

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN092319\
 Data File : BN007814.D
 Acq On : 23 Sep 2019 22:41
 Operator : HP/JU
 Sample : K4895-06
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BFDF9

Integration Parameters: LSCINT.P

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

Filtering: 5
 Min Area: 0.1 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN091619MA.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	3.152	25	28	35	rVB	85243	133051	0.84%	0.075%
2	3.281	46	50	65	rVB	4718815	5522757	35.03%	3.111%
3	3.634	107	110	116	rBV	119646	172597	1.09%	0.097%
4	3.693	116	120	126	rBV	556864	649465	4.12%	0.366%
5	3.875	146	151	156	rVV	80575	127055	0.81%	0.072%
6	3.946	159	163	170	rVB2	143784	244977	1.55%	0.138%
7	4.217	204	209	216	rBV	1146078	1400632	8.88%	0.789%
8	4.281	216	220	229	rVB	136250	204975	1.30%	0.115%
9	4.387	234	238	247	rVV	1576186	2034277	12.90%	1.146%
10	4.452	247	249	261	rVB	77562	142539	0.90%	0.080%
11	4.828	306	313	324	rBV	1876684	2589343	16.42%	1.458%
12	5.052	346	351	358	rVB	413728	596320	3.78%	0.336%
13	5.675	452	457	464	rBV	249917	346444	2.20%	0.195%
14	6.687	624	629	637	rBV	149040	248678	1.58%	0.140%
15	6.863	652	659	667	rVV3	540126	1012249	6.42%	0.570%
16	6.928	667	670	677	rVB	111957	161720	1.03%	0.091%
17	7.028	680	687	694	rBV	1403517	2291459	14.53%	1.291%
18	7.093	694	698	705	rVB	158007	256378	1.63%	0.144%
19	7.222	713	720	731	rBV	1732256	2710443	17.19%	1.527%
20	7.346	735	741	748	rVB	535978	801789	5.09%	0.452%
21	7.687	793	799	808	rBV	2100336	3337372	21.17%	1.880%
22	7.805	814	819	828	rVB	272550	465579	2.95%	0.262%
23	8.058	853	862	870	rVB	512319	946306	6.00%	0.533%
24	8.181	878	883	888	rBV	88177	137941	0.87%	0.078%
25	8.328	903	908	912	rVV2	128188	212933	1.35%	0.120%
26	8.387	912	918	927	rVV	1399187	2255074	14.30%	1.270%
27	8.493	930	936	945	rVV3	78654	173064	1.10%	0.097%
28	8.640	954	961	966	rBV	92324	148776	0.94%	0.084%
29	8.793	980	987	988	rBV	230306	328118	2.08%	0.185%
30	8.828	988	993	1007	rVV	1802774	3367629	21.36%	1.897%
31	9.305	1069	1074	1080	rVB2	205303	321051	2.04%	0.181%
32	9.369	1080	1085	1093	rVB	160095	255030	1.62%	0.144%
33	9.546	1108	1115	1123	rBV	1418663	2349715	14.90%	1.323%
34	9.722	1140	1145	1151	rVV	220237	391210	2.48%	0.220%

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35	9.869	1161	1170	1179	rVV3	385100	745702	4.73%	0.420%
36	9.963	1179	1186	1190	rVV3	89707	190712	1.21%	0.107%
37	10.081	1199	1206	1217	rVV	2524319	4078226	25.87%	2.297%
38	10.457	1261	1270	1276	rVV	2908561	4820098	30.57%	2.715%
39	10.510	1276	1279	1286	rVV	330842	546976	3.47%	0.308%
40	10.593	1286	1293	1306	rVV	1856207	3318847	21.05%	1.869%
41	11.657	1469	1474	1484	rVB2	75187	152202	0.97%	0.086%
42	11.804	1489	1499	1503	rBV2	108480	232079	1.47%	0.131%
43	12.016	1520	1535	1539	rBV6	97601	259795	1.65%	0.146%
44	12.122	1547	1553	1559	rVB	519670	836532	5.31%	0.471%
45	12.240	1567	1573	1580	rBV2	79440	164726	1.04%	0.093%
46	12.346	1580	1591	1603	rVV	1356692	2228854	14.14%	1.255%
47	13.163	1724	1730	1737	rBV5	69072	143083	0.91%	0.081%
48	13.228	1737	1741	1746	rBV	146367	230155	1.46%	0.130%
49	13.369	1755	1765	1772	rVB2	131050	328190	2.08%	0.185%
50	13.487	1778	1785	1791	rBV3	166868	416331	2.64%	0.235%
51	13.640	1801	1811	1816	rBV	370271	714350	4.53%	0.402%
52	13.693	1816	1820	1821	rVV2	137142	187601	1.19%	0.106%
53	13.728	1821	1826	1831	rVV	4950846	6714007	42.59%	3.782%
54	13.775	1831	1834	1840	rVV	892559	1165178	7.39%	0.656%
55	13.875	1847	1851	1854	rVV	277154	410864	2.61%	0.231%
56	13.910	1854	1857	1861	rVB	122589	168784	1.07%	0.095%
57	14.004	1867	1873	1886	rVB	5175494	7919139	50.23%	4.461%
58	14.316	1919	1926	1932	rVV	4469098	6575955	41.71%	3.704%
59	14.381	1932	1937	1945	rVB	1931004	2893003	18.35%	1.630%
60	14.510	1954	1959	1972	rBV2	473098	865801	5.49%	0.488%
61	14.716	1990	1994	1998	rBV2	158723	236150	1.50%	0.133%
62	14.798	2004	2008	2012	rVB	101585	142004	0.90%	0.080%
63	14.945	2028	2033	2037	rVV2	104100	177565	1.13%	0.100%
64	14.992	2037	2041	2048	rVB5	88654	159637	1.01%	0.090%
65	15.087	2053	2057	2060	rBV2	172261	257221	1.63%	0.145%
66	15.310	2089	2095	2101	rBV	6753113	9876498	62.64%	5.563%
67	15.369	2101	2105	2109	rVB	686291	988235	6.27%	0.557%
68	15.428	2109	2115	2119	rBV	1890245	2552619	16.19%	1.438%
69	15.775	2168	2174	2178	rBV5	88247	200995	1.27%	0.113%
70	15.822	2178	2182	2188	rVB2	168820	261342	1.66%	0.147%
71	15.981	2206	2209	2214	rBV	74738	116465	0.74%	0.066%

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72	16.034	2214	2218	2226	rVV4	140594	262072	1.66%	0.148%
73	16.328	2264	2268	2270	rBV	105425	145417	0.92%	0.082%
74	16.422	2280	2284	2288	rVB2	99597	124048	0.79%	0.070%
75	16.522	2297	2301	2307	rVB2	83092	134255	0.85%	0.076%
76	16.881	2358	2362	2366	rVB	131603	190412	1.21%	0.107%
77	17.063	2387	2393	2397	rBV2	5462016	7814283	49.56%	4.401%
78	17.104	2397	2400	2404	rVV	610771	812900	5.16%	0.458%
79	17.157	2404	2409	2420	rVB2	8039973	11860680	75.23%	6.681%
80	17.257	2422	2426	2431	rVB2	74499	127824	0.81%	0.072%
81	17.439	2453	2457	2458	rBV3	112201	136945	0.87%	0.077%
82	17.939	2538	2542	2544	rBV	90134	119054	0.76%	0.067%
83	17.975	2544	2548	2557	rVV2	738741	1086415	6.89%	0.612%
84	18.128	2570	2574	2579	rBV2	184187	306732	1.95%	0.173%
85	19.092	2734	2738	2740	rBV	101081	145645	0.92%	0.082%
86	19.122	2740	2743	2749	rVB	238593	338981	2.15%	0.191%
87	19.263	2763	2767	2773	rBV2	330123	481454	3.05%	0.271%
88	19.345	2779	2781	2787	rVB2	111472	126892	0.80%	0.071%
89	19.457	2794	2800	2816	rVB	10855157	15394249	97.64%	8.671%
90	20.992	3057	3061	3067	rBV	680249	969381	6.15%	0.546%
91	21.257	3100	3106	3115	rBV	7713851	9951344	63.12%	5.605%
92	21.374	3123	3126	3132	rVB	203734	269021	1.71%	0.152%
93	21.869	3207	3210	3214	rBV3	173132	224567	1.42%	0.126%
94	22.327	3285	3288	3294	rVB	274516	350502	2.22%	0.197%
95	22.833	3370	3374	3381	rVB	398184	562113	3.57%	0.317%
96	23.333	3451	3459	3465	rBV	8705210	15765856	100.00%	8.880%
97	23.392	3466	3469	3475	rVV	467381	716223	4.54%	0.403%
98	23.474	3475	3483	3493	rVB2	5383888	10286262	65.24%	5.794%
99	24.033	3573	3578	3584	rBV2	378376	651839	4.13%	0.367%
100	24.768	3698	3703	3711	rVB2	286423	570812	3.62%	0.322%

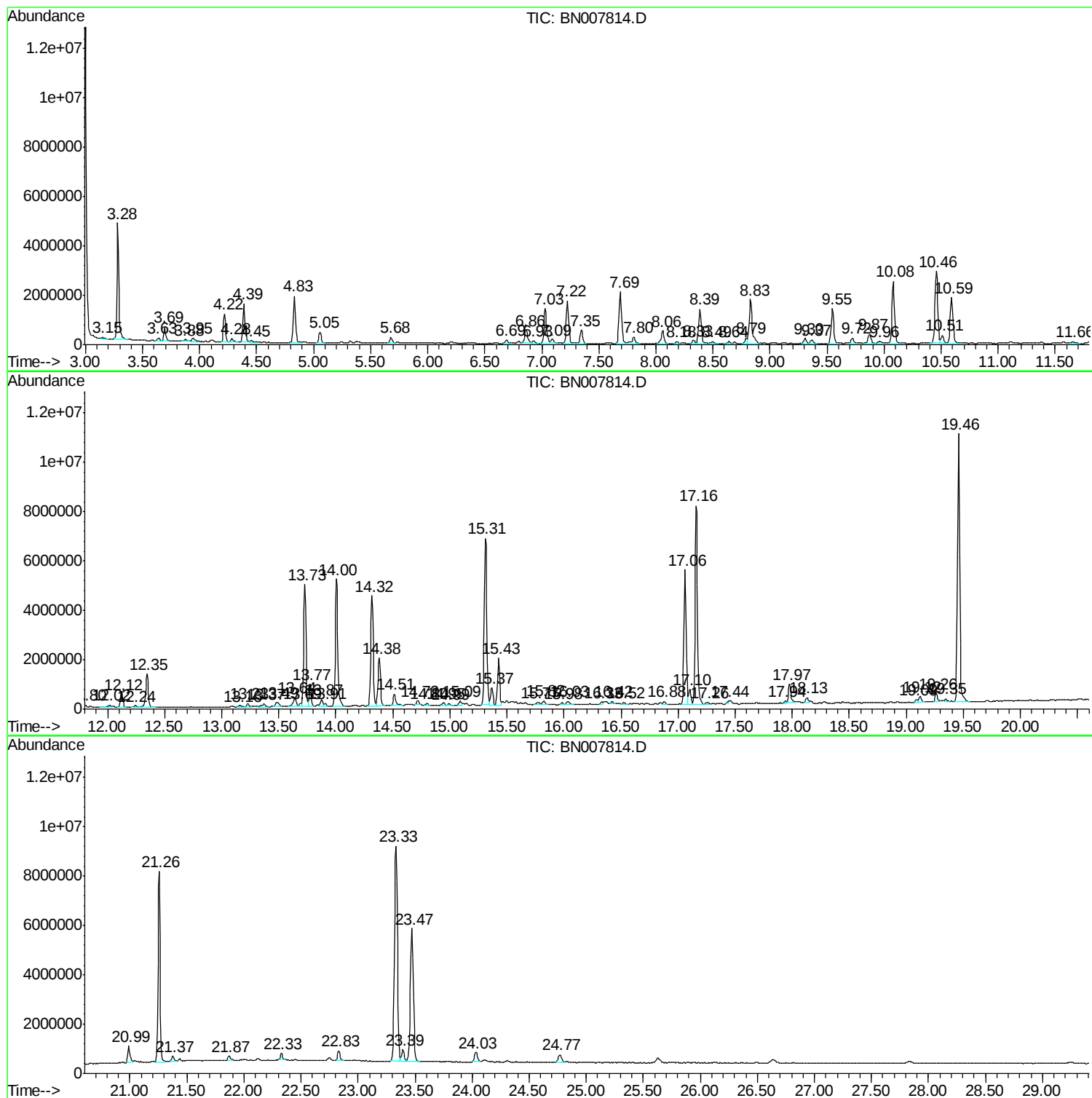
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN091619MA.M
Quant Title : SVOA CALIBRATION

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



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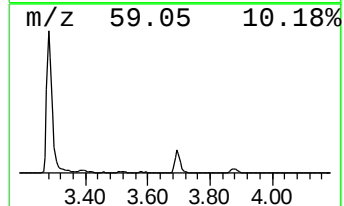
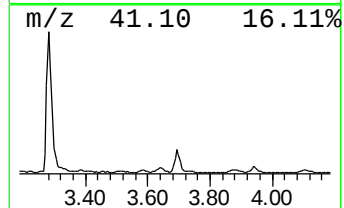
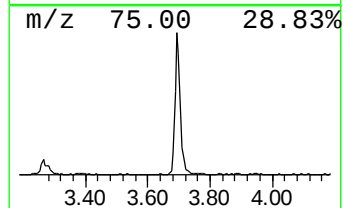
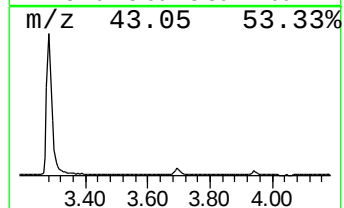
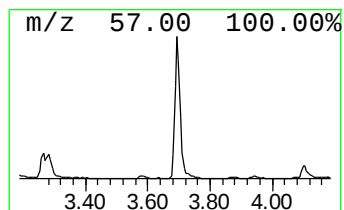
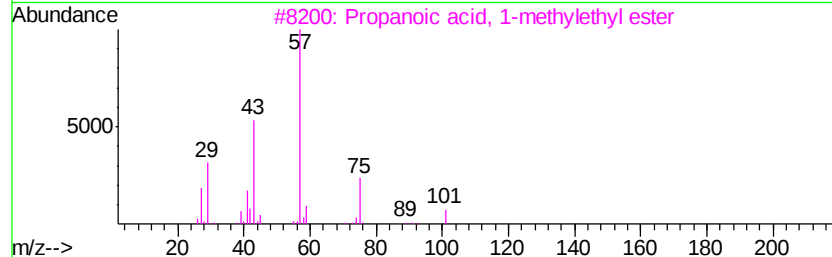
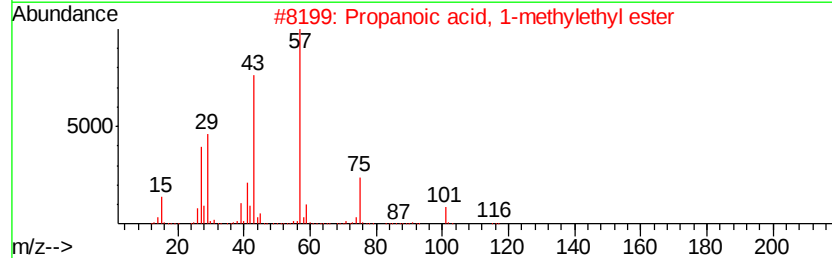
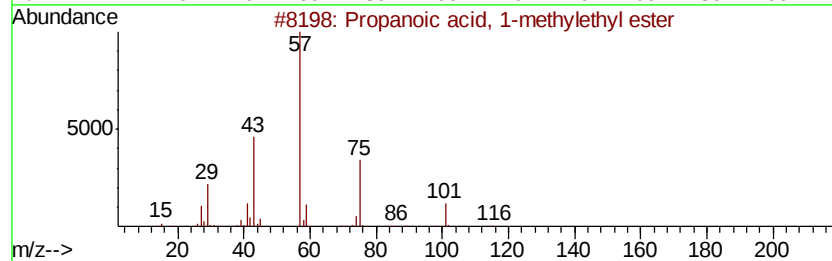
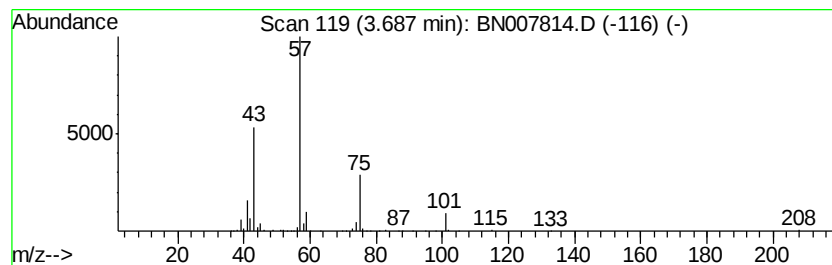
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN091619MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Propanoic acid, 1-methyleth... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.69	3.89 ng/ul	649465	1,4-Dichlorobenzene-d4	7.69

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	74
3		Propanoic acid, 1-methylethyl ester	116	C6H12O2	000637-78-5	72
4		Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	59
5		Propanoic acid, 1-methylethyl ester	116	C6H12O2	000637-78-5	50



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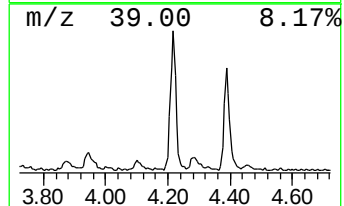
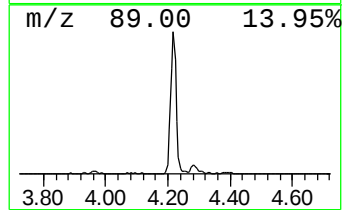
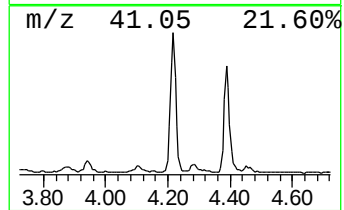
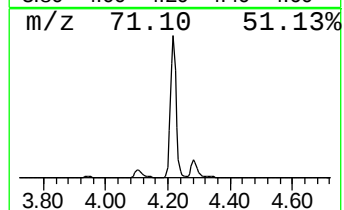
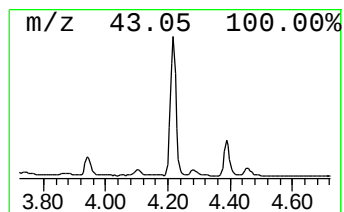
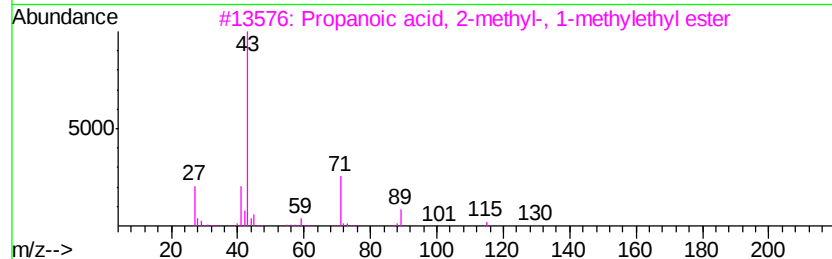
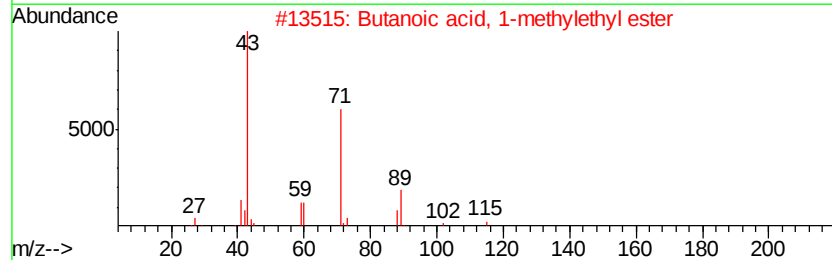
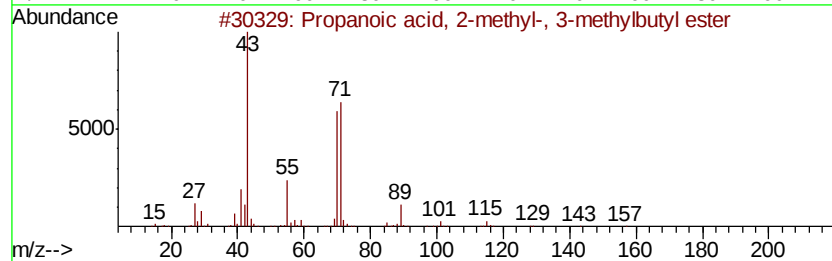
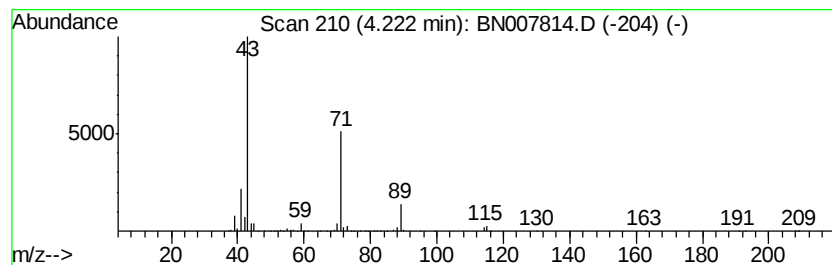
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.22	8.39 ng/ul	1400630	1,4-Dichlorobenzene-d4	7.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propanoic acid, 2-methyl-, 3-met...	158	C9H18O2	002050-01-3	56
2		Butanoic acid, 1-methylethyl ester	130	C7H14O2	000638-11-9	50
3		Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	50
4		Propanoic acid, 2-methyl-, propy...	130	C7H14O2	000644-49-5	42
5		Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	40



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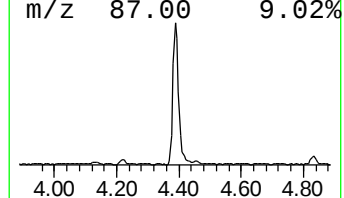
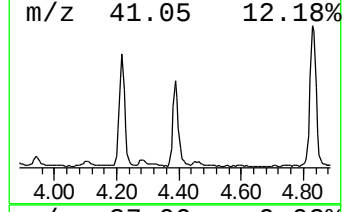
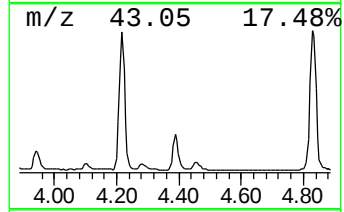
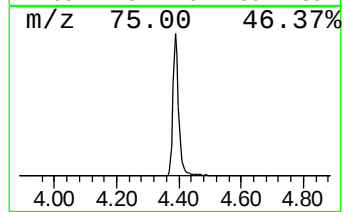
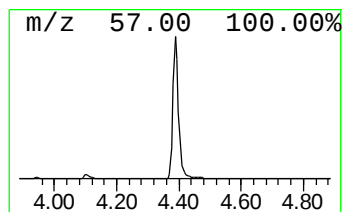
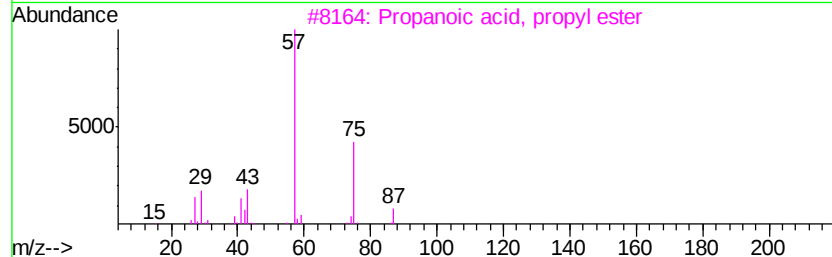
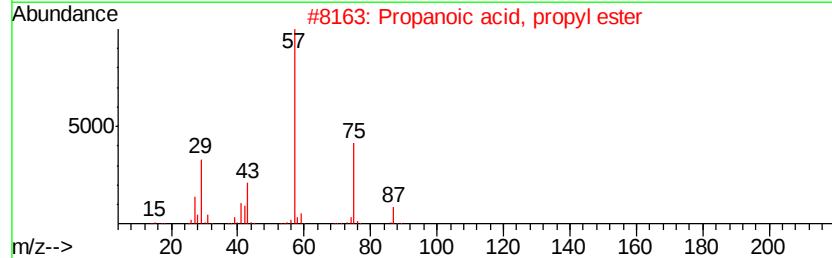
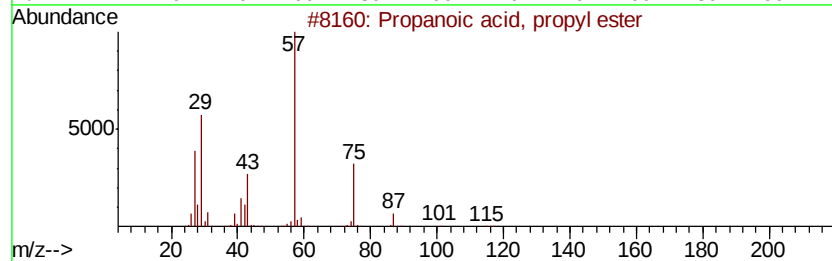
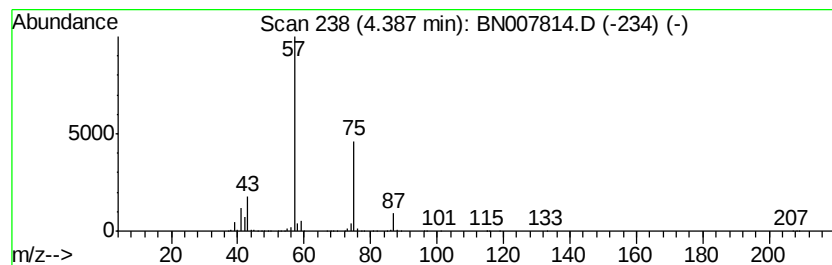
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Propanoic acid, propyl ester Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.39	12.19 ng/ul	2034280	1,4-Dichlorobenzene-d4	7.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	74
2		Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	59
3		Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	50
4		Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	37
5		Propanoic acid, butyl ester	130	C7H14O2	000590-01-2	33



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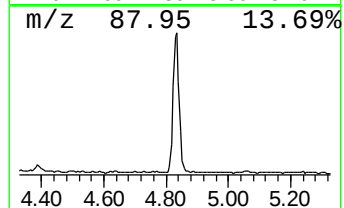
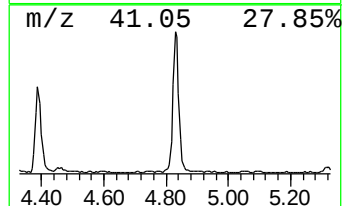
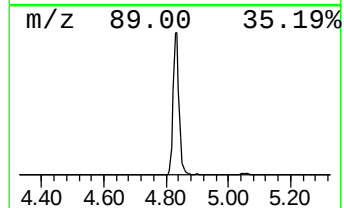
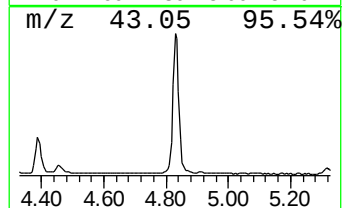
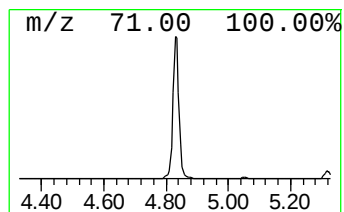
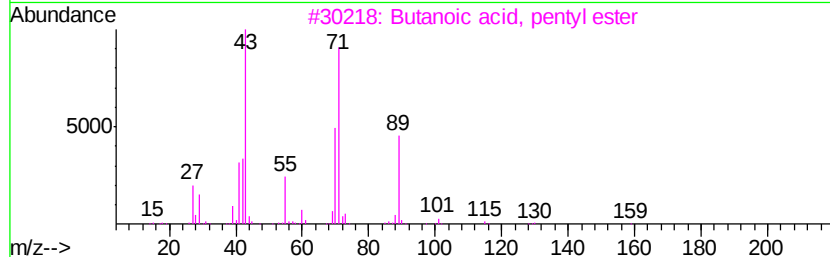
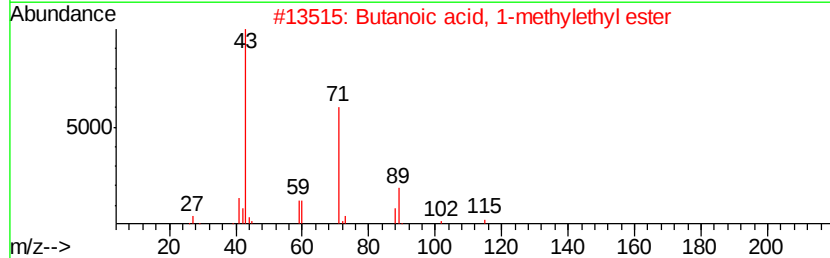
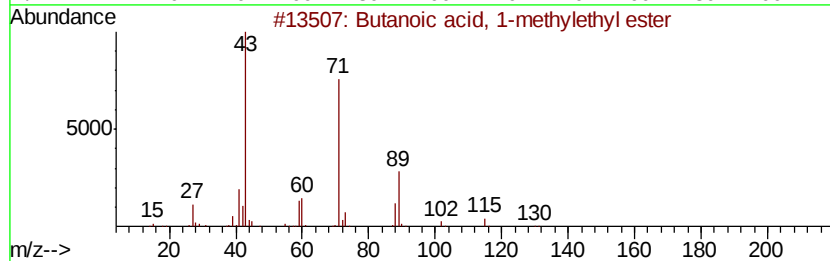
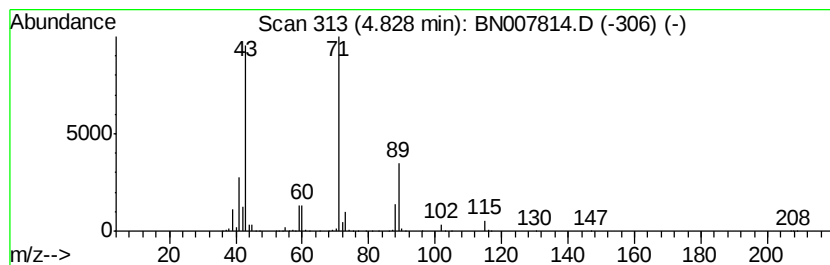
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.83	15.52 ng/ul	2589340	1,4-Dichlorobenzene-d4	7.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butanoic acid, 1-methylethyl ester	130	C7H14O2	000638-11-9	91
2		Butanoic acid, 1-methylethyl ester	130	C7H14O2	000638-11-9	83
3		Butanoic acid, pentyl ester	158	C9H18O2	000540-18-1	64
4		Butanoic acid, 1-methylethyl ester	130	C7H14O2	000638-11-9	58
5		Butanoic acid, 1-methylpropyl ester	144	C8H16O2	000819-97-6	56



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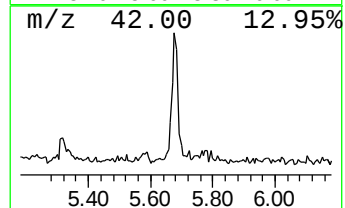
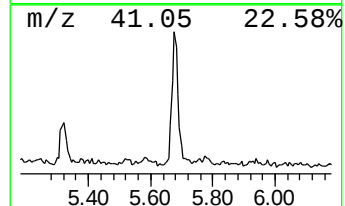
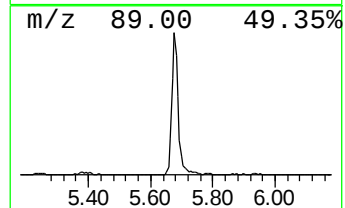
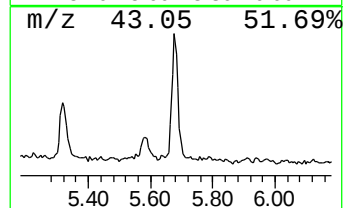
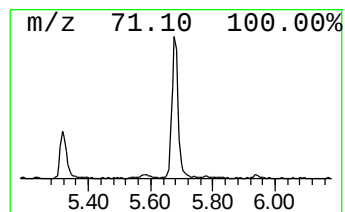
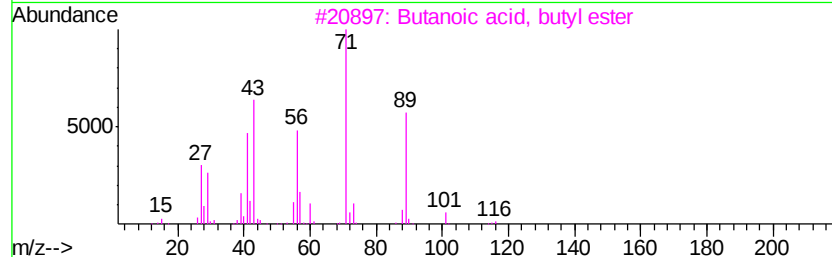
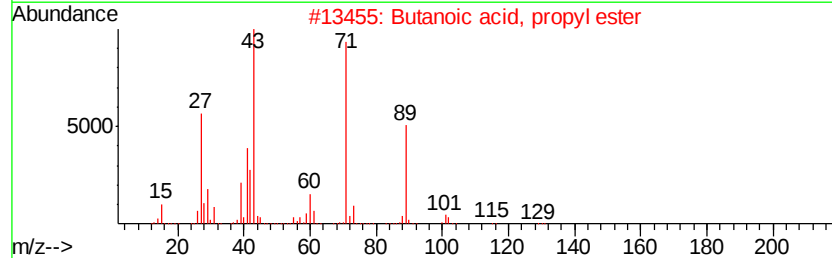
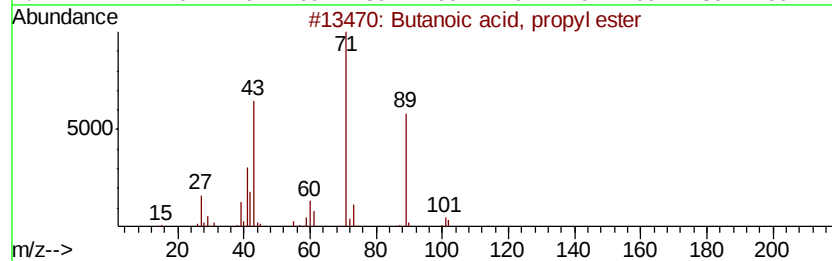
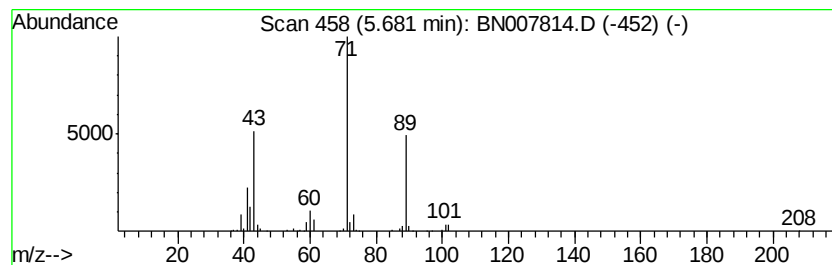
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Butanoic acid, propyl ester Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.68	2.08 ng/ul	346444	1,4-Dichlorobenzene-d4	7.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butanoic acid, propyl ester	130	C7H14O2	000105-66-8	90
2		Butanoic acid, propyl ester	130	C7H14O2	000105-66-8	83
3		Butanoic acid, butyl ester	144	C8H16O2	000109-21-7	78
4		Butanoic acid, hexyl ester	172	C10H20O2	002639-63-6	78
5		Propanoic acid, 2-methyl-, propy...	130	C7H14O2	000644-49-5	72



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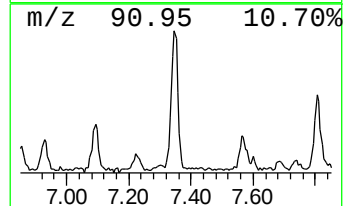
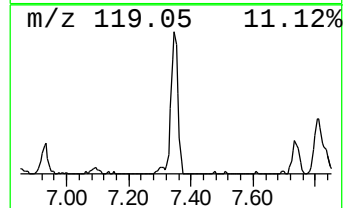
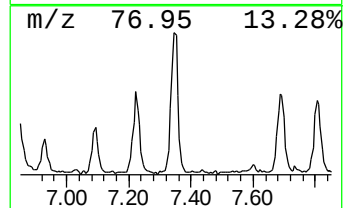
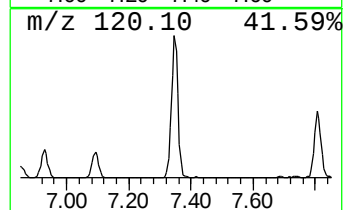
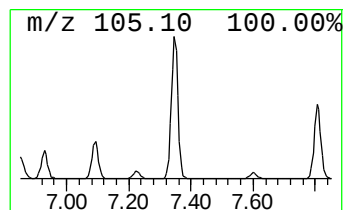
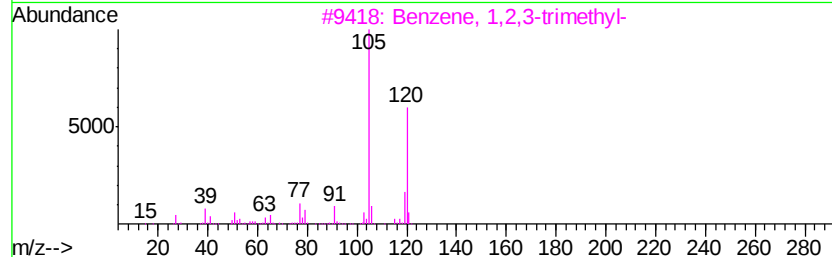
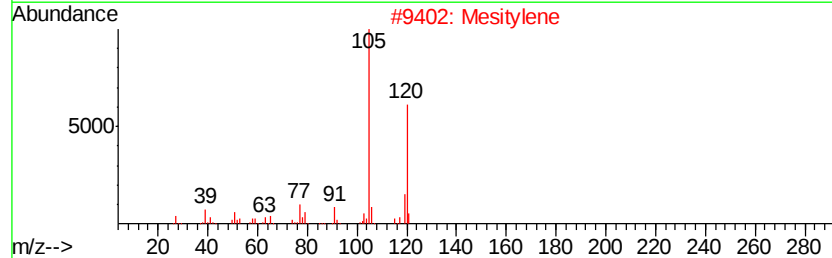
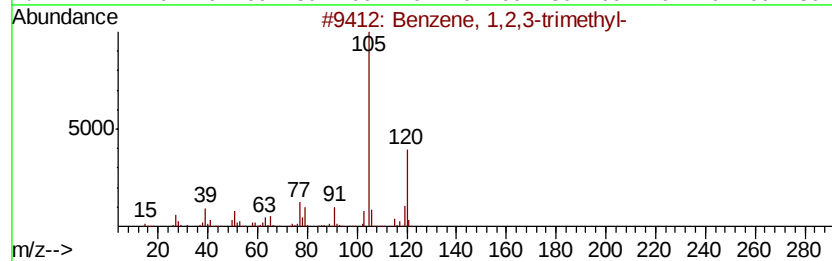
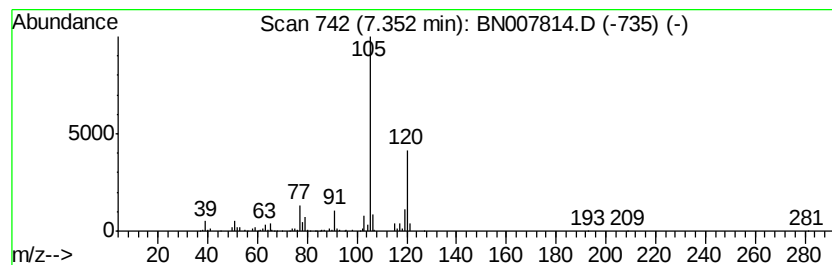
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Benzene, 1,2,3-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.35	4.80 ng/ul	801789	1,4-Dichlorobenzene-d4	7.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2		Mesitylene	120	C9H12	000108-67-8	97
3		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
4		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
5		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



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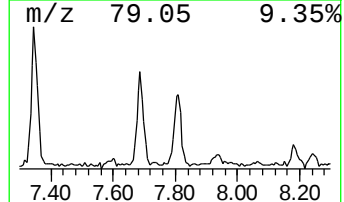
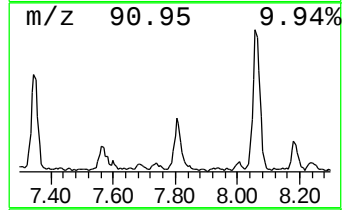
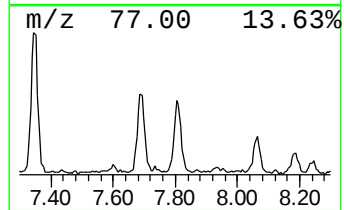
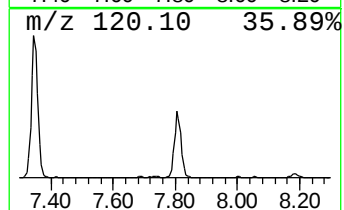
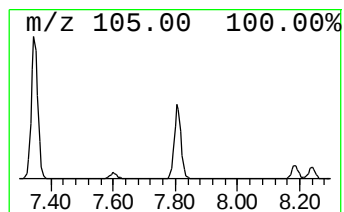
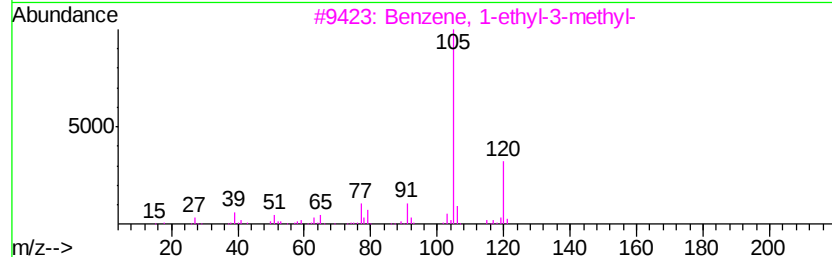
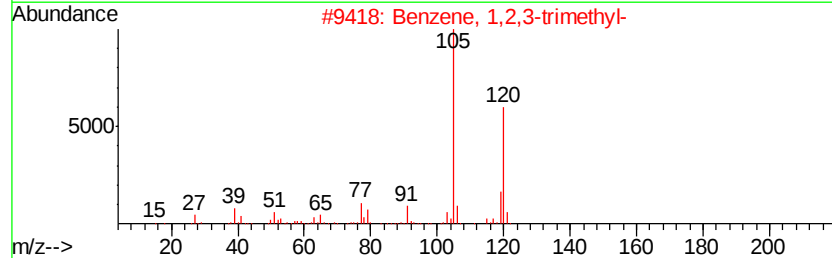
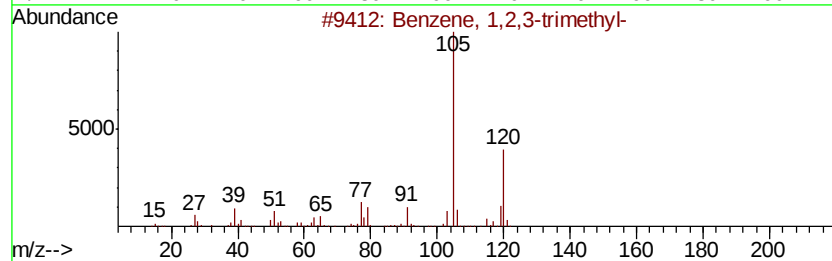
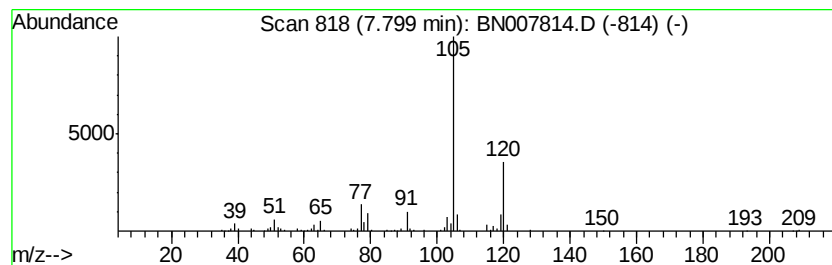
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Benzene, 1-ethyl-3-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.80	2.79 ng/ul	465579	1,4-Dichlorobenzene-d4	7.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
3		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4		Mesitylene	120	C9H12	000108-67-8	91
5		Mesitylene	120	C9H12	000108-67-8	91



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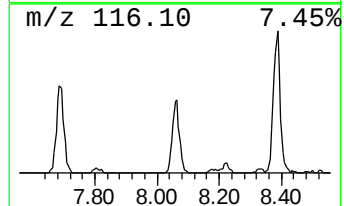
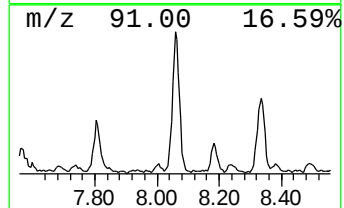
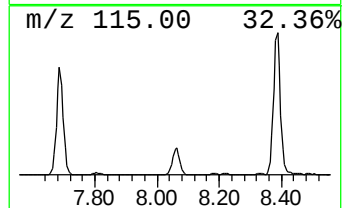
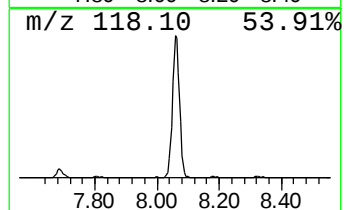
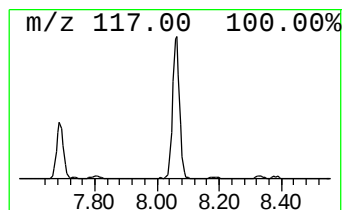
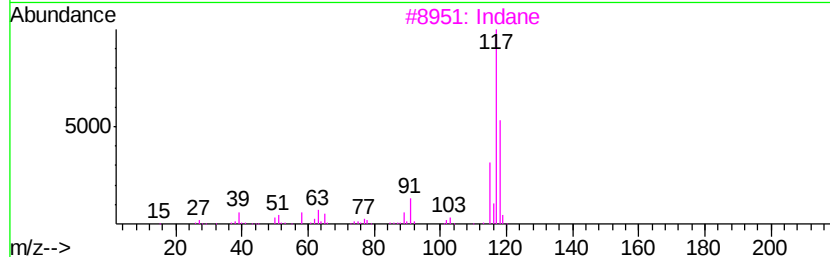
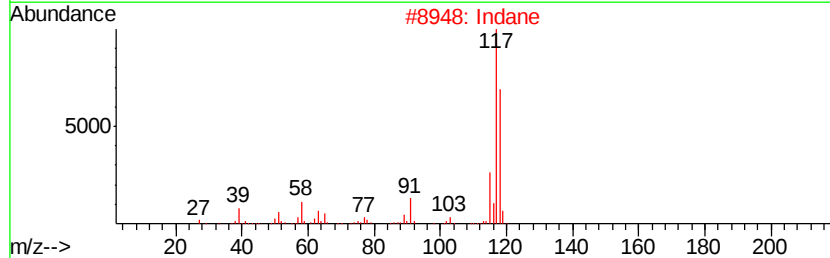
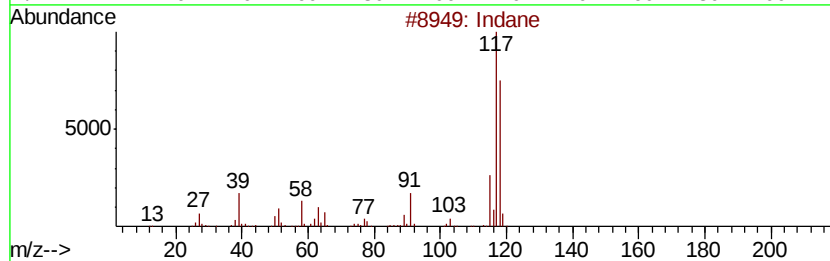
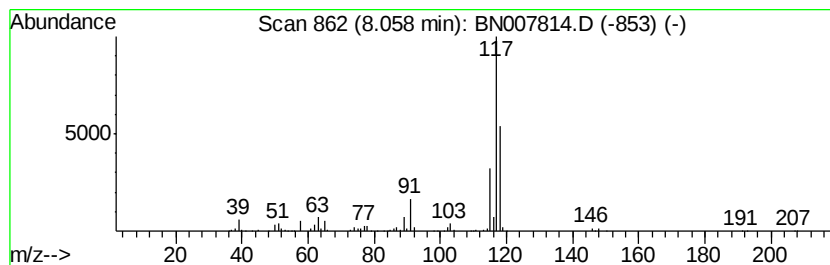
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Indane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.06	5.67 ng/ul	946306	1,4-Dichlorobenzene-d4	7.69

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane		118	C9H10	000496-11-7	83
2	Indane		118	C9H10	000496-11-7	83
3	Indane		118	C9H10	000496-11-7	81
4	Benzene, cyclopropyl-		118	C9H10	000873-49-4	72
5	Benzene, cyclopropyl-		118	C9H10	000873-49-4	68



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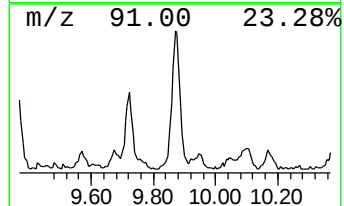
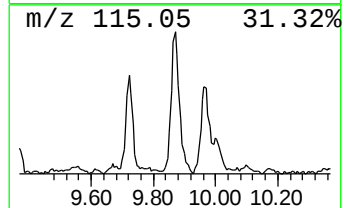
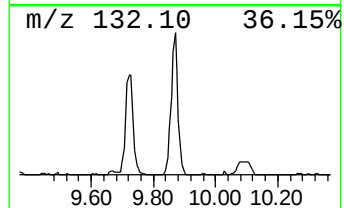
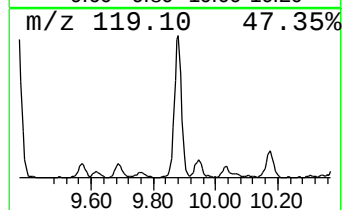
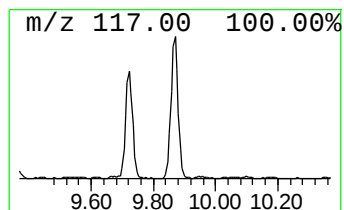
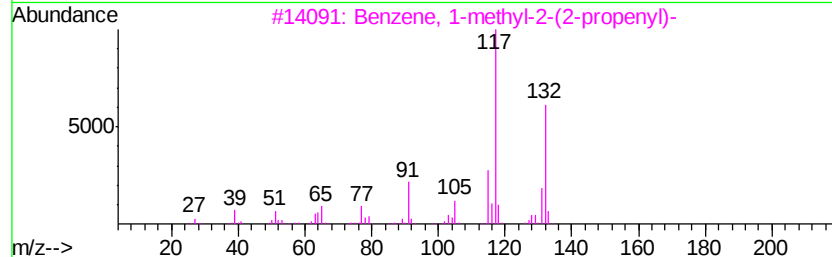
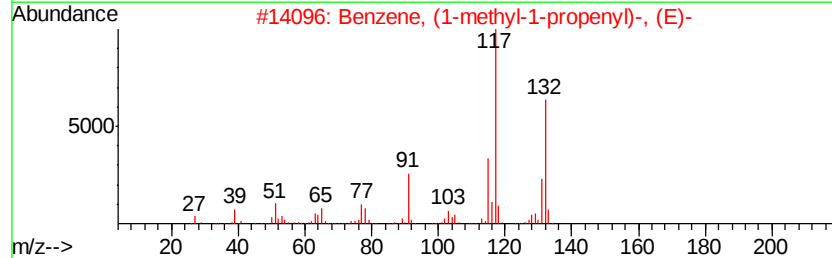
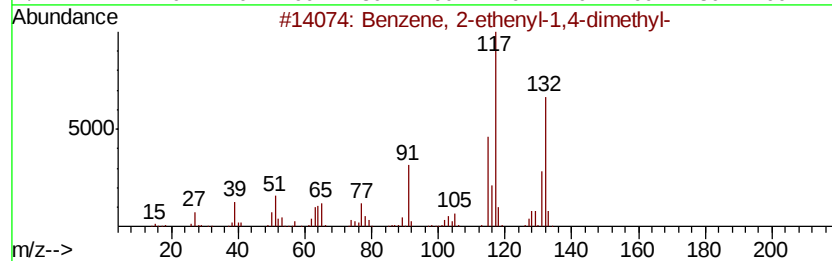
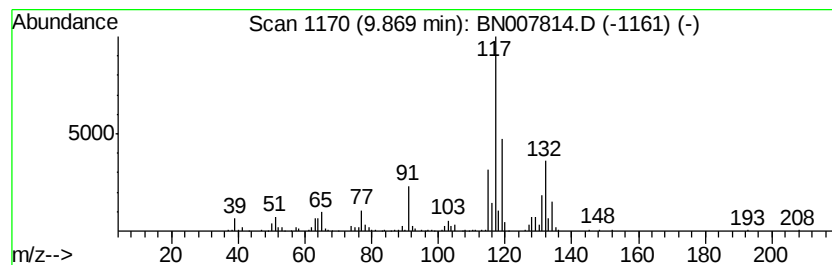
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.87	3.09 ng/ul	745702	Napthalene-d8	10.46

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
2		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	76
3		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	70
4		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	70
5		Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	70



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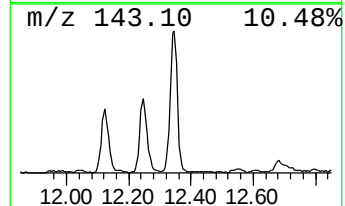
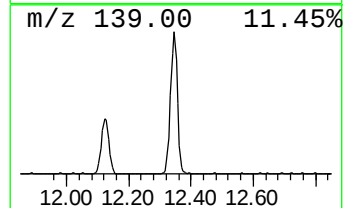
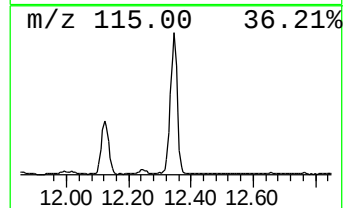
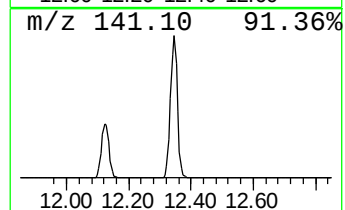
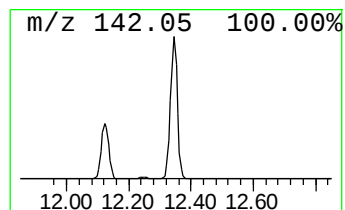
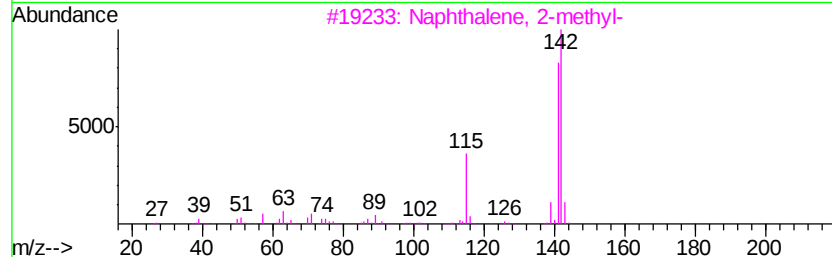
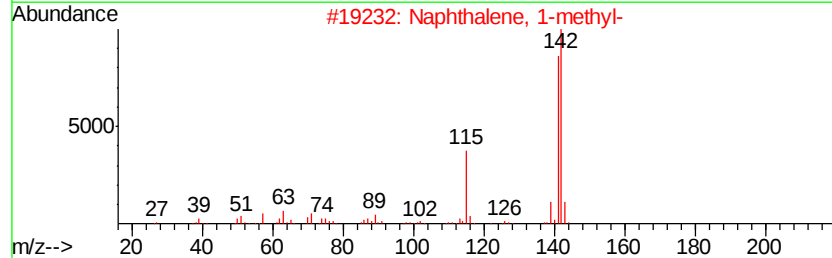
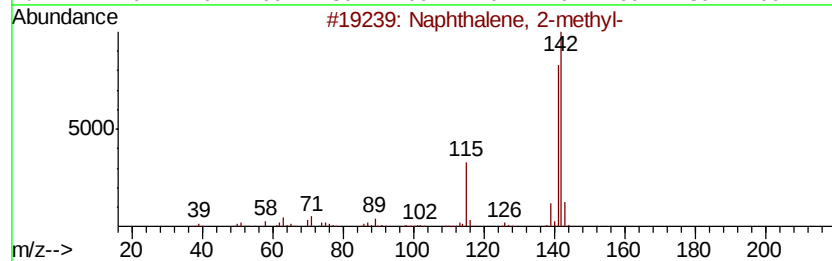
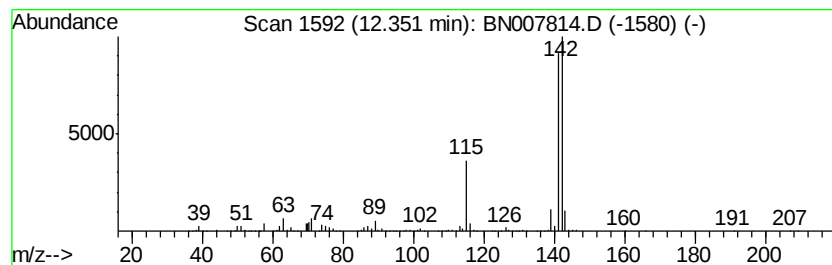
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Naphthalene, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.35	9.25 ng/ul	2228850	Naphthalene-d8	10.46

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN092319\
Data File : BN007814.D
Acq On : 23 Sep 2019 22:41
Operator : HP/JU
Sample : K4895-06
Misc :
ALS Vial : 16 Sample Multiplier: 1

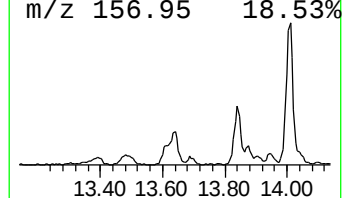
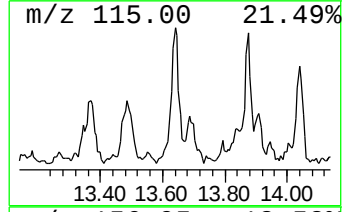
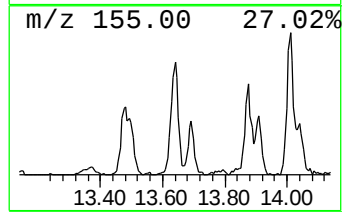
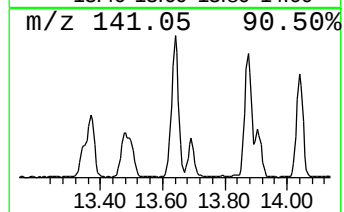
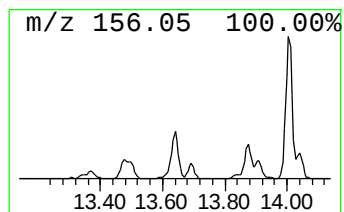
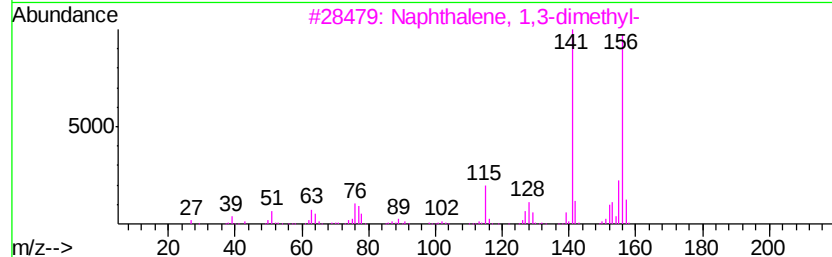
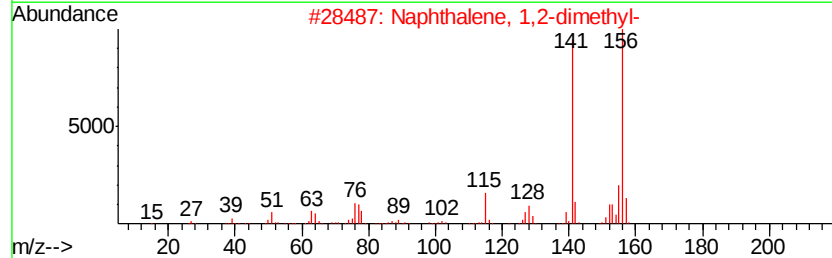
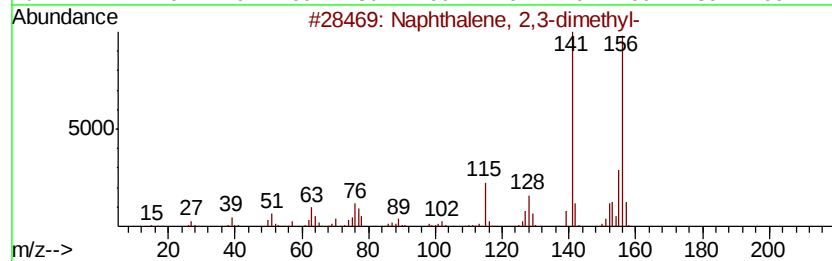
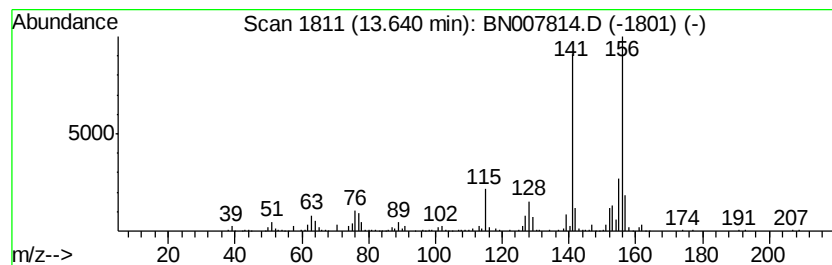
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN091619MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 12 Naphthalene, 2,3-dimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.64	2.17 ng/ul	714350	Acenaphthene-d10	14.32
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 2,3-dimethyl-	156 C12H12	000581-40-8 97
2		Naphthalene, 1,2-dimethyl-	156 C12H12	000573-98-8 97
3		Naphthalene, 1,3-dimethyl-	156 C12H12	000575-41-7 97
4		Naphthalene, 1,4-dimethyl-	156 C12H12	000571-58-4 97
5		Naphthalene, 2,3-dimethyl-	156 C12H12	000581-40-8 97



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN092319\
Data File : BN007814.D
Acq On : 23 Sep 2019 22:41
Operator : HP/JU
Sample : K4895-06
Misc :
ALS Vial : 16 Sample Multiplier: 1

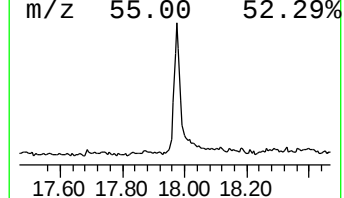
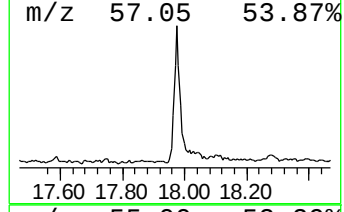
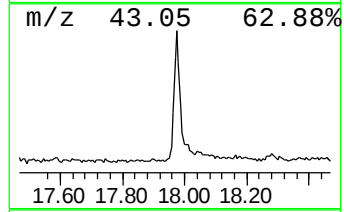
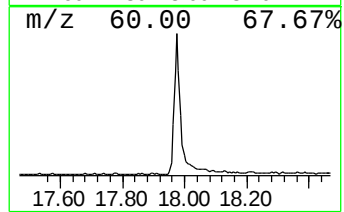
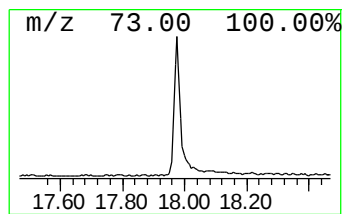
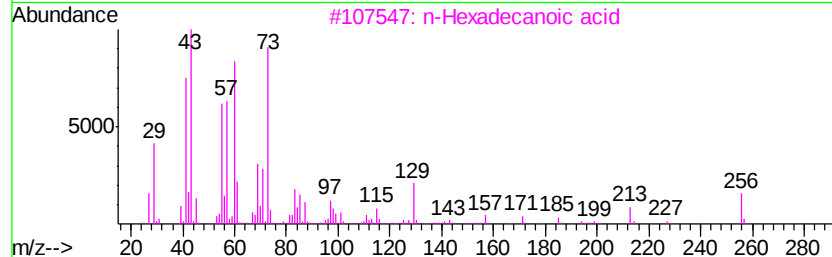
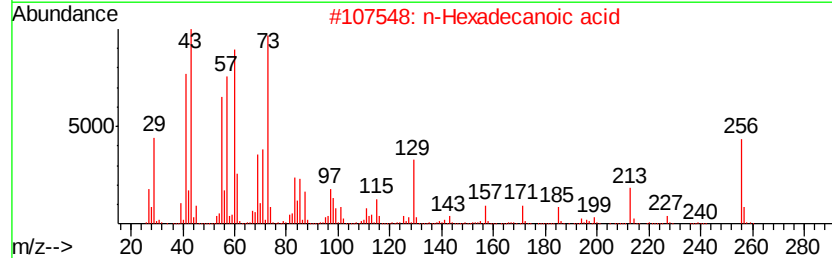
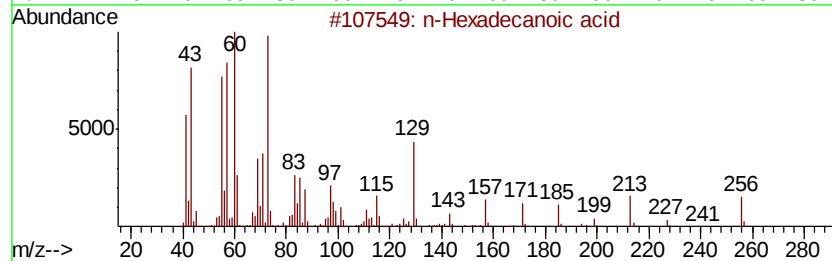
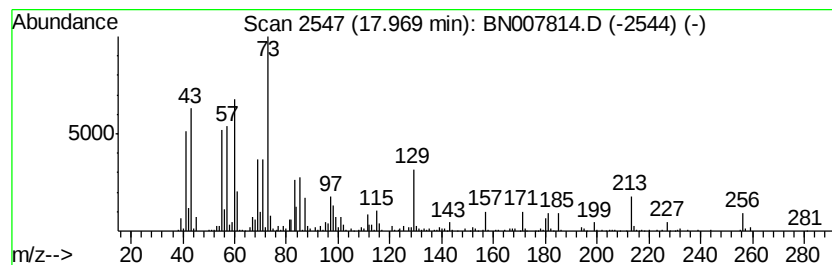
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN091619MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 13 n-Hexadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.		
17.97	2.78 ng/ul	1086420	Phenanthrene-d10		17.06		
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	98
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
4			Pentadecanoic acid	242	C15H30O2	001002-84-2	76
5			Tridecanoic acid	214	C13H26O2	000638-53-9	76



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN092319\
 Data File : BN007814.D
 Acq On : 23 Sep 2019 22:41
 Operator : HP/JU
 Sample : K4895-06
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BFDF9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN091619MA.M
 Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propanoic acid, 1...	3.69	3.9	ng/ul	649465	1	7.69	3337370	20.0
Propanoic acid, 2...	4.22	8.4	ng/ul	1400630	1	7.69	3337370	20.0
Propanoic acid, p...	4.39	12.2	ng/ul	2034280	1	7.69	3337370	20.0
Butanoic acid, 1-...	4.83	15.5	ng/ul	2589340	1	7.69	3337370	20.0
Butanoic acid, pr...	5.68	2.1	ng/ul	346444	1	7.69	3337370	20.0
Benzene, 1,2,3-tr...	7.35	4.8	ng/ul	801789	1	7.69	3337370	20.0
Benzene, 1-ethyl-...	7.80	2.8	ng/ul	465579	1	7.69	3337370	20.0
Indane	8.06	5.7	ng/ul	946306	1	7.69	3337370	20.0
Benzene, 2-etheny...	9.87	3.1	ng/ul	745702	2	10.46	4820100	20.0
Naphthalene, 1-me...	12.35	9.3	ng/ul	2228850	2	10.46	4820100	20.0
Naphthalene, 2,3-...	13.64	2.2	ng/ul	714350	3	14.32	6575960	20.0
n-Hexadecanoic acid	17.97	2.8	ng/ul	1086420	4	17.06	7814280	20.0