

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100319\
 Data File : BM022929.D
 Acq On : 03 Oct 2019 20:25
 Operator : JU
 Sample : K4983-18
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BFDL2

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-BM092719MA.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	3.158	27	29	38	rVB	31421	48873	0.79%	0.068%
2	3.269	44	48	49	rBV	75108	84745	1.37%	0.118%
3	3.299	49	53	68	rVB	1795301	2147088	34.70%	3.001%
4	3.663	111	115	121	rBV	55985	78036	1.26%	0.109%
5	3.728	121	126	131	rBV	168624	211385	3.42%	0.295%
6	3.893	150	154	165	rVV2	21523	43098	0.70%	0.060%
7	3.981	165	169	180	rVB2	60345	128091	2.07%	0.179%
8	4.146	193	197	204	rVB	27706	41346	0.67%	0.058%
9	4.263	212	217	225	rBV	364046	460566	7.44%	0.644%
10	4.334	225	229	242	rVB2	47968	78470	1.27%	0.110%
11	4.440	242	247	255	rBV	550510	704293	11.38%	0.984%
12	4.510	255	259	270	rVB	30712	52513	0.85%	0.073%
13	4.893	316	324	335	rBV	631552	829891	13.41%	1.160%
14	5.269	381	388	392	rBV	36932	66191	1.07%	0.093%
15	5.316	392	396	402	rVV	32698	50462	0.82%	0.071%
16	5.446	413	418	428	rVB	144846	280197	4.53%	0.392%
17	5.757	466	471	476	rBV	92476	128505	2.08%	0.180%
18	5.816	476	481	488	rVB	114098	163496	2.64%	0.229%
19	6.816	639	651	658	rBV2	56813	147782	2.39%	0.207%
20	6.893	658	664	669	rVV	63339	99955	1.62%	0.140%
21	6.963	669	676	684	rVV2	273789	589332	9.52%	0.824%
22	7.022	684	686	696	rVB	52369	70890	1.15%	0.099%
23	7.128	696	704	711	rBV	623896	990698	16.01%	1.385%
24	7.187	711	714	724	rVB	63808	99927	1.62%	0.140%
25	7.328	731	738	747	rVB	770219	1203332	19.45%	1.682%
26	7.446	752	758	767	rVB	224980	346803	5.61%	0.485%
27	7.793	811	817	826	rBV	880371	1384092	22.37%	1.935%
28	7.863	826	829	832	rVV	30508	45683	0.74%	0.064%
29	7.910	832	837	848	rVB	106634	182088	2.94%	0.255%
30	8.169	872	881	887	rVB2	87320	175696	2.84%	0.246%
31	8.351	906	912	913	rVV3	26957	42894	0.69%	0.060%
32	8.381	913	917	922	rVV	49274	84760	1.37%	0.118%
33	8.434	922	926	931	rVV2	38149	64120	1.04%	0.090%
34	8.498	931	937	948	rVV	682903	1099996	17.78%	1.537%

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35	8.604	950	955	963	rVB2	18275	41398	0.67%	0.058%
36	8.898	999	1005	1008	rBV	67528	109180	1.76%	0.153%
37	8.945	1008	1013	1026	rVB	803554	1361764	22.01%	1.903%
38	9.422	1090	1094	1099	rVB	47810	70201	1.13%	0.098%
39	9.481	1099	1104	1115	rVB	68219	112627	1.82%	0.157%
40	9.669	1129	1136	1144	rBV	672577	1122239	18.14%	1.569%
41	9.839	1161	1165	1172	rVB	42081	67267	1.09%	0.094%
42	9.910	1172	1177	1183	rBV	29188	45098	0.73%	0.063%
43	9.992	1183	1191	1199	rVB3	119772	245349	3.97%	0.343%
44	10.210	1220	1228	1237	rVB	1102243	1840877	29.75%	2.573%
45	10.587	1282	1292	1297	rBV	1251861	2153896	34.81%	3.011%
46	10.634	1297	1300	1307	rVB	254256	400712	6.48%	0.560%
47	10.716	1307	1314	1326	rBV	939461	1643053	26.56%	2.297%
48	11.139	1377	1386	1390	rBV7	20550	47942	0.77%	0.067%
49	11.298	1408	1413	1417	rVB	27154	40685	0.66%	0.057%
50	11.781	1490	1495	1503	rVB4	30594	53056	0.86%	0.074%
51	11.928	1510	1520	1525	rBV2	87392	157407	2.54%	0.220%
52	12.110	1545	1551	1554	rBV	32561	60684	0.98%	0.085%
53	12.251	1569	1575	1586	rVB	204352	352393	5.70%	0.493%
54	12.357	1589	1593	1603	rVB8	25478	52485	0.85%	0.073%
55	12.469	1603	1612	1618	rBV	213368	351691	5.68%	0.492%
56	12.533	1618	1623	1632	rVB10	20819	47823	0.77%	0.067%
57	13.139	1721	1726	1732	rBV6	17838	47566	0.77%	0.066%
58	13.275	1744	1749	1756	rVB2	80168	124759	2.02%	0.174%
59	13.339	1756	1760	1765	rBV4	21604	41310	0.67%	0.058%
60	13.486	1777	1785	1791	rVB2	52252	116475	1.88%	0.163%
61	13.610	1799	1806	1810	rVB5	49518	103981	1.68%	0.145%
62	13.657	1810	1814	1818	rBV5	29182	49540	0.80%	0.069%
63	13.751	1825	1830	1835	rBV2	83915	157206	2.54%	0.220%
64	13.839	1840	1845	1849	rVV	2210069	2993346	48.38%	4.184%
65	13.886	1849	1853	1865	rVB	933482	1273013	20.57%	1.779%
66	13.986	1866	1870	1873	rBV	38191	56040	0.91%	0.078%
67	14.022	1873	1876	1883	rVB6	33712	71072	1.15%	0.099%
68	14.122	1884	1893	1905	rVB	2397473	3552572	57.42%	4.965%
69	14.257	1908	1916	1919	rBV6	23137	62840	1.02%	0.088%
70	14.345	1926	1931	1936	rVB5	39908	71804	1.16%	0.100%
71	14.433	1938	1946	1953	rVV	1987959	3125893	50.52%	4.369%

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72	14.492	1953	1956	1963	rVB	102192	166626	2.69%	0.233%
73	14.633	1973	1980	1985	rBV	287836	483937	7.82%	0.676%
74	14.827	2009	2013	2015	rBV2	39314	60191	0.97%	0.084%
75	14.904	2022	2026	2030	rVB2	37281	54258	0.88%	0.076%
76	15.069	2045	2054	2058	rBV3	149831	263387	4.26%	0.368%
77	15.222	2073	2080	2090	rVV4	69889	212892	3.44%	0.298%
78	15.422	2108	2114	2120	rBV	3154337	4535590	73.30%	6.339%
79	15.469	2120	2122	2127	rVV2	92367	139878	2.26%	0.196%
80	15.539	2128	2134	2140	rVV	1020452	1474044	23.82%	2.060%
81	15.604	2142	2145	2148	rVV	37952	59262	0.96%	0.083%
82	15.657	2150	2154	2166	rVB3	72013	178458	2.88%	0.249%
83	15.933	2197	2201	2207	rVB2	75093	152400	2.46%	0.213%
84	16.098	2225	2229	2234	rBV	43976	66933	1.08%	0.094%
85	16.580	2307	2311	2315	rVB4	33954	52028	0.84%	0.073%
86	16.986	2377	2380	2386	rVB2	27131	44279	0.72%	0.062%
87	17.168	2405	2411	2416	rBV2	2341491	3366580	54.41%	4.706%
88	17.210	2416	2418	2422	rVV	102780	123311	1.99%	0.172%
89	17.268	2422	2428	2438	rVB2	3807985	5330311	86.15%	7.450%
90	17.539	2469	2474	2477	rBV2	105676	158363	2.56%	0.221%
91	17.792	2513	2517	2522	rVB2	41383	53927	0.87%	0.075%
92	18.074	2558	2565	2573	rBV	1130674	1610544	26.03%	2.251%
93	19.227	2758	2761	2765	rVB	32158	41189	0.67%	0.058%
94	19.351	2777	2782	2795	rBV	558764	855342	13.82%	1.196%
95	19.562	2812	2818	2823	rBV	4712251	6187290	100.00%	8.648%
96	19.604	2823	2825	2832	rVB	179211	209245	3.38%	0.292%
97	21.062	3069	3073	3083	rBV	539333	671346	10.85%	0.938%
98	21.351	3117	3122	3127	rBV	2355445	2959479	47.83%	4.137%
99	23.468	3475	3482	3497	rVB	2477432	4799736	77.57%	6.709%
100	23.609	3500	3506	3517	rVB	1474691	2783748	44.99%	3.891%

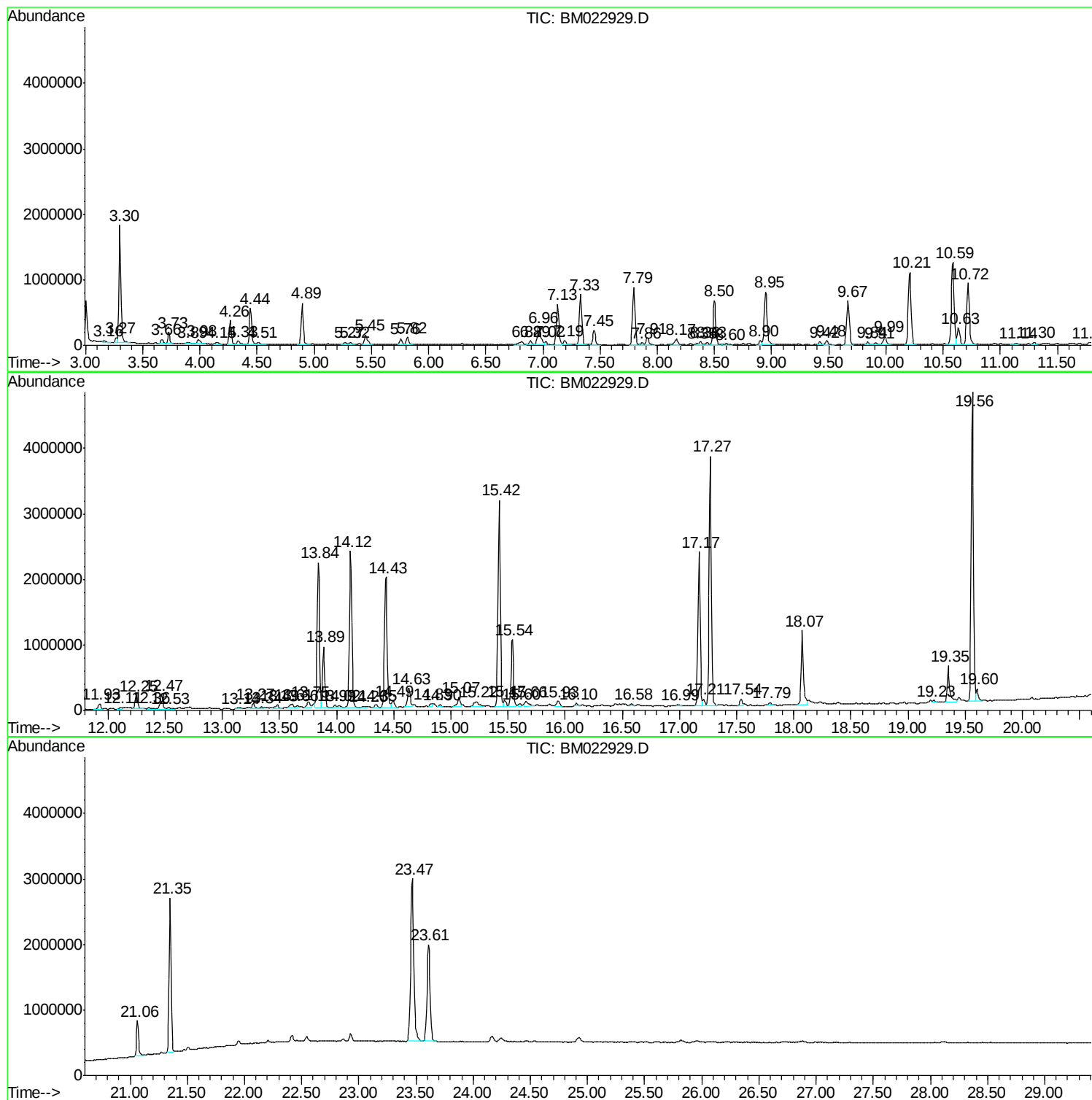
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092719MA.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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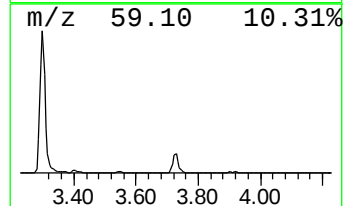
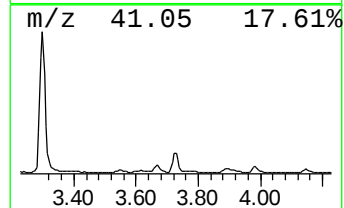
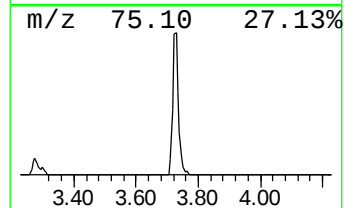
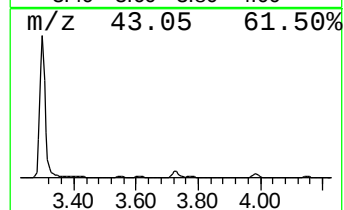
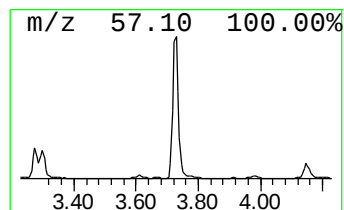
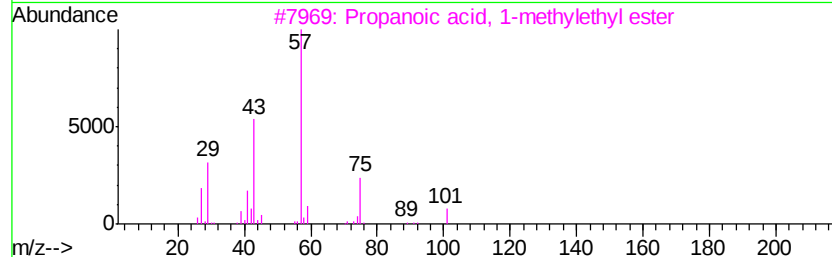
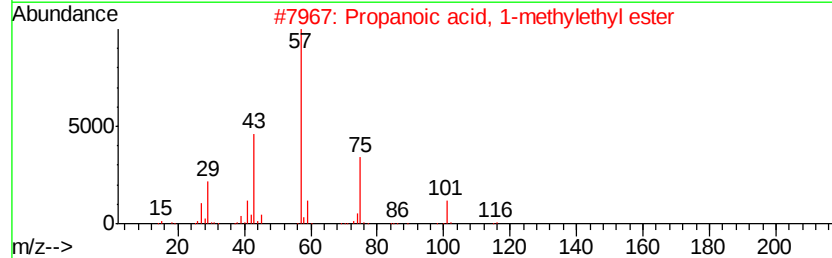
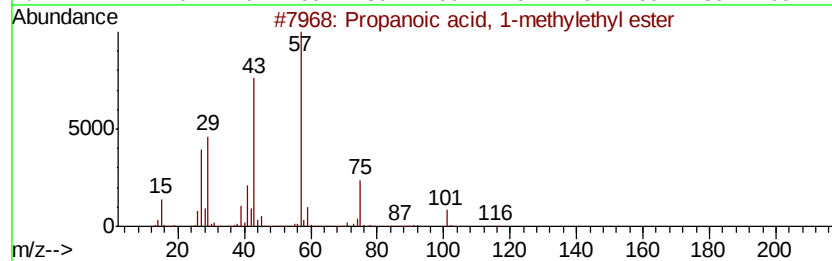
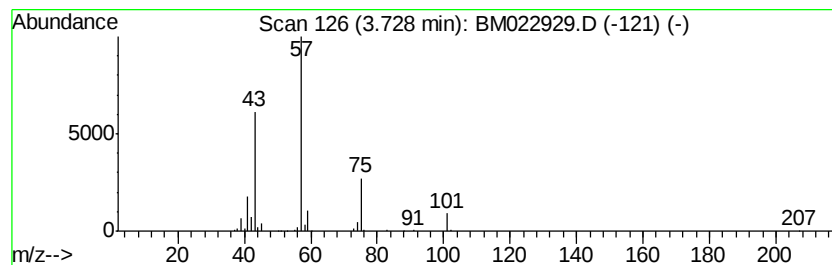
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092719MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Propanoic acid, 1-methyleth... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.73	3.05 ng/ul	211385	1,4-Dichlorobenzene-d4	7.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propanoic acid, 1-methylethyl ester	116	C6H12O2	000637-78-5	83
2		Propanoic acid, 1-methylethyl ester	116	C6H12O2	000637-78-5	83
3		Propanoic acid, 1-methylethyl ester	116	C6H12O2	000637-78-5	83
4		Propanoic acid, 1-methylethyl ester	116	C6H12O2	000637-78-5	59
5		Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	42



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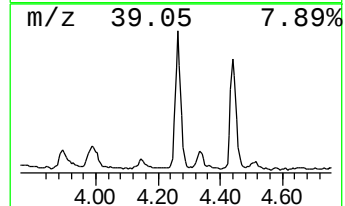
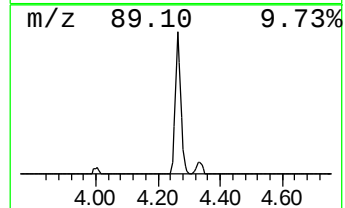
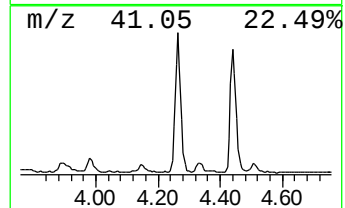
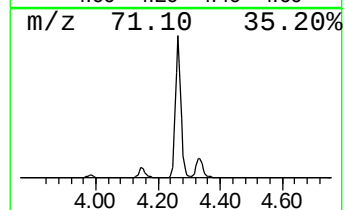
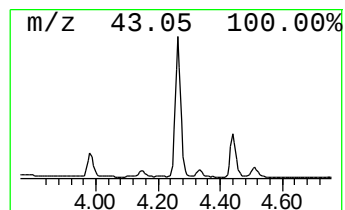
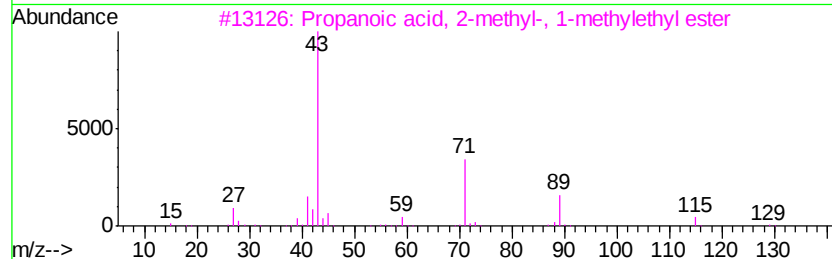
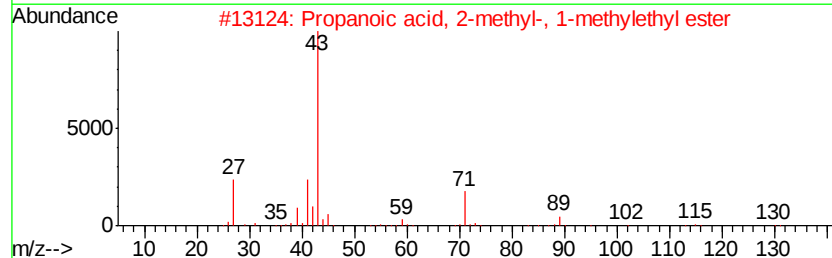
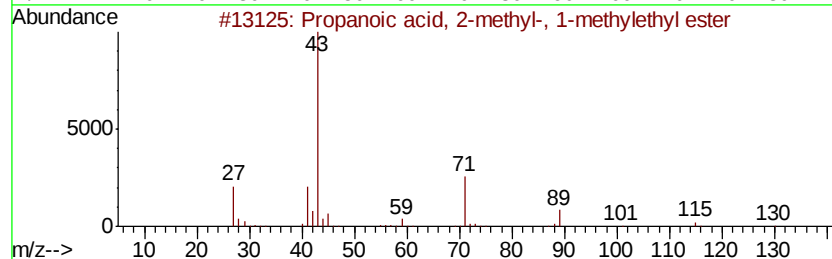
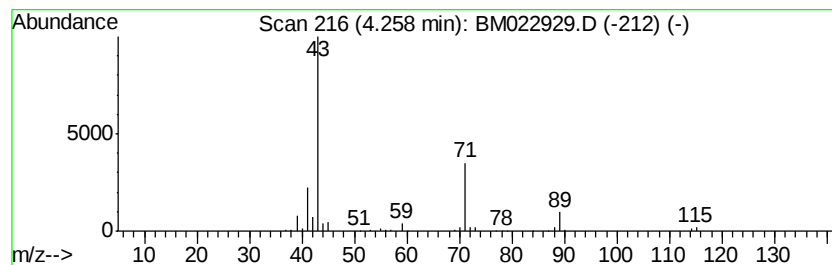
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Propanoic acid, 2-methyl-, ... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.26	6.66 ng/ul	460566	1,4-Dichlorobenzene-d4	7.79

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	78
2			Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	78
3			Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	64
4			Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	50
5			Propanoic acid, 2-methyl-, penty...	158	C9H18O2	002445-72-9	39



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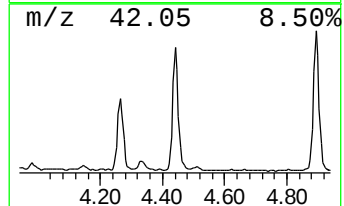
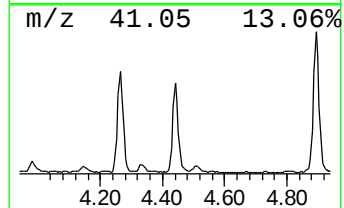
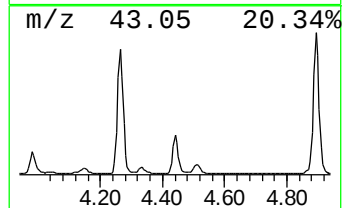
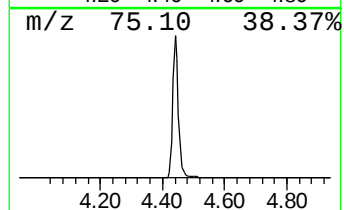
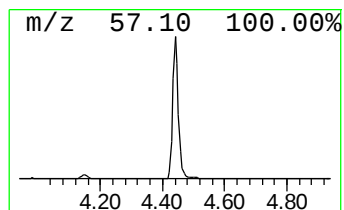
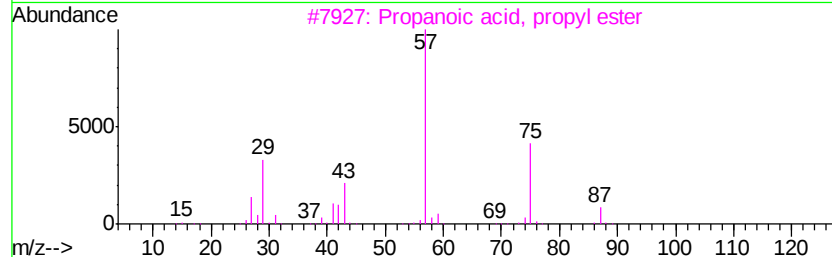
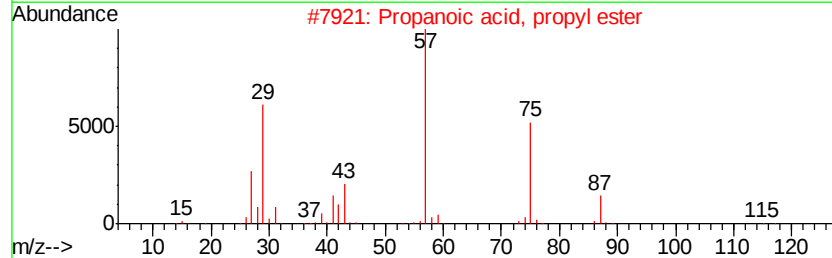
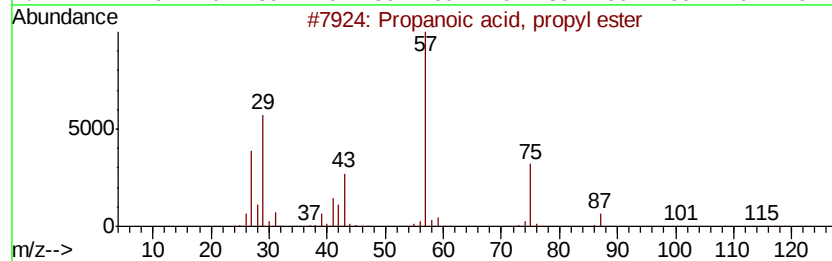
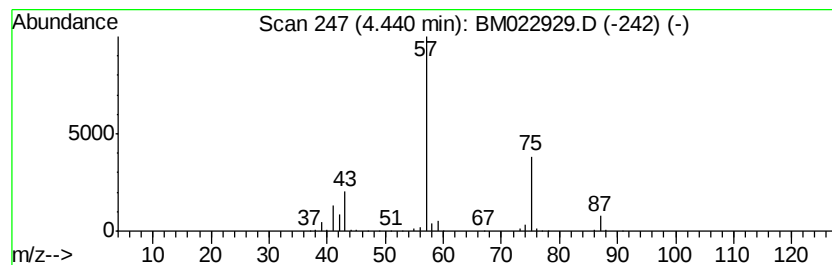
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TIC Integration Parameters: LSCINT.P

Peak Number 3 Propanoic acid, propyl ester Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.44	10.18 ng/ul	704293	1,4-Dichlorobenzene-d4	7.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	83
2		Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	83
3		Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	83
4		Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	47
5		Propanoic acid, butyl ester	130	C7H14O2	000590-01-2	39



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100319\
Data File : BM022929.D
Acq On : 03 Oct 2019 20:25
Operator : JU
Sample : K4983-18
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BFDL2

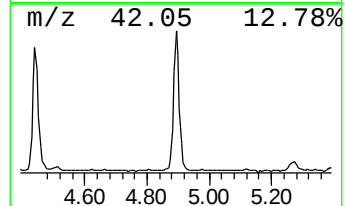
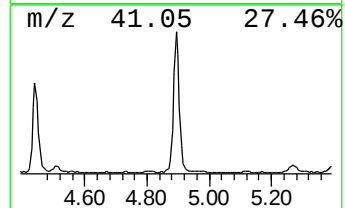
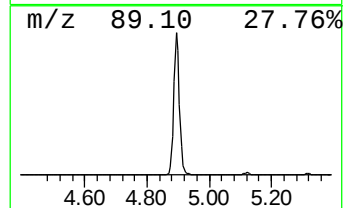
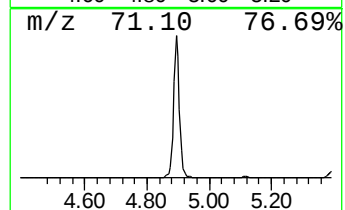
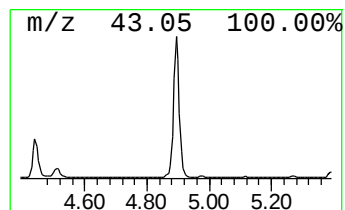
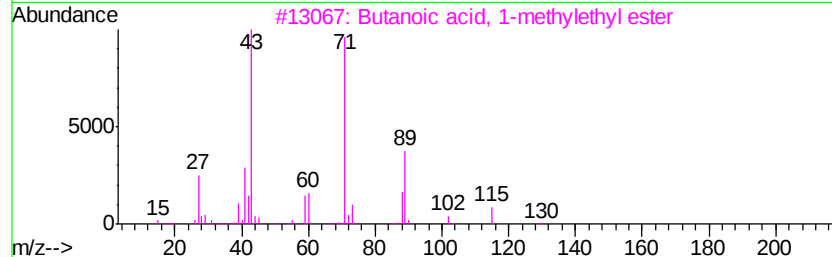
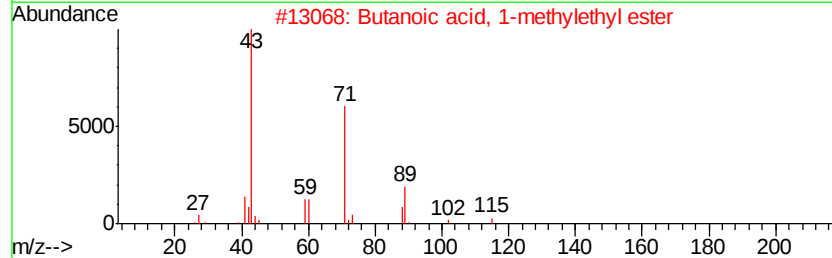
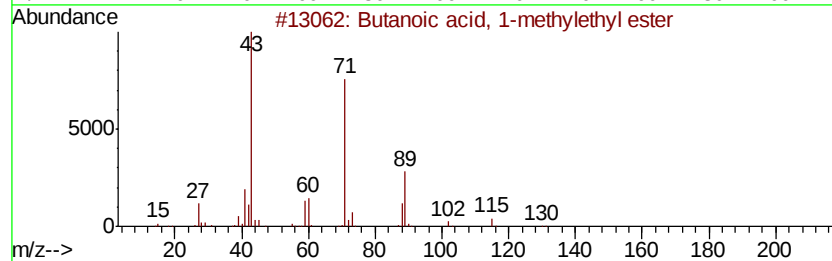
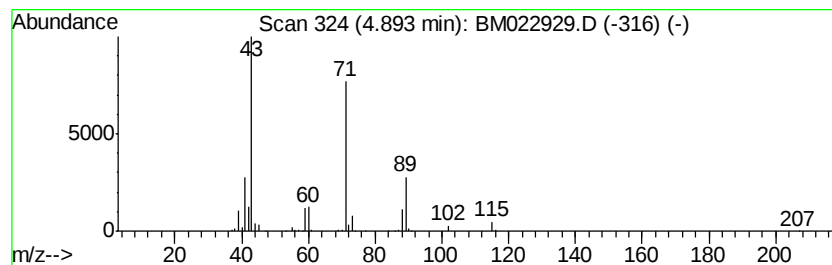
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Butanoic acid, 1-methylethy... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.89	11.99 ng/ul	829891	1,4-Dichlorobenzene-d4	7.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butanoic acid, 1-methylethyl ester	130	C7H14O2	000638-11-9	90
2		Butanoic acid, 1-methylethyl ester	130	C7H14O2	000638-11-9	90
3		Butanoic acid, 1-methylethyl ester	130	C7H14O2	000638-11-9	90
4		Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	78
5		Butanoic acid, propyl ester	130	C7H14O2	000105-66-8	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100319\
Data File : BM022929.D
Acq On : 03 Oct 2019 20:25
Operator : JU
Sample : K4983-18
Misc :
ALS Vial : 16 Sample Multiplier: 1

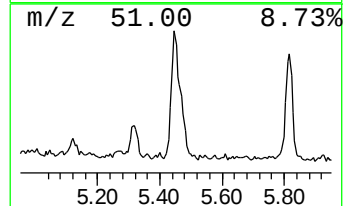
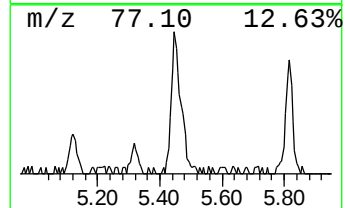
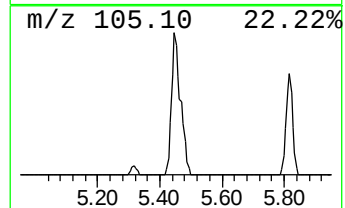
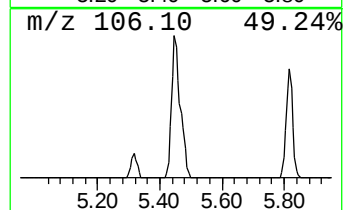
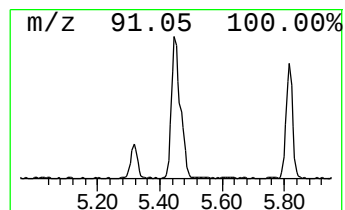
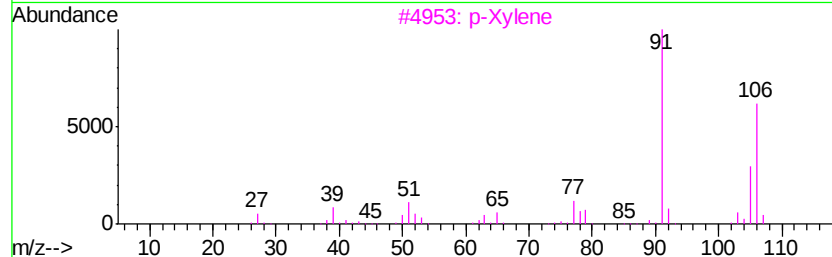
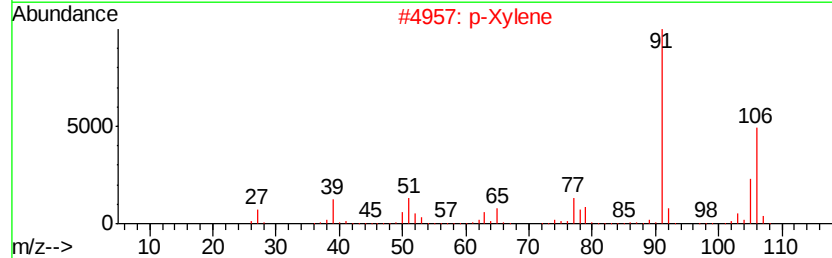
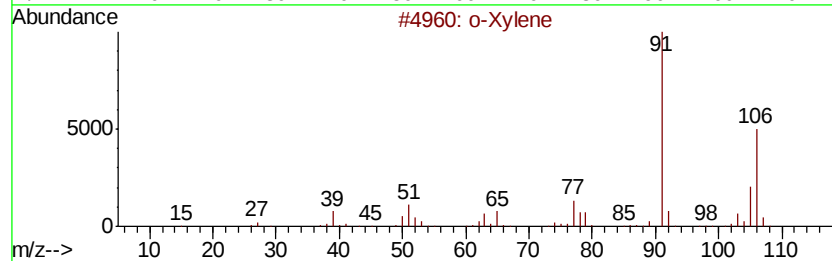
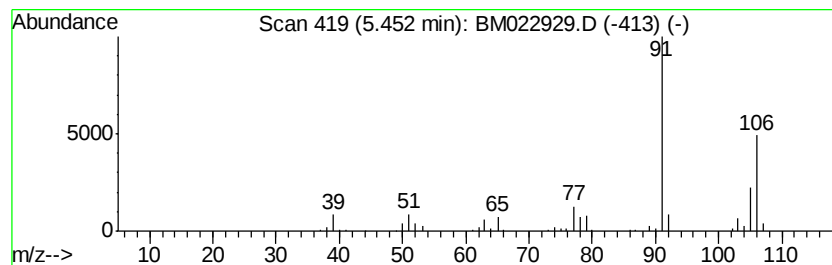
Instrument :
BNA_M
ClientSampleId :
BFDL2

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Quant Title : SVOA CALIBRATION

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

Peak Number 5 o-Xylene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
5.45	4.05 ng/ul	280197	1,4-Dichlorobenzene-d4	7.79	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Xylene	106	C8H10	000095-47-6	97
2	p-Xylene	106	C8H10	000106-42-3	97
3	p-Xylene	106	C8H10	000106-42-3	97
4	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
5	p-Xylene	106	C8H10	000106-42-3	95



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100319\
Data File : BM022929.D
Acq On : 03 Oct 2019 20:25
Operator : JU
Sample : K4983-18
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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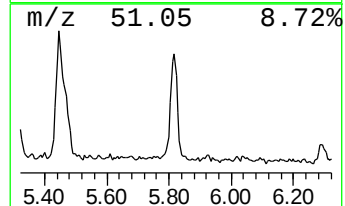
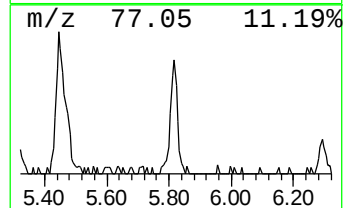
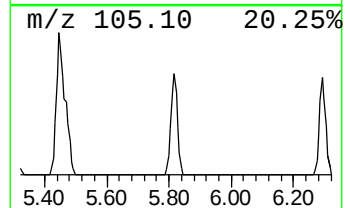
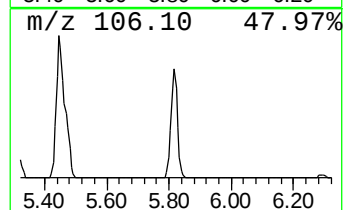
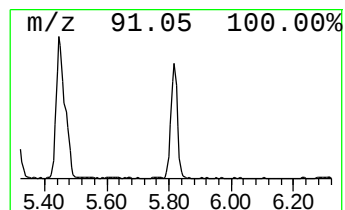
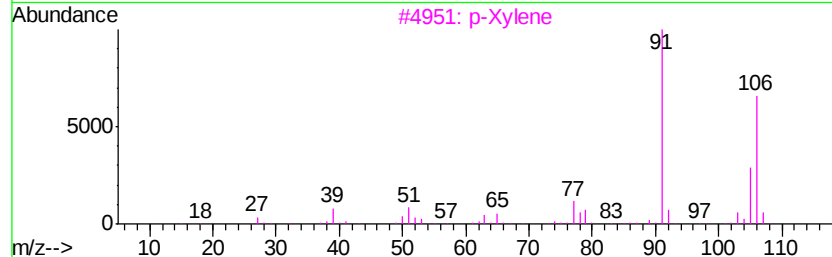
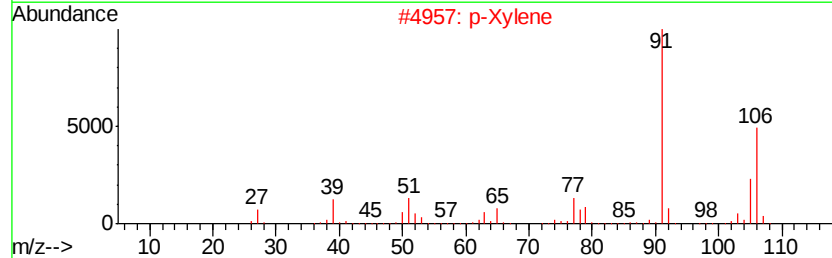
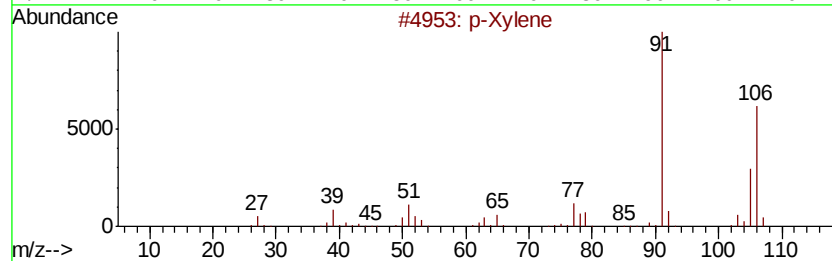
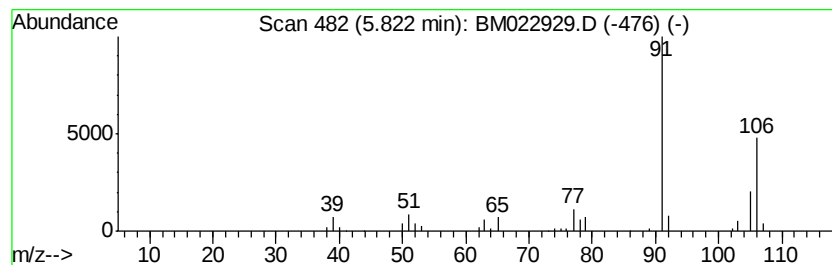
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 p-Xylene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.82	2.36 ng/ul	163496	1,4-Dichlorobenzene-d4	7.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	p-Xylene	106	C8H10	000106-42-3	95
2		p-Xylene	106	C8H10	000106-42-3	95
3		p-Xylene	106	C8H10	000106-42-3	94
4		o-Xylene	106	C8H10	000095-47-6	93
5		p-Xylene	106	C8H10	000106-42-3	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100319\
Data File : BM022929.D
Acq On : 03 Oct 2019 20:25
Operator : JU
Sample : K4983-18
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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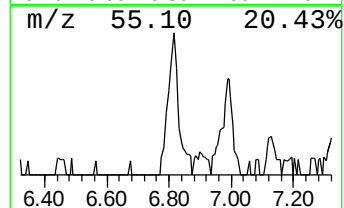
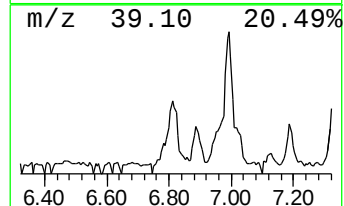
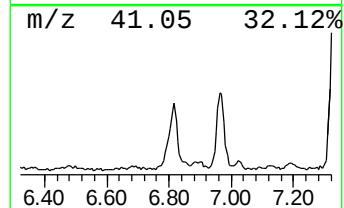
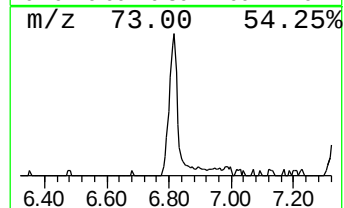
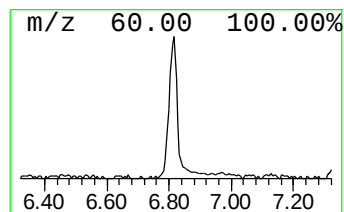
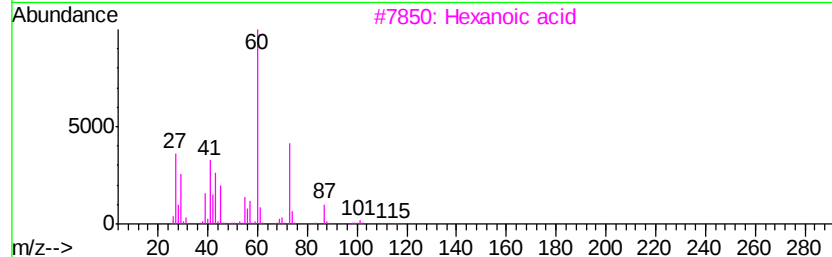
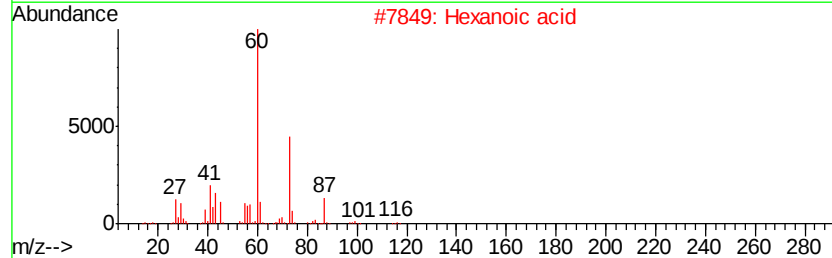
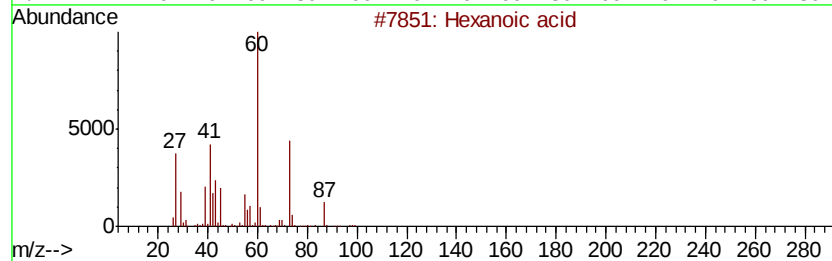
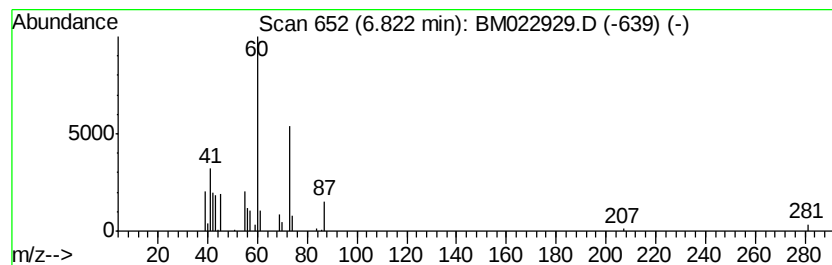
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Hexanoic acid Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.82	2.14 ng/ul	147782	1,4-Dichlorobenzene-d4	7.79

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	83
4		Propanedioic acid, propyl-	146	C6H10O4	000616-62-6	72
5		Heptanoic acid	130	C7H14O2	000111-14-8	59



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100319\
Data File : BM022929.D
Acq On : 03 Oct 2019 20:25
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Sample : K4983-18
Misc :
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Instrument :
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ClientSampleId :
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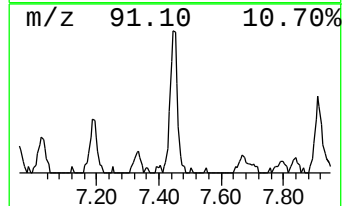
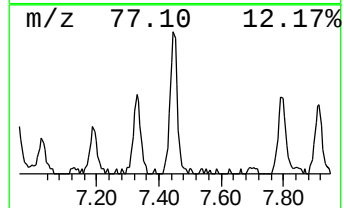
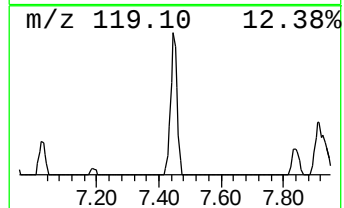
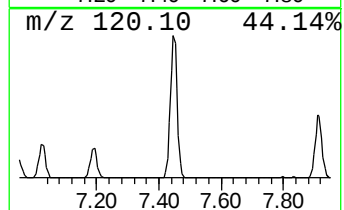
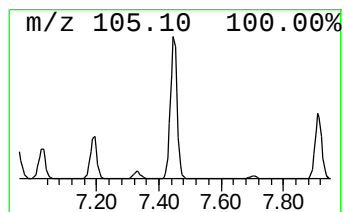
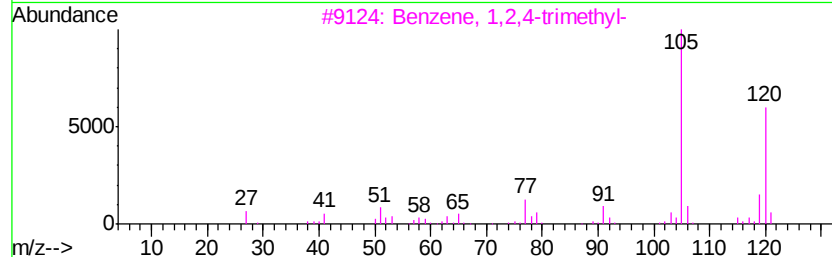
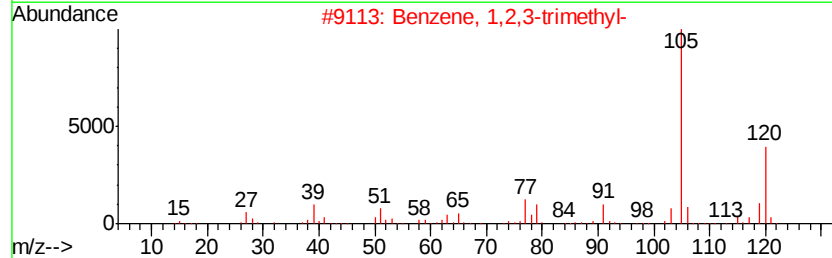
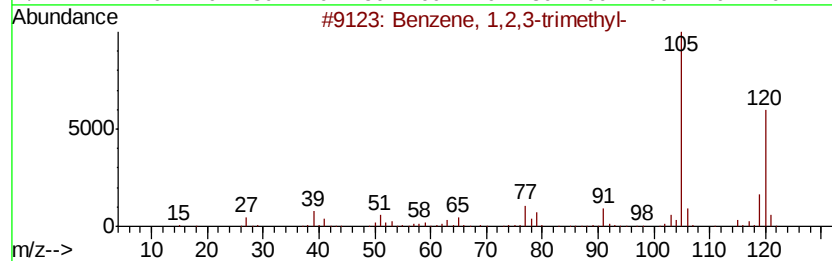
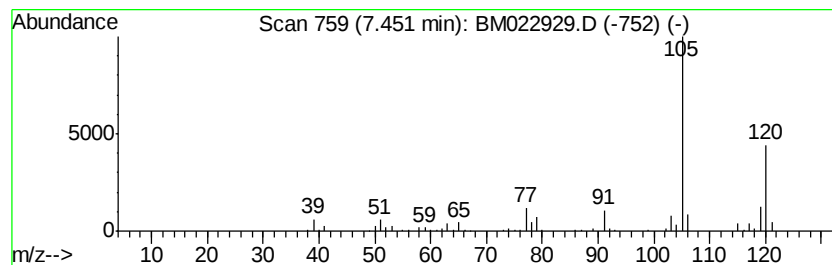
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Benzene, 1,2,3-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.45	5.01 ng/ul	346803	1,4-Dichlorobenzene-d4	7.79

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	95
4			Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	91
5			Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100319\
Data File : BM022929.D
Acq On : 03 Oct 2019 20:25
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Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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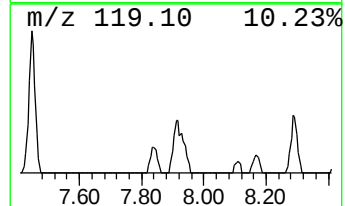
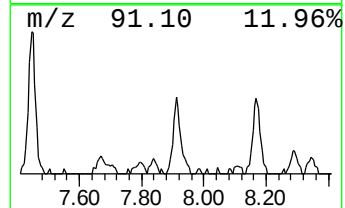
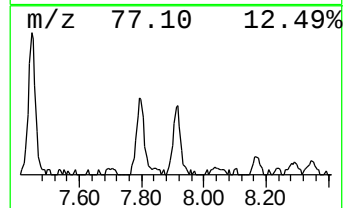
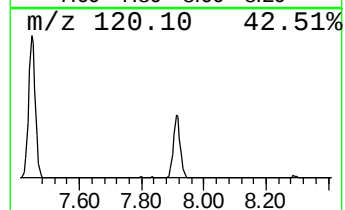
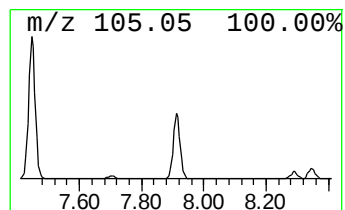
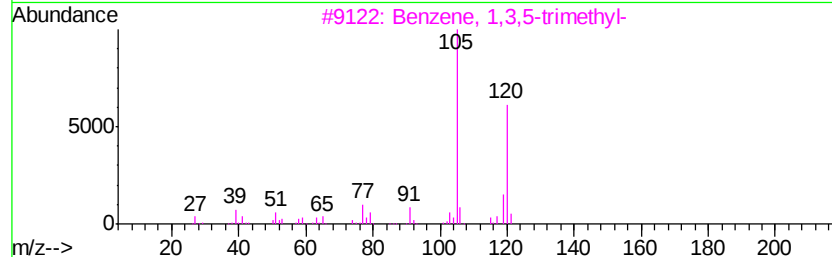
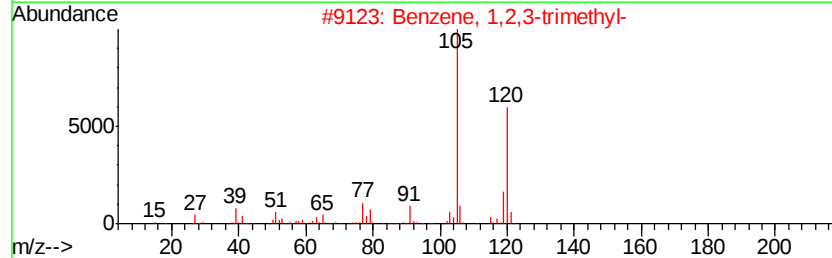
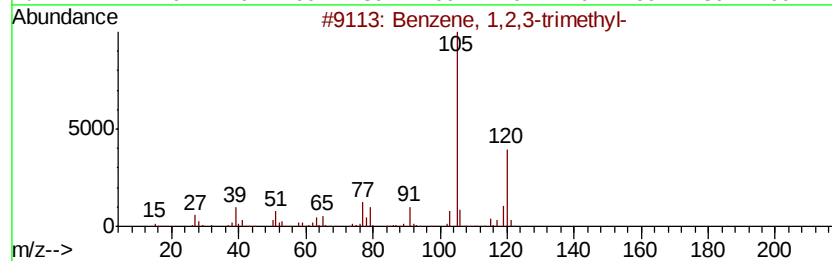
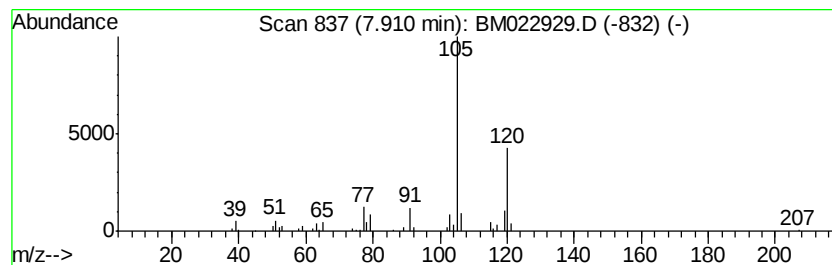
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Benzene, 1,3,5-trimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.91	2.63 ng/ul	182088	1,4-Dichlorobenzene-d4	7.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3		Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	94
4		Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	94
5		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100319\
Data File : BM022929.D
Acq On : 03 Oct 2019 20:25
Operator : JU
Sample : K4983-18
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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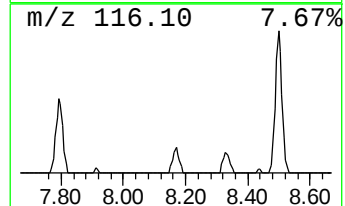
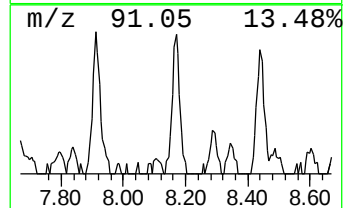
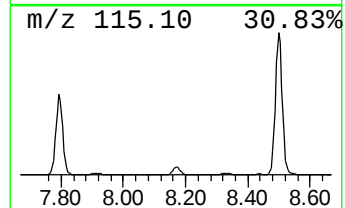
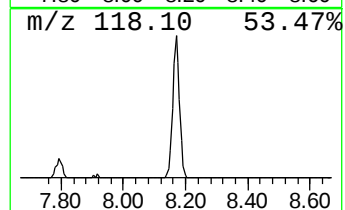
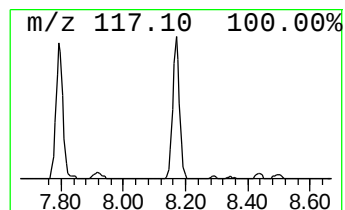
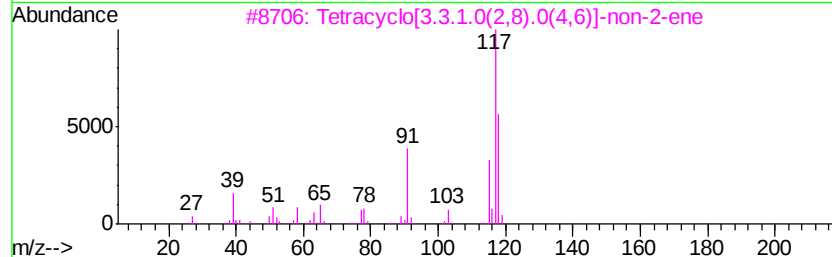
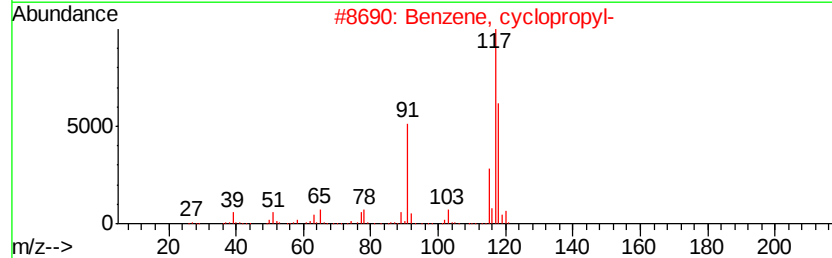
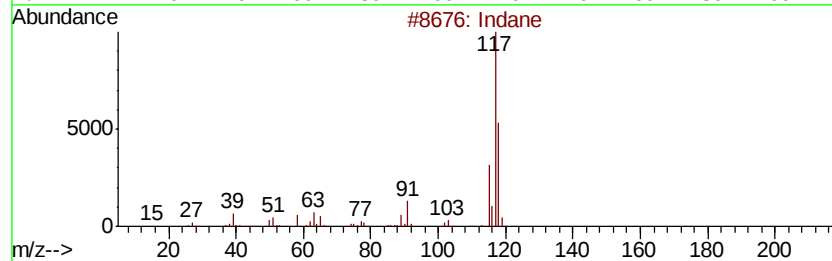
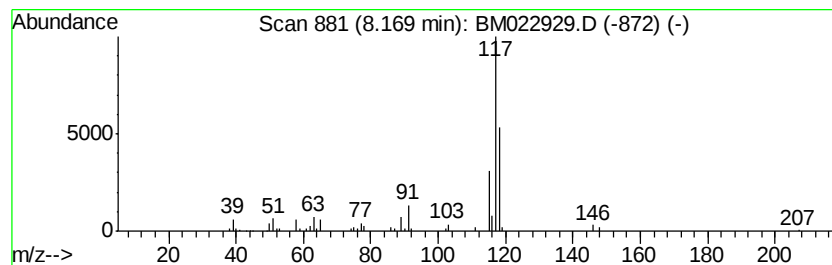
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Indane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.17	2.54 ng/ul	175696	1,4-Dichlorobenzene-d4	7.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indane	118	C9H10	000496-11-7	74
2		Benzene, cyclopropyl-	118	C9H10	000873-49-4	74
3		Tetracyclo[3.3.1.0(2,8).0(4,6)]-	118	C9H10	1000191-13-7	74
4		Benzene, cyclopropyl-	118	C9H10	000873-49-4	68
5		Indane	118	C9H10	000496-11-7	60



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100319\
Data File : BM022929.D
Acq On : 03 Oct 2019 20:25
Operator : JU
Sample : K4983-18
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BFDL2

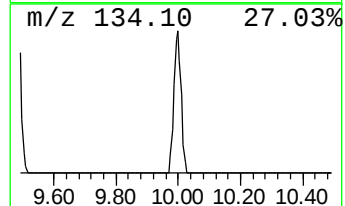
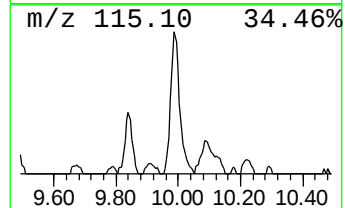
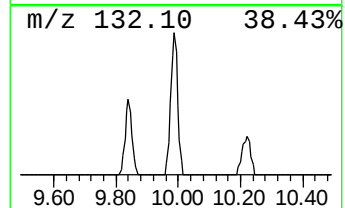
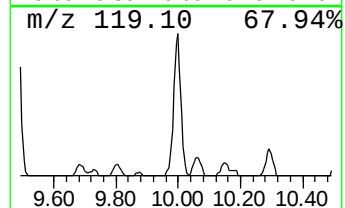
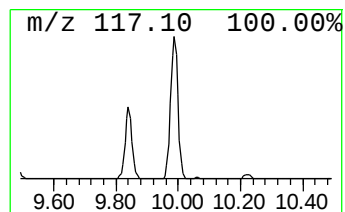
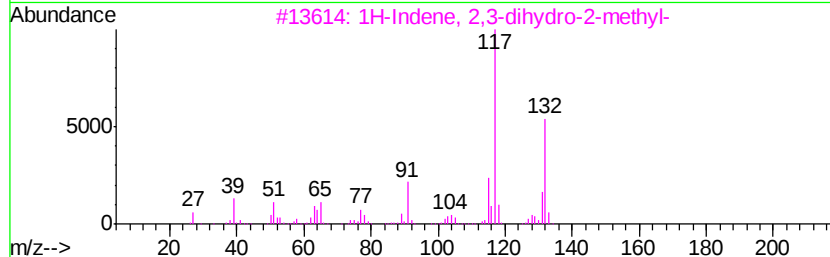
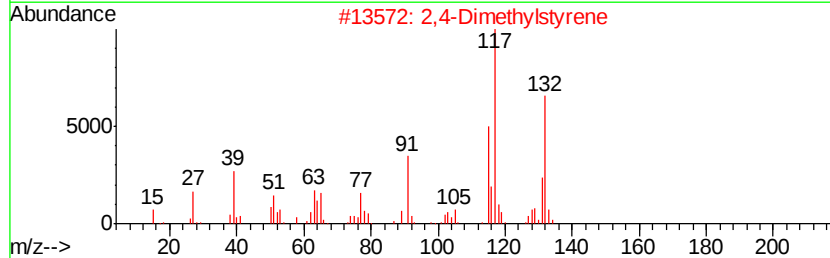
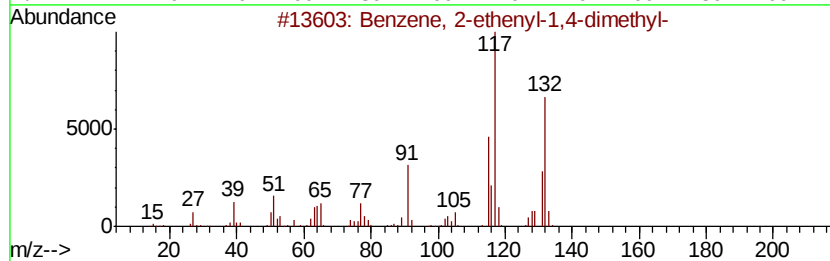
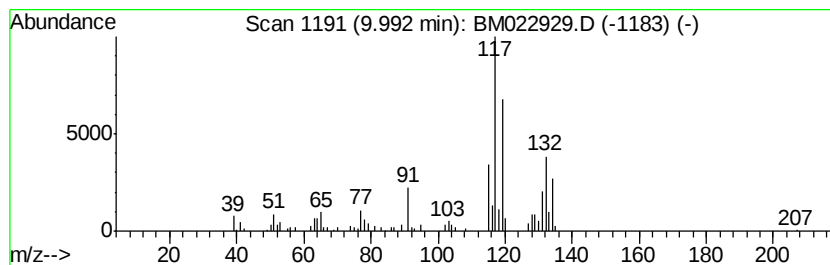
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.99	2.28 ng/ul	245349	Napthalene-d8	10.59

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
2		2,4-Dimethylstyrene	132	C10H12	002234-20-0	86
3		1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	64
4		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	64
5		1-Phenyl-1-butene	132	C10H12	000824-90-8	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100319\
Data File : BM022929.D
Acq On : 03 Oct 2019 20:25
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Instrument :
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ClientSampleId :
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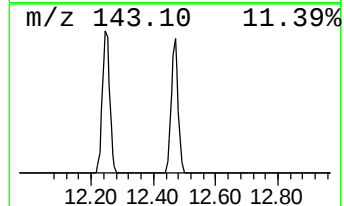
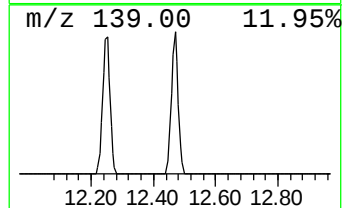
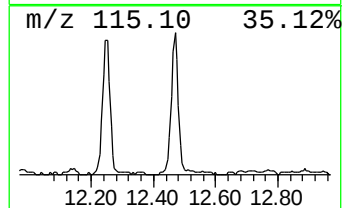
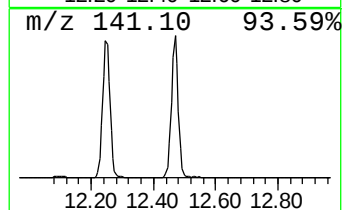
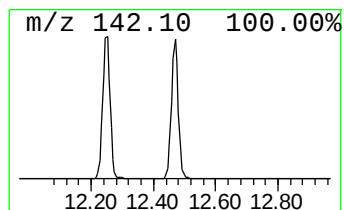
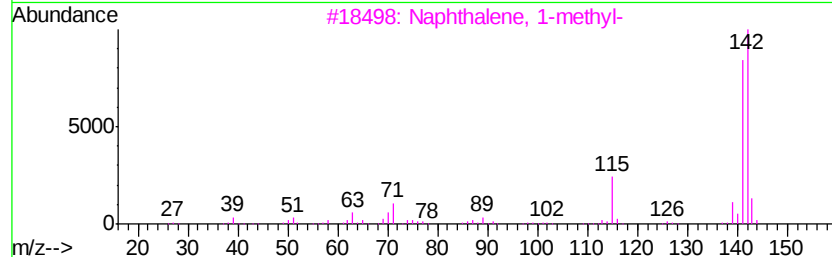
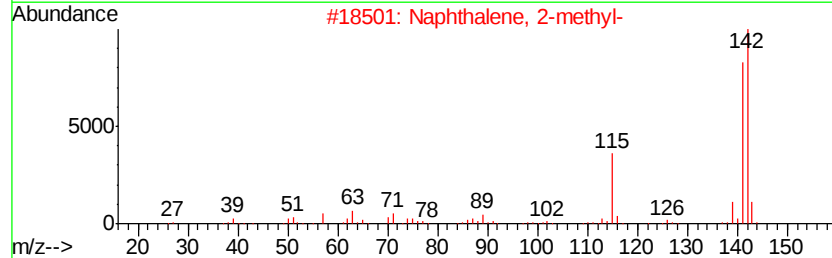
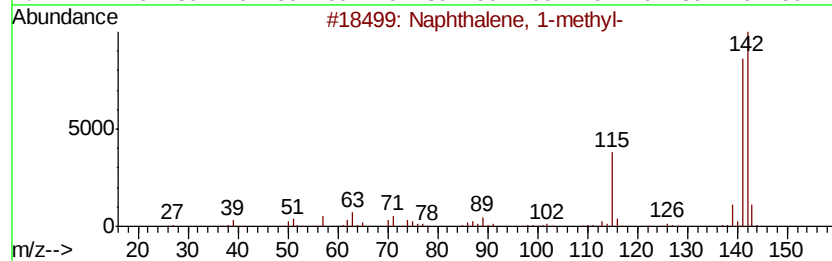
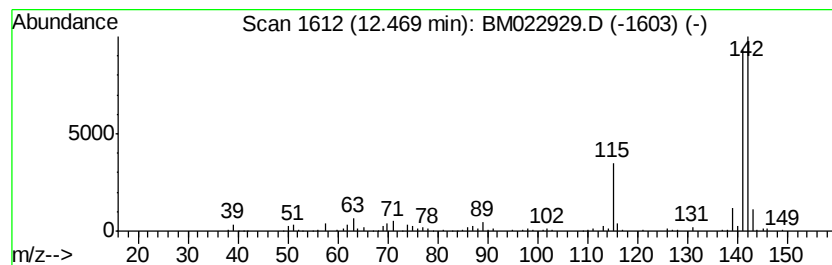
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Naphthalene, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.47	3.27 ng/ul	351691	Naphthalene-d8	10.59

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
4		Benzocycloheptatriene	142	C11H10	000264-09-5	91
5		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100319\
Data File : BM022929.D
Acq On : 03 Oct 2019 20:25
Operator : JU
Sample : K4983-18
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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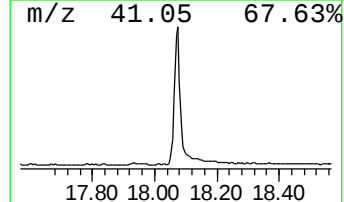
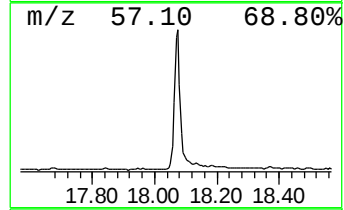
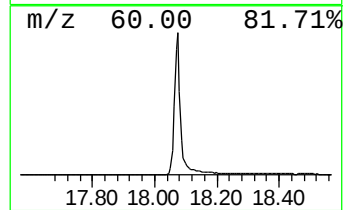
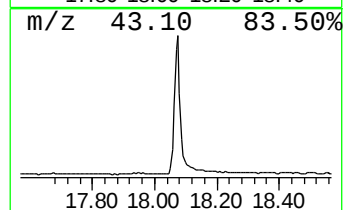
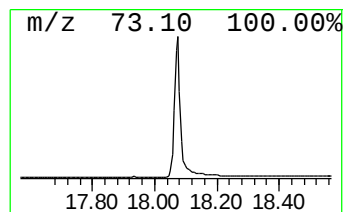
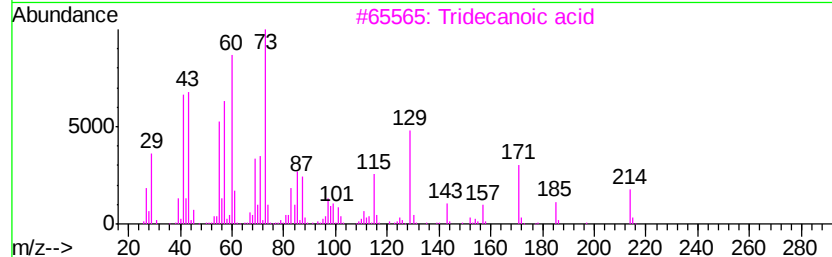
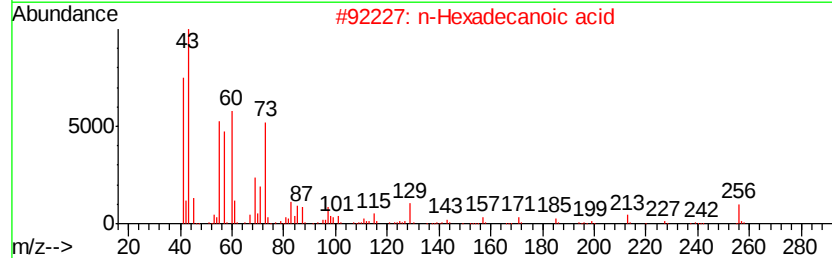
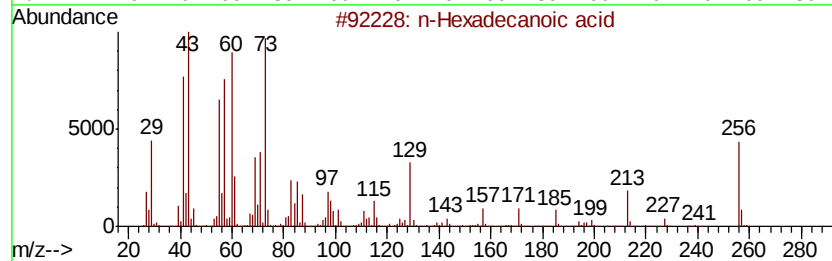
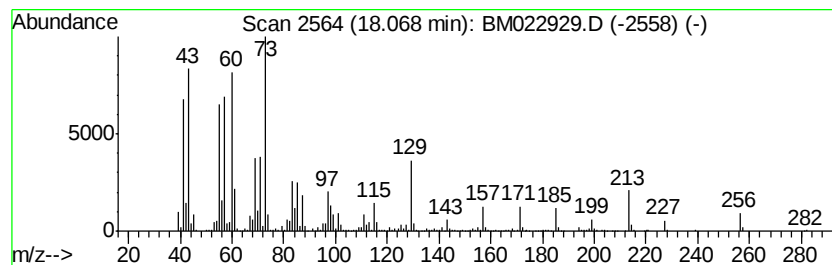
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.07	9.57 ng/ul	1610540	Phenanthrene-d10	17.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3		Tridecanoic acid	214	C13H26O2	000638-53-9	95
4		Tridecanoic acid	214	C13H26O2	000638-53-9	91
5		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100319\
Data File : BM022929.D
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Sample : K4983-18
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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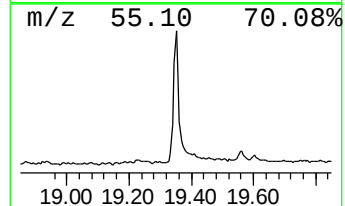
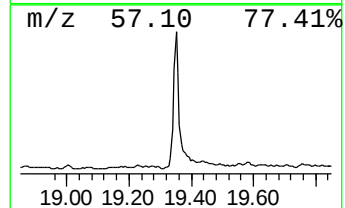
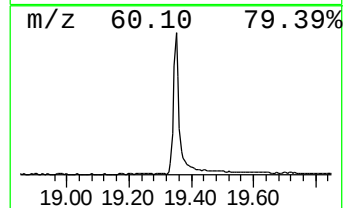
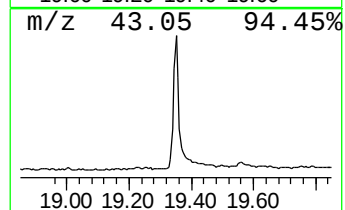
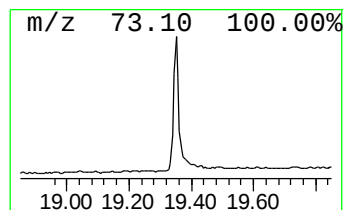
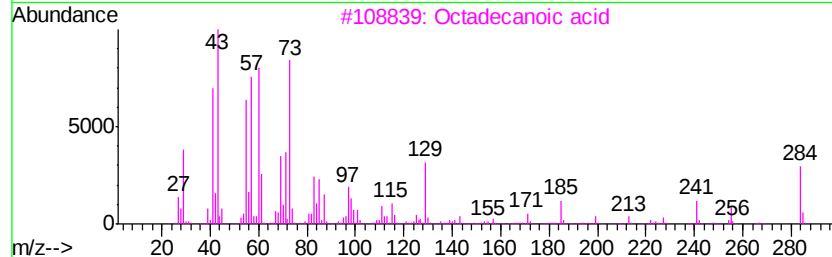
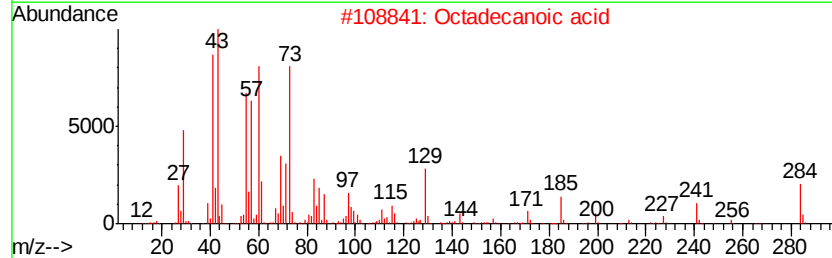
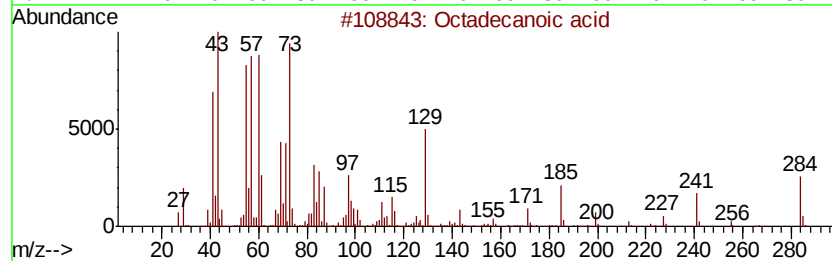
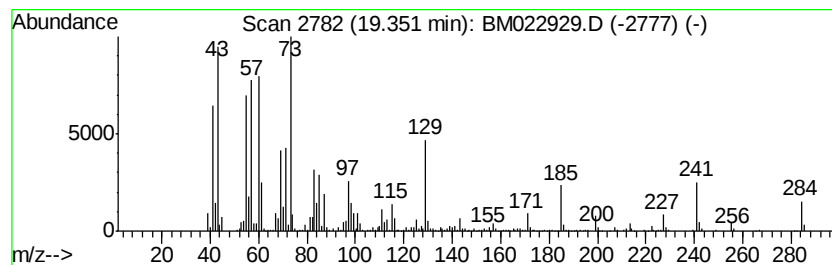
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 Octadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.35	5.78 ng/ul	855342	Chrysene-d12	21.35

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	93
4		Octadecanoic acid	284	C18H36O2	000057-11-4	93
5		Tetradecanoic acid	228	C14H28O2	000544-63-8	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100319\
Data File : BM022929.D
Acq On : 03 Oct 2019 20:25
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Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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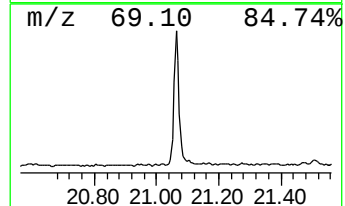
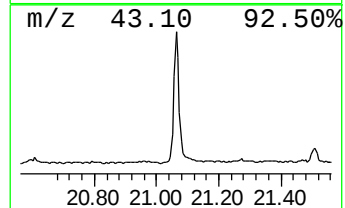
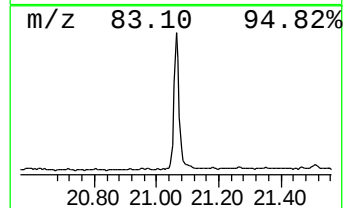
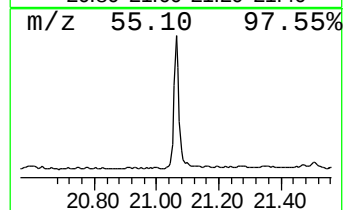
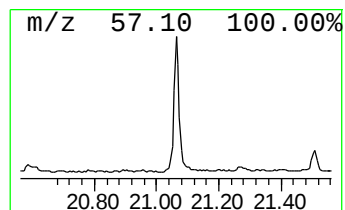
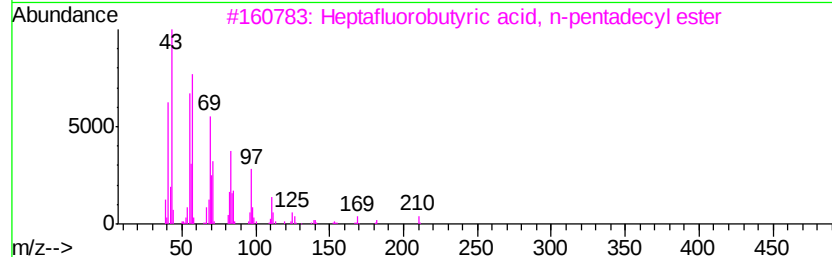
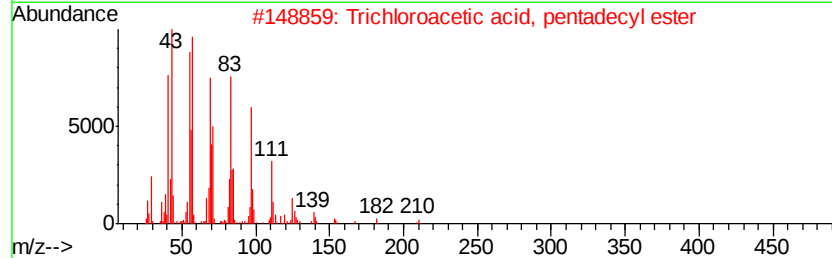
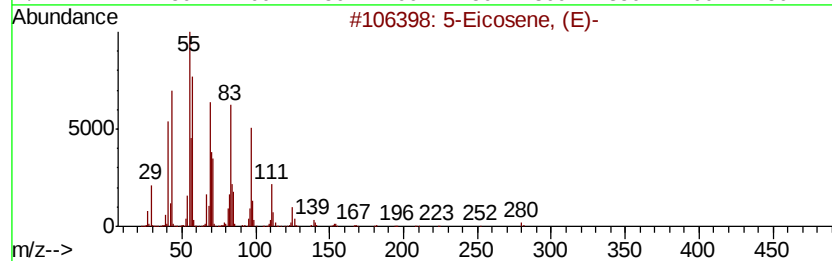
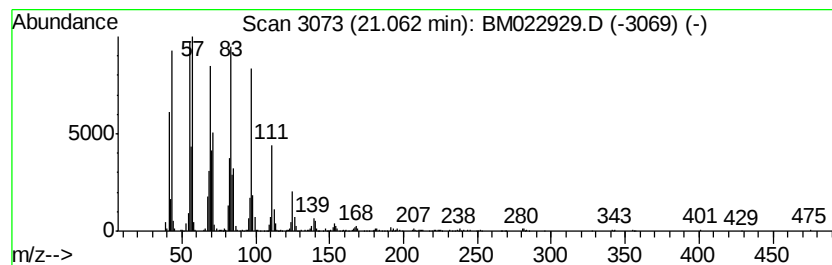
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 (DEL) Alkane: Cyclic21.06 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.06	4.54 ng/ul	671346	Chrysene-d12	21.35

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Eicosene, (E)-	280	C20H40	074685-30-6	98
2			Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	91
3			Heptafluorobutyric acid, n-penta...	424	C19H31F7O2	1000216-79-3	91
4			Trifluoroacetic acid, n-octadecy...	366	C20H37F3O2	079392-43-1	91
5			1-Heneicosyl formate	340	C22H44O2	077899-03-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM100319\
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 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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 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
Propanoic acid, 1...	3.73	3.0	ng/ul	211385	1	7.79	1384090	20.0
Propanoic acid, 2...	4.26	6.7	ng/ul	460566	1	7.79	1384090	20.0
Propanoic acid, p...	4.44	10.2	ng/ul	704293	1	7.79	1384090	20.0
Butanoic acid, 1-...	4.89	12.0	ng/ul	829891	1	7.79	1384090	20.0
o-Xylene	5.45	4.0	ng/ul	280197	1	7.79	1384090	20.0
p-Xylene	5.82	2.4	ng/ul	163496	1	7.79	1384090	20.0
Hexanoic acid	6.82	2.1	ng/ul	147782	1	7.79	1384090	20.0
Benzene, 1,2,3-tr...	7.45	5.0	ng/ul	346803	1	7.79	1384090	20.0
Benzene, 1,3,5-tr...	7.91	2.6	ng/ul	182088	1	7.79	1384090	20.0
Indane	8.17	2.5	ng/ul	175696	1	7.79	1384090	20.0
Benzene, 2-etheny...	9.99	2.3	ng/ul	245349	2	10.59	2153900	20.0
Naphthalene, 1-me...	12.47	3.3	ng/ul	351691	2	10.59	2153900	20.0
n-Hexadecanoic acid	18.07	9.6	ng/ul	1610540	4	17.17	3366580	20.0
Octadecanoic acid	19.35	5.8	ng/ul	855342	5	21.35	2959480	20.0
(DEL) Alkane: Cyc...	21.06	4.5	ng/ul	671346	5	21.35	2959480	20.0