	281	C21H15N	027869-56-3	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\
Data File : BM021278.D
Acq On : 30 Jun 2019 00:54
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Sample : K3590-01
Misc :
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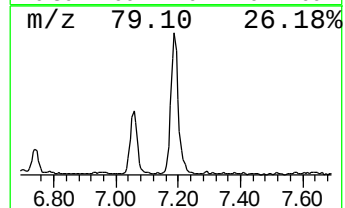
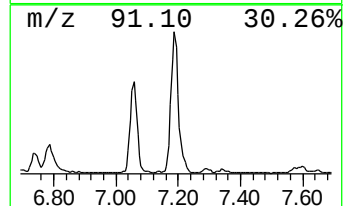
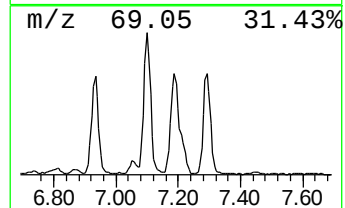
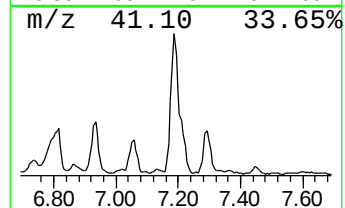
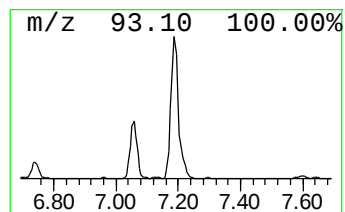
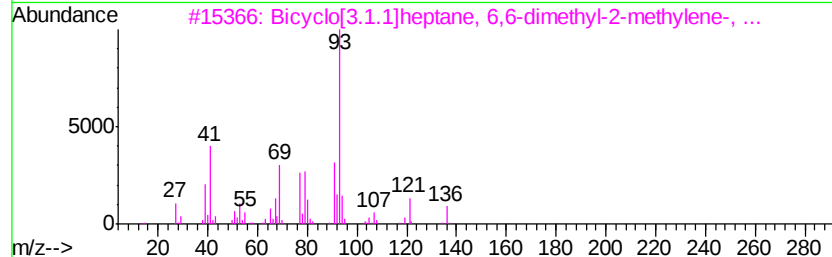
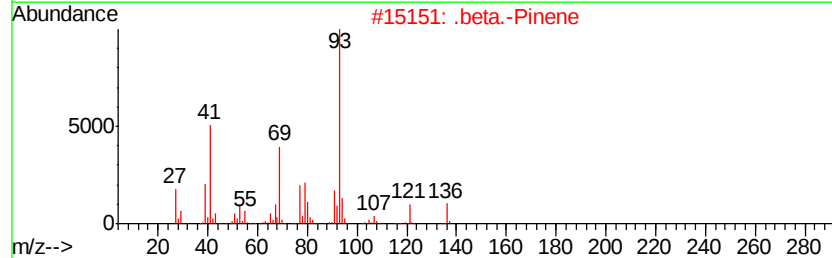
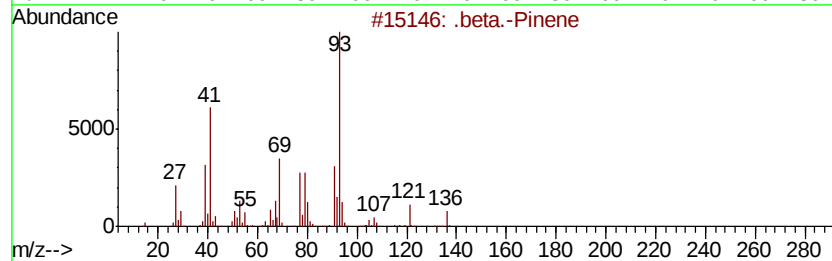
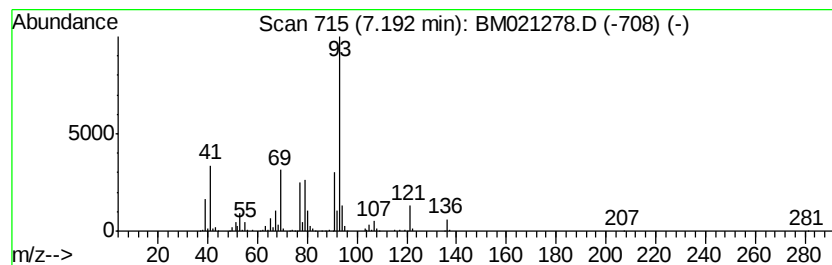
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 .beta.-Pinene Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.19	12.78 ng/ul	1190900	1,4-Dichlorobenzene-d4	7.76

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			.beta.-Pinene	136	C10H16	000127-91-3	96
2			.beta.-Pinene	136	C10H16	000127-91-3	91
3			Bicyclo[3.1.1]heptane, 6,6-dimet...	136	C10H16	018172-67-3	91
4			.beta.-Pinene	136	C10H16	000127-91-3	91
5			Cyclohexene, 4-methylene-1-(1-me...	136	C10H16	000099-84-3	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\
Data File : BM021278.D
Acq On : 30 Jun 2019 00:54
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Sample : K3590-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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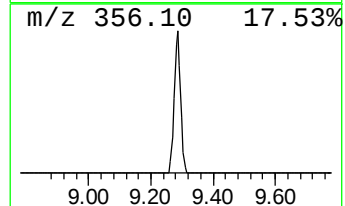
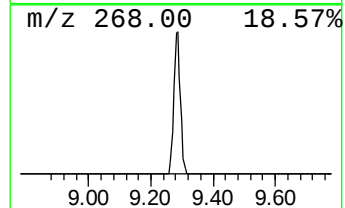
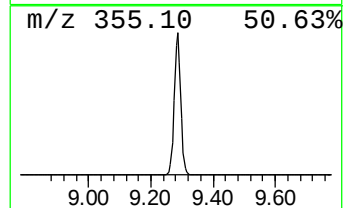
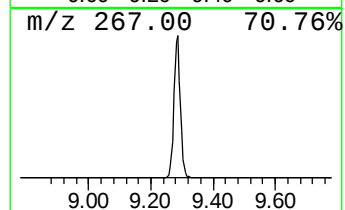
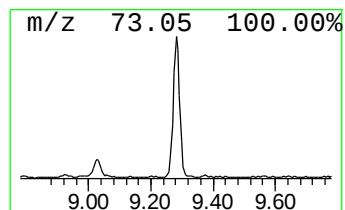
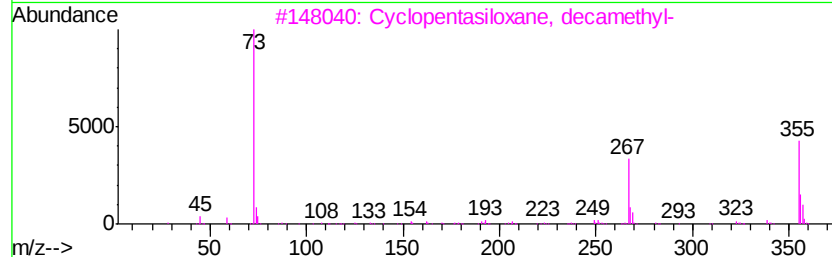
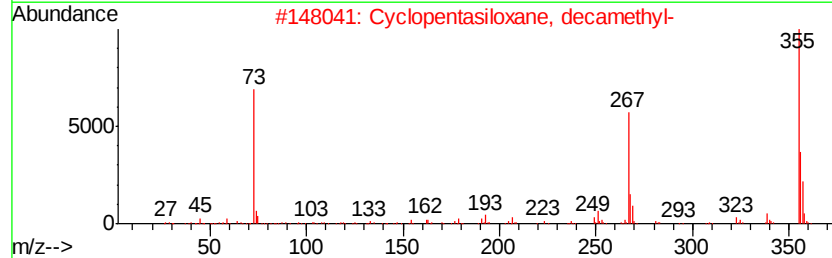
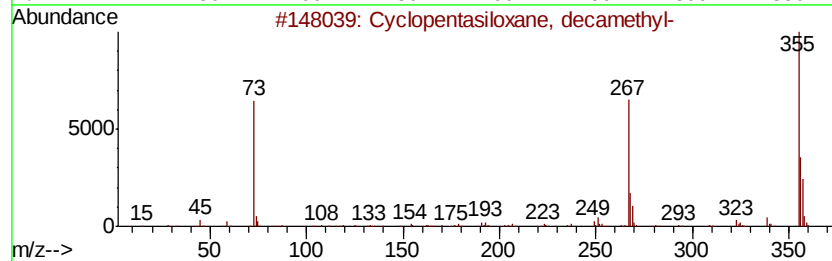
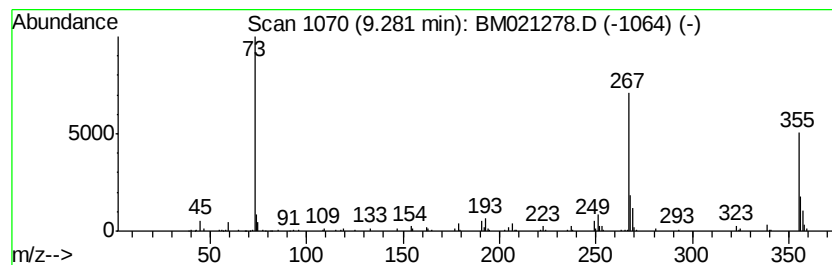
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Cyclopentasiloxane, decamet... Concentration Rank 54

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.28	2.33 ng/ul	318881	Naphthalene-d8	10.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	91
2		Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	91
3		Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	74
4		Silanamine, 1,1,1-trimethyl-N-[2...	369	C17H35N02Si3	068595-60-8	47
5		Benzoic acid, 2,5-bis(trimethyls...	370	C16H30O4Si3	003618-20-0	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\
Data File : BM021278.D
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ClientSampleId :
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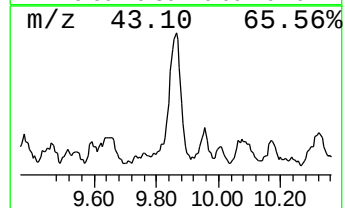
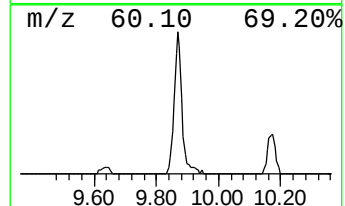
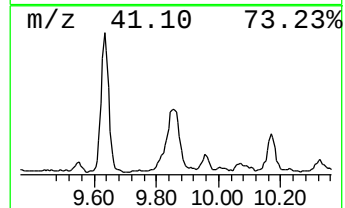
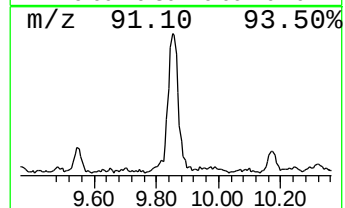
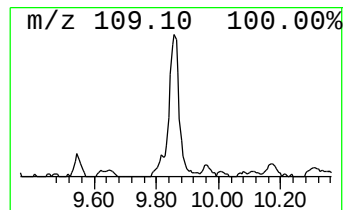
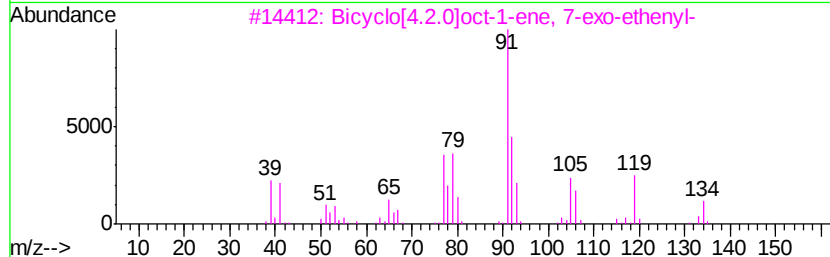
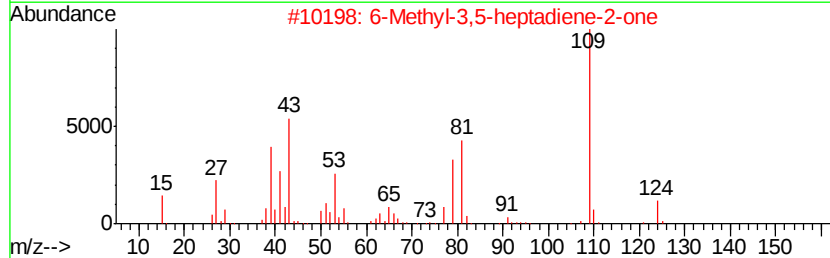
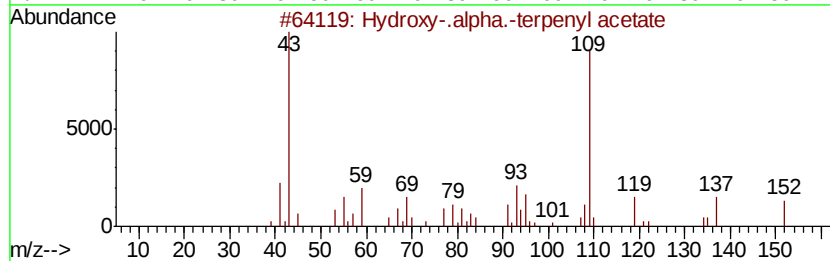
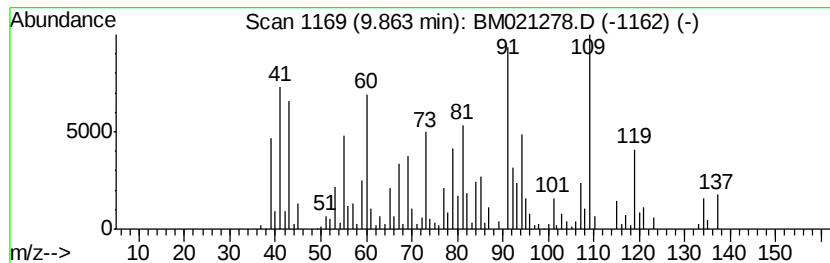
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 unknown-01 Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.86	5.46 ng/ul	746266	Naphthalene-d8	10.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hydroxy-.alpha.-terpenyl acetate	212	C12H20O3	1000293-00-9	27
2		6-Methyl-3,5-heptadiene-2-one	124	C8H12O	001604-28-0	22
3		Bicyclo[4.2.0]oct-1-ene, 7-exo-e...	134	C10H14	1000142-18-2	20
4		cis-Pinen-3-ol	152	C10H16O	1000292-85-2	15
5		Naphthalene, 1,2,3,5,8,8a-hexahy...	134	C10H14	062690-65-7	14



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\
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ALS Vial : 24 Sample Multiplier: 1

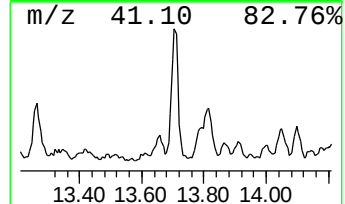
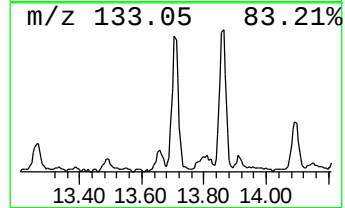
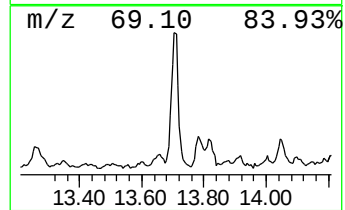
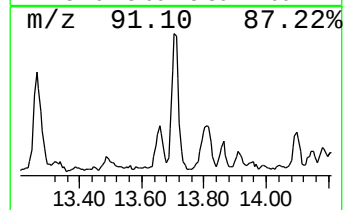
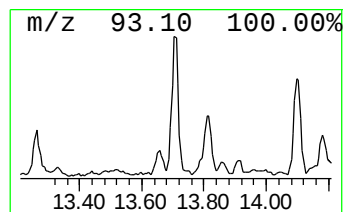
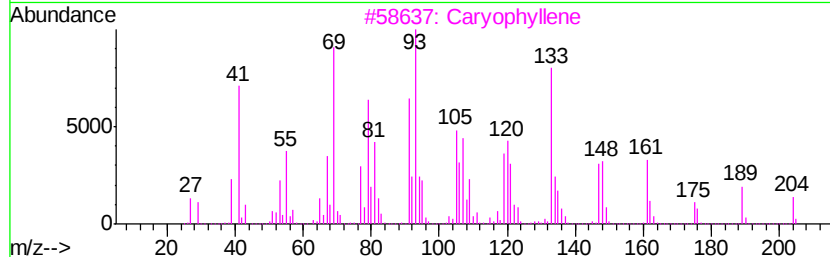
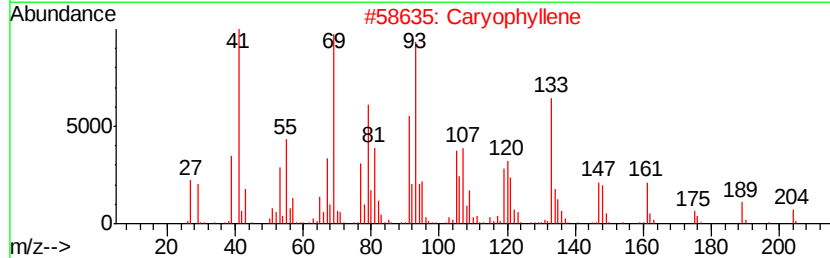
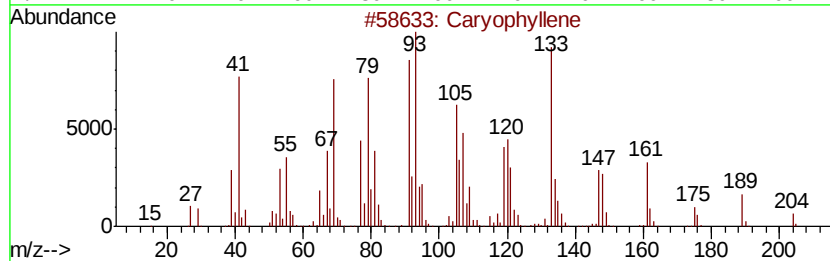
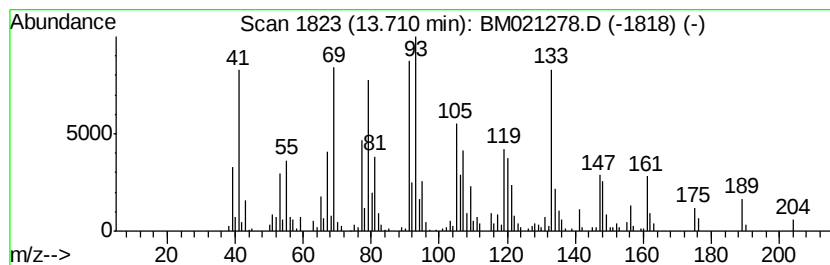
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ClientSampleId :
C5BA0

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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Caryophyllene Concentration Rank 48

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.71	3.03 ng/ul	582729	Acenaphthene-d10	14.40
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Caryophyllene	204 C15H24	000087-44-5 99
2		Caryophyllene	204 C15H24	000087-44-5 96
3		Caryophyllene	204 C15H24	000087-44-5 95
4		Bicyclo[7.2.0]undec-4-ene, 4,11,...	204 C15H24	013877-93-5 92
5		Bicyclo[7.2.0]undec-4-ene, 4,11,...	204 C15H24	000118-65-0 87



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\
Data File : BM021278.D
Acq On : 30 Jun 2019 00:54
Operator : HP/JU
Sample : K3590-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

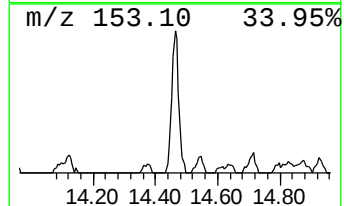
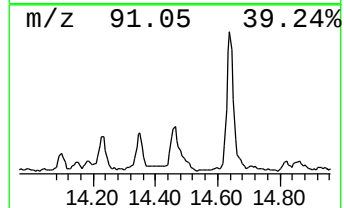
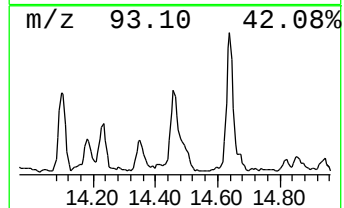
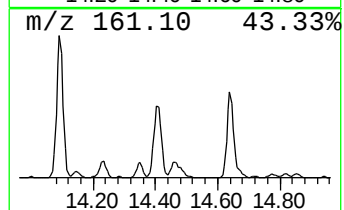
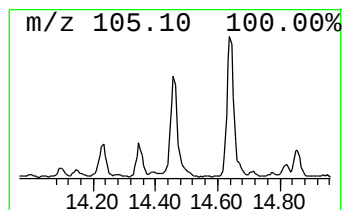
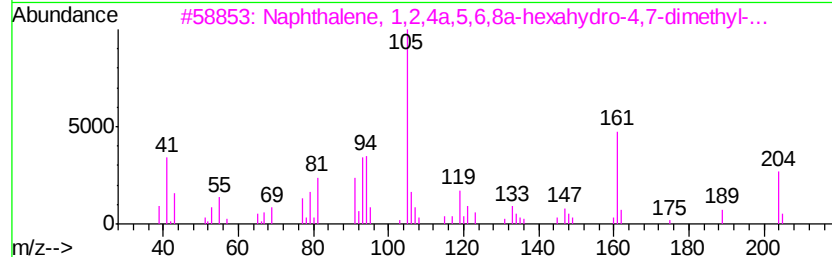
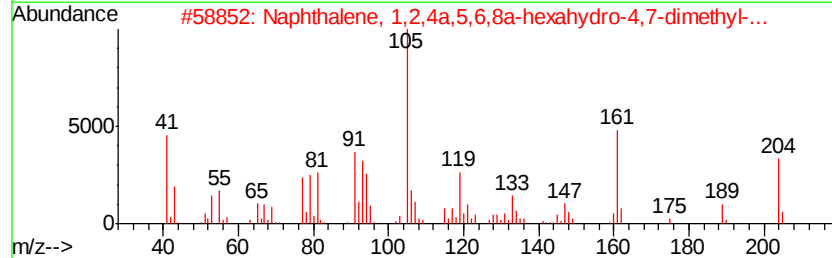
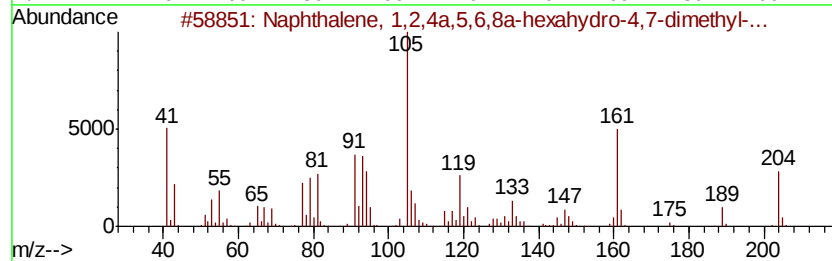
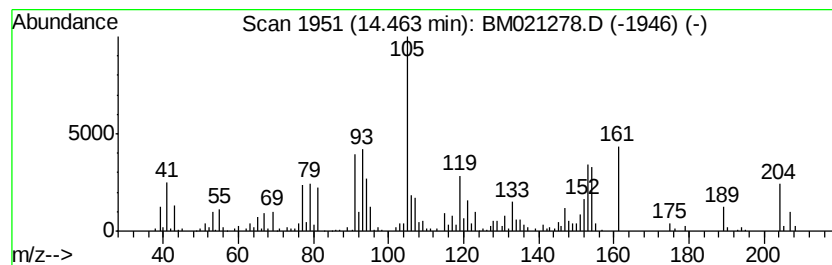
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 12 Naphthalene, 1,2,4a,5,6,8a-... Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.46	4.22 ng/ul	810262	Acenaphthene-d10	14.40
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Naphthalene, 1,2,4a,5,6,8a-hexah...	204 C15H24	031983-22-9	93
2	Naphthalene, 1,2,4a,5,6,8a-hexah...	204 C15H24	031983-22-9	91
3	Naphthalene, 1,2,4a,5,6,8a-hexah...	204 C15H24	031983-22-9	87
4	Naphthalene, 1,2,4a,5,6,8a-hexah...	204 C15H24	000483-75-0	78
5	Naphthalene, 1,2,4a,5,6,8a-hexah...	204 C15H24	000483-75-0	78



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\
Data File : BM021278.D
Acq On : 30 Jun 2019 00:54
Operator : HP/JU
Sample : K3590-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

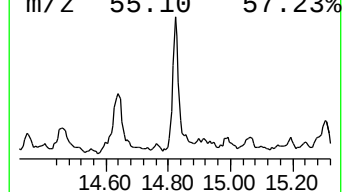
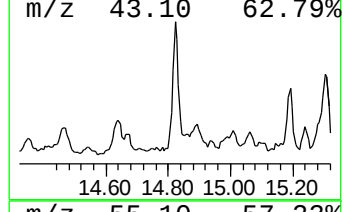
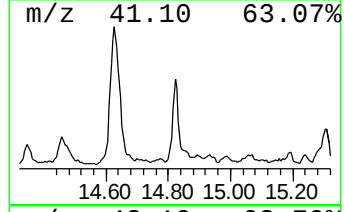
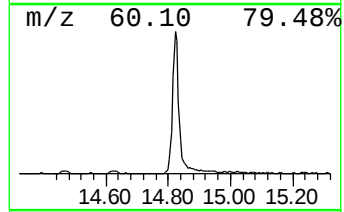
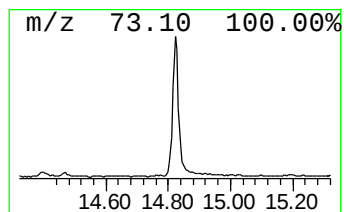
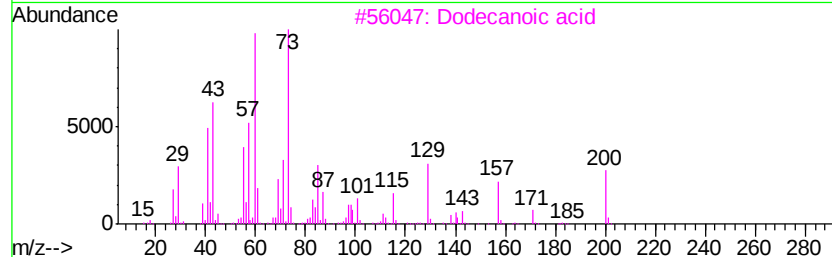
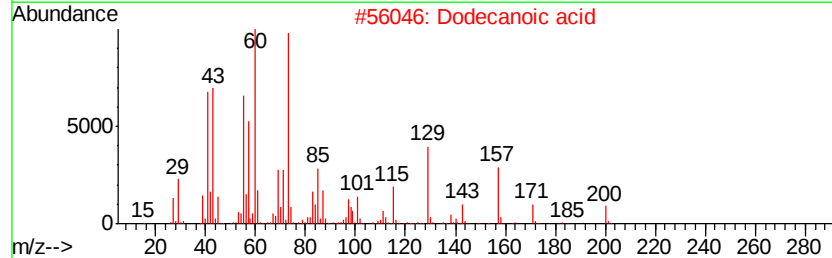
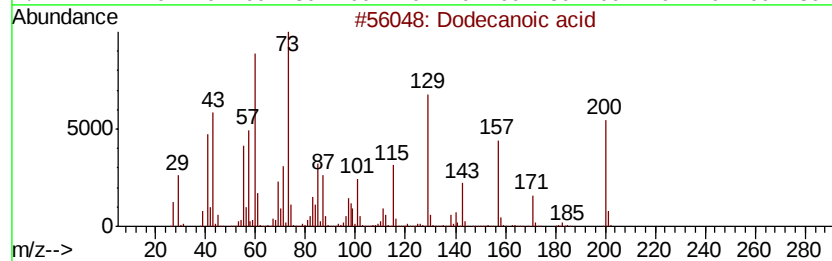
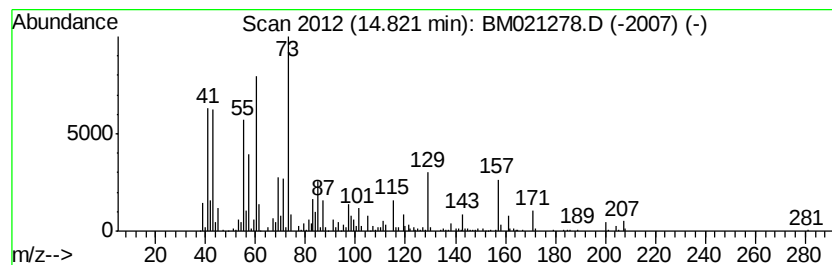
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 13 Dodecanoic acid Concentration Rank 43

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.82	3.36 ng/ul	645037	Acenaphthene-d10	14.40
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Dodecanoic acid	200 C12H24O2	000143-07-7	95
2	Dodecanoic acid	200 C12H24O2	000143-07-7	93
3	Dodecanoic acid	200 C12H24O2	000143-07-7	91
4	Dodecanoic acid	200 C12H24O2	000143-07-7	91
5	Tetradecanoic acid	228 C14H28O2	000544-63-8	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\
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Acq On : 30 Jun 2019 00:54
Operator : HP/JU
Sample : K3590-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

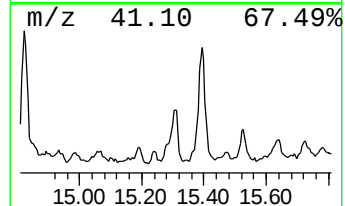
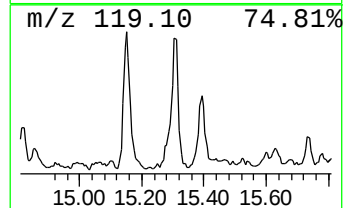
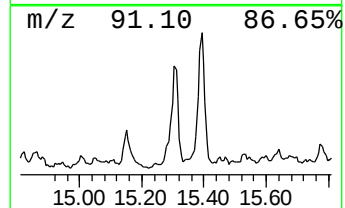
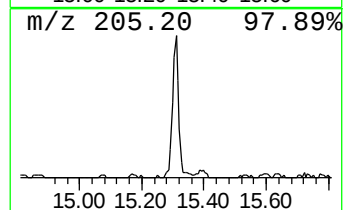
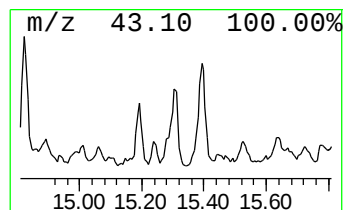
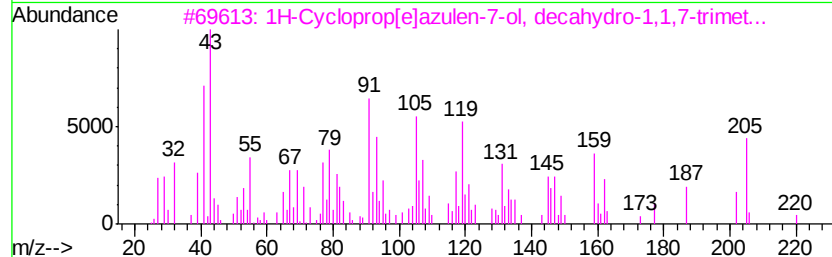
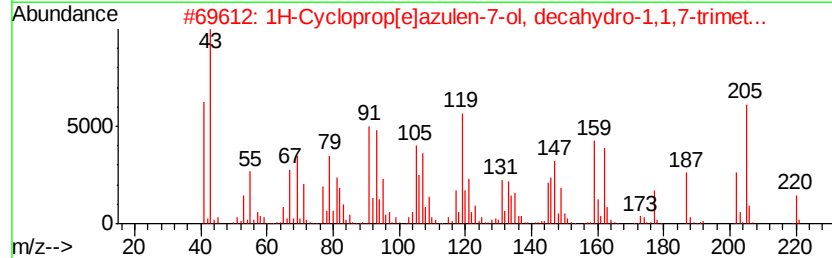
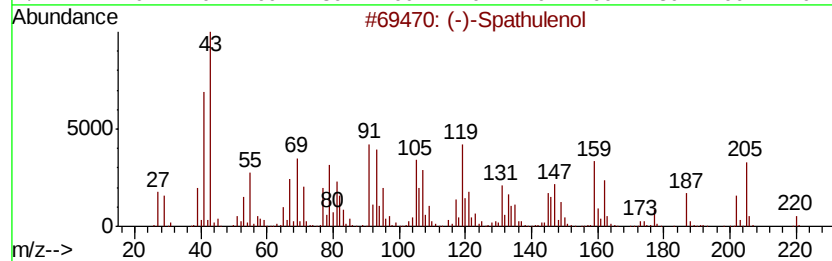
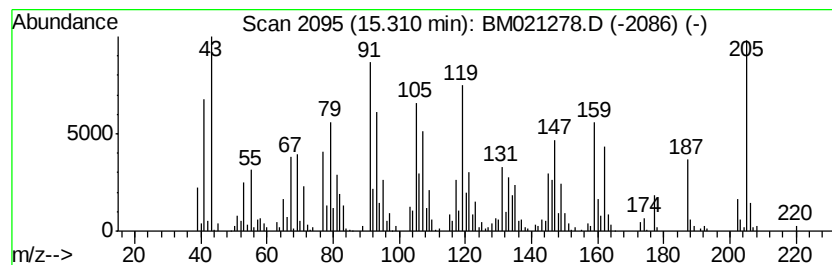
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 (-)-Spathulenol Concentration Rank 39

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.31	3.64 ng/ul	699596	Acenaphthene-d10	14.40
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		(-)-Spathulenol	220 C15H24O	077171-55-2 91
2		1H-Cycloprop[elazulen-7-ol, deca...	220 C15H24O	006750-60-3 87
3		1H-Cycloprop[elazulen-7-ol, deca...	220 C15H24O	006750-60-3 43
4		1,3,6,10-Dodecatetraene, 3,7,11-...	204 C15H24	026560-14-5 35
5		1H-1,5-Benzodiazepine, 2,3,4,5-t...	162 C10H14N2	040358-34-7 30



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\
Data File : BM021278.D
Acq On : 30 Jun 2019 00:54
Operator : HP/JU
Sample : K3590-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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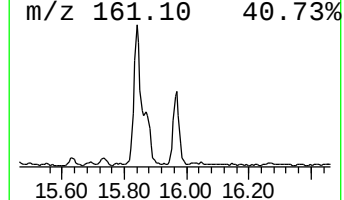
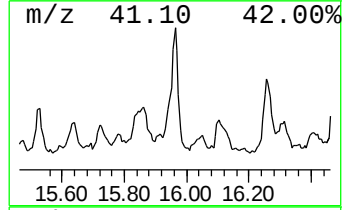
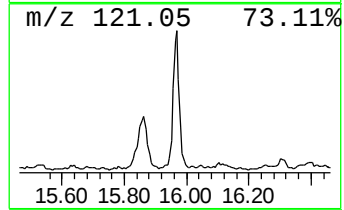
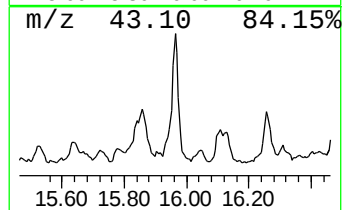
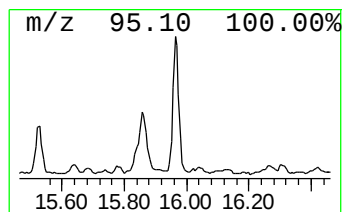
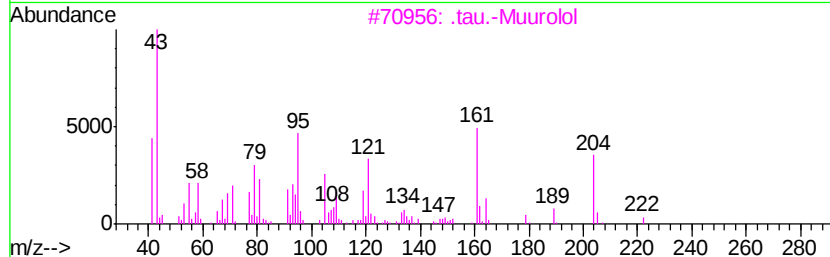
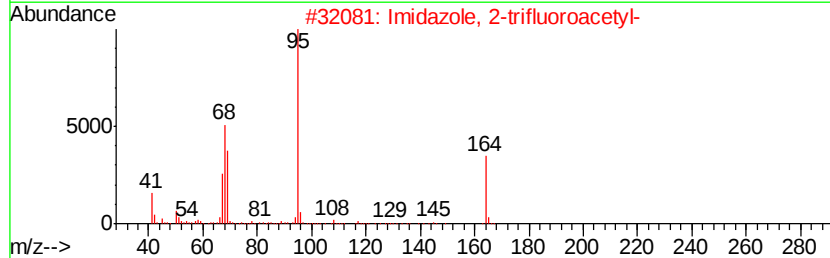
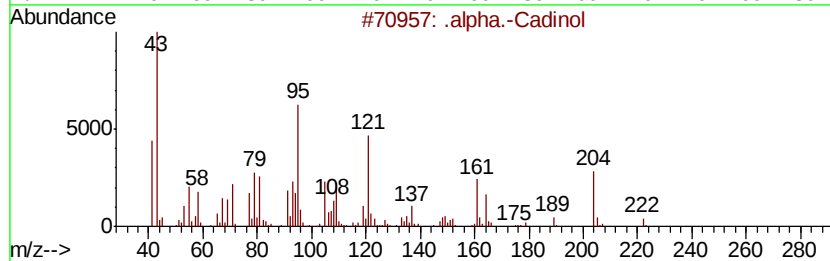
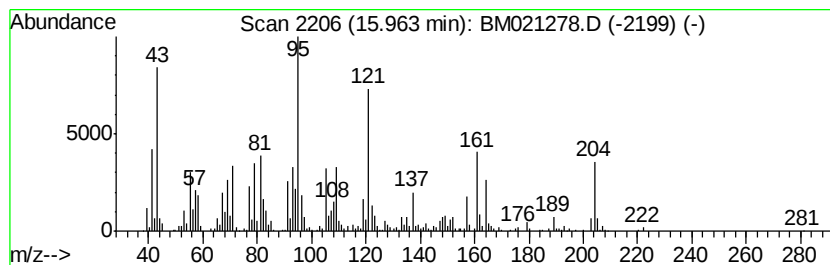
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 .alpha.-Cadinol Concentration Rank 35

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.96	4.89 ng/ul	1056270	Phenanthrene-d10	17.15

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			.alpha.-Cadinol	222	C15H26O	000481-34-5	87
2			Imidazole, 2-trifluoroacetyl-	164	C5H3F3N2O	105480-29-3	30
3			.tau.-Muuroiol	222	C15H26O	019912-62-0	30
4			Cyclopentanecarboxylic acid, 2-m...	154	C9H14O2	074764-25-3	27
5			4-Methylimidazole-5-butyric acid	168	C8H12N2O2	1000129-14-7	22



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\
Data File : BM021278.D
Acq On : 30 Jun 2019 00:54
Operator : HP/JU
Sample : K3590-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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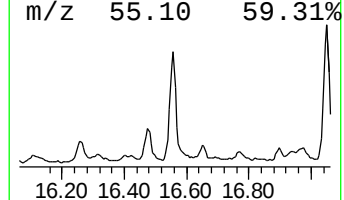
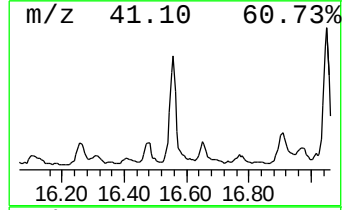
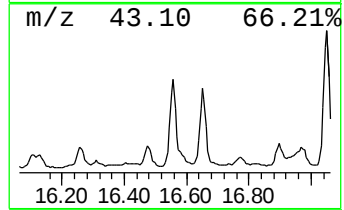
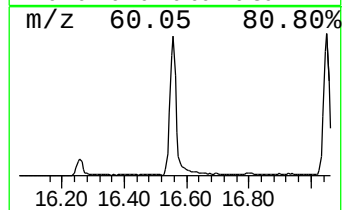
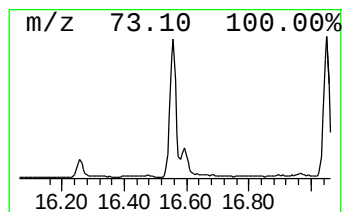
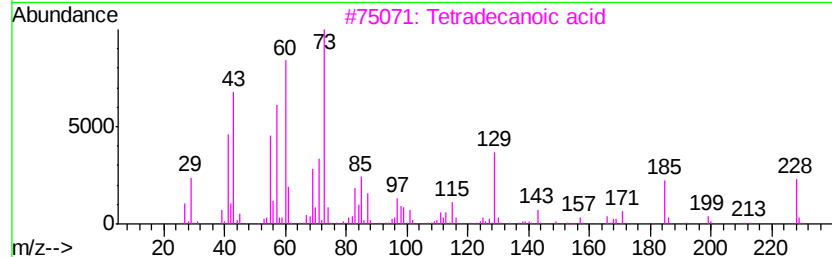
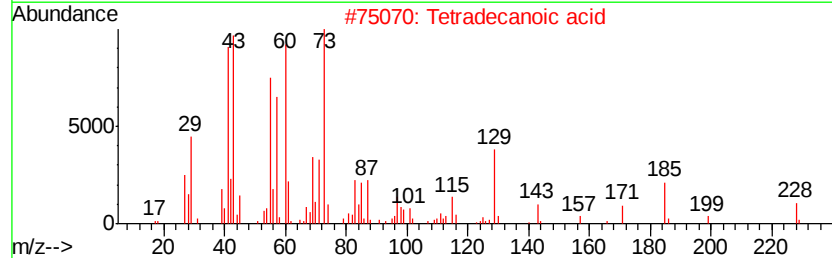
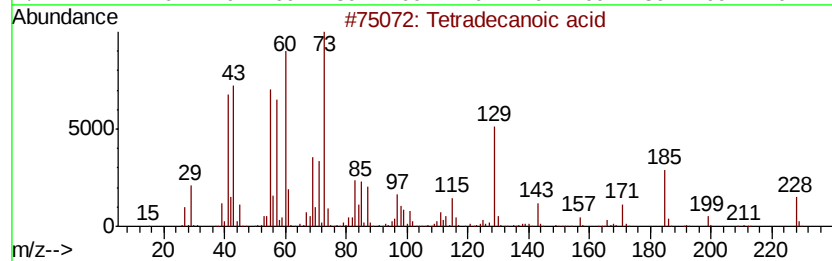
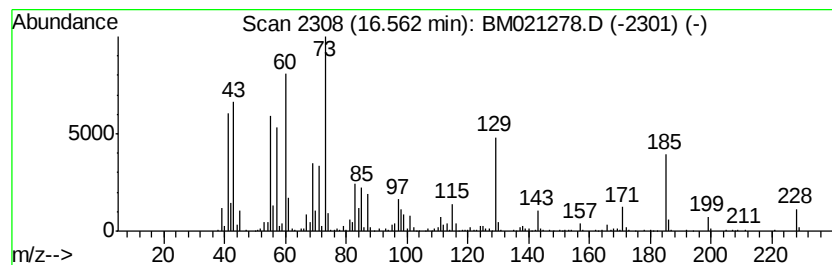
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 Tetradecanoic acid Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.56	6.48 ng/ul	1397920	Phenanthrene-d10	17.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetradecanoic acid	228	C14H28O2	000544-63-8	99
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	96
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	83
4		n-Decanoic acid	172	C10H20O2	000334-48-5	70
5		Dodecanoic acid	200	C12H24O2	000143-07-7	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\
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Acq On : 30 Jun 2019 00:54
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Sample : K3590-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

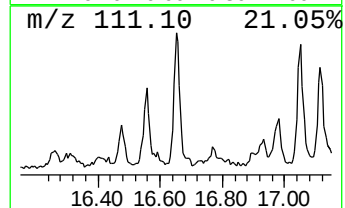
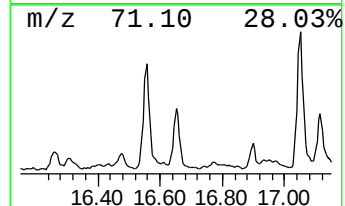
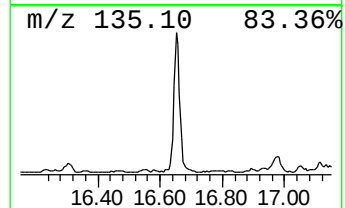
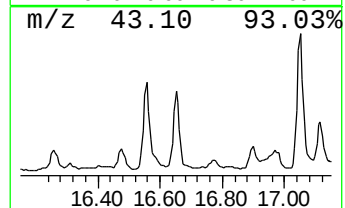
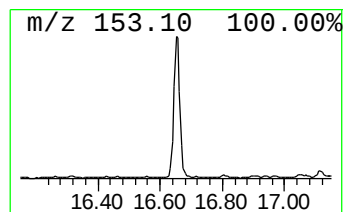
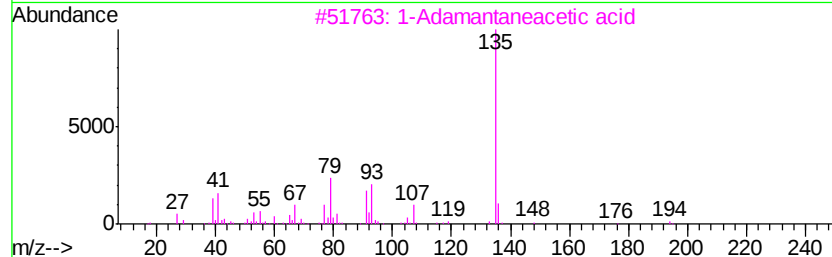
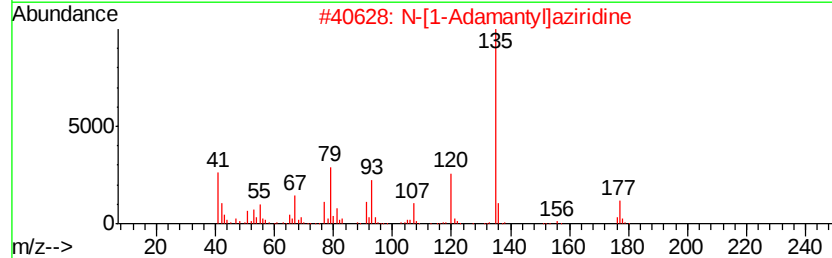
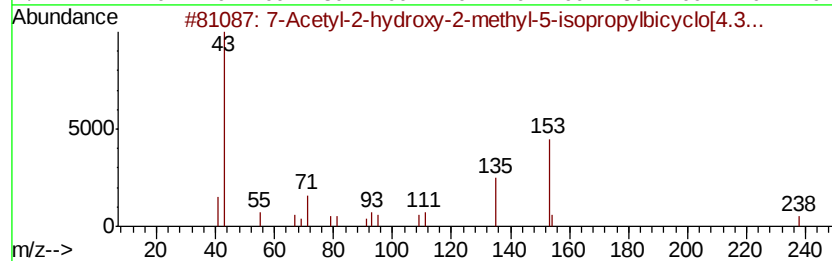
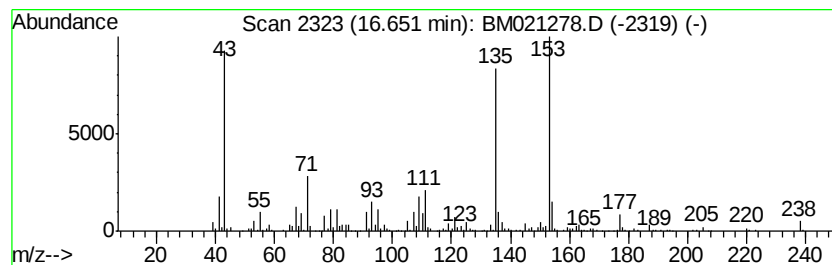
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 17 7-Acetyl-2-hydroxy-2-methyl... Concentration Rank 44

R.T.	EstConc	Area	Relative to ISTD	R.T.		
16.65	3.24 ng/ul	700014	Phenanthrene-d10	17.15		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	7-Acetyl-2-hydroxy-2-methyl-5-is...	238	C15H26O2	096093-81-1	53	
2	N-[1-Adamantyl]aziridine	177	C12H19N	1000256-62-9	38	
3	1-Adamantaneacetic acid	194	C12H18O2	004942-47-6	30	
4	Tricyclo[3.3.1.1(3,7)]decane, 1-...	181	C10H15NO2	007575-82-8	27	
5	3-Adamantan-1-yl-3-oxo-propionit...	203	C13H17NO	1000296-74-6	27	



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\
Data File : BM021278.D
Acq On : 30 Jun 2019 00:54
Operator : HP/JU
Sample : K3590-01
Misc :
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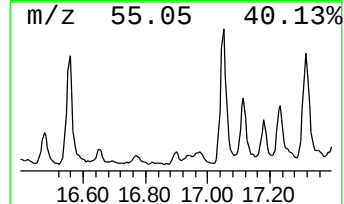
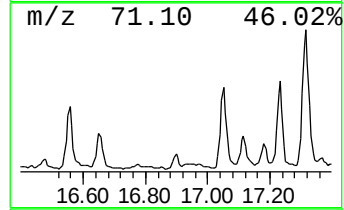
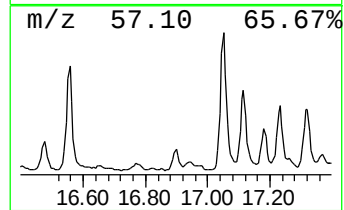
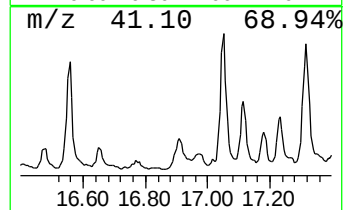
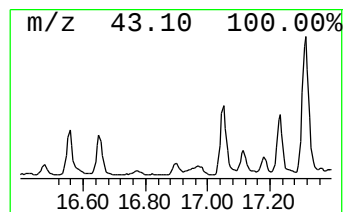
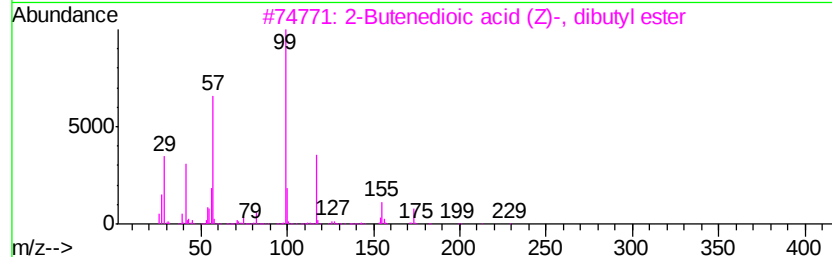
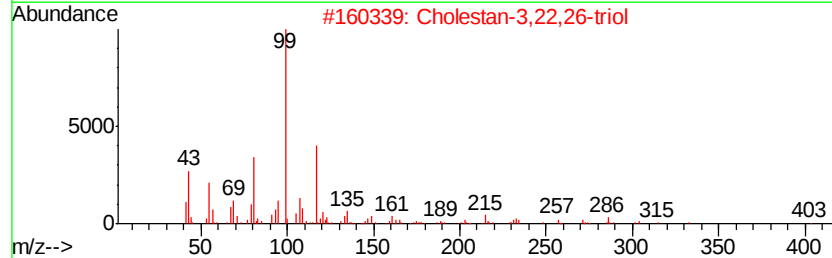
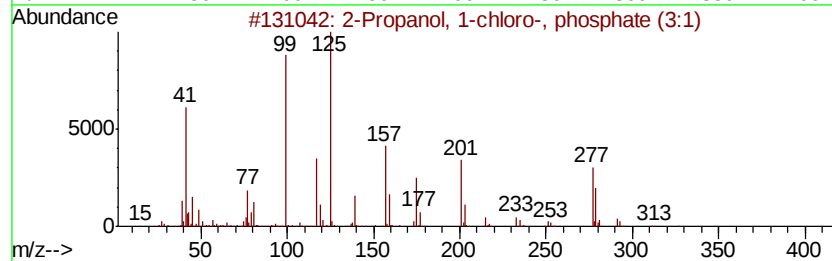
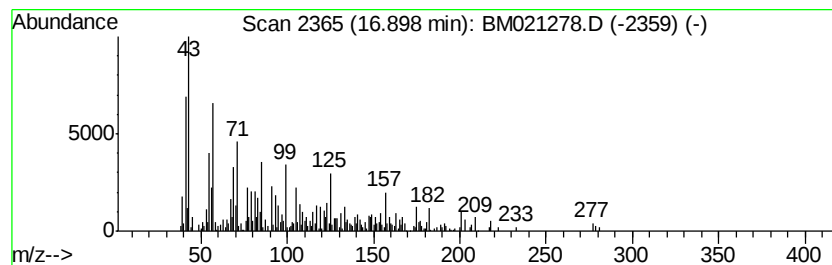
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 2-Propanol, 1-chloro-, phos... Concentration Rank 41

R.T.	EstConc	Area	Relative to ISTD	R.T.		
16.90	3.43 ng/ul	740025	Phenanthrene-d10	17.15		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-chloro-, phosphate...	326	C9H18Cl3O4P	013674-84-5	87	
2	Cholestan-3,22,26-triol	420	C27H48O3	1000252-39-8	27	
3	2-Butenedioic acid (Z)-, dibutyl...	228	C12H20O4	000105-76-0	27	
4	2-Heptyl methylphosphonofluoridate	196	C8H18FO2P	000674-95-3	25	
5	Hexanohydroxamic acid	131	C6H13NO2	004312-93-0	18	



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\
Data File : BM021278.D
Acq On : 30 Jun 2019 00:54
Operator : HP/JU
Sample : K3590-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C5BA0

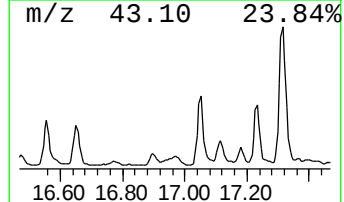
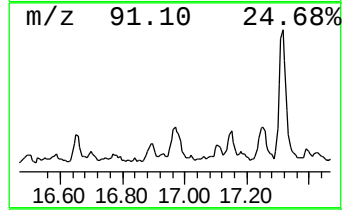
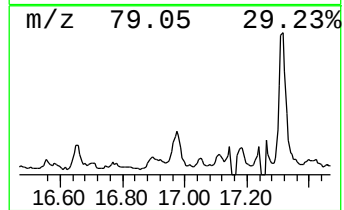
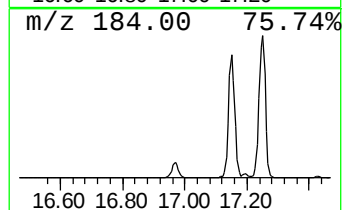
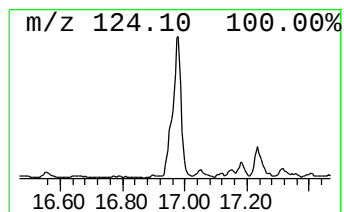
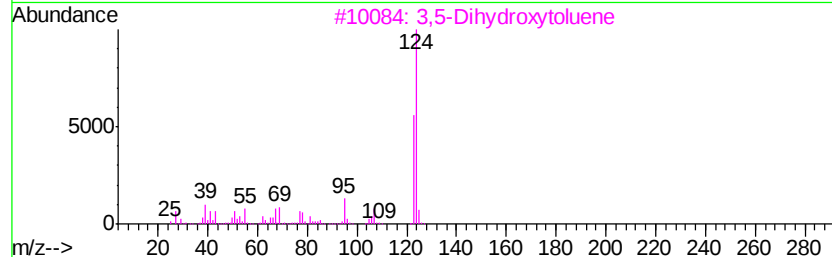
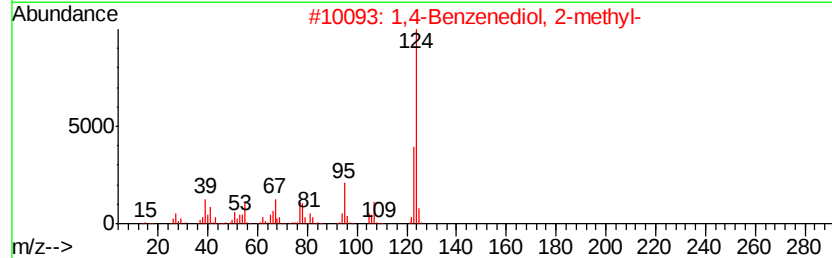
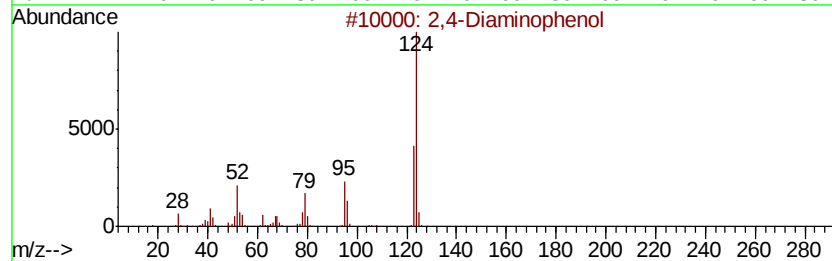
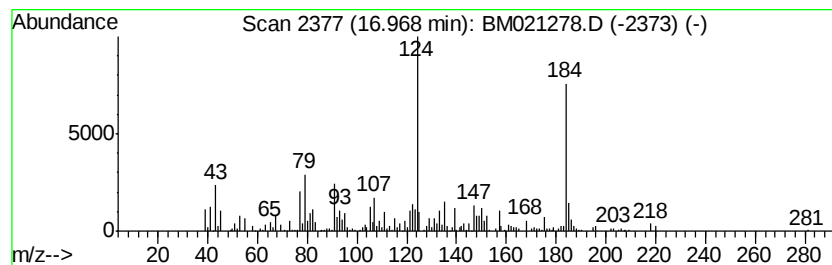
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 unknown-02 Concentration Rank 50

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.97	2.85 ng/ul	615160	Phenanthrene-d10	17.15

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,4-Diaminophenol	124	C6H8N2O	000095-86-3	43
2			1,4-Benzenediol, 2-methyl-	124	C7H8O2	000095-71-6	43
3			3,5-Dihydroxytoluene	124	C7H8O2	000504-15-4	38
4			Benzeneacetic acid, .alpha.-hydr...	291	C17H25NO3	000100-91-4	37
5			1,4-Benzenediol, 2-methyl-	124	C7H8O2	000095-71-6	37



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\
Data File : BM021278.D
Acq On : 30 Jun 2019 00:54
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Sample : K3590-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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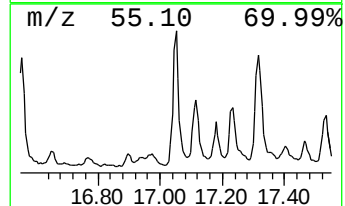
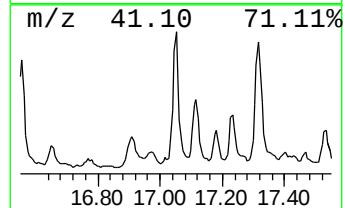
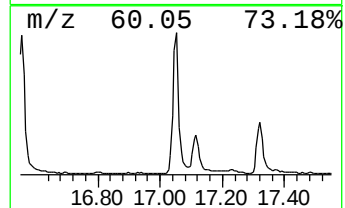
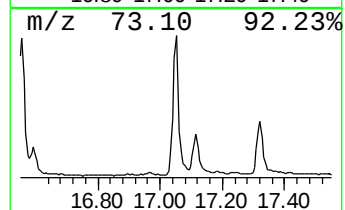
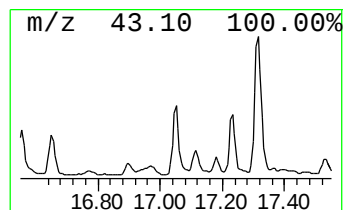
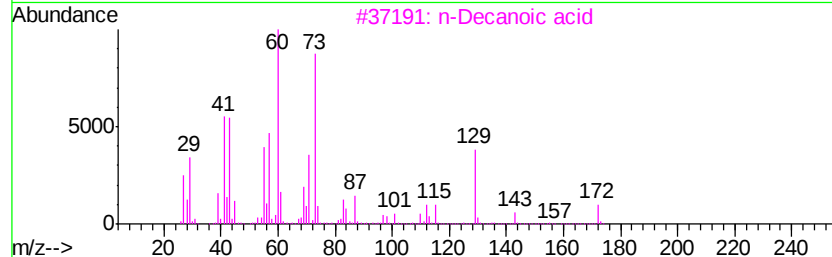
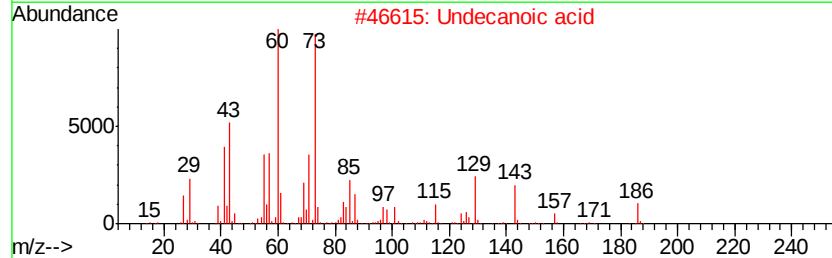
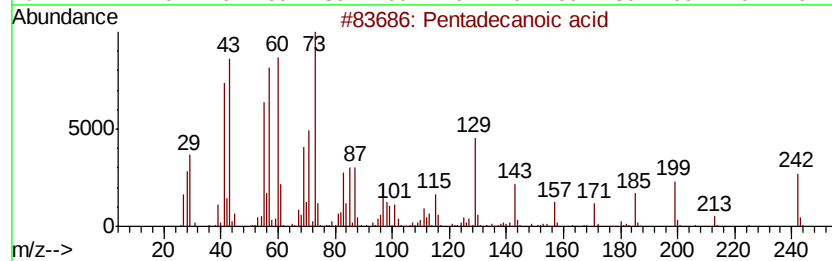
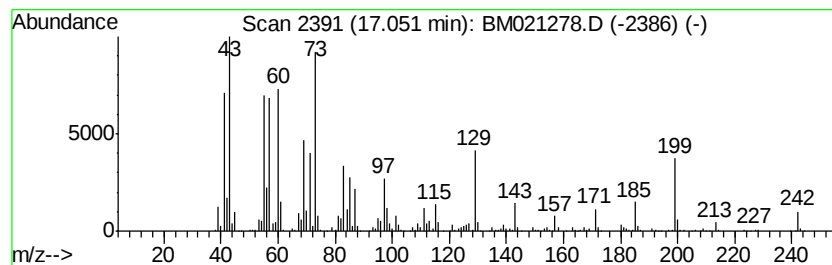
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 Pentadecanoic acid Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.05	7.69 ng/ul	1658740	Phenanthrene-d10	17.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecanoic acid	242	C15H30O2	001002-84-2	87
2		Undecanoic acid	186	C11H22O2	000112-37-8	74
3		n-Decanoic acid	172	C10H20O2	000334-48-5	70
4		n-Decanoic acid	172	C10H20O2	000334-48-5	70
5		Dodecanoic acid	200	C12H24O2	000143-07-7	68



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\
Data File : BM021278.D
Acq On : 30 Jun 2019 00:54
Operator : HP/JU
Sample : K3590-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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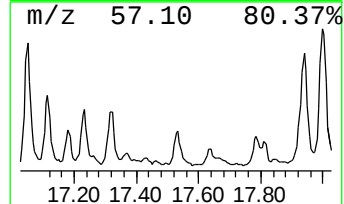
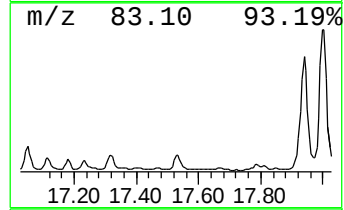
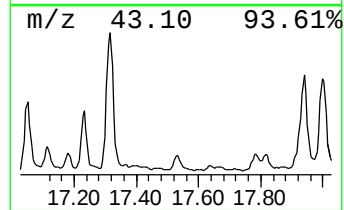
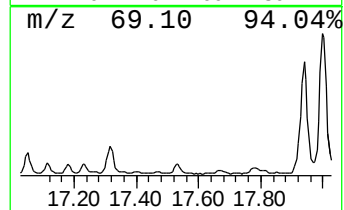
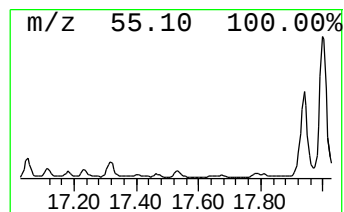
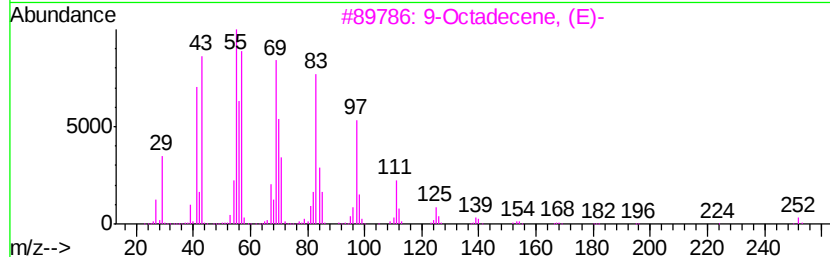
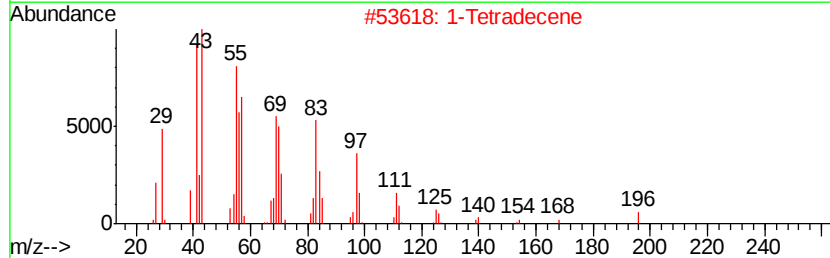
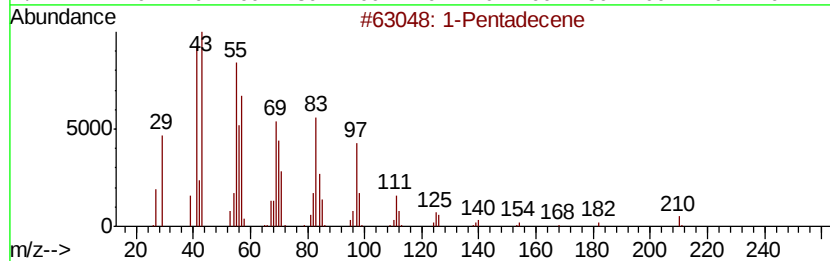
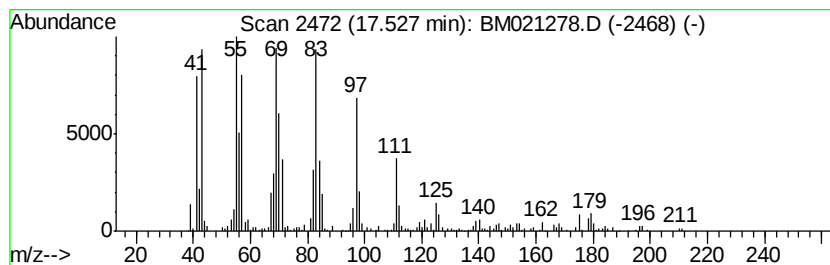
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 (DEL) Alkane: Cyclic17.53 Concentration Rank 53

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.53	2.37 ng/ul	510592	Phenanthrene-d10	17.15

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Pentadecene	210	C15H30	013360-61-7	98
2			1-Tetradecene	196	C14H28	001120-36-1	95
3			9-Octadecene, (E)-	252	C18H36	007206-25-9	91
4			1-Pentadecanol	228	C15H32O	000629-76-5	91
5			1-Tetradecanol	214	C14H30O	000112-72-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\
Data File : BM021278.D
Acq On : 30 Jun 2019 00:54
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Sample : K3590-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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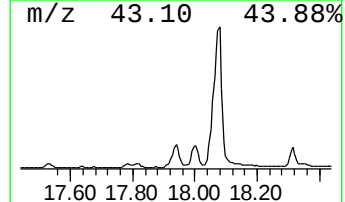
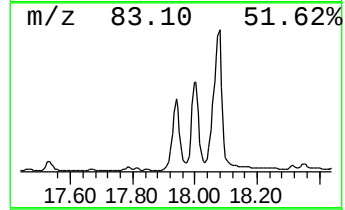
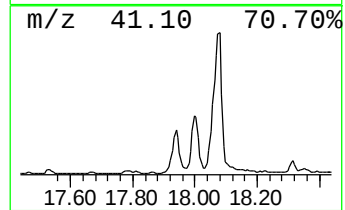
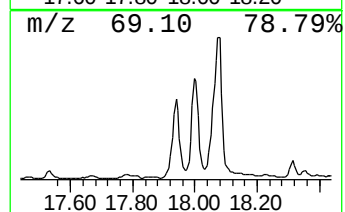
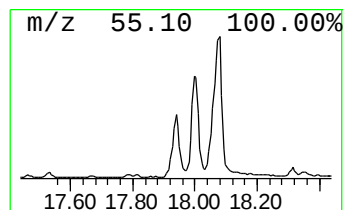
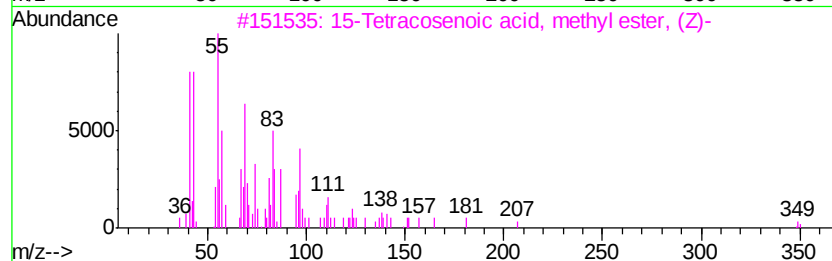
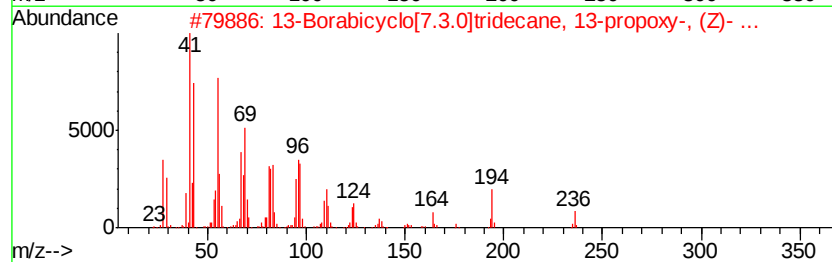
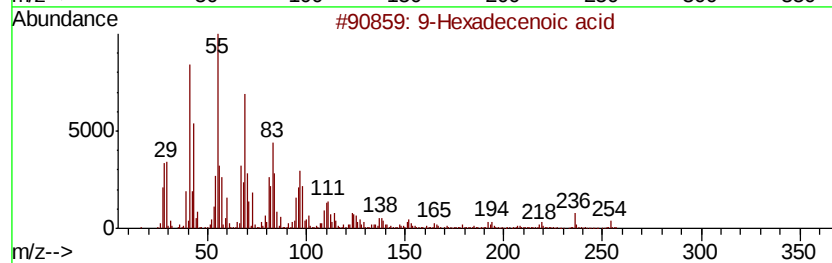
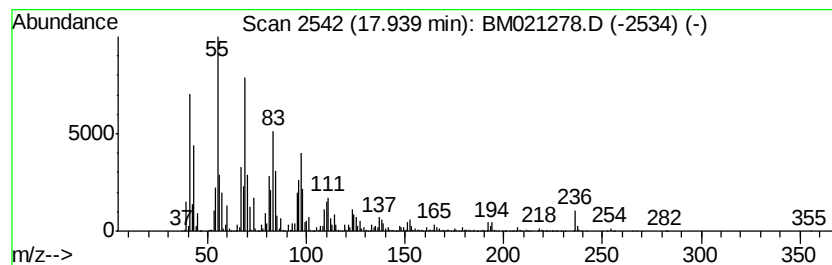
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 9-Hexadecenoic acid Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.94	29.48 ng/ul	6363750	Phenanthrene-d10	17.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Hexadecenoic acid	254	C16H30O2	002091-29-4	78
2		13-Borabicyclo[7.3.0]tridecane, ...	236	C15H29B0	1000156-41-7	55
3		15-Tetracosenoic acid, methyl es...	380	C25H48O2	002733-88-2	55
4		1-Cyclohexyl-2-methyl-prop-2-en-...	152	C10H16O	025183-82-8	55
5		Hexadecenoic acid, Z-11-	254	C16H30O2	002416-20-8	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\
Data File : BM021278.D
Acq On : 30 Jun 2019 00:54
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Sample : K3590-01
Misc :
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ClientSampleId :
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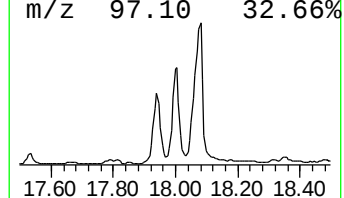
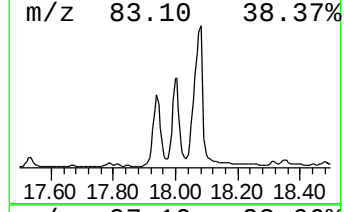
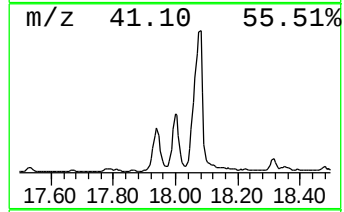
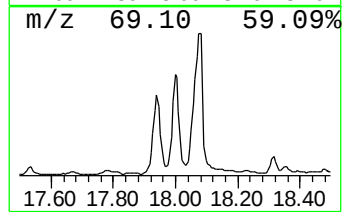
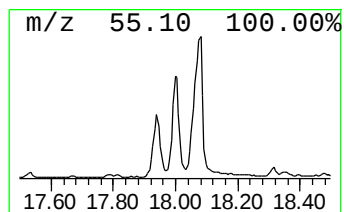
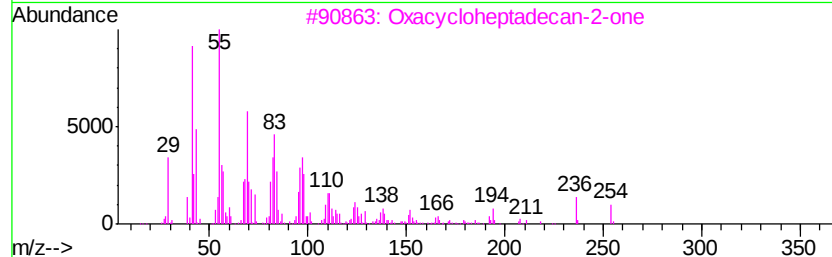
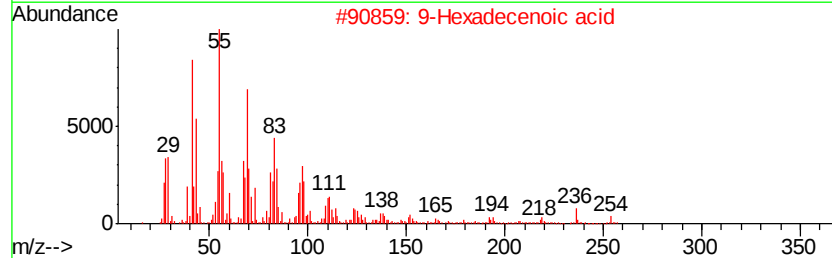
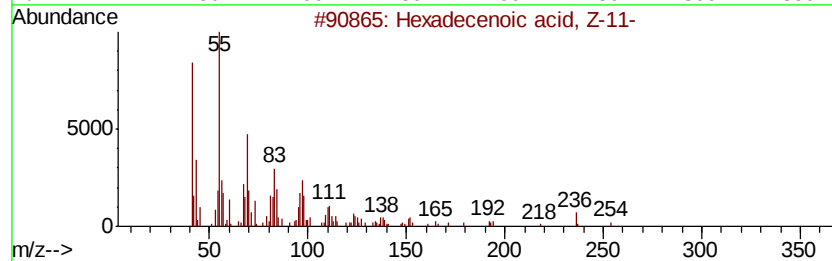
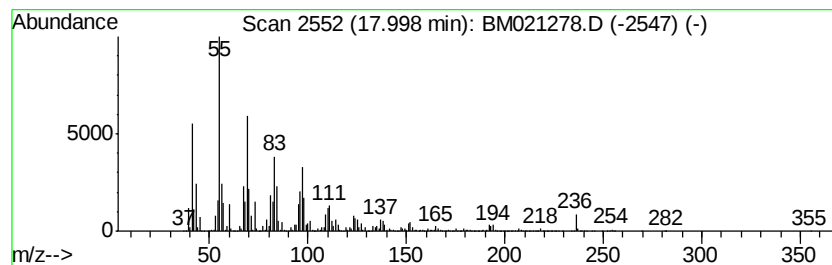
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Quant Title : SVOA CALIBRATION

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

Peak Number 23 Hexadecenoic acid, Z-11- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.00	34.13 ng/ul	7366070	Phenanthrene-d10	17.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecenoic acid, Z-11-	254	C16H30O2	002416-20-8	94
2		9-Hexadecenoic acid	254	C16H30O2	002091-29-4	91
3		Oxacycloheptadecan-2-one	254	C16H30O2	000109-29-5	91
4		15-Tetracosenoic acid, methyl es...	380	C25H48O2	002733-88-2	90
5		cis-2-Methyl-4-n-pentylthiane, S...	218	C11H22O2S	1000215-67-9	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\
Data File : BM021278.D
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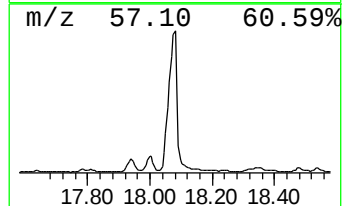
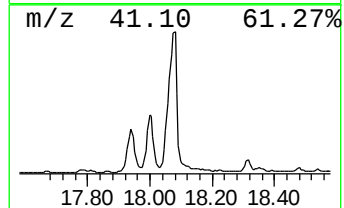
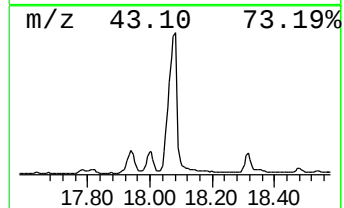
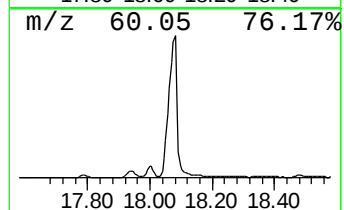
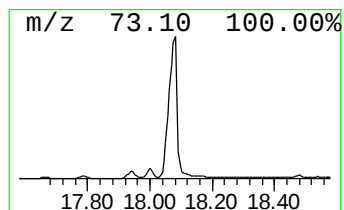
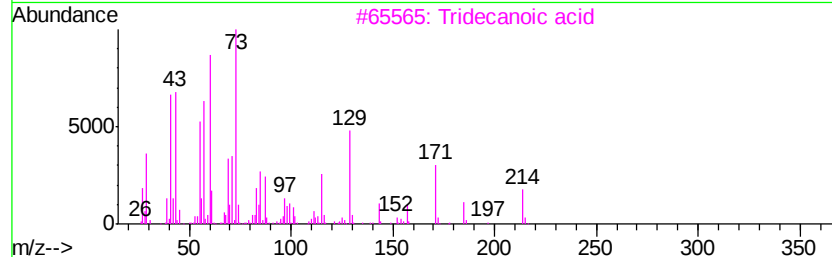
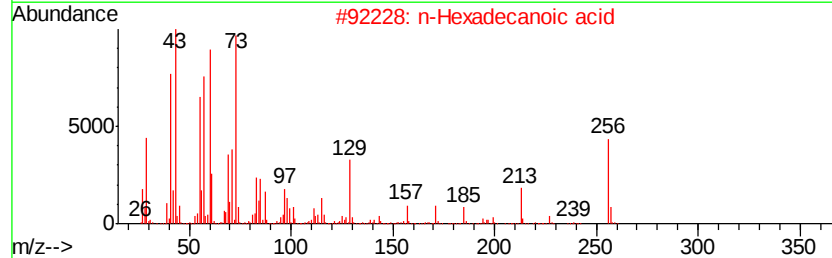
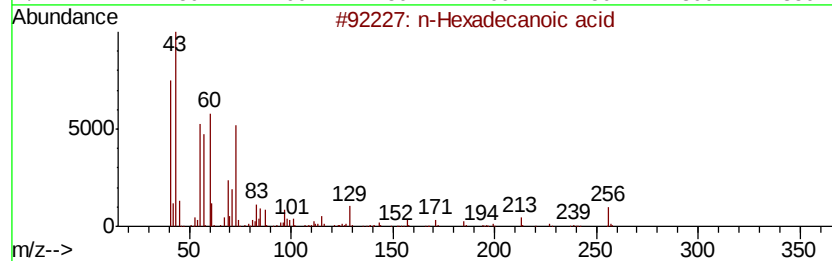
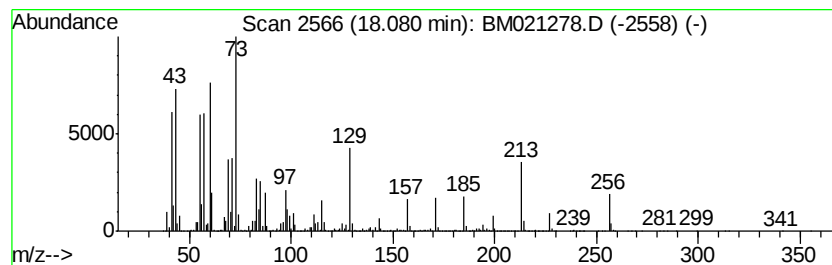
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Quant Title : SVOA CALIBRATION

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

Peak Number 24 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.08	114.38 ng/ul	24686800	Phenanthrene-d10	17.15

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	97
3		Tridecanoic acid	214	C13H26O2	000638-53-9	89
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	86
5		Tridecanoic acid	214	C13H26O2	000638-53-9	76



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\
Data File : BM021278.D
Acq On : 30 Jun 2019 00:54
Operator : HP/JU
Sample : K3590-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

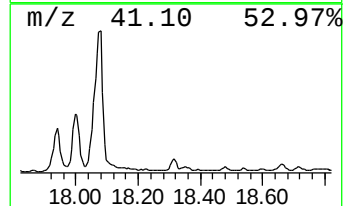
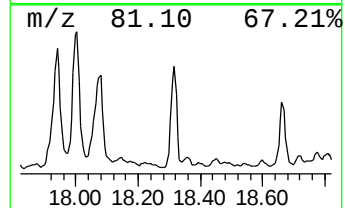
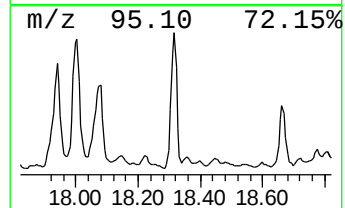
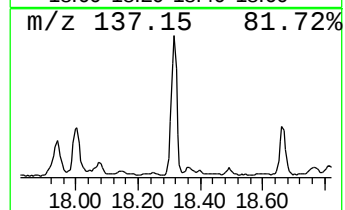
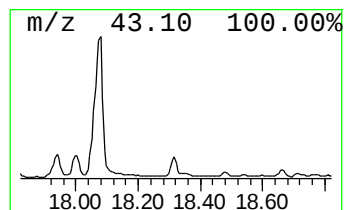
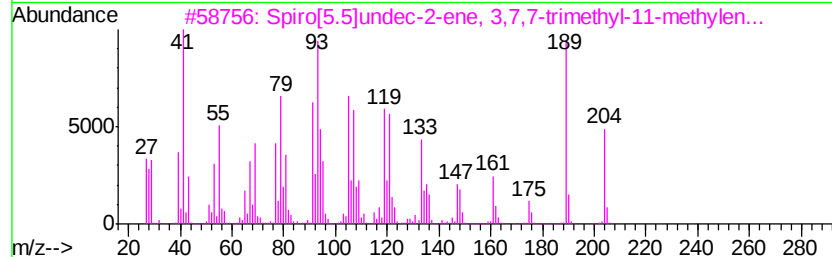
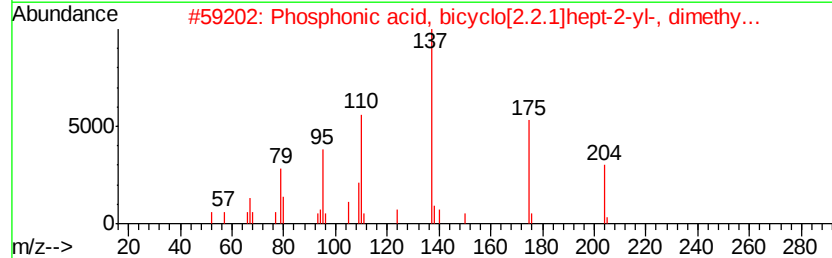
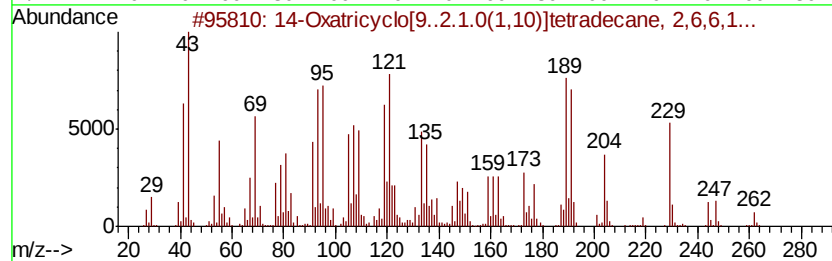
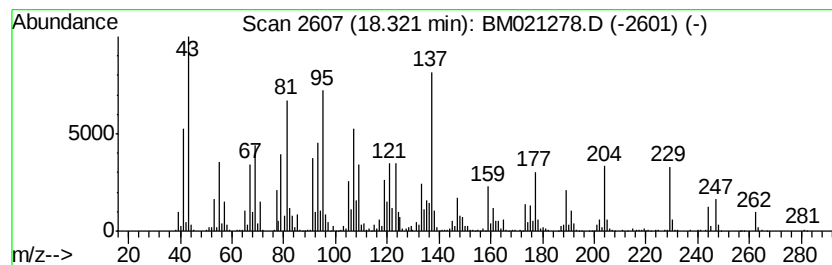
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 25 unknown-03 Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.		
18.32	11.77 ng/ul	2540470	Phenanthrene-d10	17.15		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	14-Oxatricyclo[9..2.1.0(1,10)]te...	262	C18H30O	1000193-06-8	46	
2	Phosphonic acid, bicyclo[2.2.1]h...	204	C9H17O3P	031061-88-8	45	
3	Spiro[5.5]undec-2-ene, 3,7,7-tri...	204	C15H24	018431-82-8	41	
4	N-Cyclohexanecarbonylanthranilic...	247	C14H17NO3	025437-73-4	35	
5	Naphthalene, 2-butyldecahydro-	194	C14H26	006305-52-8	27	



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\
Data File : BM021278.D
Acq On : 30 Jun 2019 00:54
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Sample : K3590-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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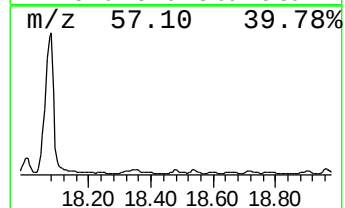
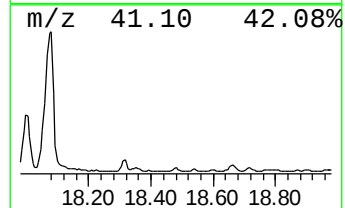
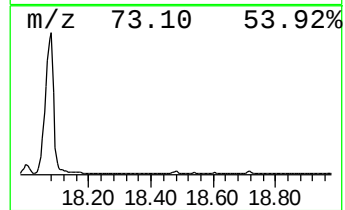
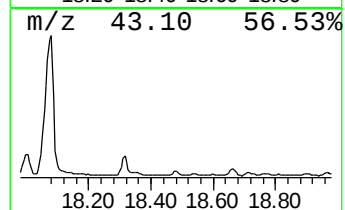
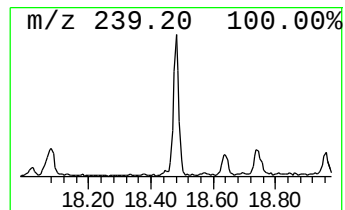
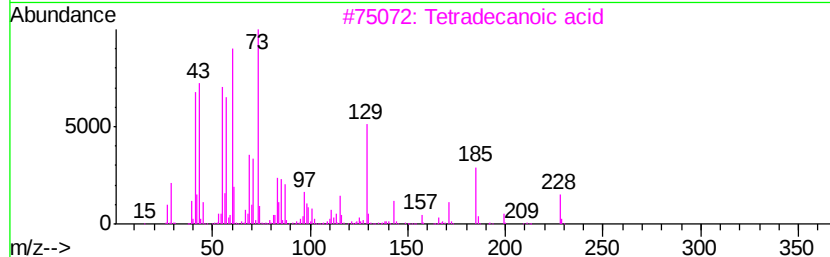
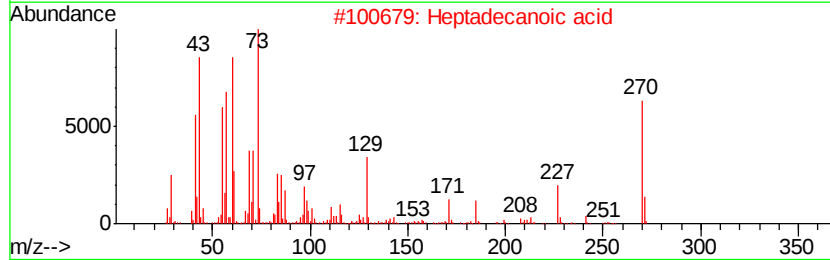
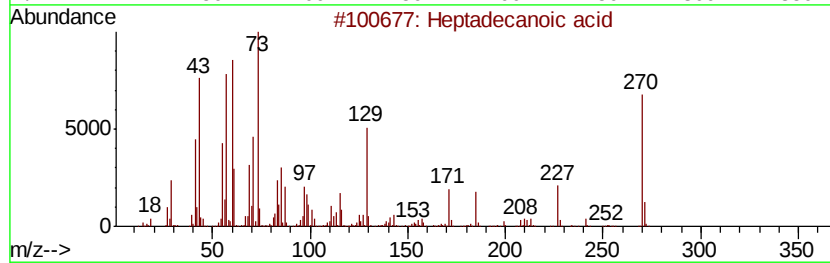
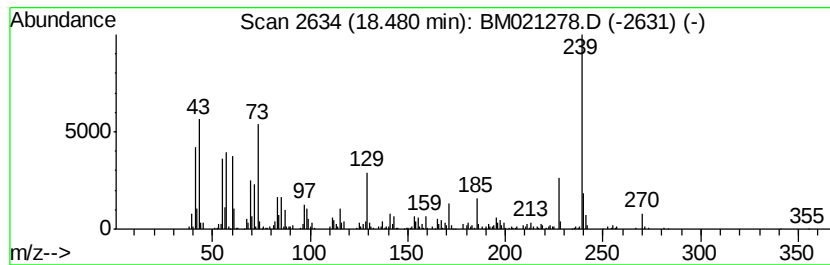
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Quant Title : SVOA CALIBRATION

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

Peak Number 26 Heptadecanoic acid Concentration Rank 55

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.48	2.32 ng/ul	501103	Phenanthrene-d10	17.15

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecanoic acid	270	C17H34O2	000506-12-7	95
2			Heptadecanoic acid	270	C17H34O2	000506-12-7	90
3			Tetradecanoic acid	228	C14H28O2	000544-63-8	84
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	78
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\
Data File : BM021278.D
Acq On : 30 Jun 2019 00:54
Operator : HP/JU
Sample : K3590-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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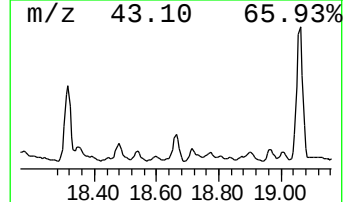
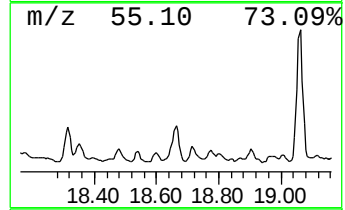
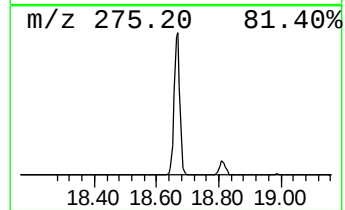
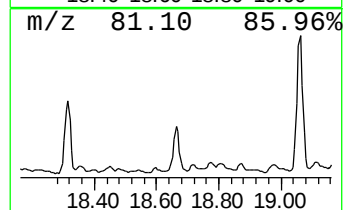
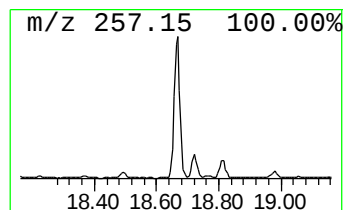
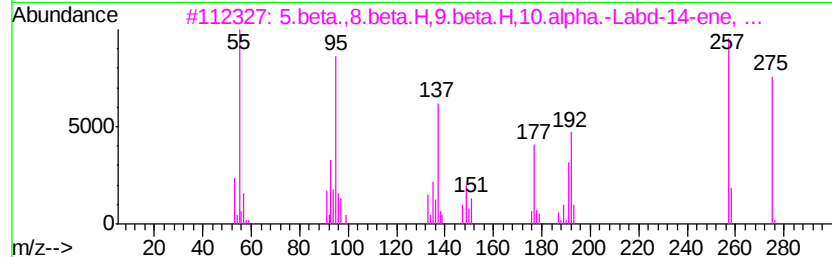
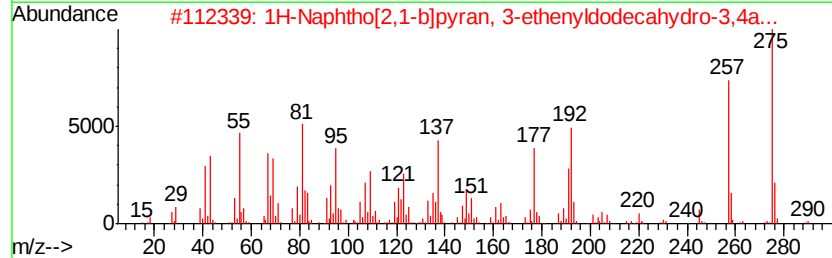
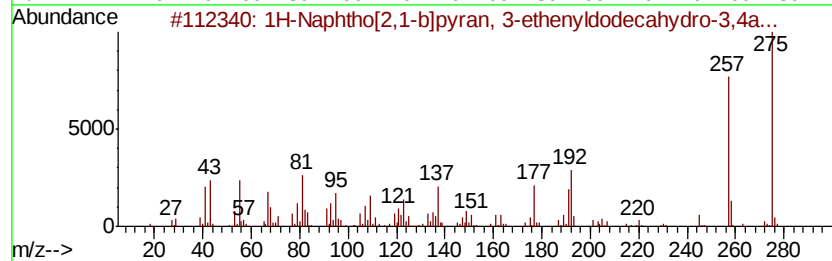
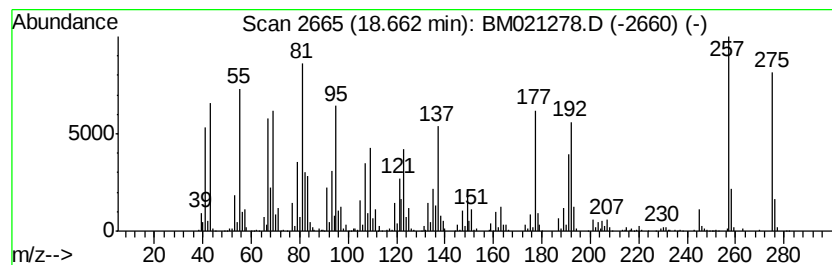
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Quant Title : SVOA CALIBRATION

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

Peak Number 27 1H-Naphtho[2,1-b]pyran, 3-e... Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.66	7.34 ng/ul	1584020	Phenanthrene-d10	17.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Naphtho[2,1-b]pyran, 3-etheny...	290	C20H34O	001227-93-6	80
2		1H-Naphtho[2,1-b]pyran, 3-etheny...	290	C20H34O	000596-84-9	74
3		5.beta.,8.beta.H,9.beta.H,10.alpha...	290	C20H34O	019123-29-6	64
4		4a,7,7,10a-Tetramethyl-dodecahyd...	264	C17H28O2	1000192-27-9	38
5		4H-1-Benzopyran-4-one, 2,3-dihyd...	192	C11H12O3	017771-33-4	25



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\
Data File : BM021278.D
Acq On : 30 Jun 2019 00:54
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Sample : K3590-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampled :
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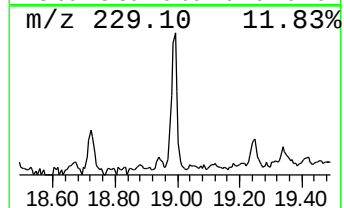
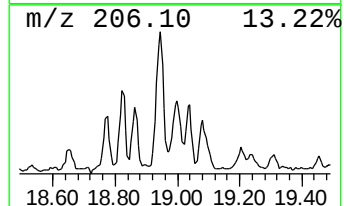
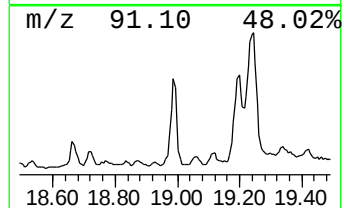
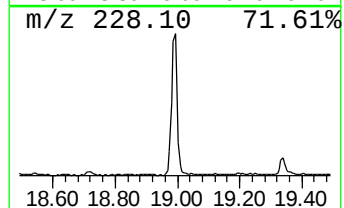
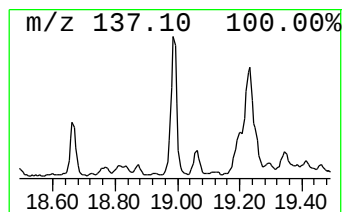
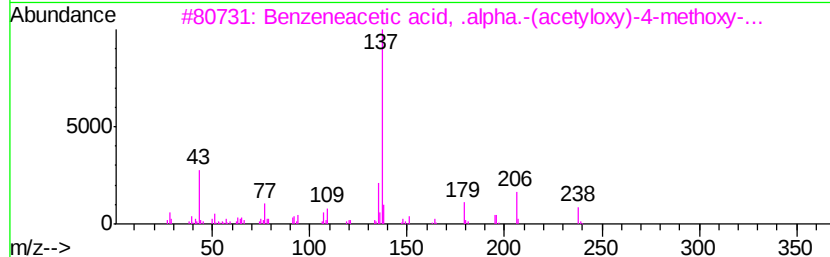
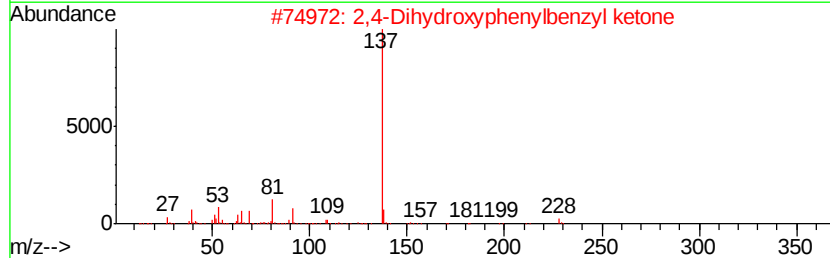
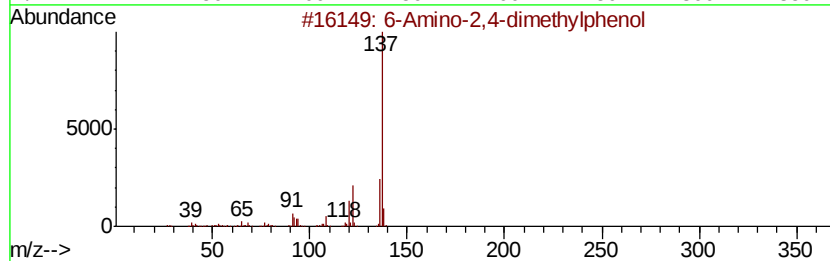
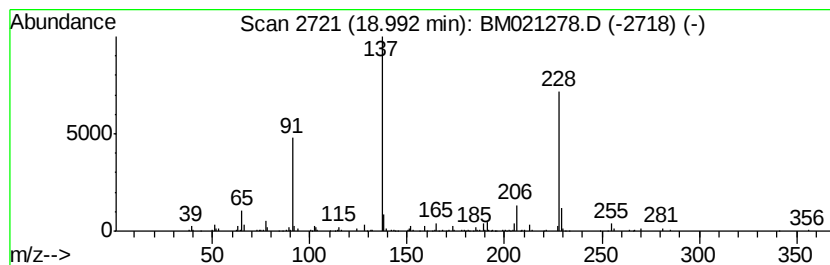
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Quant Title : SVOA CALIBRATION

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

Peak Number 29 unknown-04 Concentration Rank 40

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.99	3.64 ng/ul	785704	Phenanthrene-d10	17.15

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			6-Amino-2,4-dimethylphenol	137	C8H11NO	041458-65-5	43
2			2,4-Dihydroxyphenylbenzyl ketone	228	C14H12O3	003669-41-8	25
3			Benzeneacetic acid, .alpha.-(ace...	238	C12H14O5	055538-79-9	25
4			Benzeneacetic acid, .alpha.-(ace...	238	C12H14O5	055538-77-7	25
5			6-Methylnicotinic acid	137	C7H7NO2	003222-47-7	23



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\
Data File : BM021278.D
Acq On : 30 Jun 2019 00:54
Operator : HP/JU
Sample : K3590-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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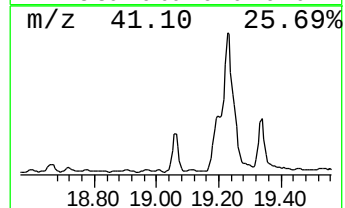
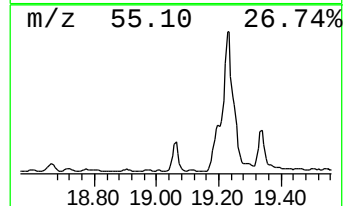
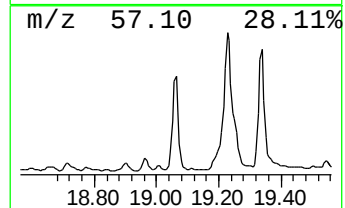
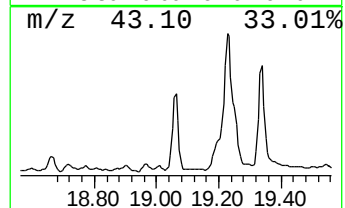
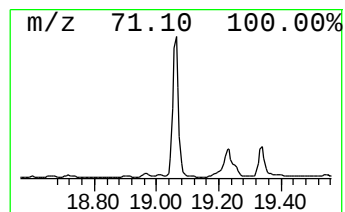
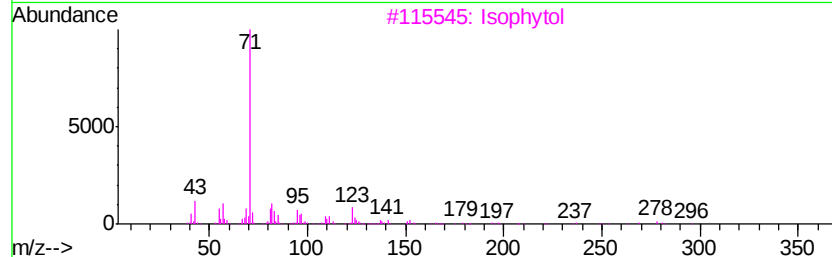
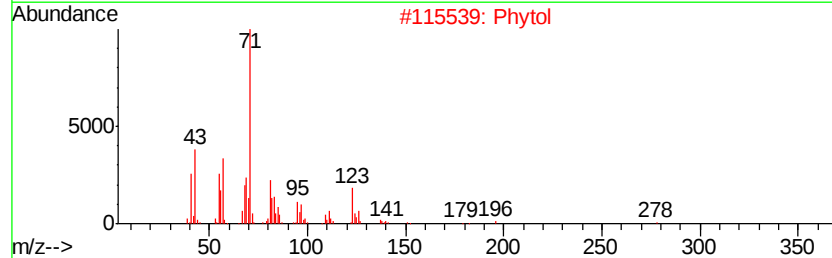
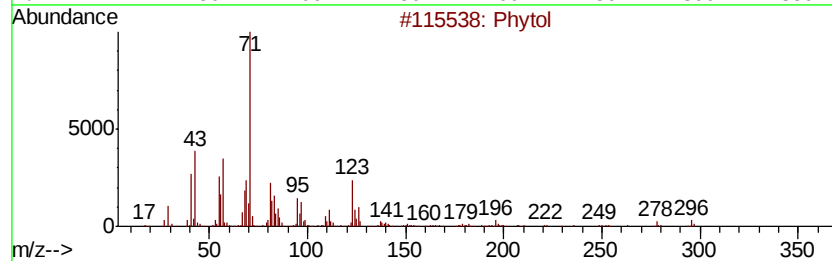
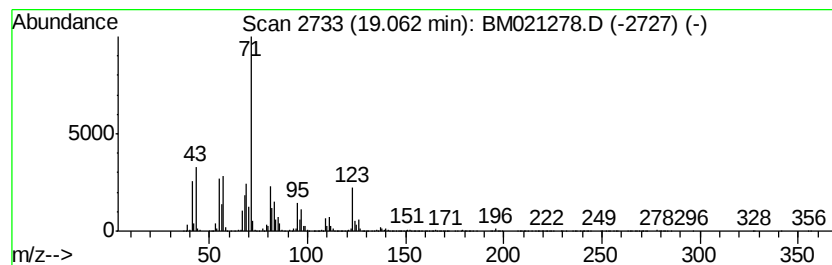
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Quant Title : SVOA CALIBRATION

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

Peak Number 30 Phytol Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.06	21.66 ng/ul	4675430	Phenanthrene-d10	17.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phytol	296	C20H40O	000150-86-7	93
2		Phytol	296	C20H40O	000150-86-7	90
3		Isophytol	296	C20H40O	000505-32-8	87
4		Hexadecane, 1,1-dimethoxy-	286	C18H38O2	002791-29-9	53
5		Phytol	296	C20H40O	000150-86-7	52



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\
Data File : BM021278.D
Acq On : 30 Jun 2019 00:54
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Sample : K3590-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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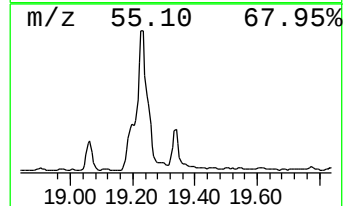
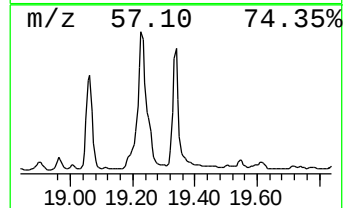
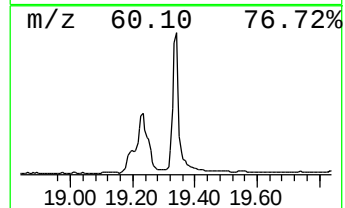
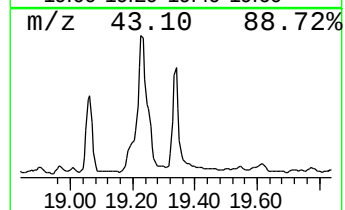
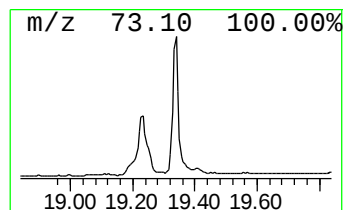
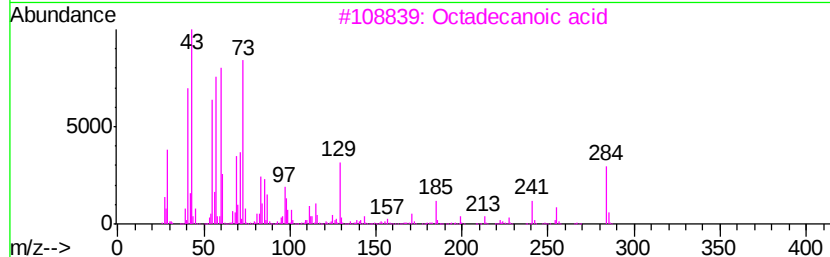
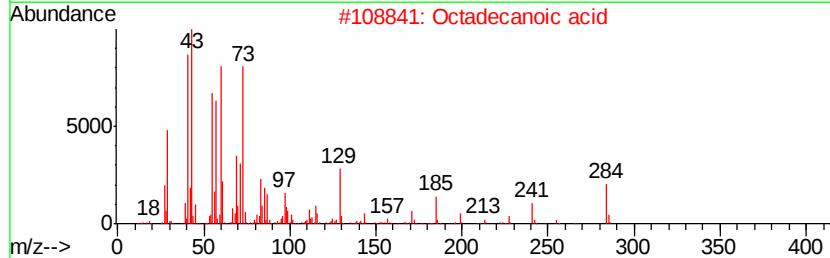
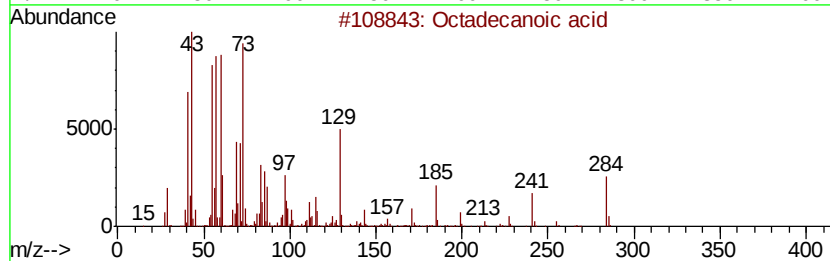
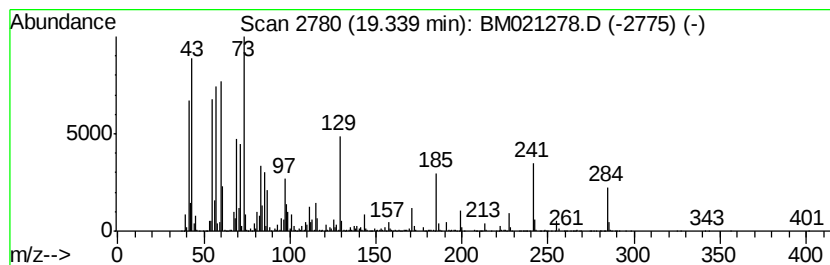
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Quant Title : SVOA CALIBRATION

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

Peak Number 31 Octadecanoic acid Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.34	17.39 ng/ul	5742540	Chrysene-d12	21.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		Octadecanoic acid	284	C18H36O2	000057-11-4	98
3		Octadecanoic acid	284	C18H36O2	000057-11-4	94
4		Octadecanoic acid	284	C18H36O2	000057-11-4	93
5		Pentadecanoic acid	242	C15H30O2	001002-84-2	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\
Data File : BM021278.D
Acq On : 30 Jun 2019 00:54
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Sample : K3590-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

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ClientSampleId :
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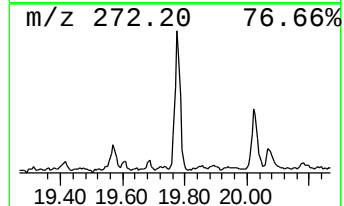
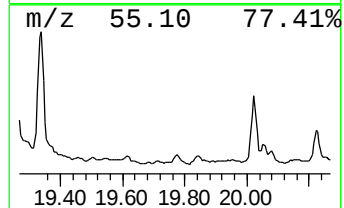
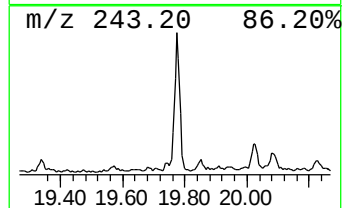
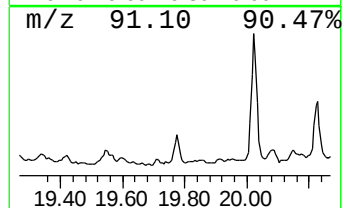
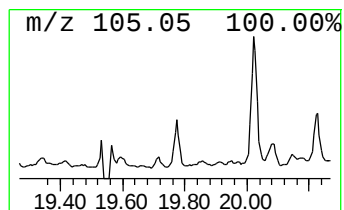
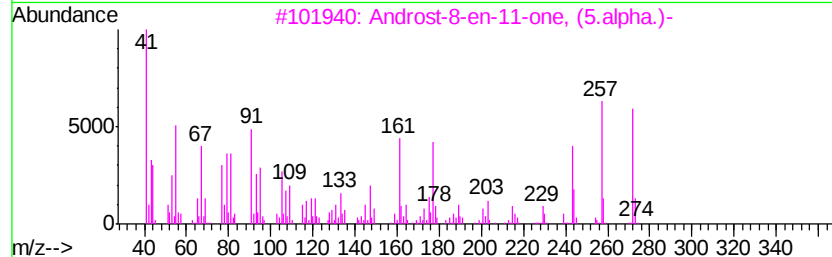
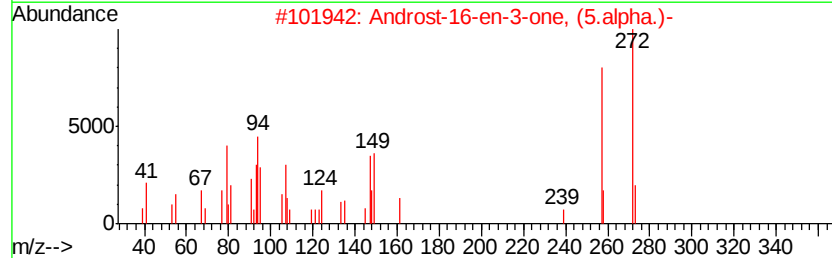
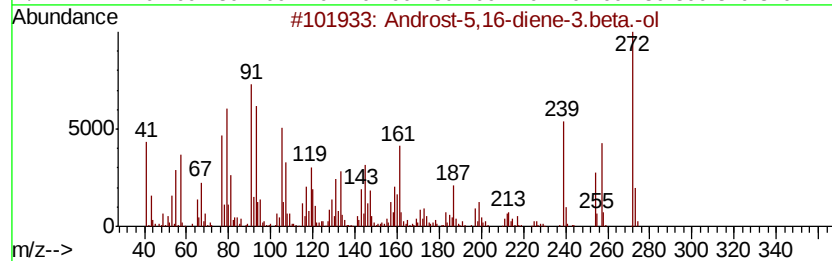
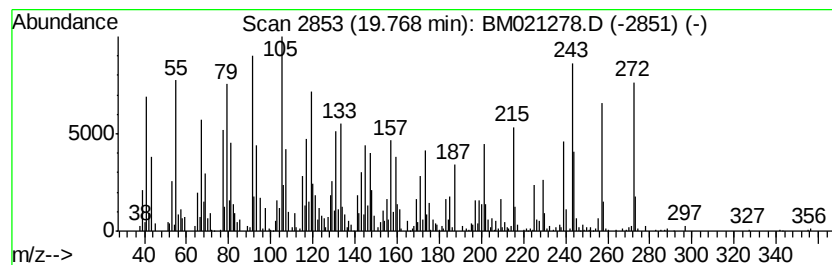
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Quant Title : SVOA CALIBRATION

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

Peak Number 32 Androst-5,16-diene-3.beta.-ol Concentration Rank 51

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.77	2.55 ng/ul	841610	Chrysene-d12	21.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Androst-5,16-diene-3.beta.-ol	272	C19H28O	001224-94-8	50
2		Androst-16-en-3-one, (5.alpha.)-	272	C19H28O	018339-16-7	49
3		Androst-8-en-11-one, (5.alpha.)-	272	C19H28O	054498-82-7	41
4		Benzo[fg]naphthacene, 9a,10,11,1...	272	C21H20	1000155-18-3	40
5		Benz[a]anthracen-12-ol, 7,12-dih...	272	C20H16O	077573-42-3	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\
Data File : BM021278.D
Acq On : 30 Jun 2019 00:54
Operator : HP/JU
Sample : K3590-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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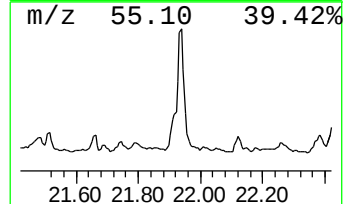
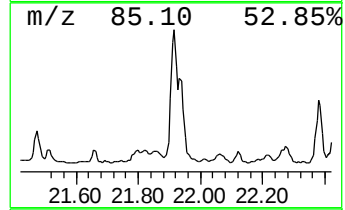
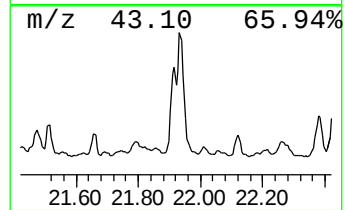
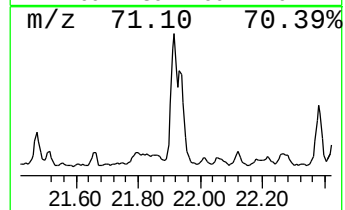
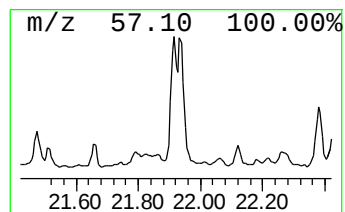
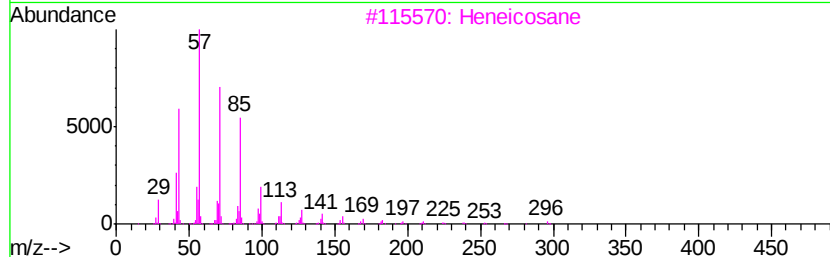
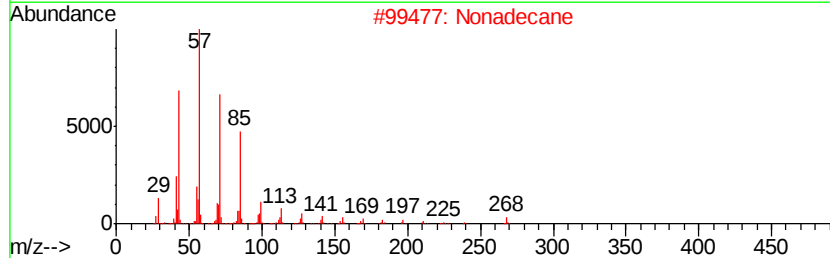
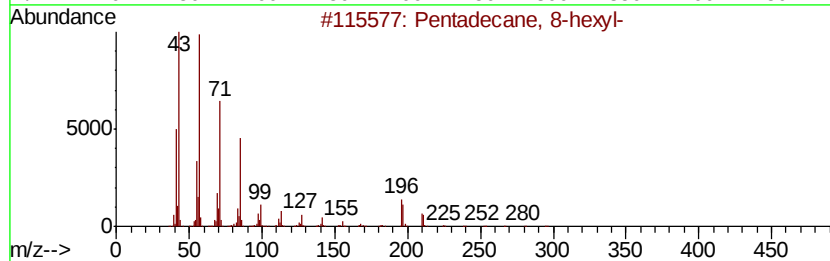
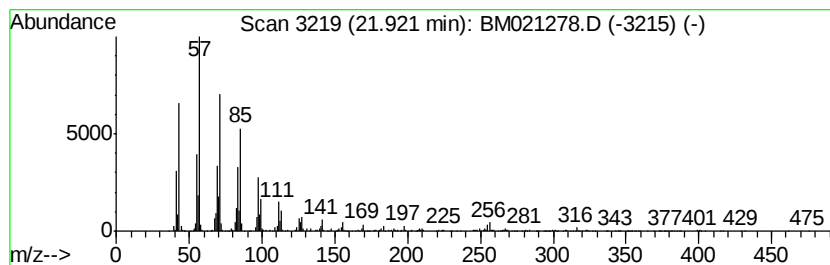
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Quant Title : SVOA CALIBRATION

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

Peak Number 40 (DEL) Alkane: Straight-Chai... Concentration Rank 47

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.92	3.03 ng/ul	1002210	Chrysene-d12	21.34

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	94
2			Nonadecane	268	C19H40	000629-92-5	93
3			Heneicosane	296	C21H44	000629-94-7	91
4			Tetratriacontane	479	C34H70	014167-59-0	91
5			Hexadecane	226	C16H34	000544-76-3	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\
 Data File : BM021278.D
 Acq On : 30 Jun 2019 00:54
 Operator : HP/JU
 Sample : K3590-01
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
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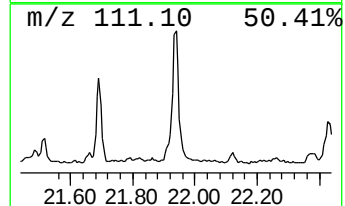
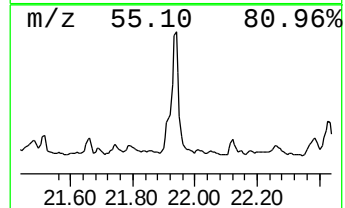
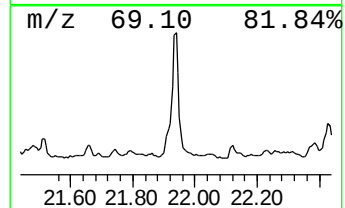
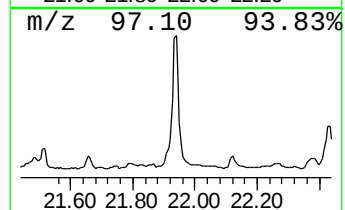
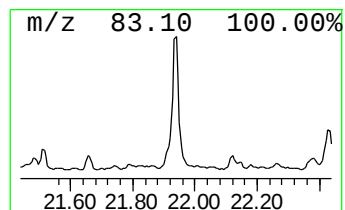
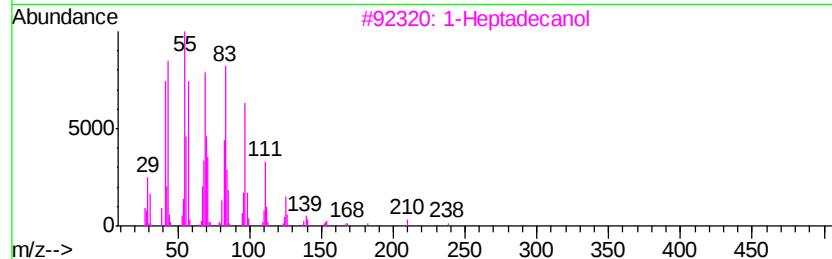
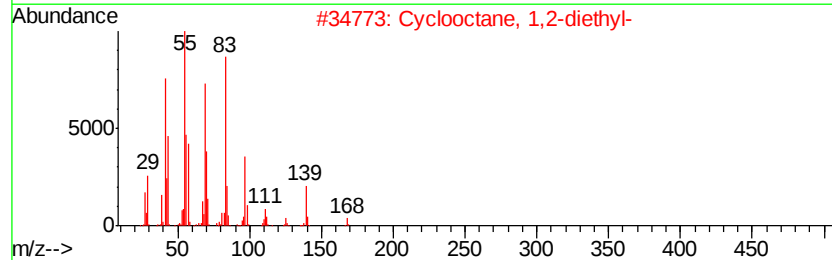
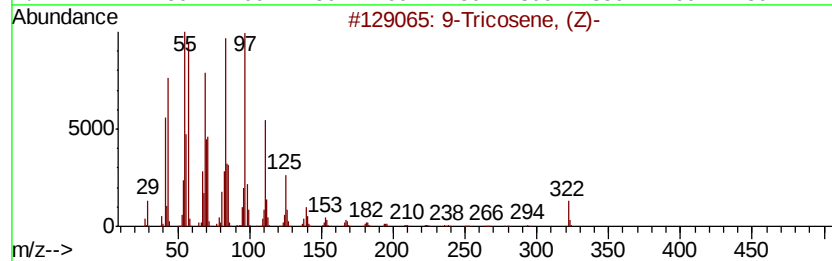
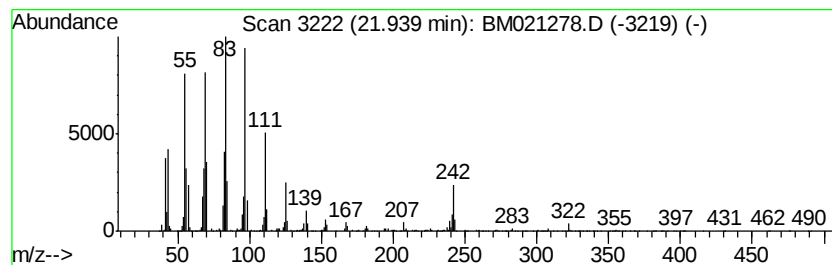
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 Quant Title : SVOA CALIBRATION

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

 Peak Number 41 (DEL) Alkane: Cyclic21.94 Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.94	11.02 ng/ul	3640570	Chrysene-d12	21.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Tricosene, (Z)-	322	C23H46	027519-02-4	78
2		Cyclooctane, 1,2-diethyl-	168	C12H24	023609-46-3	55
3		1-Heptadecanol	256	C17H36O	001454-85-9	49
4		1-Pentadecanol	228	C15H32O	000629-76-5	47
5		9-Tricosene, (Z)-	322	C23H46	027519-02-4	46



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\
Data File : BM021278.D
Acq On : 30 Jun 2019 00:54
Operator : HP/JU
Sample : K3590-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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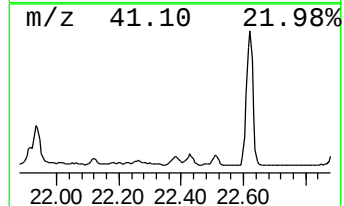
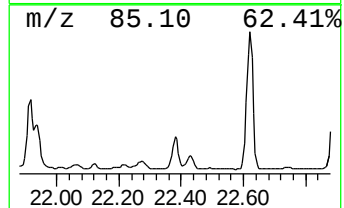
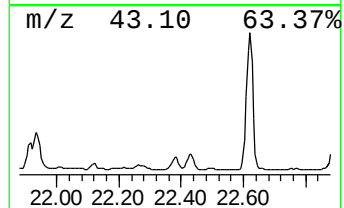
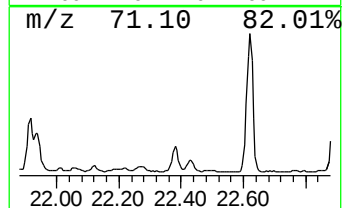
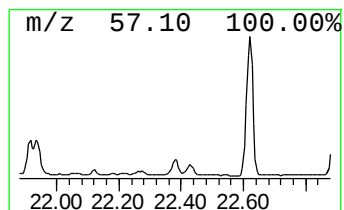
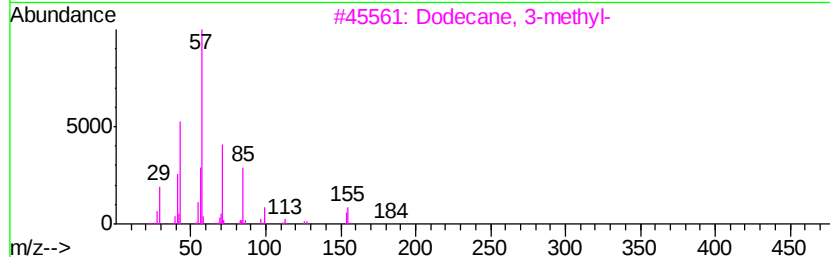
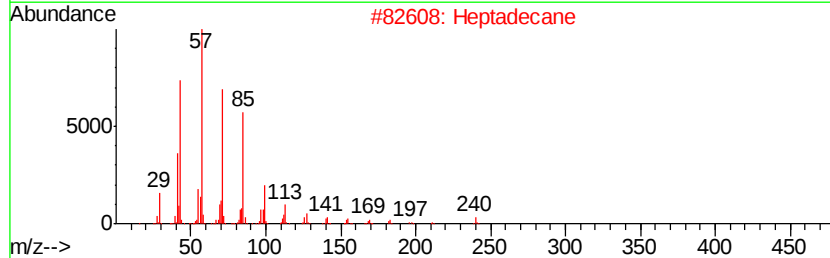
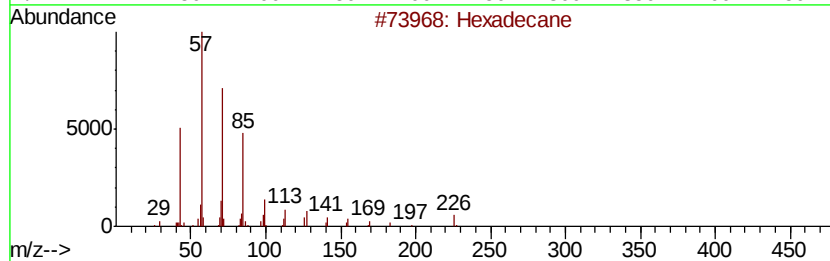
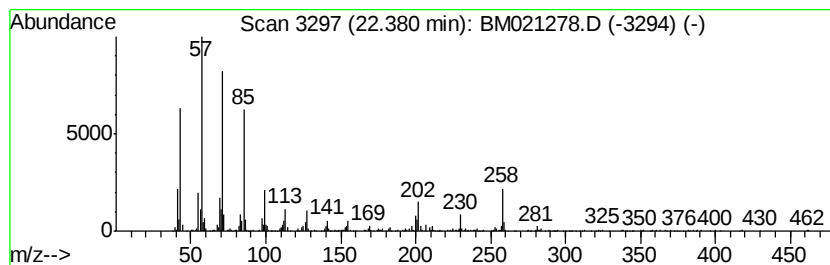
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Quant Title : SVOA CALIBRATION

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

Peak Number 42 (DEL) Alkane: Straight-Chai... Concentration Rank 52

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.38	2.38 ng/ul	786627	Chrysene-d12	21.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	81
2		Heptadecane	240	C17H36	000629-78-7	72
3		Dodecane, 3-methyl-	184	C13H28	017312-57-1	72
4		Octadecane, 1-iodo-	380	C18H37I	000629-93-6	72
5		Heneicosane	296	C21H44	000629-94-7	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\
Data File : BM021278.D
Acq On : 30 Jun 2019 00:54
Operator : HP/JU
Sample : K3590-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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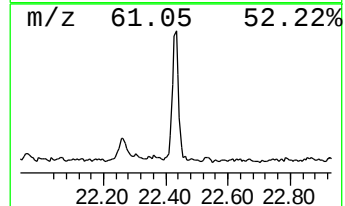
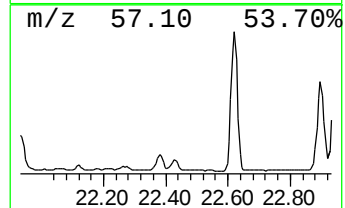
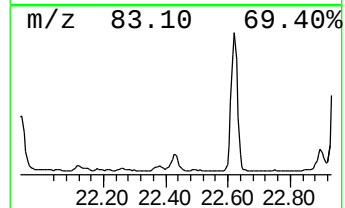
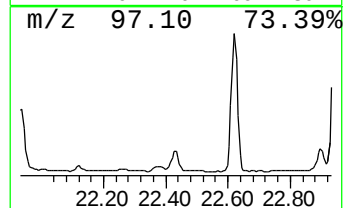
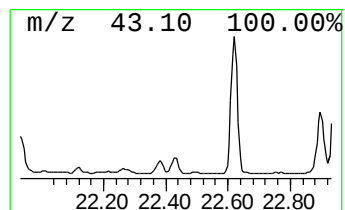
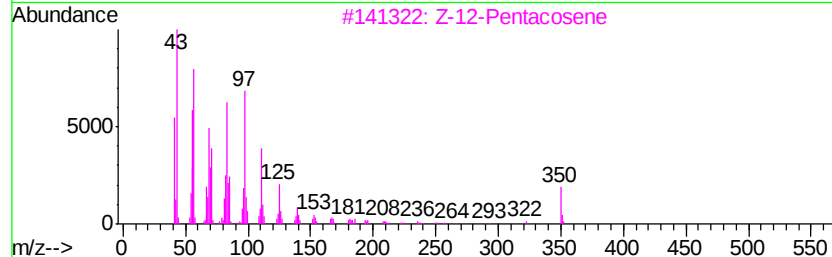
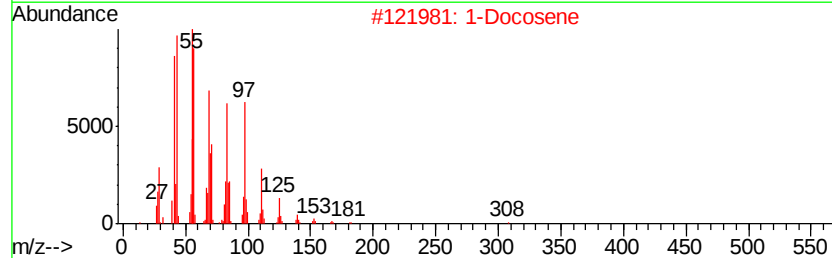
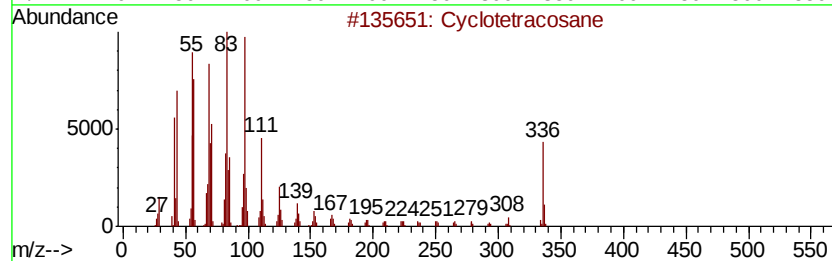
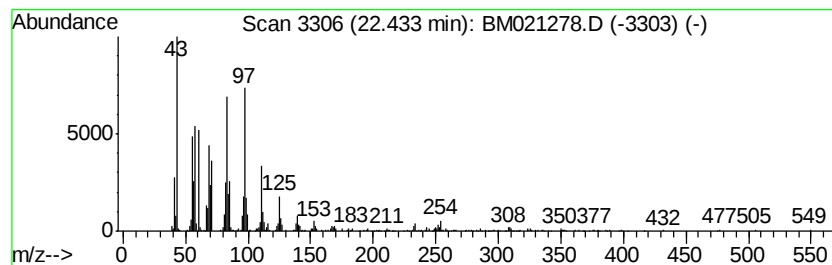
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Quant Title : SVOA CALIBRATION

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

Peak Number 43 (DEL) Alkane: Cyclic22.43 Concentration Rank 42

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.43	3.42 ng/ul	1129410	Chrysene-d12	21.34

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclotetracosane	336	C24H48	000297-03-0	98
2			1-Docosene	308	C22H44	001599-67-3	95
3			Z-12-Pentacosene	350	C25H50	1000131-09-4	93
4			1-Acetoxynonadecane	326	C21H42O2	001577-43-1	93
5			Acetic acid, octadecyl ester	312	C20H40O2	000822-23-1	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\
Data File : BM021278.D
Acq On : 30 Jun 2019 00:54
Operator : HP/JU
Sample : K3590-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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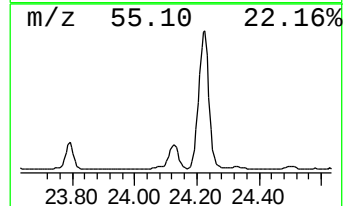
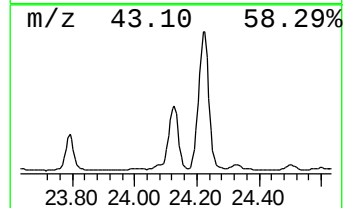
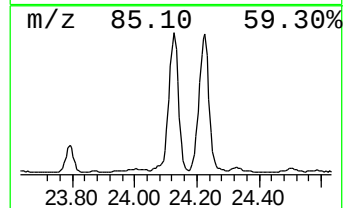
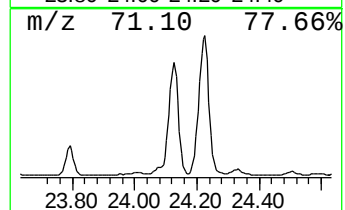
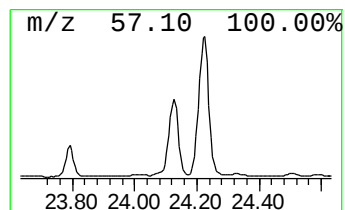
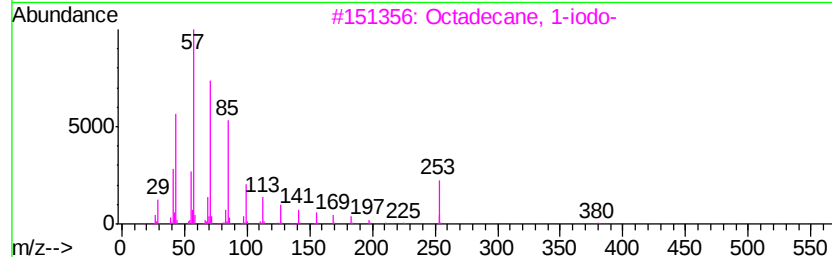
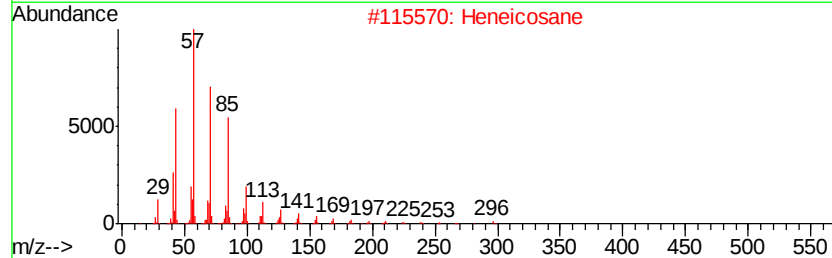
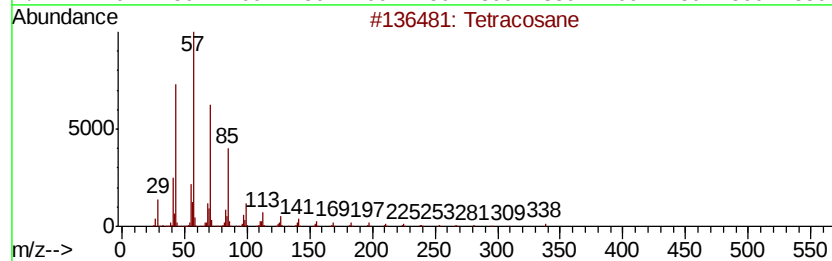
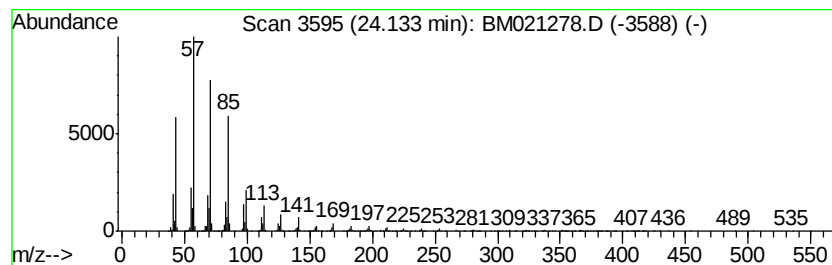
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Quant Title : SVOA CALIBRATION

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

Peak Number 47 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.13	47.12 ng/ul	8383000	Perylene-d12	23.61

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosane	338	C24H50	000646-31-1	97
2		Heneicosane	296	C21H44	000629-94-7	91
3		Octadecane, 1-iodo-	380	C18H37I	000629-93-6	91
4		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	91
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\
Data File : BM021278.D
Acq On : 30 Jun 2019 00:54
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Sample : K3590-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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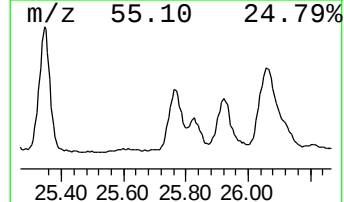
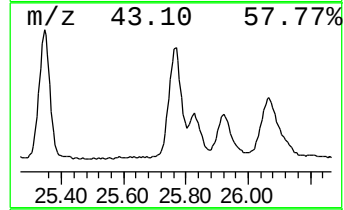
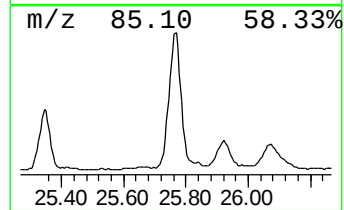
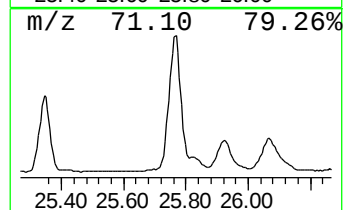
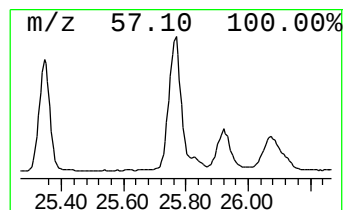
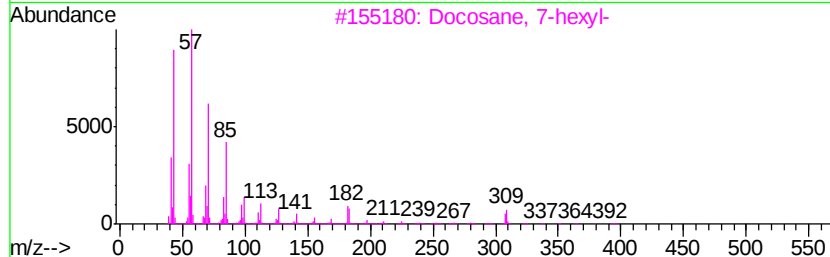
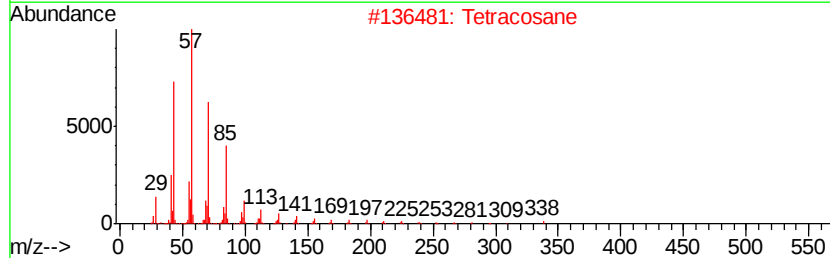
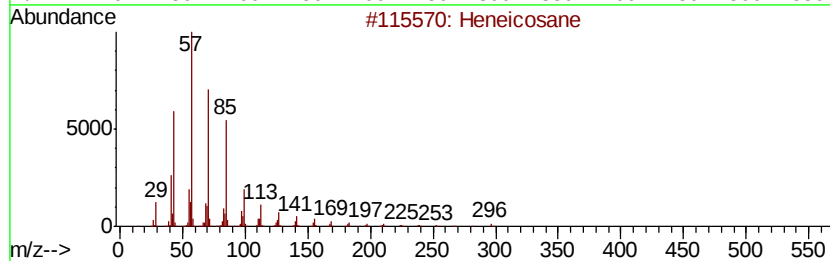
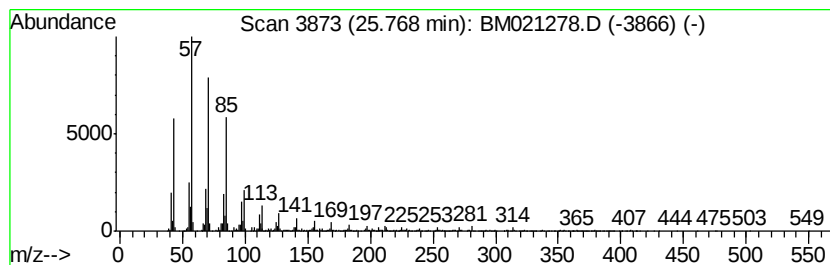
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Quant Title : SVOA CALIBRATION

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

Peak Number 51 (DEL) Alkane: Straight-Chai... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.77	27.46 ng/ul	4884740	Perylene-d12	23.61

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	97
2		Tetracosane	338	C24H50	000646-31-1	97
3		Docosane, 7-hexyl-	394	C28H58	055373-86-9	94
4		Tetratriacontane	479	C34H70	014167-59-0	94
5		triacontane	422	C30H62	000638-68-6	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM062919\
 Data File : BM021278.D
 Acq On : 30 Jun 2019 00:54
 Operator : HP/JU
 Sample : K3590-01
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C5BA0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.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	33.5	ng/ul	3117240	1	7.76	1863820	20.0
Cyclotrisiloxane,...	4.40	5.4	ng/ul	500264	1	7.76	1863820	20.0
Bicyclo[3.1.0]hex...	6.27	4.0	ng/ul	368440	1	7.76	1863820	20.0
1R-.alpha.-Pinene	6.43	24.6	ng/ul	2292210	1	7.76	1863820	20.0
Hexanoic acid	6.82	6.1	ng/ul	564510	1	7.76	1863820	20.0
Cyclotetrasiloxan...	6.86	4.2	ng/ul	386384	1	7.76	1863820	20.0
.beta.-Pinene	7.19	12.8	ng/ul	1190900	1	7.76	1863820	20.0
Cyclopentasiloxan...	9.28	2.3	ng/ul	318881	2	10.55	2735610	20.0
unknown-01	9.86	5.5	ng/ul	746266	2	10.55	2735610	20.0
Bicyclo[4.4.0]dec...	13.26	2.1	ng/ul	412280	3	14.40	3842380	20.0
Caryophyllene	13.71	3.0	ng/ul	582729	3	14.40	3842380	20.0
Naphthalene, 1,2,...	14.46	4.2	ng/ul	810262	3	14.40	3842380	20.0
Dodecanoic acid	14.82	3.4	ng/ul	645037	3	14.40	3842380	20.0
(-)-Spathulenol	15.31	3.6	ng/ul	699596	3	14.40	3842380	20.0
.alpha.-Cadinol	15.96	4.9	ng/ul	1056270	4	17.15	4316670	20.0
Tetradecanoic acid	16.56	6.5	ng/ul	1397920	4	17.15	4316670	20.0
7-Acetyl-2-hydrox...	16.65	3.2	ng/ul	700014	4	17.15	4316670	20.0
2-Propanol, 1-chl...	16.90	3.4	ng/ul	740025	4	17.15	4316670	20.0
unknown-02	16.97	2.9	ng/ul	615160	4	17.15	4316670	20.0
Pentadecanoic acid	17.05	7.7	ng/ul	1658740	4	17.15	4316670	20.0
(DEL) Alkane: Cyc...	17.53	2.4	ng/ul	510592	4	17.15	4316670	20.0
9-Hexadecenoic acid	17.94	29.5	ng/ul	6363750	4	17.15	4316670	20.0
Hexadecenoic acid...	18.00	34.1	ng/ul	7366070	4	17.15	4316670	20.0
n-Hexadecanoic acid	18.08	114.4	ng/ul	24686800	4	17.15	4316670	20.0
unknown-03	18.32	11.8	ng/ul	2540470	4	17.15	4316670	20.0
Heptadecanoic acid	18.48	2.3	ng/ul	501103	4	17.15	4316670	20.0
1H-Naphtho[2,1-b]...	18.66	7.3	ng/ul	1584020	4	17.15	4316670	20.0
unknown-04	18.99	3.6	ng/ul	785704	4	17.15	4316670	20.0
Phytol	19.06	21.7	ng/ul	4675430	4	17.15	4316670	20.0
Octadecanoic acid	19.34	17.4	ng/ul	5742540	5	21.34	6604780	20.0
Androst-5,16-dien...	19.77	2.5	ng/ul	841610	5	21.34	6604780	20.0
(DEL) Alkane: Str...	21.92	3.0	ng/ul	1002210	5	21.34	6604780	20.0
(DEL) Alkane: Cyc...	21.94	11.0	ng/ul	3640570	5	21.34	6604780	20.0
(DEL) Alkane: Str...	22.38	2.4	ng/ul	786627	5	21.34	6604780	20.0
(DEL) Alkane: Cyc...	22.43	3.4	ng/ul	1129410	5	21.34	6604780	20.0
(DEL) Alkane: Str...	24.13	47.1	ng/ul	8383000	6	23.61	3557800	20.0
(DEL) Alkane: Str...	25.77	27.5	ng/ul	4884740	6	23.61	3557800	20.0