

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN092319\
 Data File : BN007810.D
 Acq On : 23 Sep 2019 20:18
 Operator : HP/JU
 Sample : K4895-01
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BFD5

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	3.281	46	50	66	rVB	5662163	6798040	35.75%	2.538%
2	3.693	116	120	127	rVB	639254	754423	3.97%	0.282%
3	4.217	204	209	216	rBV	1405087	1723907	9.07%	0.644%
4	4.281	216	220	234	rVB	182545	346545	1.82%	0.129%
5	4.387	234	238	246	rBV	2009143	2603973	13.69%	0.972%
6	4.475	246	253	262	rVB3	294810	603525	3.17%	0.225%
7	4.834	307	314	323	rVV	2441196	3303209	17.37%	1.233%
8	5.375	401	406	423	rVB	601100	1234305	6.49%	0.461%
9	5.675	452	457	462	rVV	376070	537696	2.83%	0.201%
10	5.734	462	467	480	rVB	1208450	1845241	9.70%	0.689%
11	6.799	642	648	653	rVV	1508345	2212735	11.64%	0.826%
12	6.858	653	658	665	rVV2	1143733	1947669	10.24%	0.727%
13	6.928	665	670	678	rVB	1714791	2494090	13.12%	0.931%
14	7.028	681	687	693	rBV	1585059	2555374	13.44%	0.954%
15	7.093	693	698	706	rVB	1338620	2013257	10.59%	0.752%
16	7.222	713	720	730	rBV	1858583	2941444	15.47%	1.098%
17	7.346	734	741	748	rVB	2740767	4090241	21.51%	1.527%
18	7.687	793	799	804	rBV	2054377	3216710	16.92%	1.201%
19	7.734	804	807	814	rVV2	262523	506642	2.66%	0.189%
20	7.805	814	819	829	rVV	2164269	3543152	18.63%	1.323%
21	8.058	843	862	869	rVB2	897513	2120898	11.15%	0.792%
22	8.187	878	884	888	rBV	365233	566538	2.98%	0.212%
23	8.240	888	893	899	rVV	607185	918965	4.83%	0.343%
24	8.328	902	908	913	rVV	1391043	2078044	10.93%	0.776%
25	8.387	913	918	925	rVV	1559107	2485929	13.07%	0.928%
26	8.493	931	936	945	rVV	412486	675156	3.55%	0.252%
27	8.640	953	961	966	rBV	611971	942589	4.96%	0.352%
28	8.693	966	970	977	rVB	676737	1010850	5.32%	0.377%
29	8.793	977	987	990	rBV	1570267	2593254	13.64%	0.968%
30	8.828	990	993	997	rVV	1988360	3228664	16.98%	1.206%
31	8.863	997	999	1007	rVV2	774054	1286148	6.76%	0.480%
32	9.040	1020	1029	1036	rVB4	130469	326224	1.72%	0.122%
33	9.122	1036	1043	1048	rBV	482141	825501	4.34%	0.308%
34	9.310	1069	1075	1079	rVV	929260	1428104	7.51%	0.533%

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35	9.369	1079	1085	1092	rVB	1462705	2275890	11.97%	0.850%
36	9.546	1108	1115	1123	rBV	1511167	2644503	13.91%	0.987%
37	9.681	1132	1138	1141	rVV4	206288	441100	2.32%	0.165%
38	9.722	1141	1145	1157	rVB	855926	1419883	7.47%	0.530%
39	9.869	1159	1170	1178	rBV3	2174812	4189321	22.03%	1.564%
40	9.946	1178	1183	1192	rVV3	225835	501109	2.64%	0.187%
41	10.081	1193	1206	1216	rVB2	2730629	5336873	28.07%	1.993%
42	10.457	1261	1270	1275	rVV	3068519	5561001	29.25%	2.076%
43	10.510	1275	1279	1287	rVV	2323674	4075844	21.44%	1.522%
44	10.587	1287	1292	1305	rVV	2433487	4418715	23.24%	1.650%
45	11.122	1380	1383	1389	rVB	242192	359584	1.89%	0.134%
46	11.381	1420	1427	1435	rBV	370075	640109	3.37%	0.239%
47	11.569	1452	1459	1464	rVV	321846	598551	3.15%	0.223%
48	11.657	1468	1474	1483	rVB	281995	585735	3.08%	0.219%
49	11.851	1502	1507	1516	rVV2	254695	443245	2.33%	0.165%
50	12.016	1530	1535	1545	rVV2	323927	631925	3.32%	0.236%
51	12.122	1545	1553	1563	rVB	3097242	5046293	26.54%	1.884%
52	12.346	1580	1591	1602	rVV	4819993	7810538	41.08%	2.916%
53	13.146	1722	1727	1733	rBV	242450	363452	1.91%	0.136%
54	13.228	1735	1741	1749	rVV2	171178	428134	2.25%	0.160%
55	13.345	1755	1761	1763	rVV2	274313	489149	2.57%	0.183%
56	13.369	1763	1765	1771	rVB	284701	378185	1.99%	0.141%
57	13.493	1773	1786	1793	rBV2	617460	1603870	8.43%	0.599%
58	13.640	1801	1811	1816	rBV	1220785	2207897	11.61%	0.824%
59	13.693	1816	1820	1822	rVV	711874	1059069	5.57%	0.395%
60	13.728	1822	1826	1831	rVV	5754837	7751631	40.77%	2.894%
61	13.775	1831	1834	1840	rVV	643288	870666	4.58%	0.325%
62	13.875	1842	1851	1854	rVV	691718	1025603	5.39%	0.383%
63	13.904	1854	1856	1862	rVB	320884	435219	2.29%	0.163%
64	14.004	1867	1873	1878	rBV2	5799515	8891277	46.76%	3.320%
65	14.316	1916	1926	1932	rBV	4468589	6758964	35.55%	2.524%
66	14.381	1932	1937	1946	rVV	3359033	4834958	25.43%	1.805%
67	14.510	1954	1959	1971	rVV2	504169	1069892	5.63%	0.399%
68	14.716	1984	1994	2003	rVV2	581100	1123859	5.91%	0.420%
69	14.798	2003	2008	2015	rVB	247355	343481	1.81%	0.128%
70	14.940	2024	2032	2037	rBV2	246205	481907	2.53%	0.180%
71	15.310	2082	2095	2101	rVV	8092794	11994566	63.08%	4.478%

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72	15.363	2101	2104	2109	rVV2	637640	982847	5.17%	0.367%
73	15.428	2109	2115	2119	rVV	2147917	2915036	15.33%	1.088%
74	15.492	2123	2126	2129	rVV	313753	438344	2.31%	0.164%
75	15.528	2129	2132	2138	rVV4	193863	456296	2.40%	0.170%
76	15.775	2167	2174	2177	rVV2	191870	354709	1.87%	0.132%
77	16.369	2266	2275	2280	rBV2	255425	592032	3.11%	0.221%
78	16.881	2354	2362	2369	rVB	348381	571512	3.01%	0.213%
79	17.063	2387	2393	2396	rBV2	5335300	7868278	41.38%	2.938%
80	17.104	2396	2400	2404	rVB	3342902	4498829	23.66%	1.680%
81	17.157	2404	2409	2414	rBV2	9556260	13941384	73.32%	5.205%
82	17.457	2456	2460	2467	rVB	459734	632797	3.33%	0.236%
83	17.686	2494	2499	2504	rVB	531683	700684	3.68%	0.262%
84	17.933	2537	2541	2544	rBV	330292	444492	2.34%	0.166%
85	17.975	2544	2548	2559	rVV3	1148022	2014259	10.59%	0.752%
86	18.133	2569	2575	2578	rVV2	566835	1030295	5.42%	0.385%
87	18.163	2578	2580	2587	rVB	272796	372989	1.96%	0.139%
88	19.122	2739	2743	2749	rVB	717803	969063	5.10%	0.362%
89	19.263	2763	2767	2774	rBV3	517804	803479	4.23%	0.300%
90	19.457	2794	2800	2817	rVB	13077279	18884768	99.32%	7.051%
91	20.992	3057	3061	3073	rBV2	860893	1589671	8.36%	0.594%
92	21.257	3100	3106	3116	rBV	7966020	10039167	52.80%	3.748%
93	21.869	3206	3210	3222	rBV	268754	437760	2.30%	0.163%
94	22.327	3284	3288	3293	rVB	356395	451879	2.38%	0.169%
95	22.833	3369	3374	3387	rVB	514752	831443	4.37%	0.310%
96	23.333	3451	3459	3466	rBV	10615057	19014717	100.00%	7.100%
97	23.392	3466	3469	3475	rVV	539701	841920	4.43%	0.314%
98	23.474	3475	3483	3498	rVB	5439192	10147046	53.36%	3.789%
99	24.033	3574	3578	3584	rVB	374416	626888	3.30%	0.234%
100	24.768	3698	3703	3709	rBV2	261392	522878	2.75%	0.195%

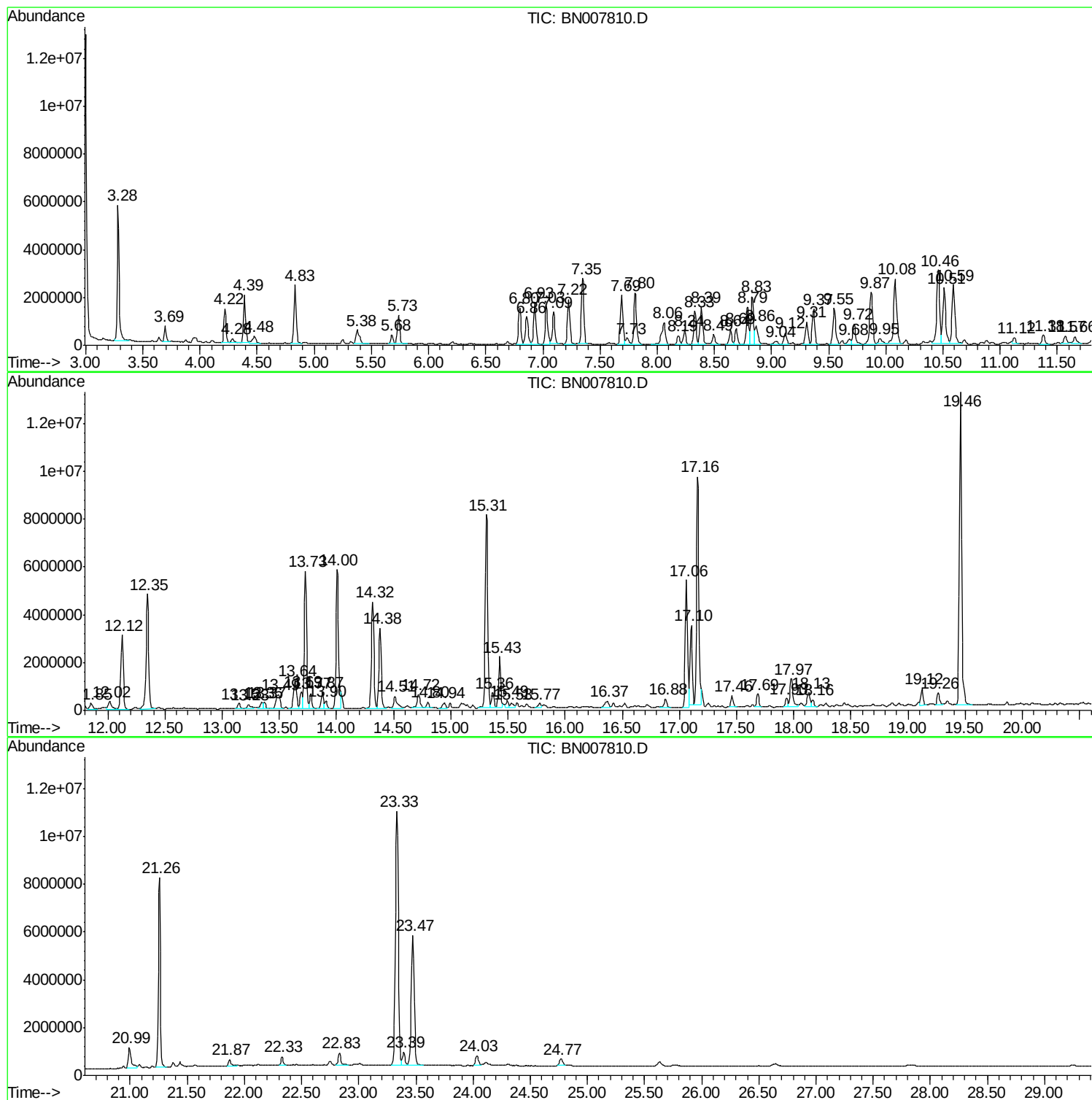
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN091619MA.M
Quant Title : SVOA CALIBRATION

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



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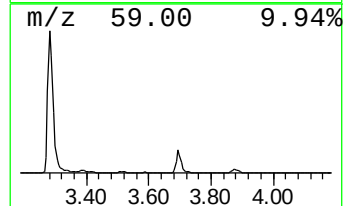
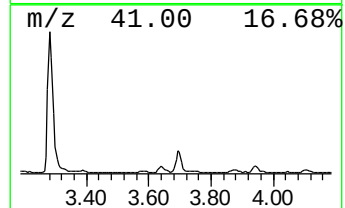
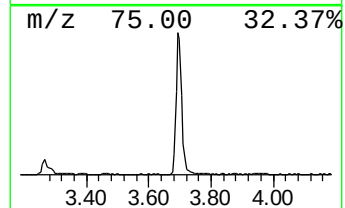
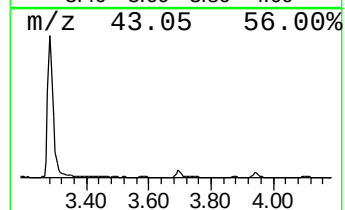
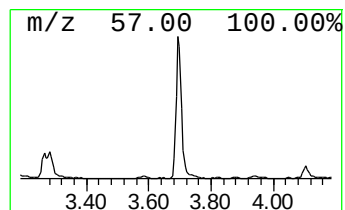
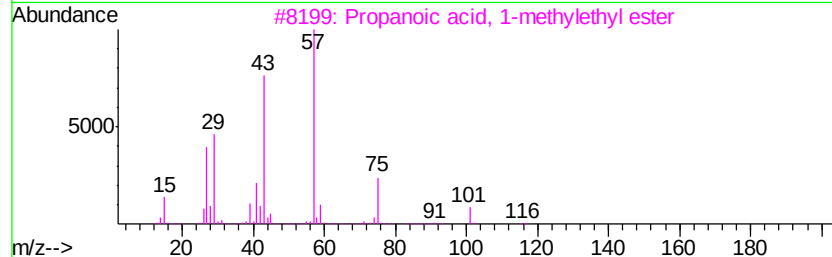
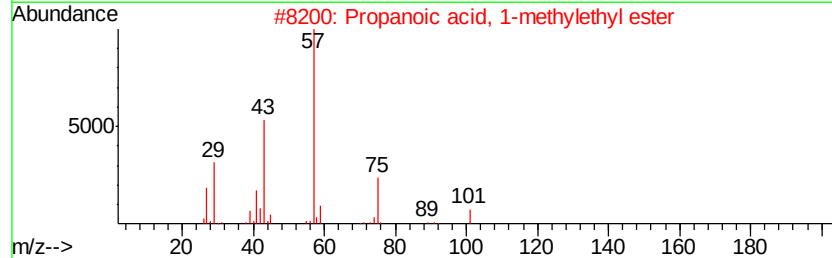
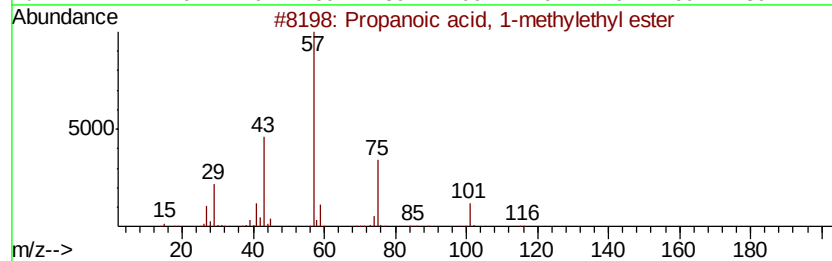
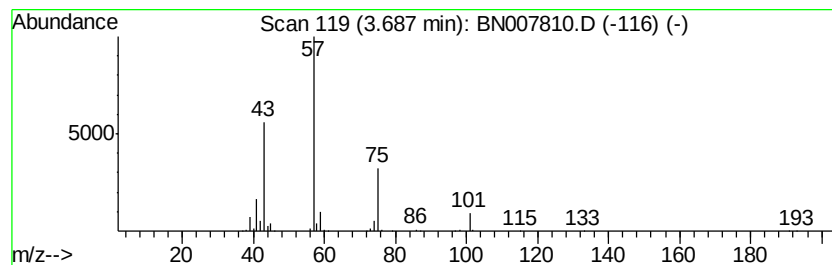
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN091619MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Propanoic acid, 1-methyleth... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.69	4.69 ng/ul	754423	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	78
2		Propanoic acid, 1-methylethyl ester	116	C6H12O2	000637-78-5	78
3		Propanoic acid, 1-methylethyl ester	116	C6H12O2	000637-78-5	74
4		Propanoic acid, 1-methylethyl ester	116	C6H12O2	000637-78-5	53
5		Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	53



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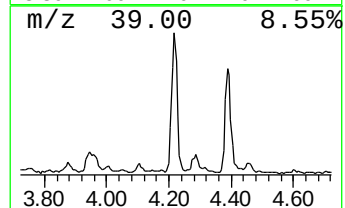
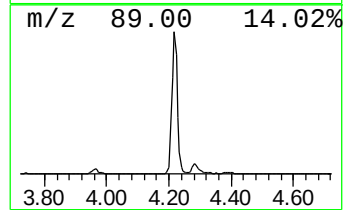
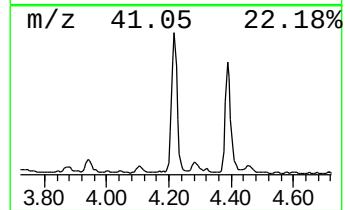
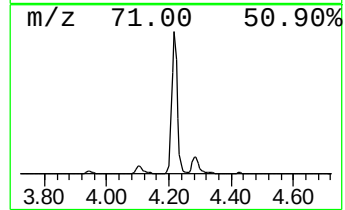
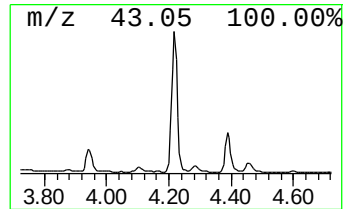
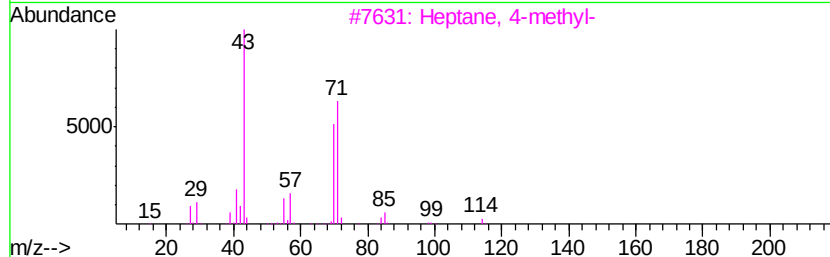
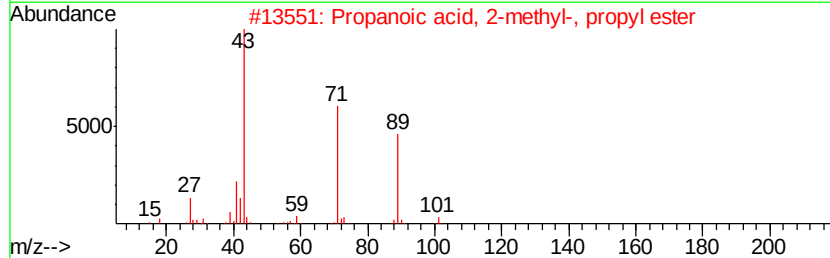
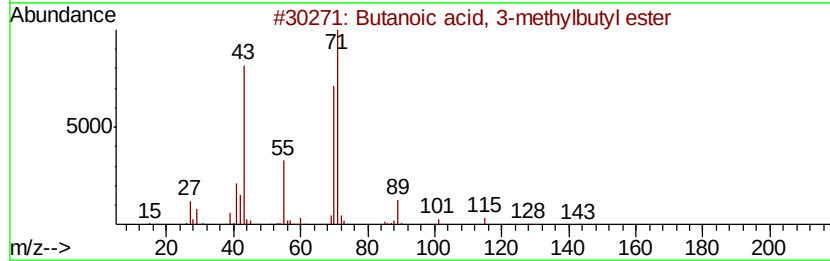
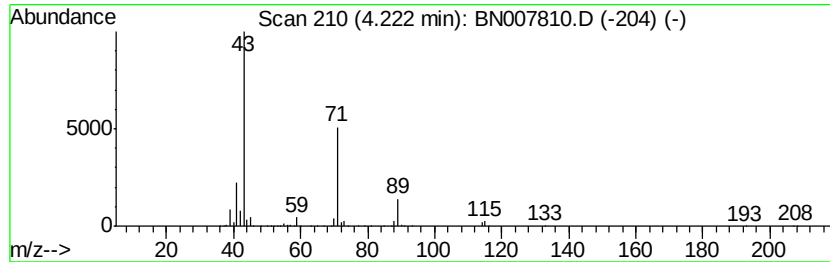
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Butanoic acid, 3-methylbuty... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.22	10.72 ng/ul	1723910	1,4-Dichlorobenzene-d4	7.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butanoic acid, 3-methylbutyl ester	158	C9H18O2	000106-27-4	53
2		Propanoic acid, 2-methyl-, propyl...	130	C7H14O2	000644-49-5	53
3		Heptane, 4-methyl-	114	C8H18	000589-53-7	50
4		Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	50
5		Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	50



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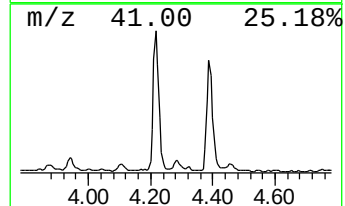
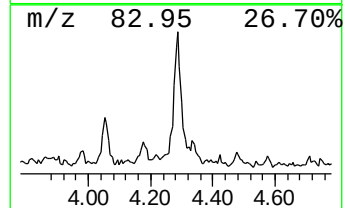
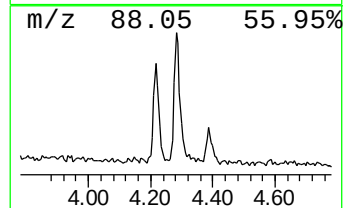
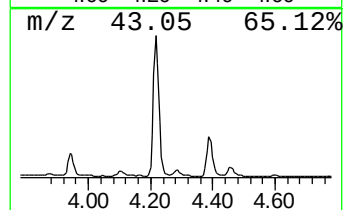
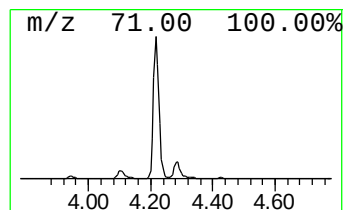
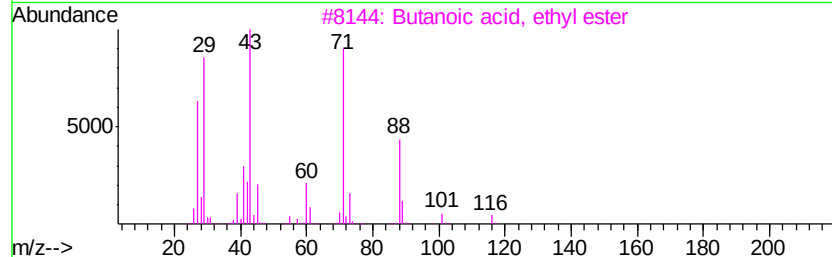
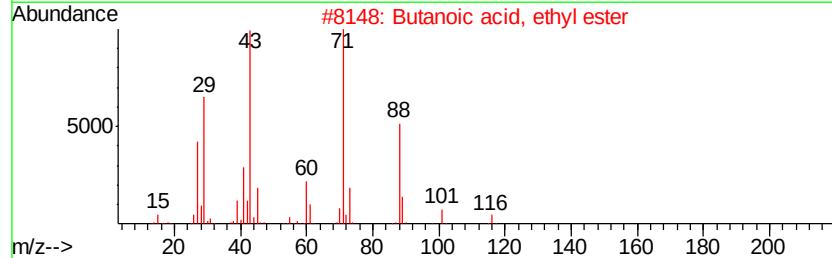
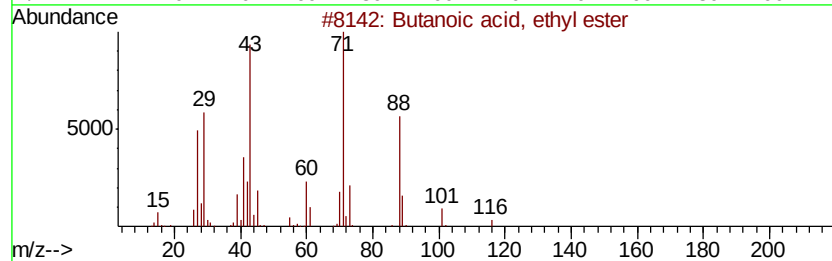
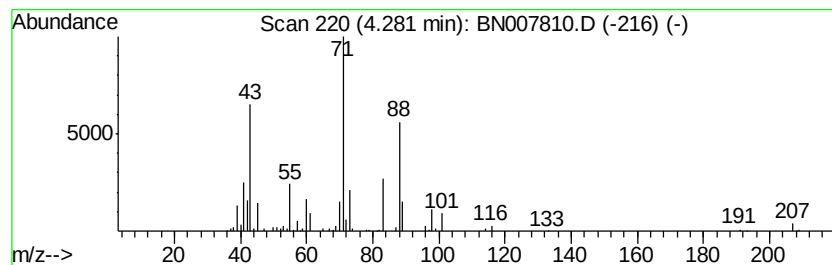
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Peak Number 3 Butanoic acid, ethyl ester Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.28	2.15 ng/ul	346545	1,4-Dichlorobenzene-d4	7.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butanoic acid, ethyl ester	116	C6H12O2	000105-54-4	96
2		Butanoic acid, ethyl ester	116	C6H12O2	000105-54-4	81
3		Butanoic acid, ethyl ester	116	C6H12O2	000105-54-4	81
4		Butanoic acid, ethyl ester	116	C6H12O2	000105-54-4	64
5		Butanoic acid, propyl ester	130	C7H14O2	000105-66-8	43



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Sample : K4895-01
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ClientSampleId :
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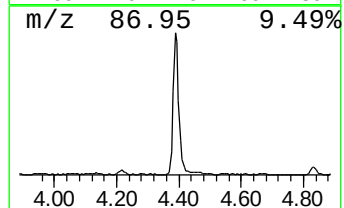
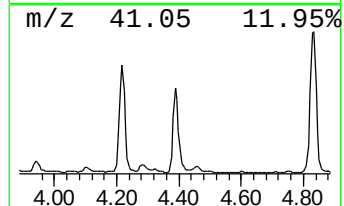
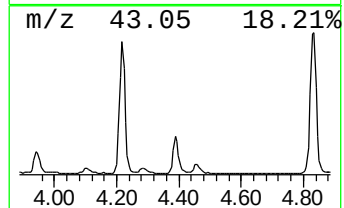
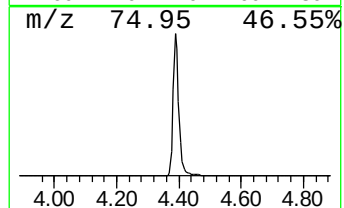
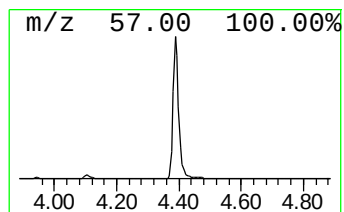
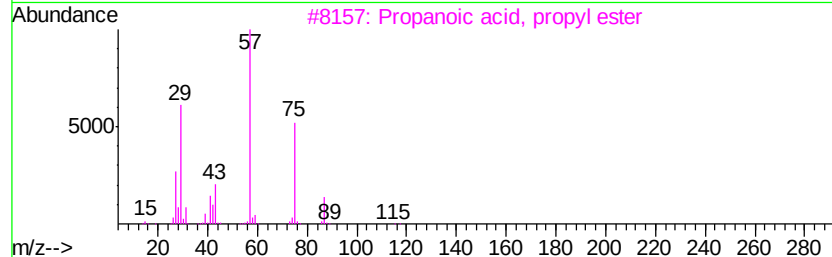
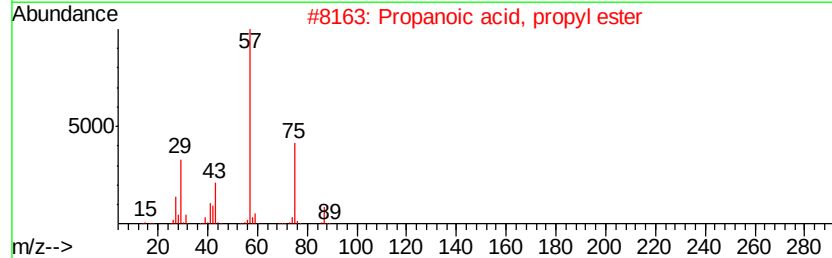
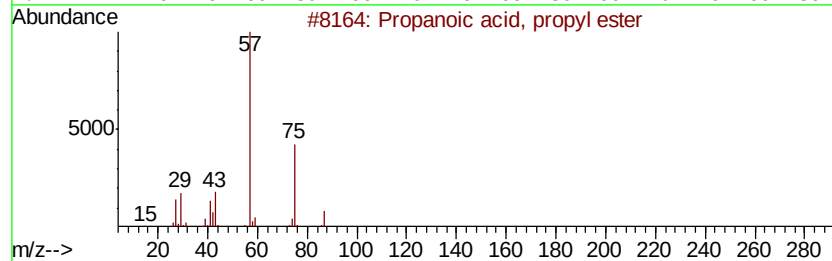
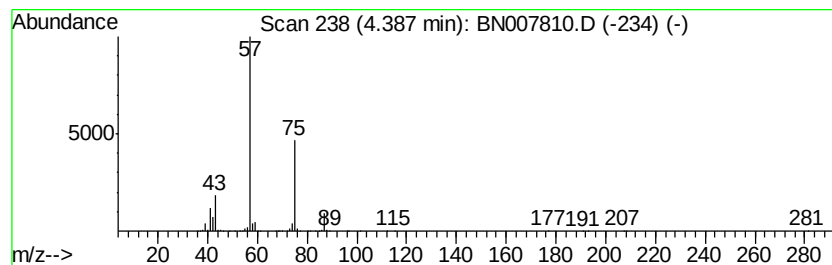
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Propanoic acid, propyl ester Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.39	16.19 ng/ul	2603970	1,4-Dichlorobenzene-d4	7.69

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN092319\
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Instrument :
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ClientSampleId :
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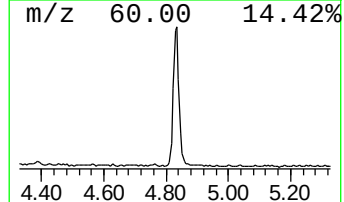
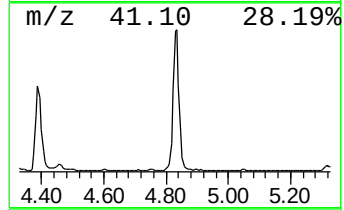
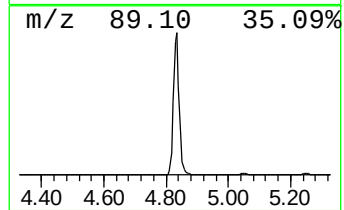
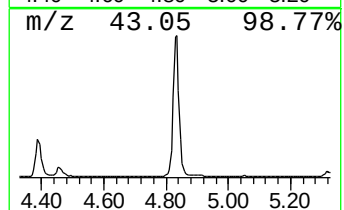
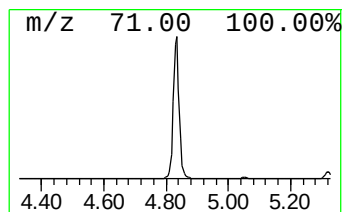
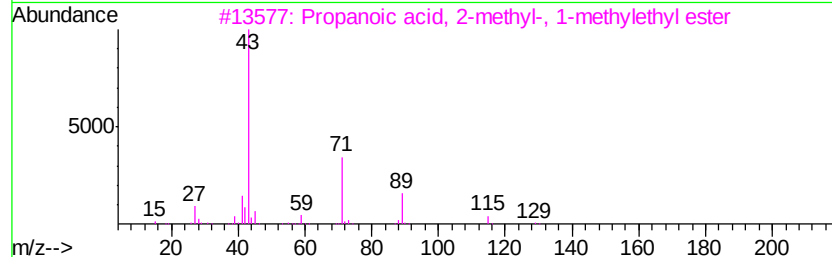
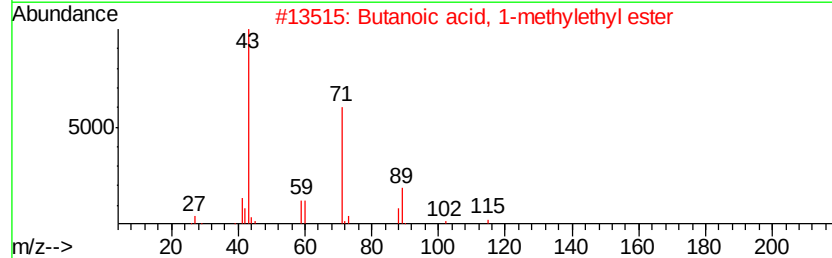
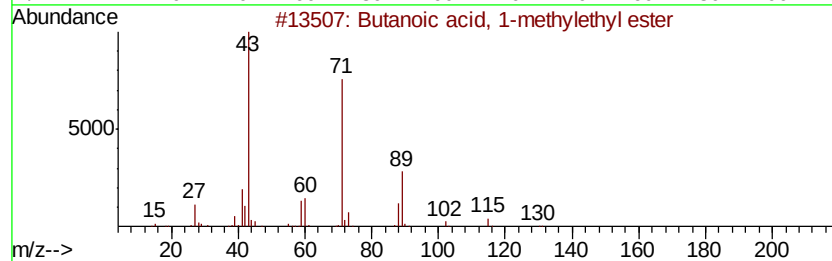
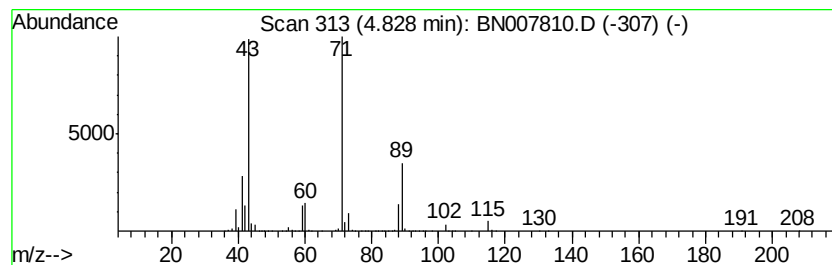
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Butanoic acid, 1-methylethy... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.83	20.54 ng/ul	3303210	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		Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	72
4		Butanoic acid, pentyl ester	158	C9H18O2	000540-18-1	64
5		Butanoic acid, 1-methylethyl ester	130	C7H14O2	000638-11-9	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN092319\
Data File : BN007810.D
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Sample : K4895-01
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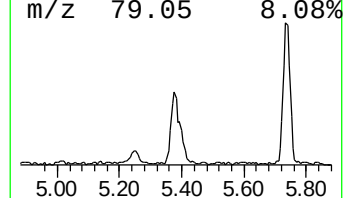
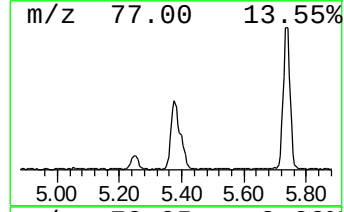
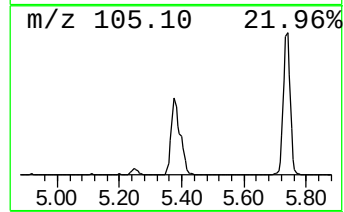
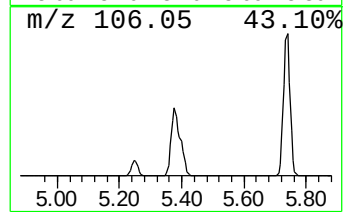
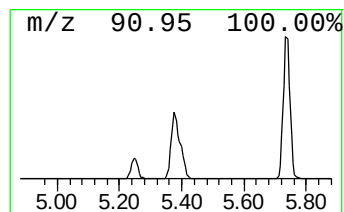
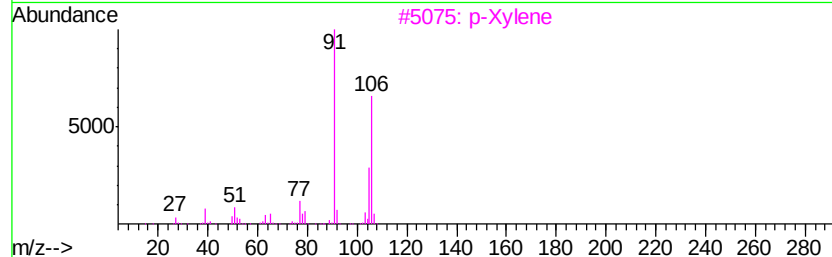
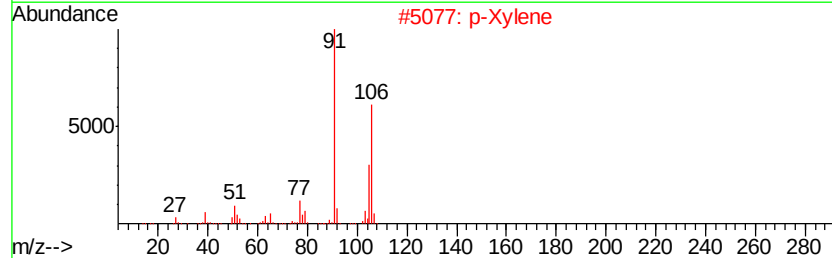
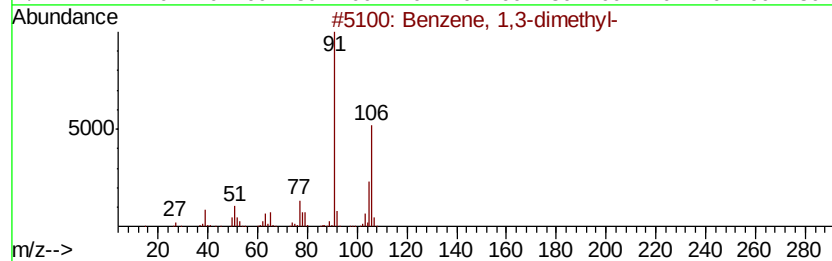
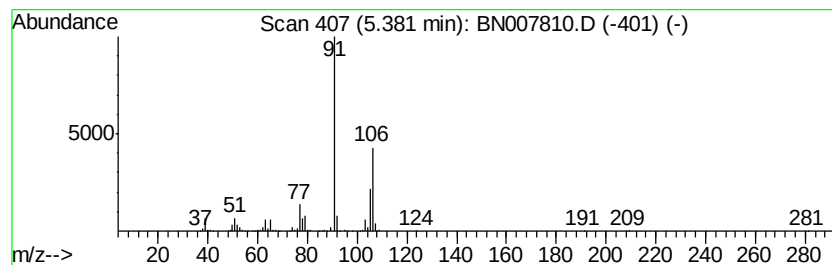
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzene, 1,3-dimethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.38	7.67 ng/ul	1234310	1,4-Dichlorobenzene-d4	7.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
2		p-Xylene	106	C8H10	000106-42-3	95
3		p-Xylene	106	C8H10	000106-42-3	95
4		p-Xylene	106	C8H10	000106-42-3	95
5		o-Xylene	106	C8H10	000095-47-6	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN092319\
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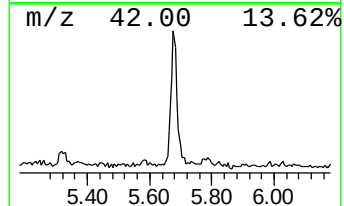
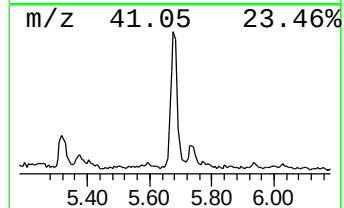
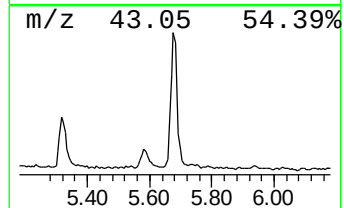
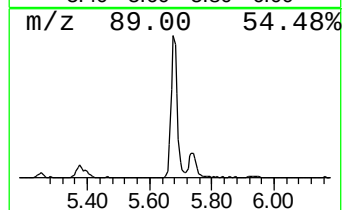
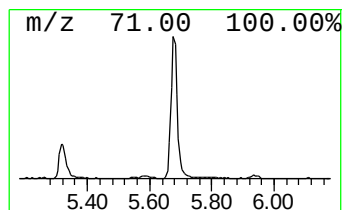
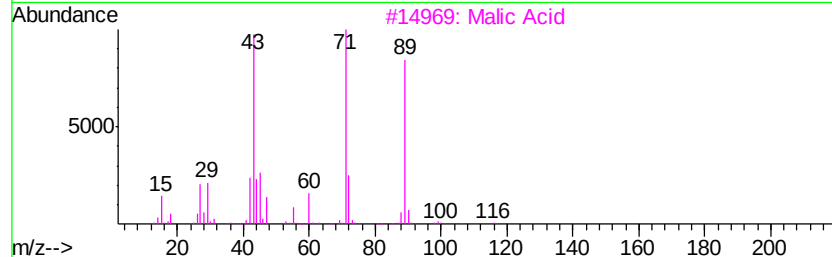
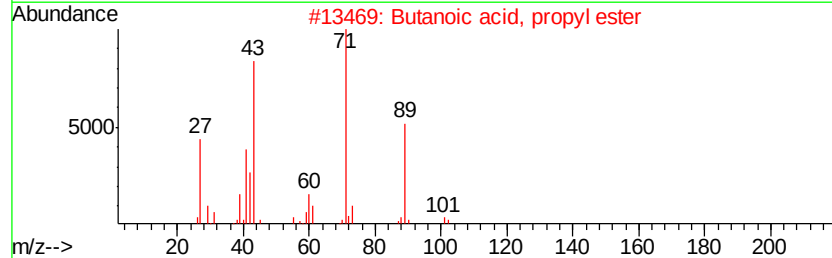
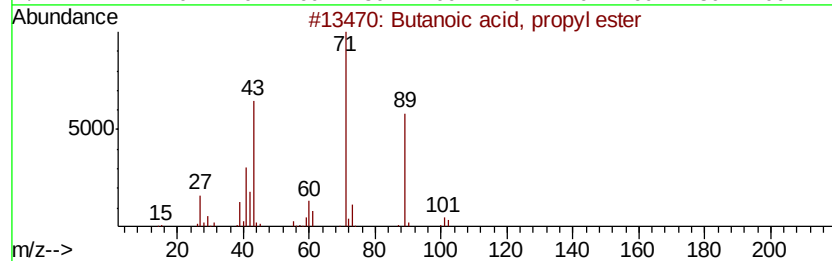
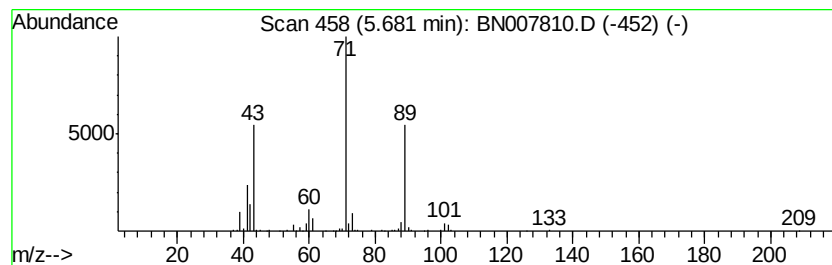
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Butanoic acid, propyl ester Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.68	3.34 ng/ul	537696	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	72
3		Malic Acid	134	C4H6O5	006915-15-7	64
4		Butanoic acid, hexyl ester	172	C10H20O2	002639-63-6	64
5		Butanoic acid, butyl ester	144	C8H16O2	000109-21-7	64



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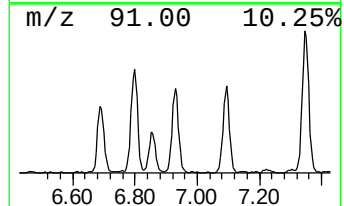
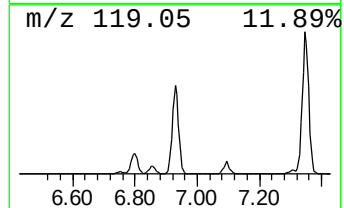
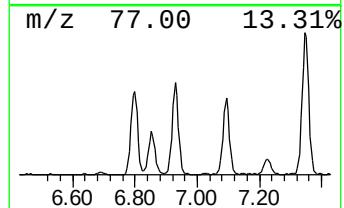
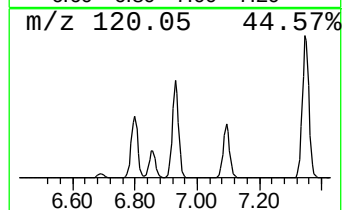
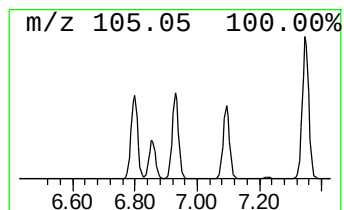
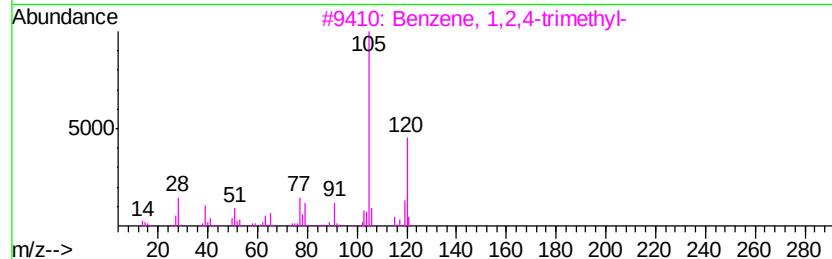
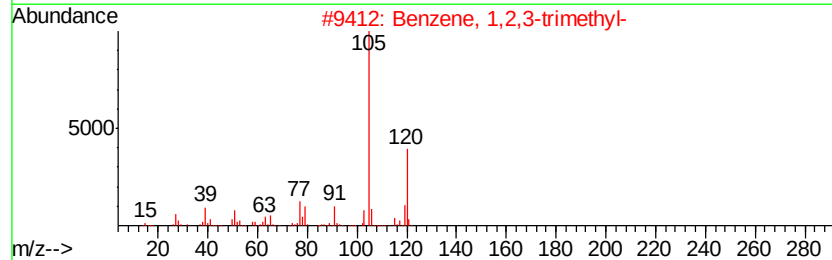
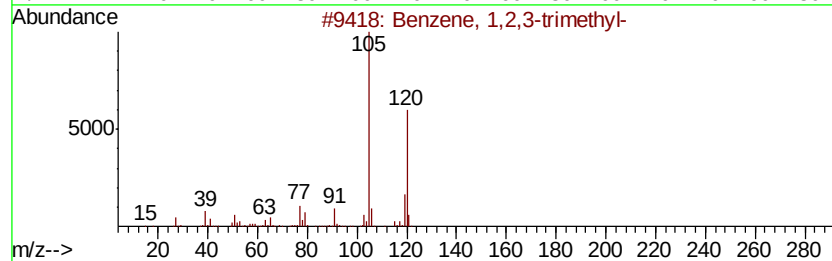
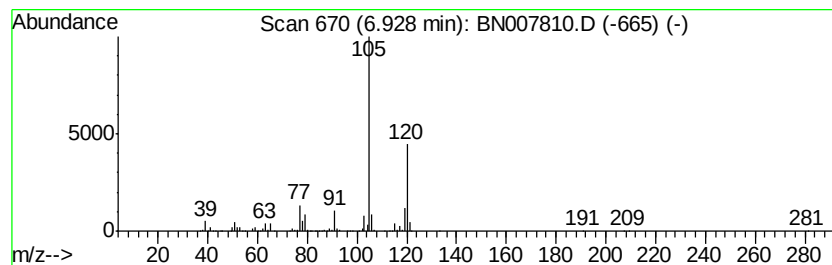
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Benzene, 1,2,3-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.93	15.51 ng/ul	2494090	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	95
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
5		Mesitylene	120	C9H12	000108-67-8	91



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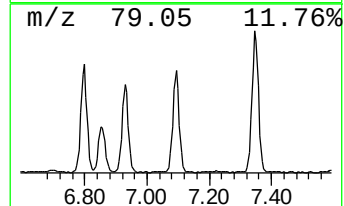
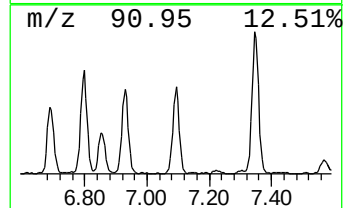
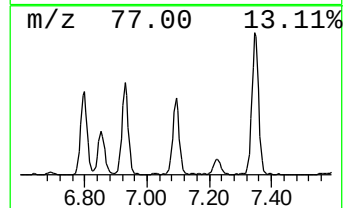
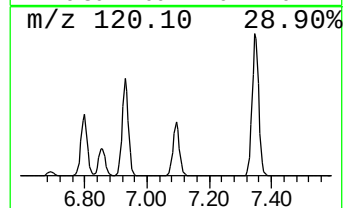
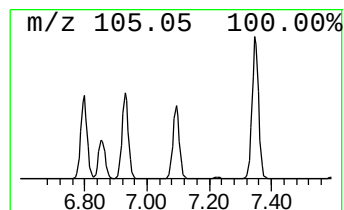
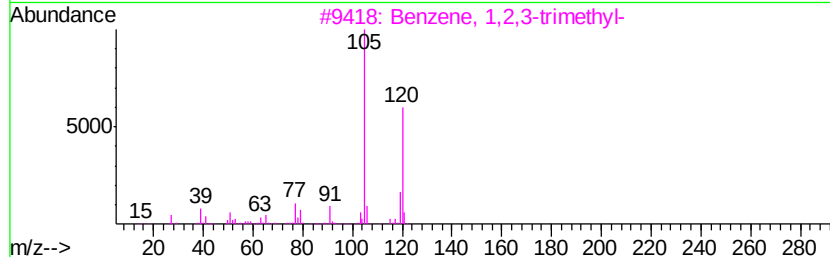
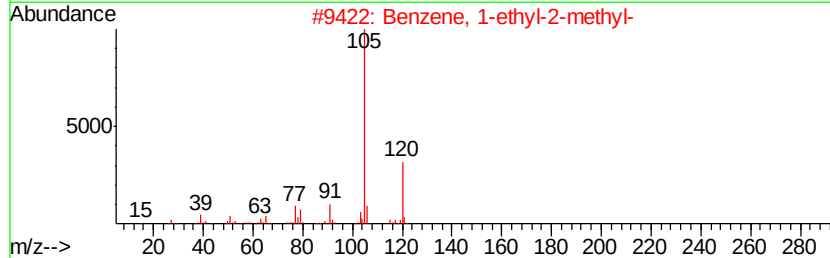
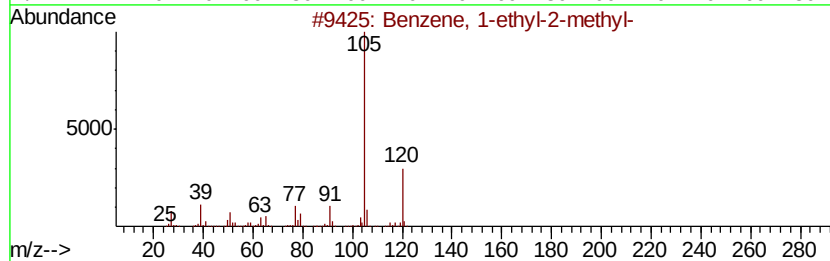
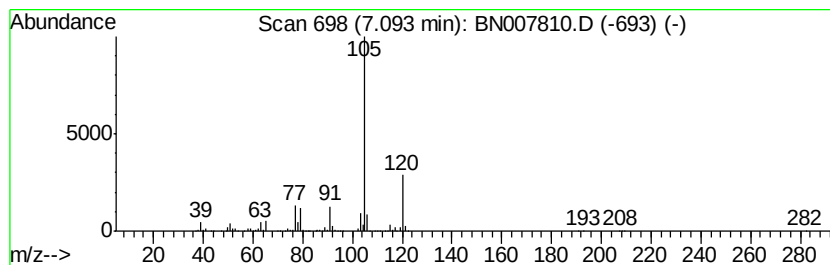
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Benzene, 1-ethyl-2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.09	12.52 ng/ul	2013260	1,4-Dichlorobenzene-d4	7.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
4		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91
5		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN092319\
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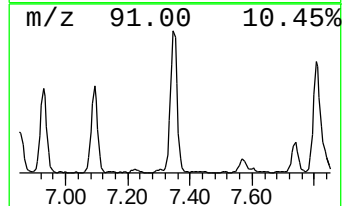
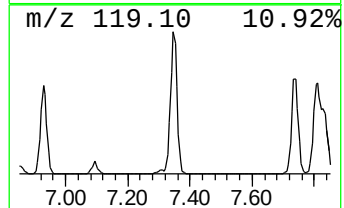
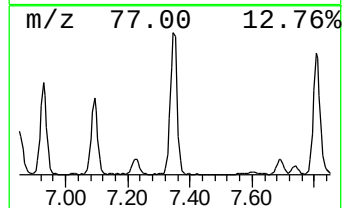
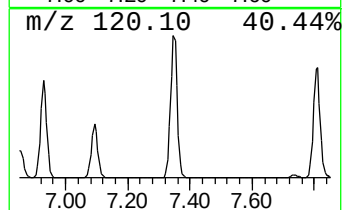
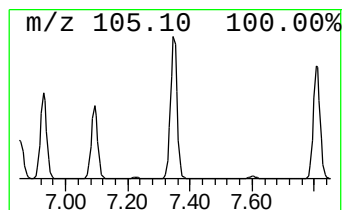
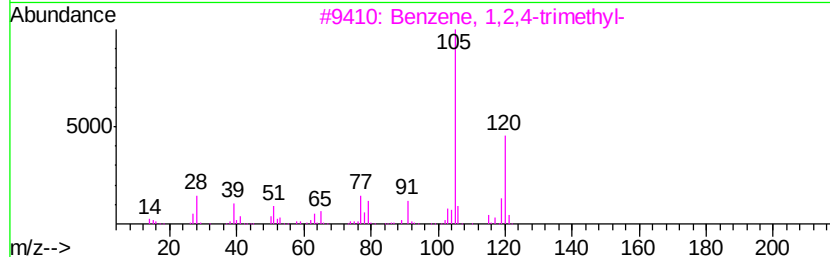
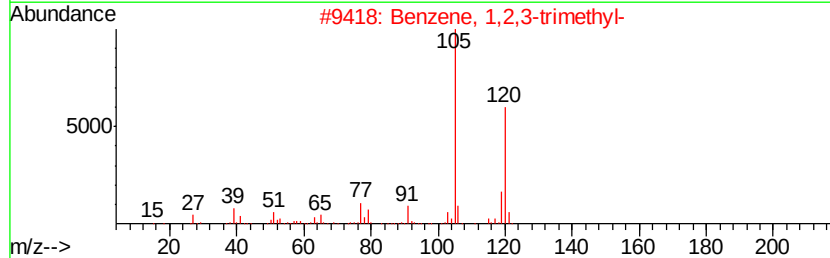
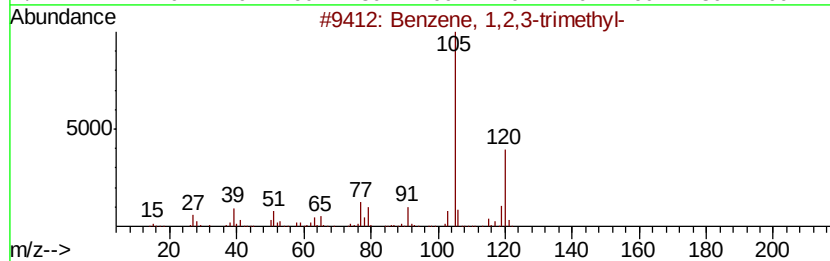
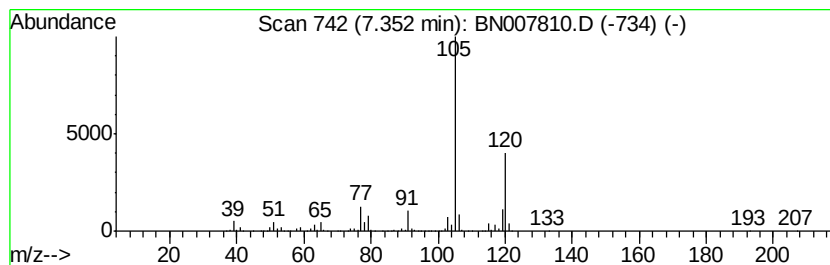
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 Benzene, 1,2,4-trimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.35	25.43 ng/ul	4090240	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	95
2		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN092319\
Data File : BN007810.D
Acq On : 23 Sep 2019 20:18
Operator : HP/JU
Sample : K4895-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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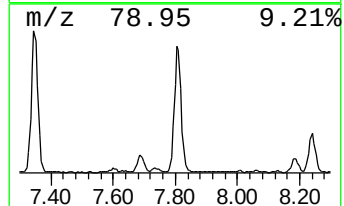
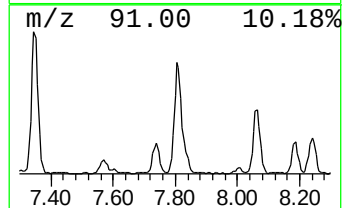
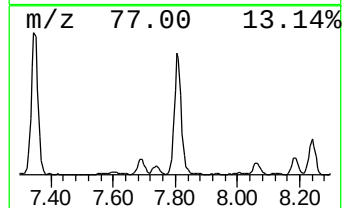
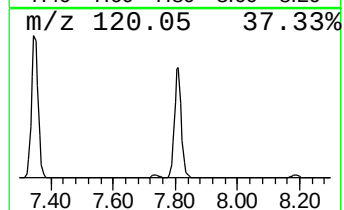
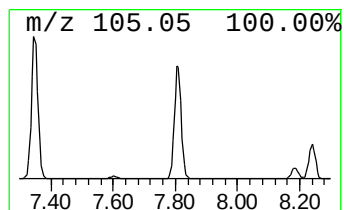
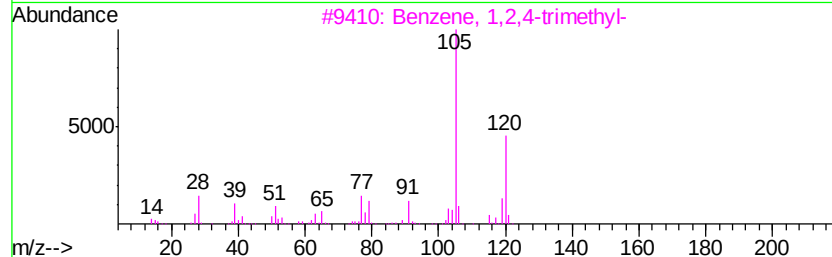
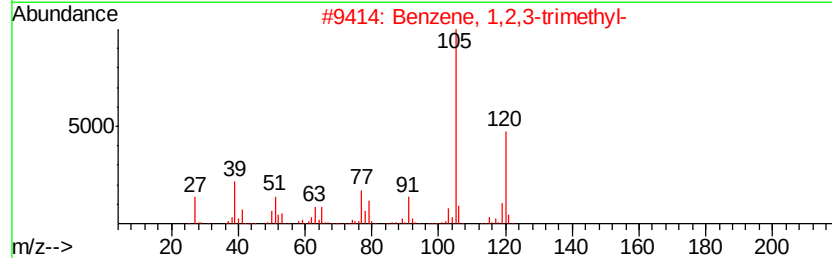
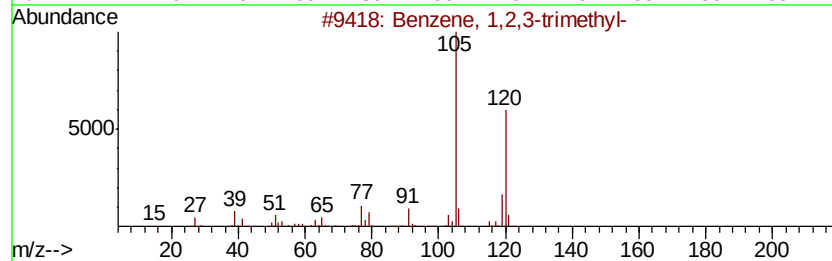
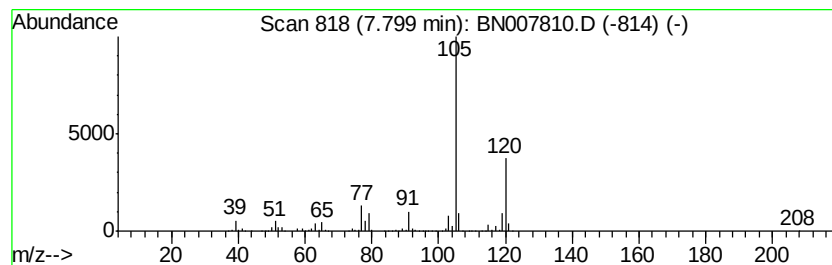
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 Mesitylene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.80	22.03 ng/ul	3543150	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	94
2		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4		Mesitylene	120	C9H12	000108-67-8	91
5		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN092319\
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Acq On : 23 Sep 2019 20:18
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Sample : K4895-01
Misc :
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Instrument :
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ClientSampleId :
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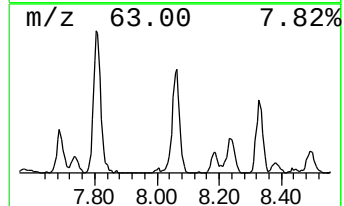
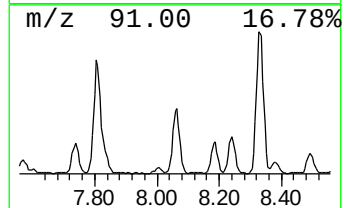
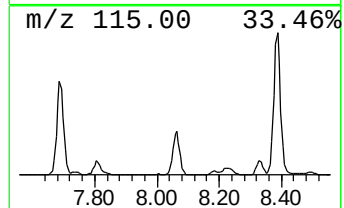
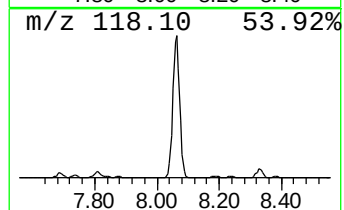
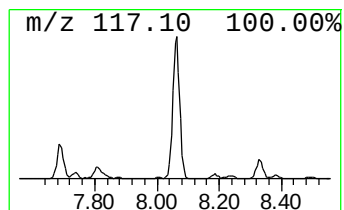
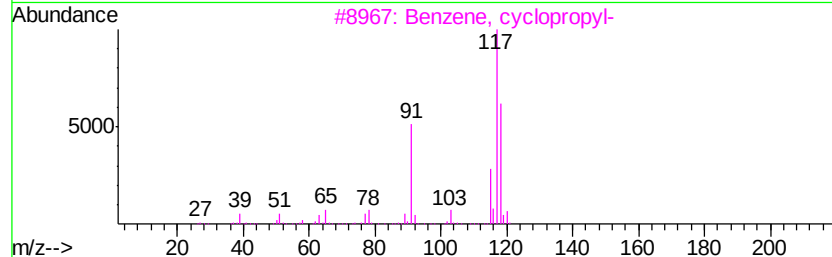
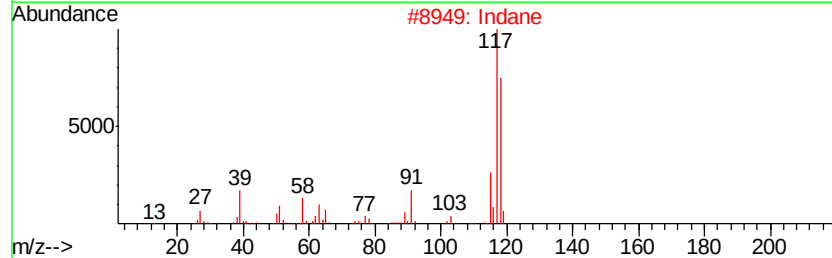
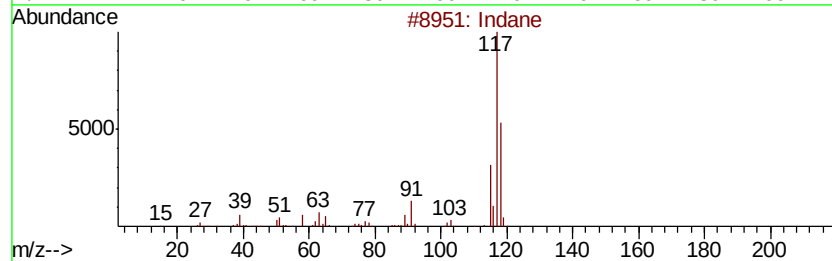
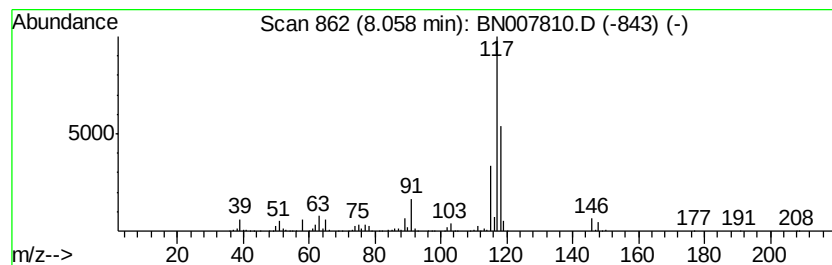
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 Indane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.06	13.19 ng/ul	2120900	1,4-Dichlorobenzene-d4	7.69

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane		118	C9H10	000496-11-7	93
2	Indane		118	C9H10	000496-11-7	90
3	Benzene, cyclopropyl-		118	C9H10	000873-49-4	74
4	Deltacyclene		118	C9H10	007785-10-6	72
5	Benzene, cyclopropyl-		118	C9H10	000873-49-4	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN092319\
Data File : BN007810.D
Acq On : 23 Sep 2019 20:18
Operator : HP/JU
Sample : K4895-01
Misc :
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Instrument :
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ClientSampleId :
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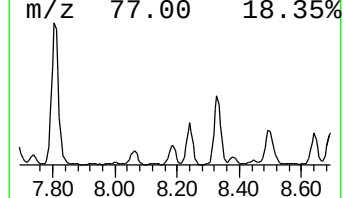
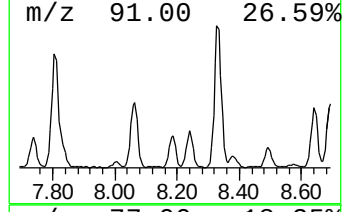
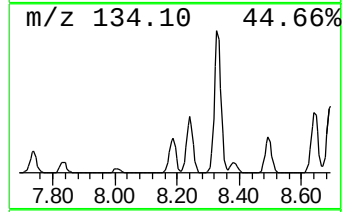
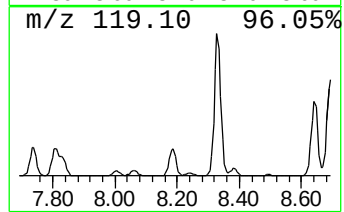
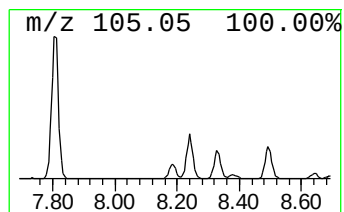
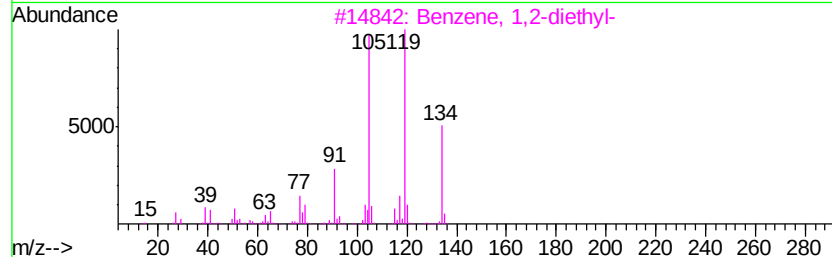
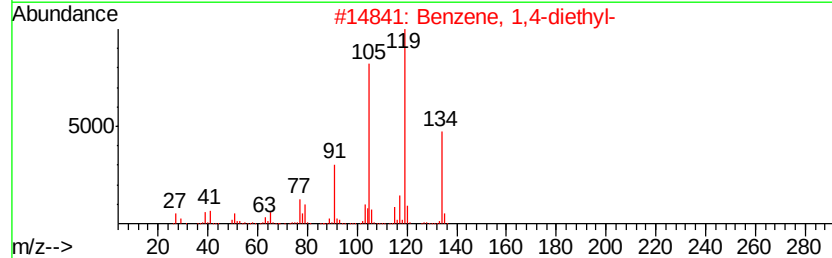
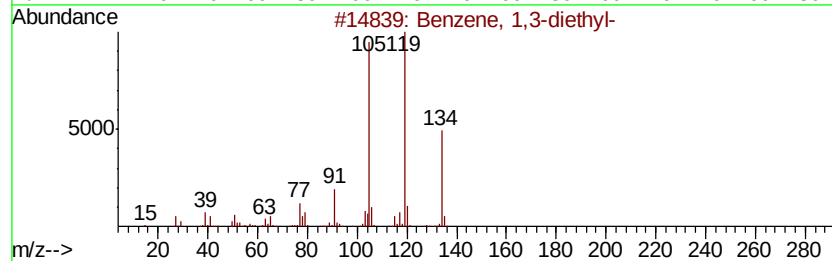
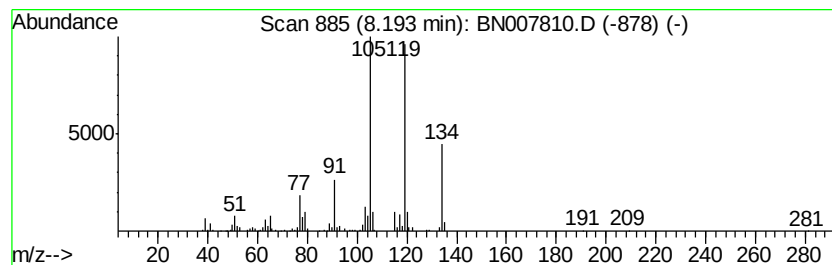
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 Benzene, 1,3-diethyl- Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.19	3.52 ng/ul	566538	1,4-Dichlorobenzene-d4	7.69

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	97
2			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	97
3			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
4			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	95
5			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	95



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN092319\
Data File : BN007810.D
Acq On : 23 Sep 2019 20:18
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ClientSampleId :
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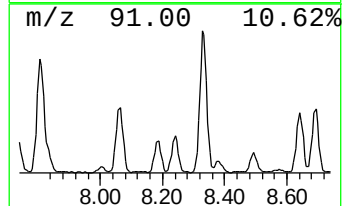
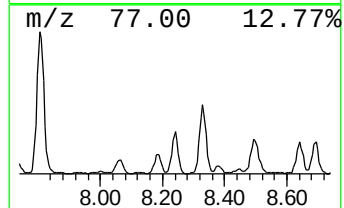
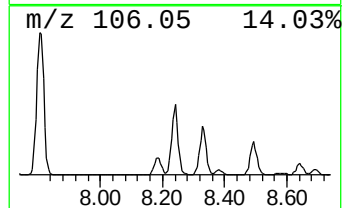
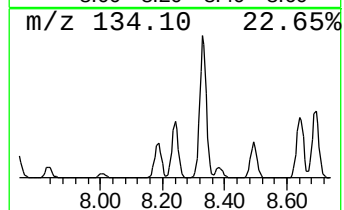
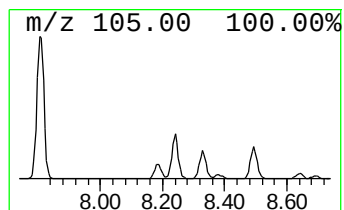
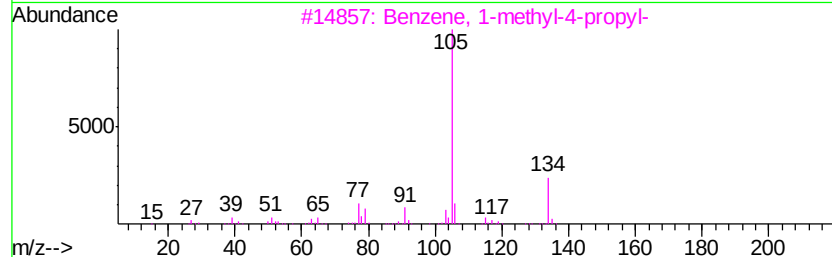
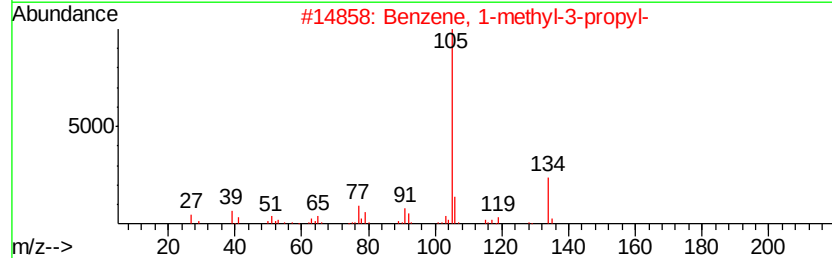
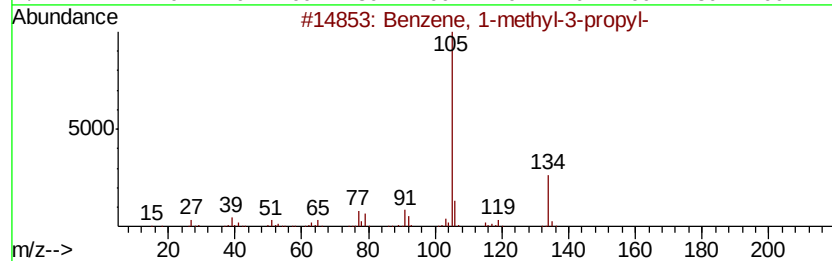
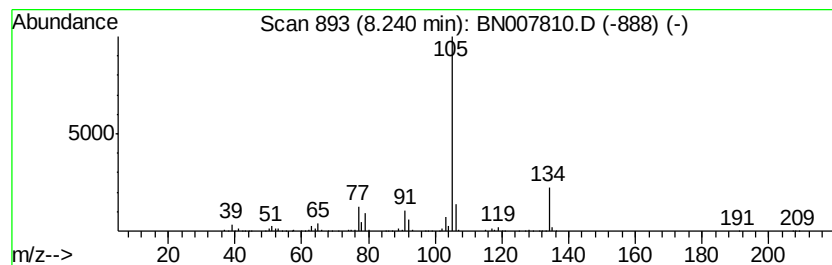
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 Benzene, 1-methyl-3-propyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.24	5.71 ng/ul	918965	1,4-Dichlorobenzene-d4	7.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	91
2		Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	91
3		Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
4		Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	91
5		Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	91



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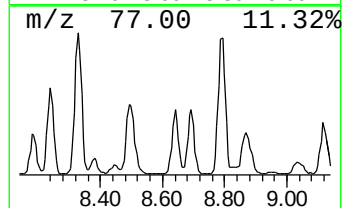
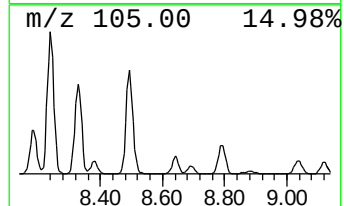
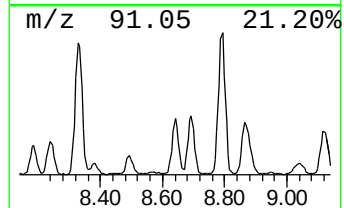
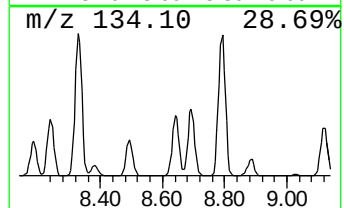
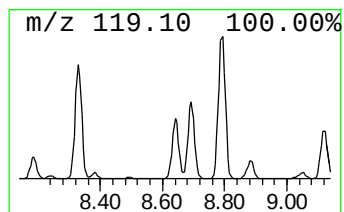
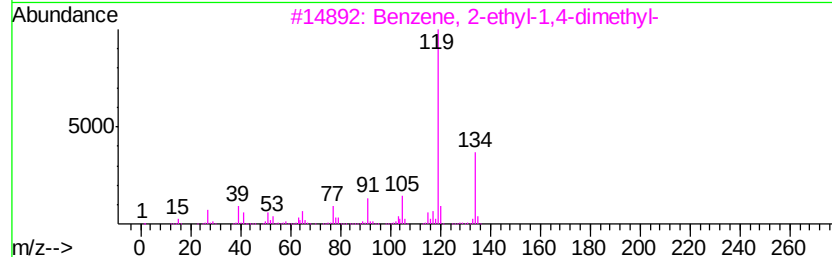
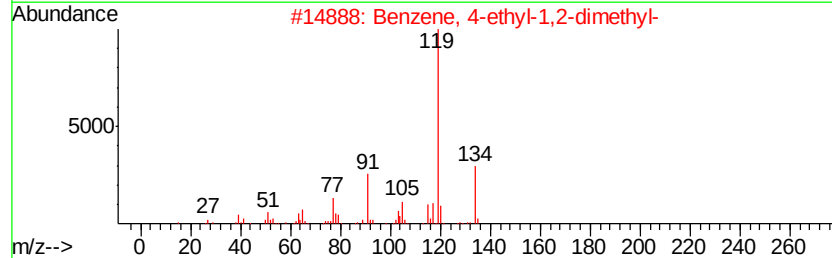
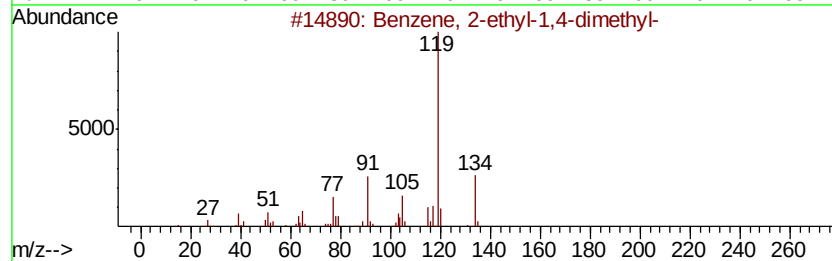
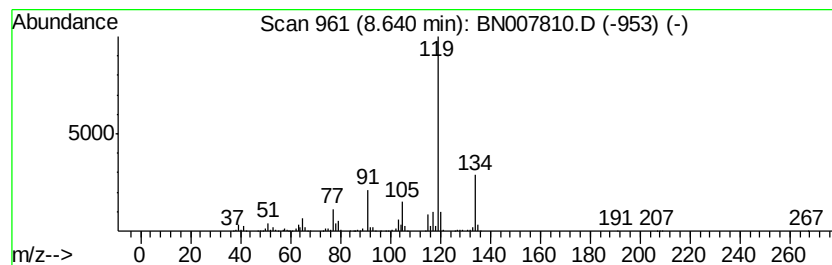
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.64	5.86 ng/ul	942589	1,4-Dichlorobenzene-d4	7.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
3		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
4		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
5		o-Cymene	134	C10H14	000527-84-4	95



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN092319\
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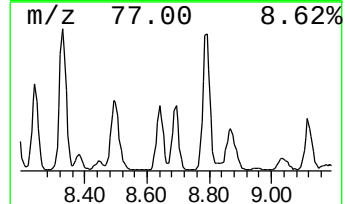
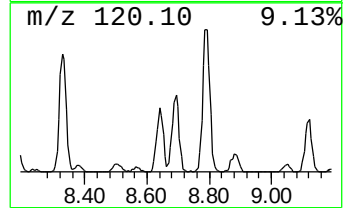
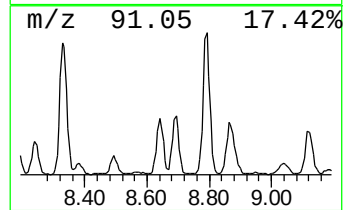
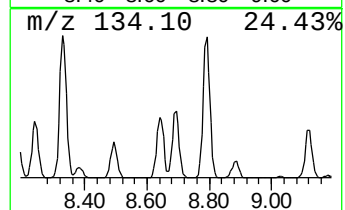
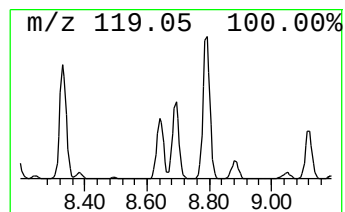
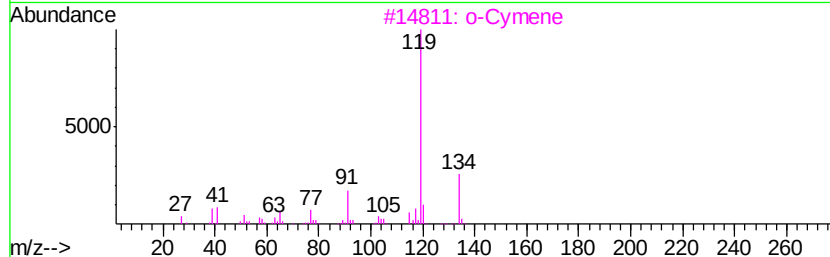
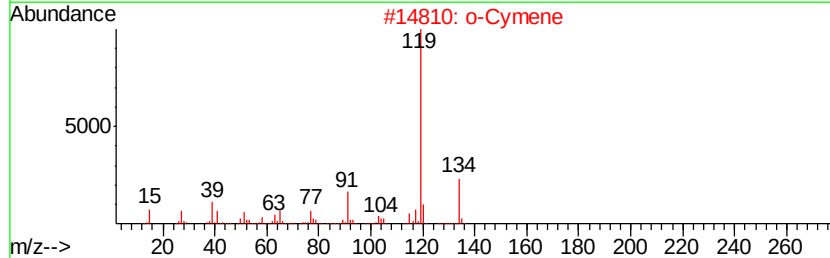
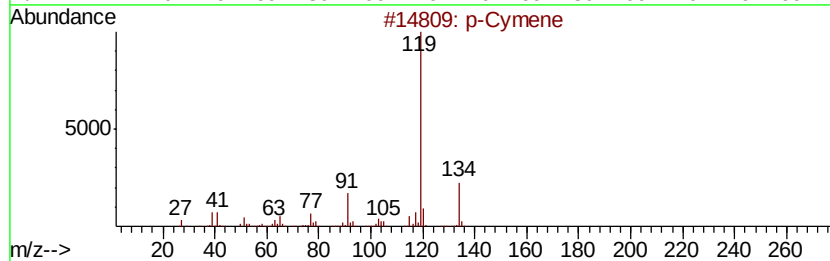
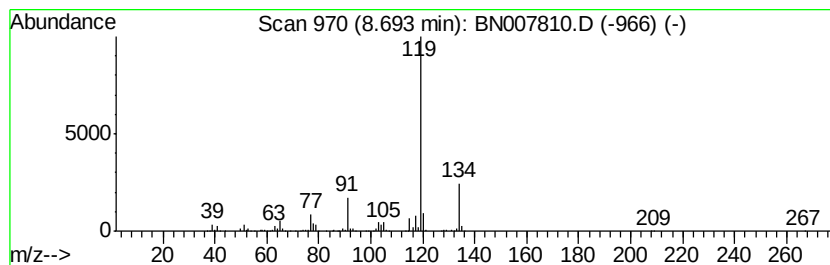
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 p-Cymene Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.69	6.28 ng/ul	1010850	1,4-Dichlorobenzene-d4	7.69

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			p-Cymene	134	C10H14	000099-87-6	97
2			o-Cymene	134	C10H14	000527-84-4	97
3			o-Cymene	134	C10H14	000527-84-4	97
4			o-Cymene	134	C10H14	000527-84-4	95
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN092319\
Data File : BN007810.D
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ClientSampleId :
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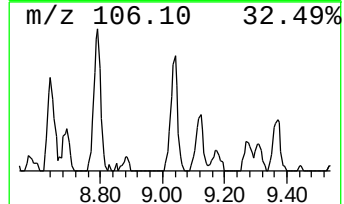
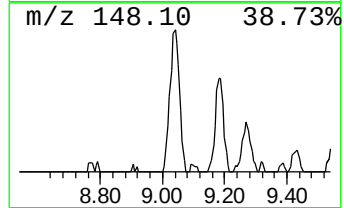
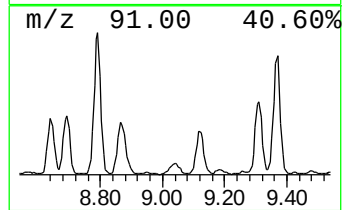
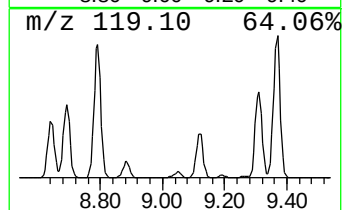
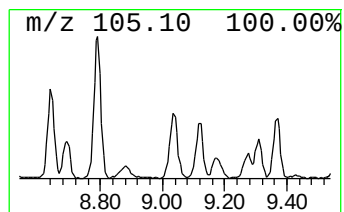
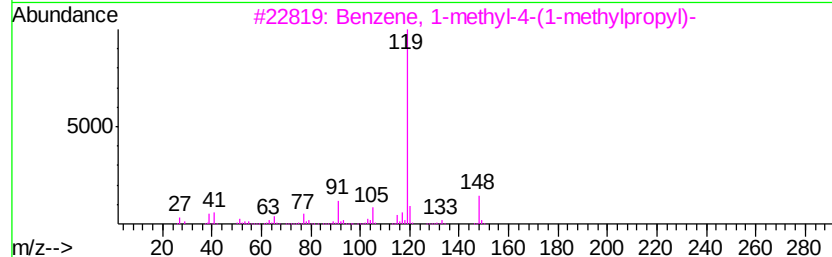
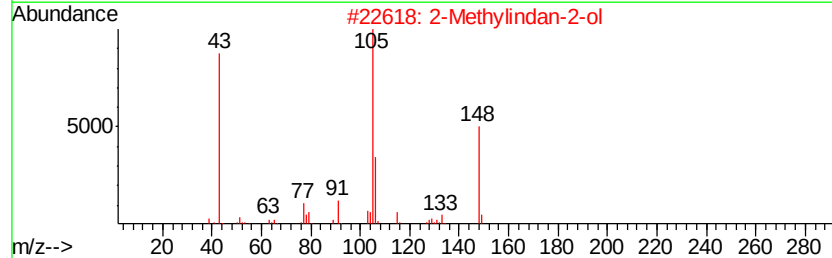
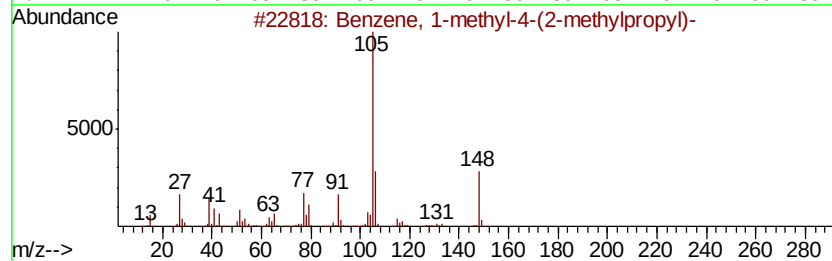
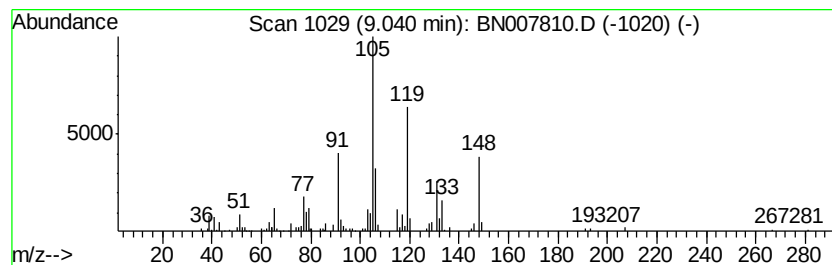
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 Benzene, 1-methyl-4-(2-meth... Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.04	2.03 ng/ul	326224	1,4-Dichlorobenzene-d4	7.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-4-(2-methylpro...	148	C11H16	005161-04-6	60
2		2-Methylindan-2-ol	148	C10H12O	033223-84-6	53
3		Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	52
4		Benzene, 1-methyl-4-butyl	148	C11H16	001595-05-7	50
5		3-Buten-2-ol, 4-phenyl-, (E)-	148	C10H12O	1000163-70-9	46



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN092319\
Data File : BN007810.D
Acq On : 23 Sep 2019 20:18
Operator : HP/JU
Sample : K4895-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BFDF5

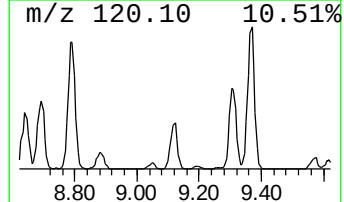
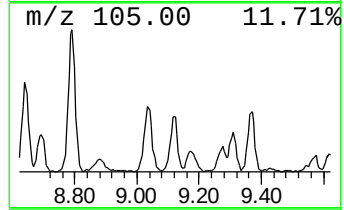
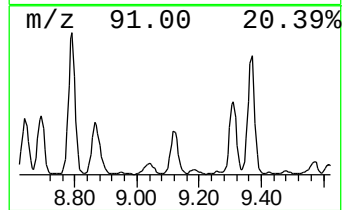
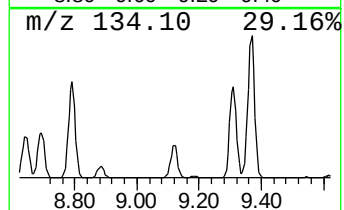
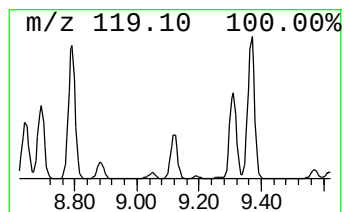
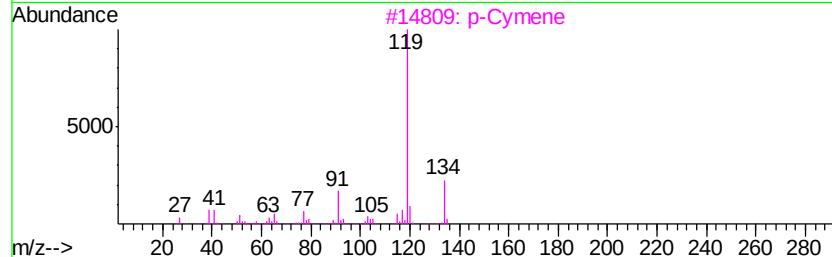
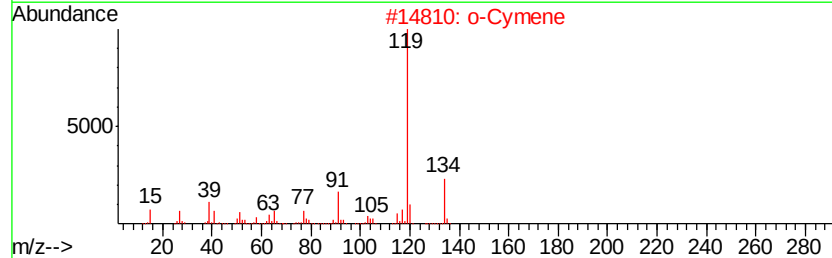
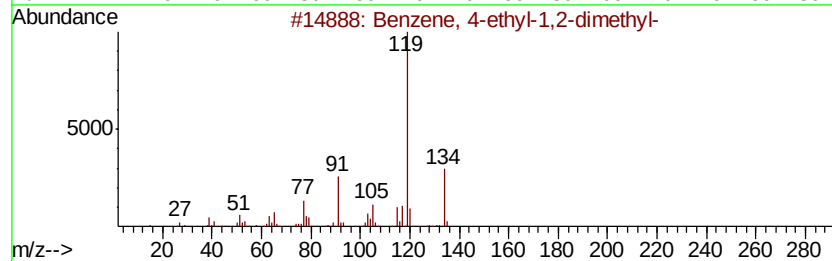
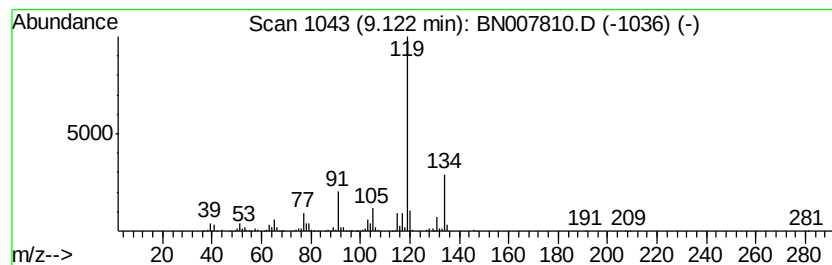
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.12	2.97 ng/ul	825501	Napthalene-d8	10.46

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
2		o-Cymene	134	C10H14	000527-84-4	95
3		p-Cymene	134	C10H14	000099-87-6	95
4		o-Cymene	134	C10H14	000527-84-4	95
5		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN092319\
Data File : BN007810.D
Acq On : 23 Sep 2019 20:18
Operator : HP/JU
Sample : K4895-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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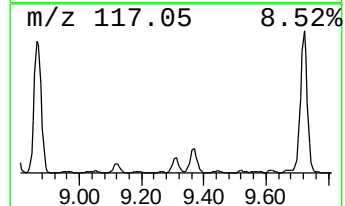
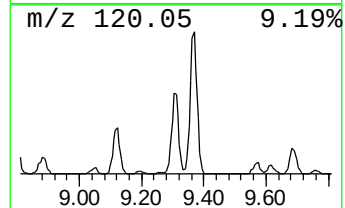
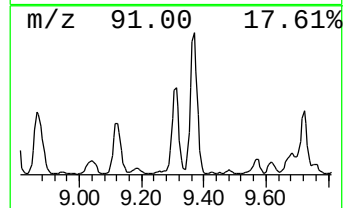
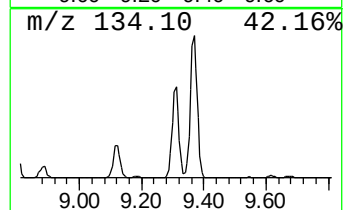
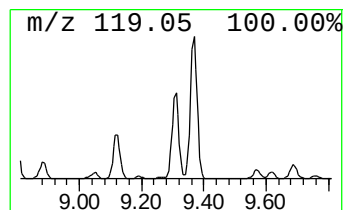
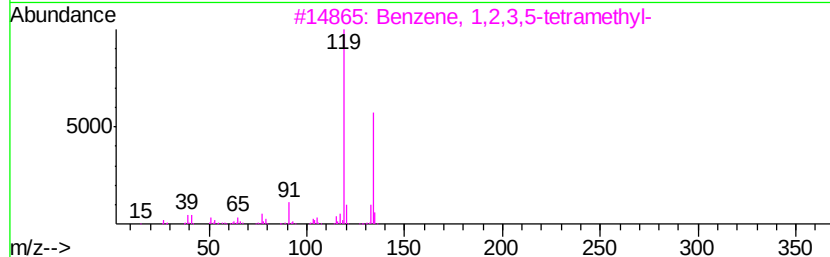
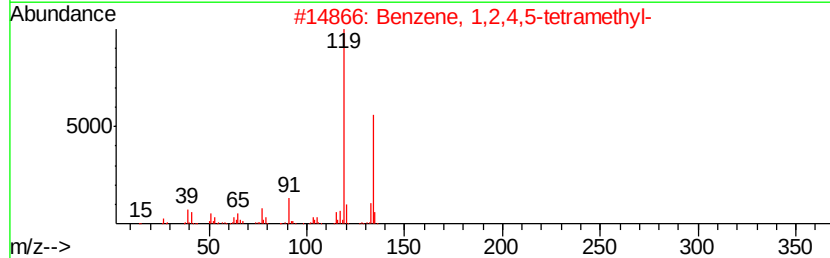
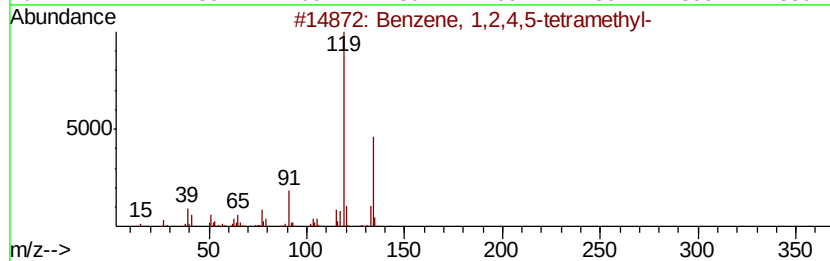
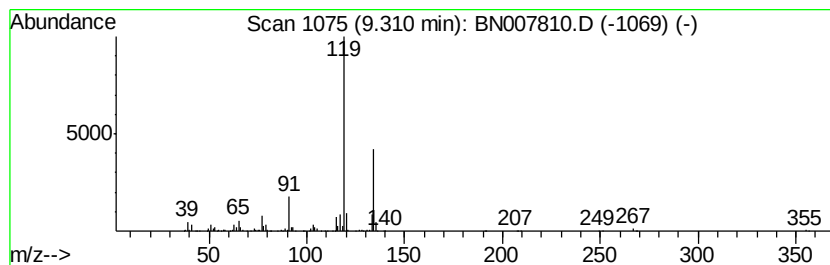
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Quant Title : SVOA CALIBRATION

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

Peak Number 23 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.31	5.14 ng/ul	1428100	Napthalene-d8	10.46

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
2		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
3		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
4		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
5		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN092319\
Data File : BN007810.D
Acq On : 23 Sep 2019 20:18
Operator : HP/JU
Sample : K4895-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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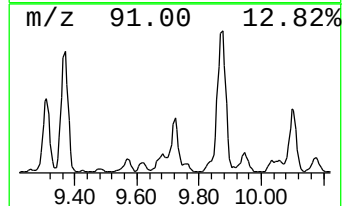
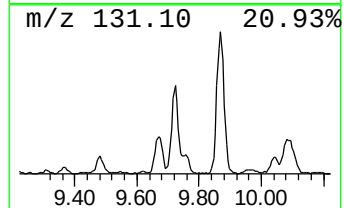
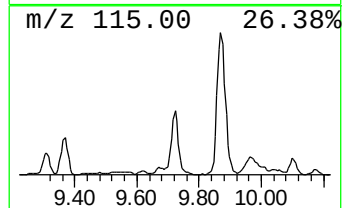
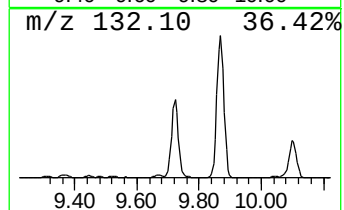
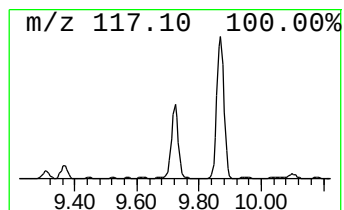
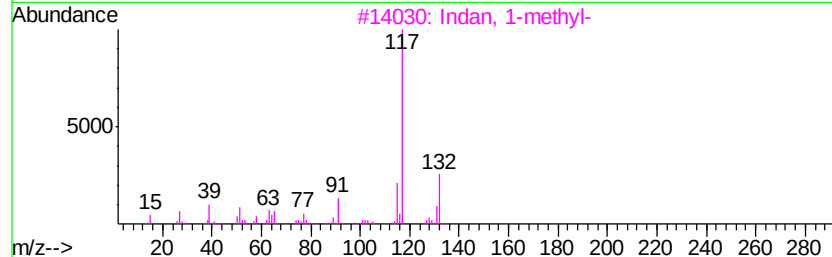
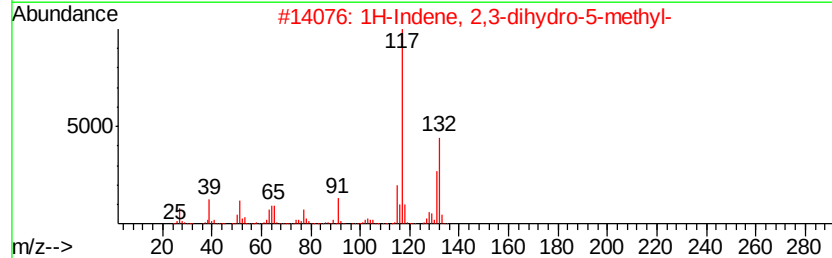
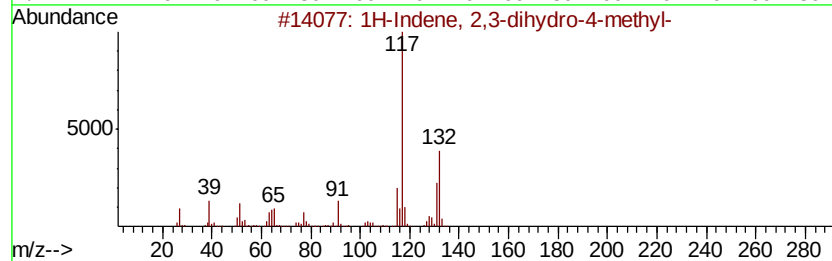
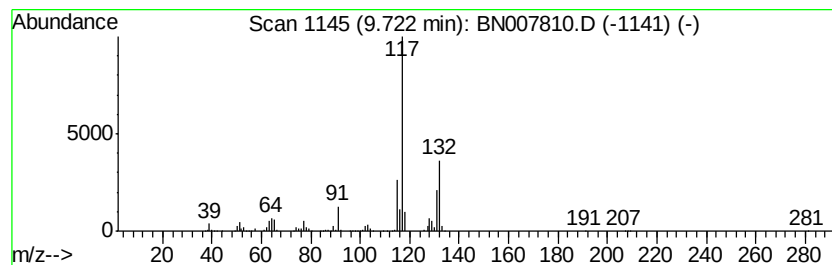
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Quant Title : SVOA CALIBRATION

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

Peak Number 25 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.72	5.11 ng/ul	1419880	Napthalene-d8	10.46

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	94
2			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
3			Indan, 1-methyl-	132	C10H12	000767-58-8	87
4			Indan, 1-methyl-	132	C10H12	000767-58-8	81
5			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN092319\
Data File : BN007810.D
Acq On : 23 Sep 2019 20:18
Operator : HP/JU
Sample : K4895-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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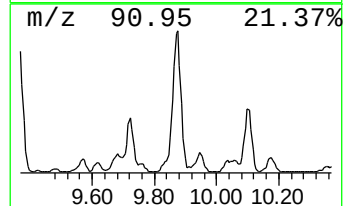
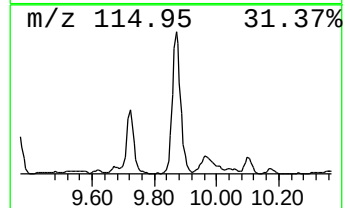
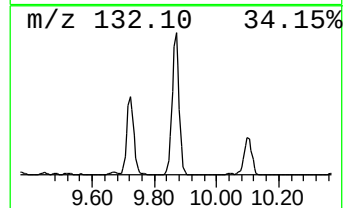
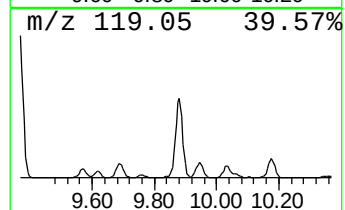
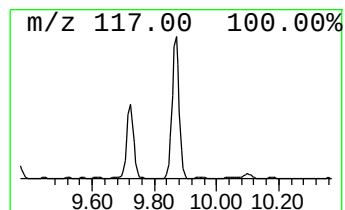
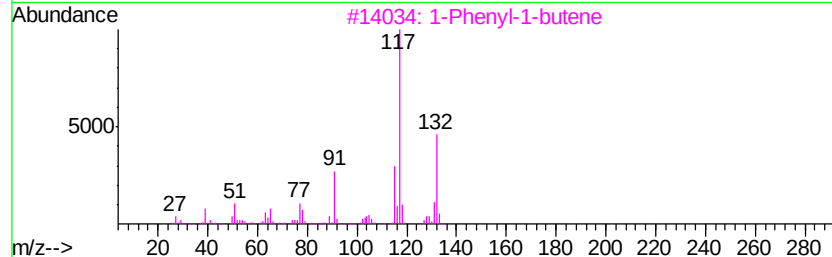
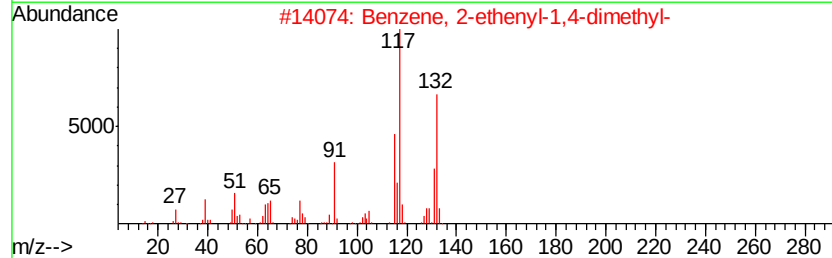
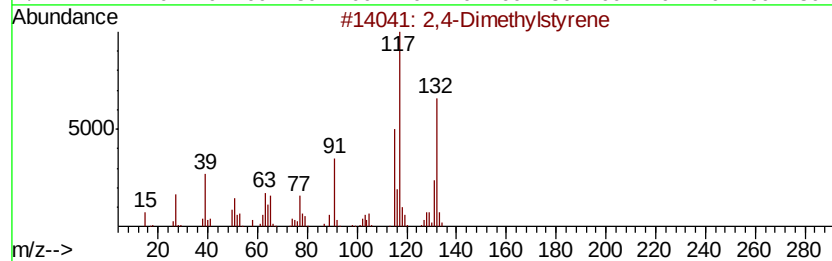
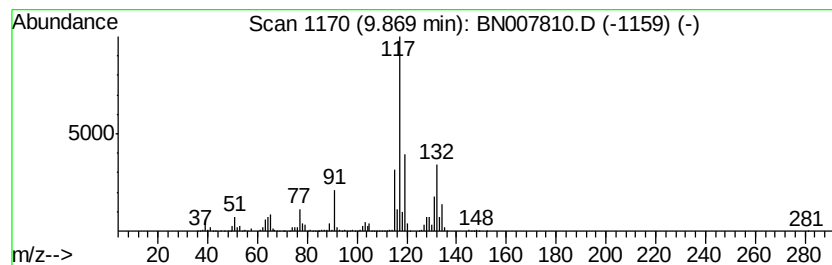
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Quant Title : SVOA CALIBRATION

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

Peak Number 26 2,4-Dimethylstyrene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.87	15.07 ng/ul	4189320	Napthalene-d8	10.46

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,4-Dimethylstvrene	132	C10H12	002234-20-0	95
2		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
3		1-Phenvl-1-butene	132	C10H12	000824-90-8	76
4		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	70
5		Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	70



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN092319\
Data File : BN007810.D
Acq On : 23 Sep 2019 20:18
Operator : HP/JU
Sample : K4895-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

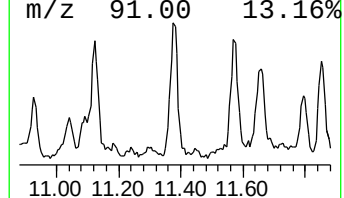
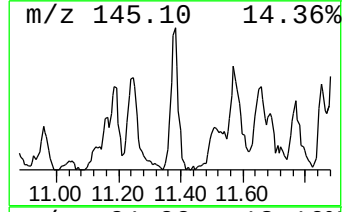
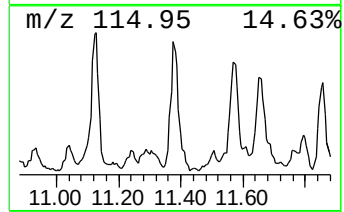
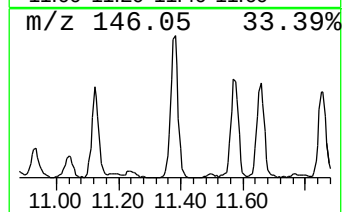
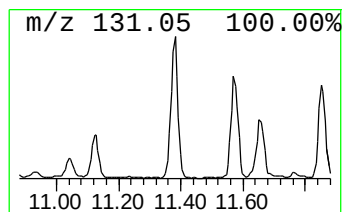
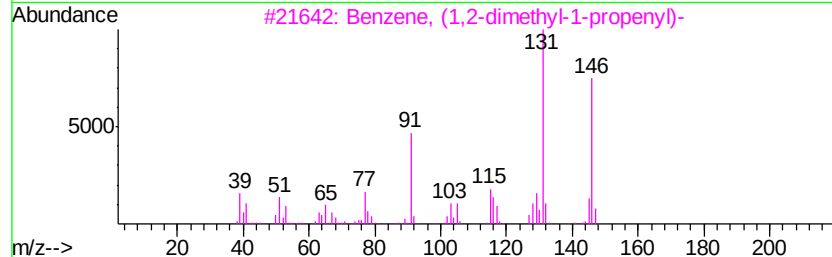
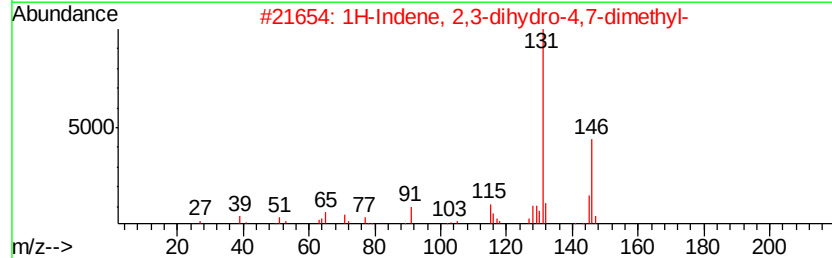
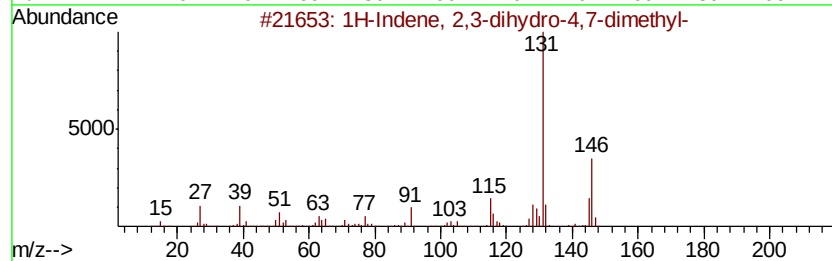
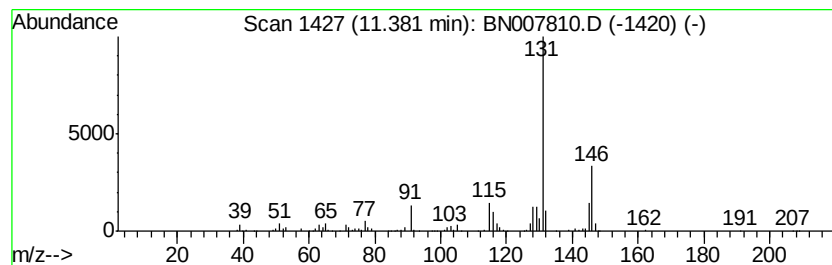
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 27 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.38	2.30 ng/ul	640109	Napthalene-d8	10.46
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1H-Indene, 2,3-dihydro-4,7-dimet...	146 C11H14	006682-71-9	96
2	1H-Indene, 2,3-dihydro-4,7-dimet...	146 C11H14	006682-71-9	95
3	Benzene, (1,2-dimethyl-1-propenyl)-	146 C11H14	000769-57-3	94
4	1H-Indene, 2,3-dihydro-4,7-dimet...	146 C11H14	006682-71-9	94
5	1H-Indene, 2,3-dihydro-5,6-dimet...	146 C11H14	001075-22-5	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN092319\
Data File : BN007810.D
Acq On : 23 Sep 2019 20:18
Operator : HP/JU
Sample : K4895-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BFDF5

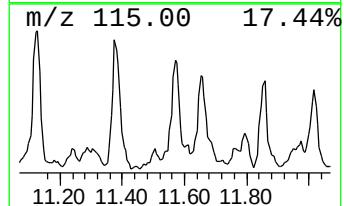
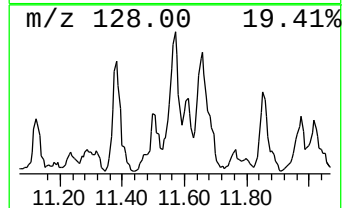
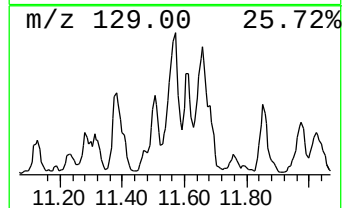
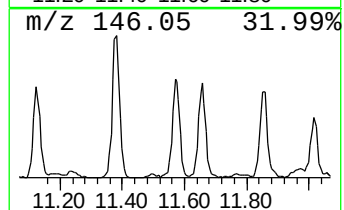
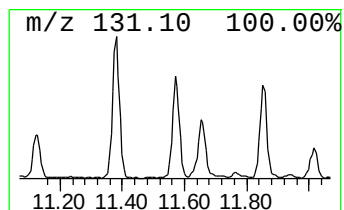
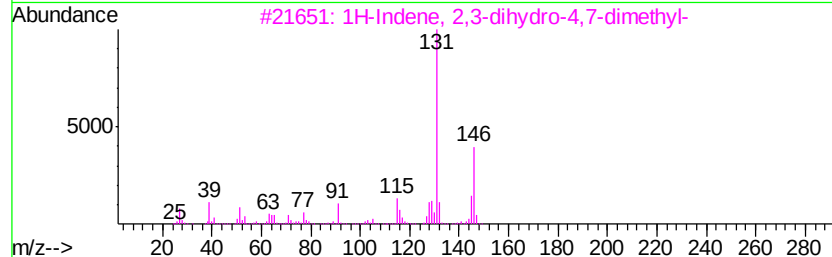
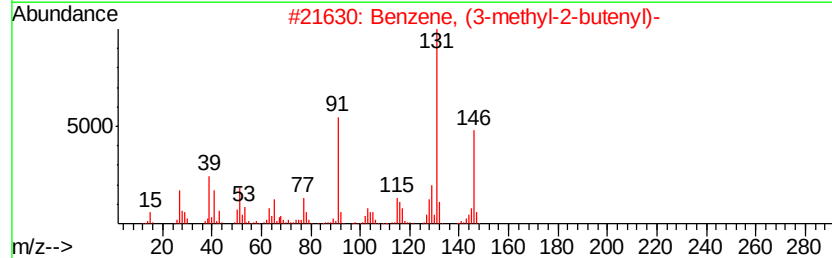
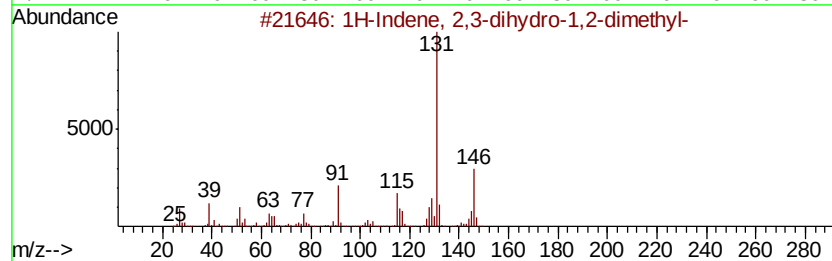
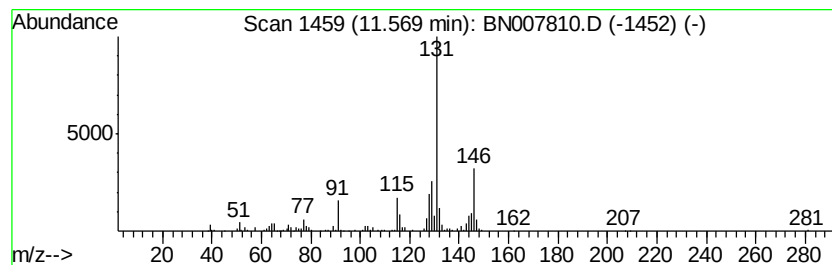
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Quant Title : SVOA CALIBRATION

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

Peak Number 28 1H-Indene, 2,3-dihydro-1,2-... Concentration Rank 35

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.57	2.15 ng/ul	598551	Napthalene-d8	10.46

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	87
2		Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	83
3		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	83
4		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	81
5		1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	81



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN092319\
Data File : BN007810.D
Acq On : 23 Sep 2019 20:18
Operator : HP/JU
Sample : K4895-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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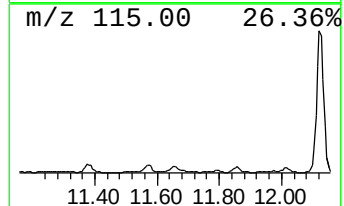
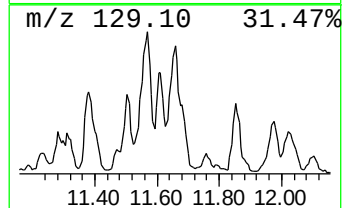
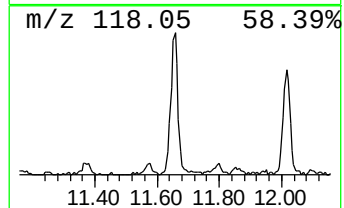
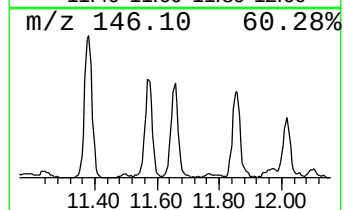
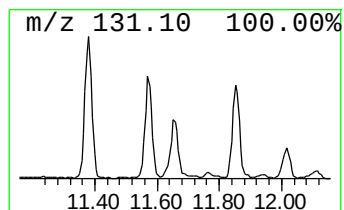
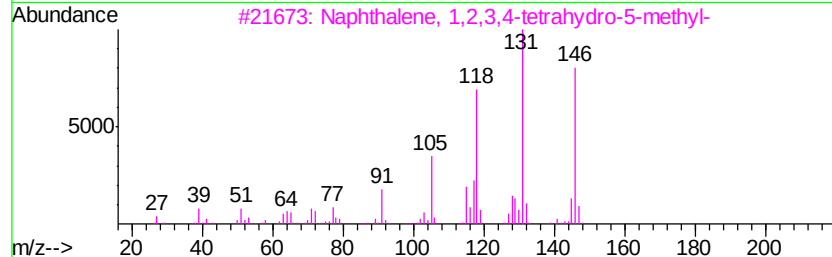
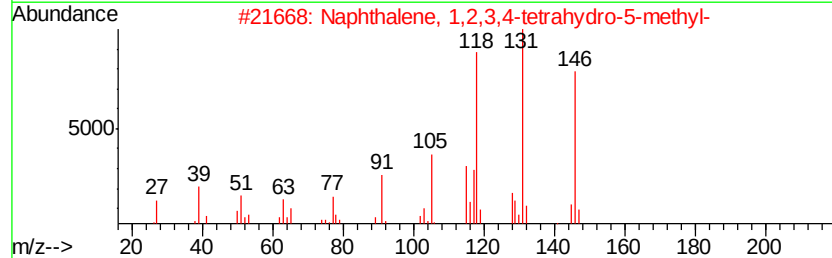
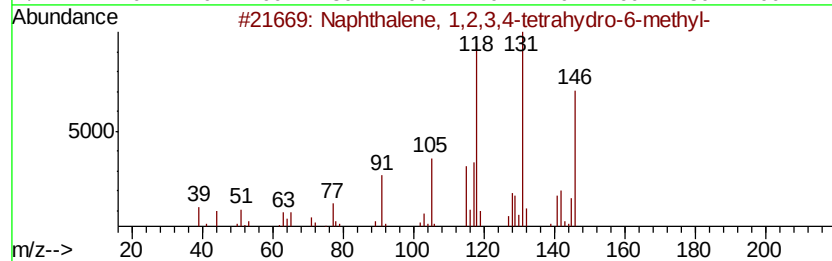
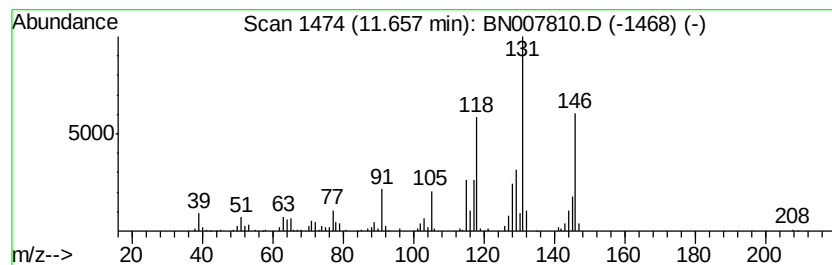
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Quant Title : SVOA CALIBRATION

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

Peak Number 29 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.66	2.11 ng/ul	585735	Naphthalene-d8	10.46

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	91
2		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	87
3		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	87
4		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	80
5		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN092319\
Data File : BN007810.D
Acq On : 23 Sep 2019 20:18
Operator : HP/JU
Sample : K4895-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

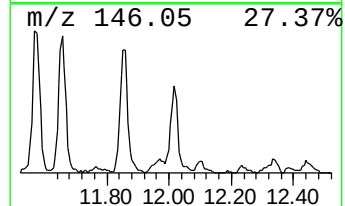
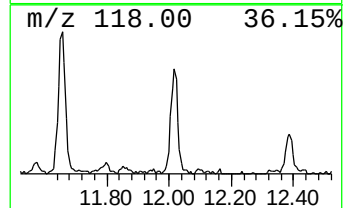
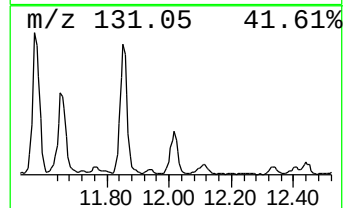
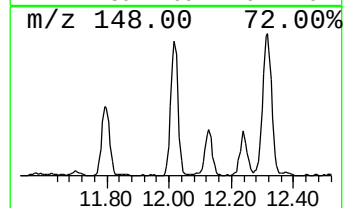
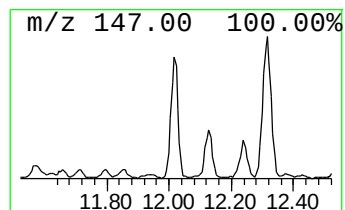
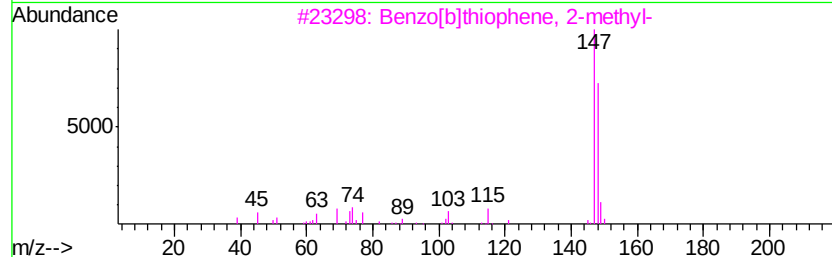
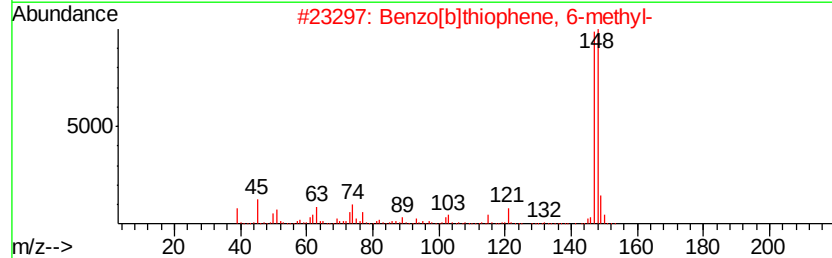
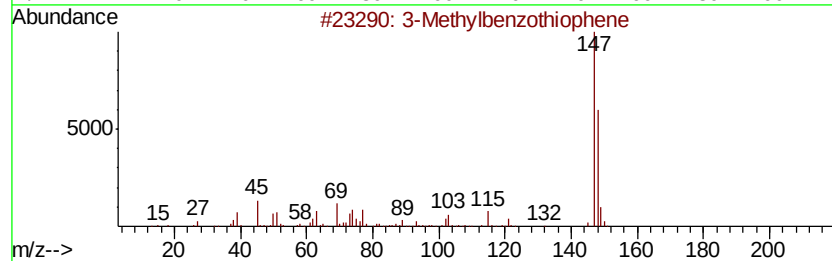
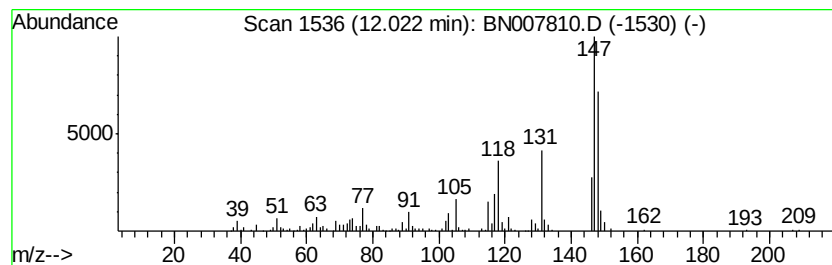
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN091619MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 30 3-Methylbenzothiophene Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.02	2.27 ng/ul	631925	Napthalene-d8	10.46
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		3-Methylbenzothiophene	148 C9H8S	001455-18-1 64
2		Benzo[b]thiophene, 6-methyl-	148 C9H8S	016587-47-6 64
3		Benzo[b]thiophene, 2-methyl-	148 C9H8S	001195-14-8 58
4		Benzaldehyde, 2,4,5-trimethyl-	148 C10H12O	005779-72-6 58
5		Benzo[b]thiophene, 2-methyl-	148 C9H8S	001195-14-8 58



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN092319\
Data File : BN007810.D
Acq On : 23 Sep 2019 20:18
Operator : HP/JU
Sample : K4895-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BDF5

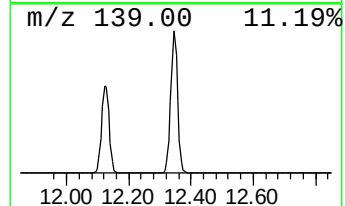
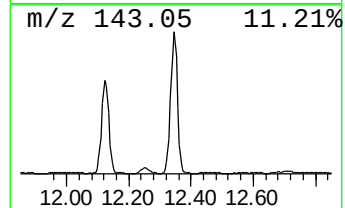
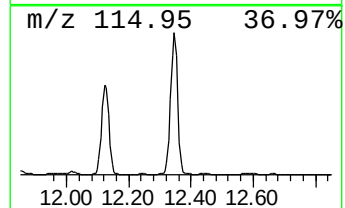
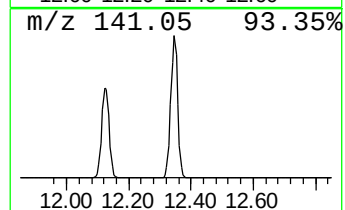
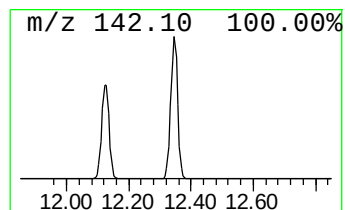
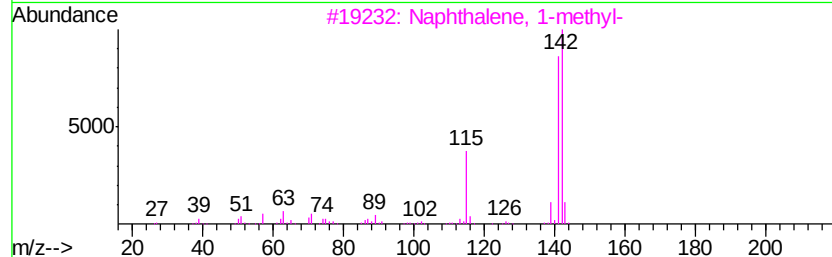
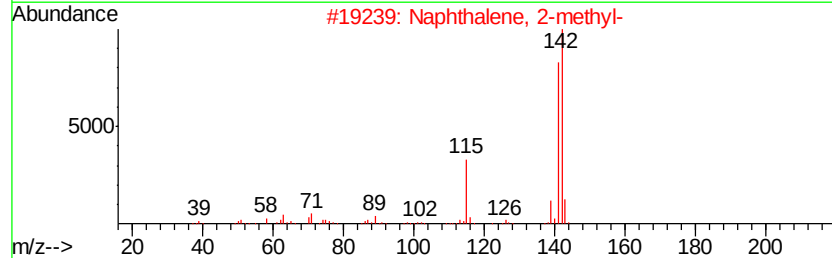
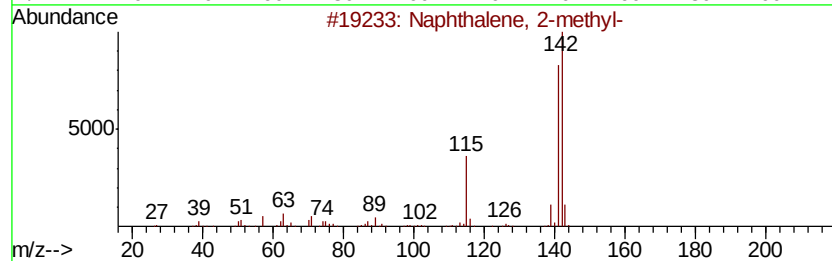
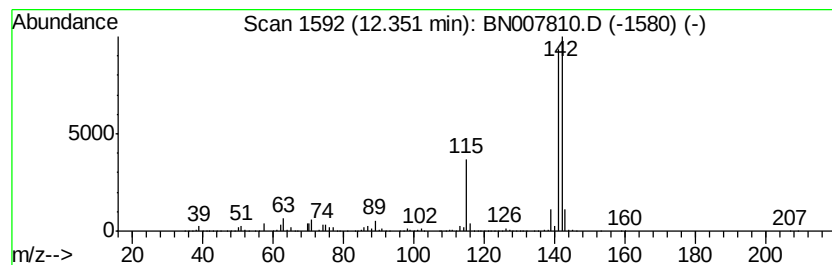
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN091619MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 31 Naphthalene, 1-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.35	28.09 ng/ul	7810540	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, 2-methyl-	142	C11H10	000091-57-6	96
3		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
4		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
5		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN092319\
 Data File : BN007810.D
 Acq On : 23 Sep 2019 20:18
 Operator : HP/JU
 Sample : K4895-01
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BFDF5

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	4.7	ng/ul	754423	1	7.69	3216710	20.0
Butanoic acid, 3-...	4.22	10.7	ng/ul	1723910	1	7.69	3216710	20.0
Butanoic acid, et...	4.28	2.1	ng/ul	346545	1	7.69	3216710	20.0
Propanoic acid, p...	4.39	16.2	ng/ul	2603970	1	7.69	3216710	20.0
Butanoic acid, 1-...	4.83	20.5	ng/ul	3303210	1	7.69	3216710	20.0
Benzene, 1,3-dime...	5.38	7.7	ng/ul	1234310	1	7.69	3216710	20.0
Butanoic acid, pr...	5.68	3.3	ng/ul	537696	1	7.69	3216710	20.0
Benzene, 1,2,3-tr...	6.93	15.5	ng/ul	2494090	1	7.69	3216710	20.0
Benzene, 1-ethyl-...	7.09	12.5	ng/ul	2013260	1	7.69	3216710	20.0
Benzene, 1,2,4-tr...	7.35	25.4	ng/ul	4090240	1	7.69	3216710	20.0
Mesitylene	7.80	22.0	ng/ul	3543150	1	7.69	3216710	20.0
Indane	8.06	13.2	ng/ul	2120900	1	7.69	3216710	20.0
Benzene, 1,3-diet...	8.19	3.5	ng/ul	566538	1	7.69	3216710	20.0
Benzene, 1-methyl...	8.24	5.7	ng/ul	918965	1	7.69	3216710	20.0
Benzene, 2-ethyl-...	8.64	5.9	ng/ul	942589	1	7.69	3216710	20.0
p-Cymene	8.69	6.3	ng/ul	1010850	1	7.69	3216710	20.0
Benzene, 1-methyl...	9.04	2.0	ng/ul	326224	1	7.69	3216710	20.0
Benzene, 4-ethyl-...	9.12	3.0	ng/ul	825501	2	10.46	5561000	20.0
Benzene, 1,2,4,5-...	9.31	5.1	ng/ul	1428100	2	10.46	5561000	20.0
1H-Indene, 2,3-di...	9.72	5.1	ng/ul	1419880	2	10.46	5561000	20.0
2,4-Dimethylstyrene	9.87	15.1	ng/ul	4189320	2	10.46	5561000	20.0
1H-Indene, 2,3-di...	11.38	2.3	ng/ul	640109	2	10.46	5561000	20.0
1H-Indene, 2,3-di...	11.57	2.1	ng/ul	598551	2	10.46	5561000	20.0
Naphthalene, 1,2,...	11.66	2.1	ng/ul	585735	2	10.46	5561000	20.0
3-Methylbenzothio...	12.02	2.3	ng/ul	631925	2	10.46	5561000	20.0
Naphthalene, 1-me...	12.35	28.1	ng/ul	7810540	2	10.46	5561000	20.0