

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN092319\
 Data File : BN007816.D
 Acq On : 23 Sep 2019 23:52
 Operator : HP/JU
 Sample : K4895-08
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BFDH8

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	64	rVB	5619947	6667541	37.62%	2.788%
2	3.634	106	110	116	rBV	180903	262054	1.48%	0.110%
3	3.693	116	120	127	rVB	689475	793921	4.48%	0.332%
4	3.940	158	162	175	rVV2	202234	387824	2.19%	0.162%
5	4.216	204	209	216	rBV	1507629	1846075	10.42%	0.772%
6	4.281	216	220	234	rVV	181362	285389	1.61%	0.119%
7	4.387	234	238	246	rVV	1986094	2582693	14.57%	1.080%
8	4.834	305	314	324	rBV	2517459	3466585	19.56%	1.450%
9	5.675	452	457	463	rBV	368182	524446	2.96%	0.219%
10	6.405	573	581	589	rVB	169205	266703	1.50%	0.112%
11	6.799	643	648	652	rBV	184569	270958	1.53%	0.113%
12	6.863	652	659	666	rVV3	624972	1235906	6.97%	0.517%
13	6.928	666	670	677	rVB	193173	291197	1.64%	0.122%
14	7.028	681	687	694	rBV	1495485	2460481	13.88%	1.029%
15	7.093	694	698	703	rVV	272335	451703	2.55%	0.189%
16	7.222	713	720	731	rVV	1829532	2914330	16.44%	1.219%
17	7.346	735	741	748	rVV	731097	1121792	6.33%	0.469%
18	7.687	792	799	805	rBV	2143236	3356899	18.94%	1.404%
19	7.805	815	819	828	rVB	368903	640006	3.61%	0.268%
20	8.063	855	863	869	rVB	226454	378820	2.14%	0.158%
21	8.328	901	908	912	rBV	216912	328255	1.85%	0.137%
22	8.387	912	918	924	rBV	1457244	2242152	12.65%	0.938%
23	8.793	979	987	989	rBV	361797	591527	3.34%	0.247%
24	8.828	989	993	998	rBV	1719464	2518388	14.21%	1.053%
25	9.310	1070	1075	1079	rVV2	173034	264434	1.49%	0.111%
26	9.363	1079	1084	1093	rVB	293984	491886	2.78%	0.206%
27	9.546	1108	1115	1123	rBV	1533914	2570997	14.51%	1.075%
28	9.640	1123	1131	1134	rVV2	136516	286278	1.62%	0.120%
29	9.722	1141	1145	1150	rVV	212075	328691	1.85%	0.137%
30	9.875	1159	1171	1180	rBV6	597047	1419881	8.01%	0.594%
31	9.963	1180	1186	1191	rVV2	599055	1271068	7.17%	0.532%
32	10.004	1191	1193	1199	rVV3	234450	350686	1.98%	0.147%
33	10.081	1199	1206	1217	rVB	2710440	4550661	25.67%	1.903%
34	10.457	1260	1270	1275	rVV	2981641	5262122	29.69%	2.200%

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35	10.510	1275	1279	1286	rVV	1969611	3348014	18.89%	1.400%
36	10.593	1286	1293	1306	rVV	2167066	3987480	22.50%	1.667%
37	11.287	1407	1411	1421	rVV4	116883	336152	1.90%	0.141%
38	11.504	1439	1448	1451	rBV3	131126	264914	1.49%	0.111%
39	11.569	1451	1459	1463	rVV4	147960	402691	2.27%	0.168%
40	11.657	1469	1474	1487	rVB	204863	441770	2.49%	0.185%
41	12.016	1531	1535	1546	rVV4	182479	419293	2.37%	0.175%
42	12.122	1546	1553	1564	rVB	1226515	1972730	11.13%	0.825%
43	12.345	1580	1591	1600	rVB	5470722	8640744	48.75%	3.613%
44	13.151	1723	1728	1736	rBV3	172432	330216	1.86%	0.138%
45	13.345	1755	1761	1762	rVV2	193040	283767	1.60%	0.119%
46	13.369	1763	1765	1774	rVB3	282001	395606	2.23%	0.165%
47	13.492	1775	1786	1793	rBV3	433254	998475	5.63%	0.418%
48	13.640	1803	1811	1816	rBV	1292067	2078635	11.73%	0.869%
49	13.692	1816	1820	1821	rVV	280128	350310	1.98%	0.146%
50	13.728	1821	1826	1831	rVV	5413492	7502167	42.33%	3.137%
51	13.775	1831	1834	1840	rVB	744982	958298	5.41%	0.401%
52	13.875	1847	1851	1854	rVV	485917	707753	3.99%	0.296%
53	13.904	1854	1856	1861	rVB	233016	294006	1.66%	0.123%
54	14.004	1867	1873	1878	rBV	5517357	8615840	48.61%	3.603%
55	14.316	1918	1926	1932	rVV2	4527101	7040338	39.72%	2.944%
56	14.381	1932	1937	1945	rVB	4714116	6891805	38.88%	2.882%
57	14.510	1954	1959	1971	rBV2	544118	1050230	5.93%	0.439%
58	14.716	1989	1994	1999	rBV2	774498	1163490	6.56%	0.487%
59	14.945	2025	2033	2038	rBV4	218179	479133	2.70%	0.200%
60	15.092	2053	2058	2060	rVV3	202276	331700	1.87%	0.139%
61	15.310	2089	2095	2101	rBV	8001493	11503303	64.90%	4.810%
62	15.369	2101	2105	2110	rVV	1718290	2326305	13.12%	0.973%
63	15.428	2110	2115	2119	rVV	2222450	2851131	16.09%	1.192%
64	15.463	2119	2121	2124	rVV	223928	262065	1.48%	0.110%
65	15.581	2138	2141	2145	rVB2	211200	287224	1.62%	0.120%
66	15.775	2168	2174	2178	rBV	510404	808776	4.56%	0.338%
67	16.369	2269	2275	2280	rBV2	274917	586778	3.31%	0.245%
68	16.522	2297	2301	2308	rVB2	222958	339621	1.92%	0.142%
69	16.875	2357	2361	2370	rVB	493901	748347	4.22%	0.313%
70	17.063	2387	2393	2396	rBV2	5685046	8026961	45.29%	3.357%
71	17.104	2396	2400	2404	rVV	5253346	7188995	40.56%	3.006%

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72	17.163	2404	2410	2414	rVV	8758579	13348720	75.31%	5.582%
73	17.192	2414	2415	2420	rVV	1127031	977369	5.51%	0.409%
74	17.251	2420	2425	2432	rVB	479520	716851	4.04%	0.300%
75	17.457	2452	2460	2466	rBV	964939	1489115	8.40%	0.623%
76	17.686	2495	2499	2503	rBV	429803	522516	2.95%	0.218%
77	17.939	2538	2542	2544	rBV	275276	332026	1.87%	0.139%
78	17.980	2544	2549	2557	rVB3	2192702	3222806	18.18%	1.348%
79	18.128	2569	2574	2578	rBV2	663622	1168933	6.60%	0.489%
80	18.169	2578	2581	2586	rVB2	321982	427909	2.41%	0.179%
81	18.280	2596	2600	2612	rVB5	264138	455236	2.57%	0.190%
82	19.122	2739	2743	2750	rVB	903968	1215524	6.86%	0.508%
83	19.263	2763	2767	2776	rBV2	1111517	1689331	9.53%	0.706%
84	19.345	2778	2781	2789	rVB2	197701	333981	1.88%	0.140%
85	19.457	2795	2800	2816	rVB	12046922	17493873	98.70%	7.315%
86	20.510	2975	2979	2982	rBV	352156	439021	2.48%	0.184%
87	20.945	3050	3053	3057	rBV2	519520	629212	3.55%	0.263%
88	20.986	3057	3060	3068	rVB	1611995	2001622	11.29%	0.837%
89	21.257	3100	3106	3115	rVB	7790907	10368467	58.50%	4.336%
90	21.374	3123	3126	3132	rBV	230823	300676	1.70%	0.126%
91	21.433	3133	3136	3145	rVB2	229279	384273	2.17%	0.161%
92	21.868	3207	3210	3221	rVB2	345809	508455	2.87%	0.213%
93	22.327	3284	3288	3293	rVB	533463	663119	3.74%	0.277%
94	22.827	3369	3373	3379	rVB	739506	971314	5.48%	0.406%
95	23.333	3451	3459	3465	rBV	9822324	17724258	100.00%	7.412%
96	23.392	3465	3469	3475	rVV	783957	1271957	7.18%	0.532%
97	23.468	3475	3482	3492	rVB2	5828117	10693830	60.33%	4.472%
98	24.033	3572	3578	3585	rBV	675866	1227013	6.92%	0.513%
99	24.768	3697	3703	3711	rBV2	461484	1014820	5.73%	0.424%
100	25.627	3845	3849	3858	rVB4	277040	648344	3.66%	0.271%

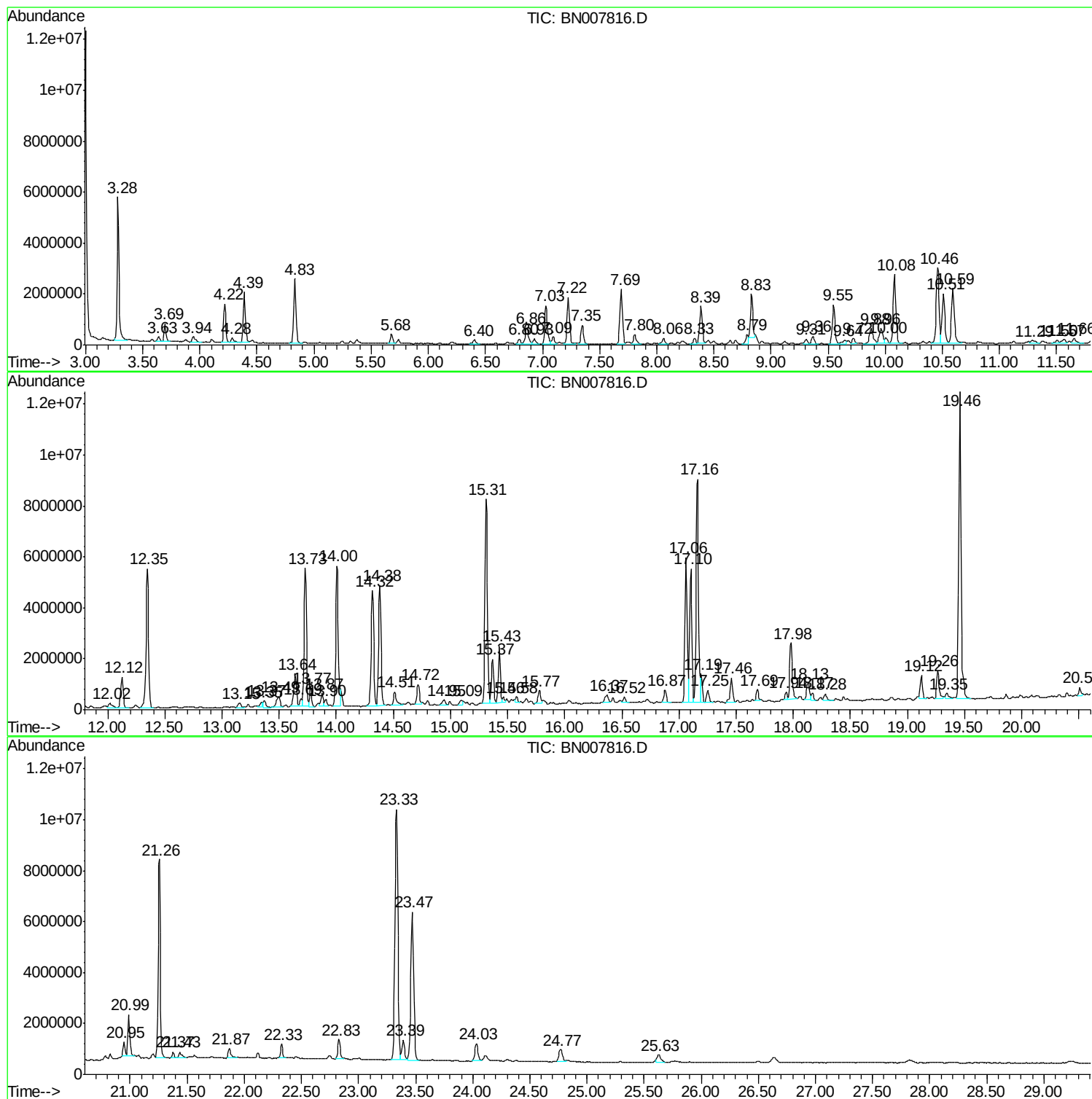
Sum of corrected areas: 239137983

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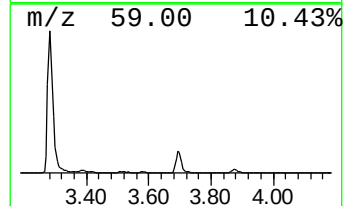
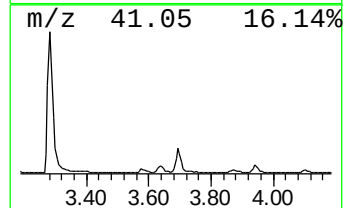
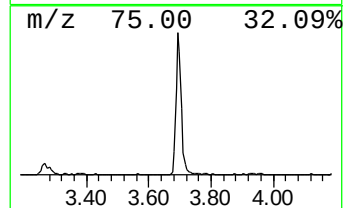
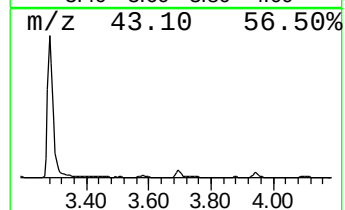
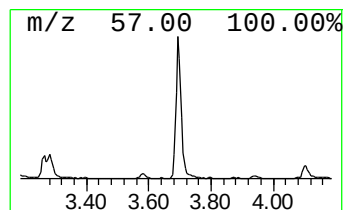
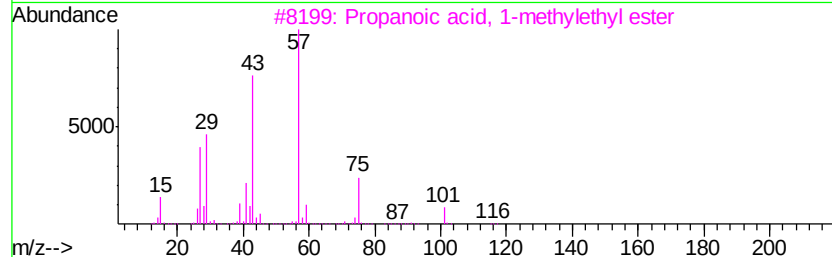
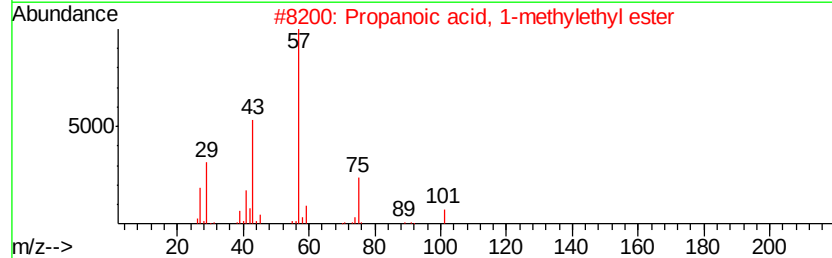
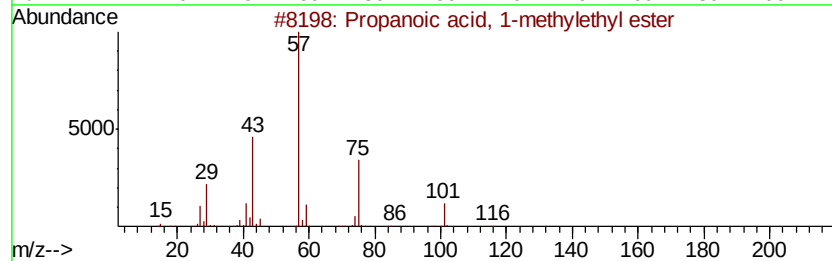
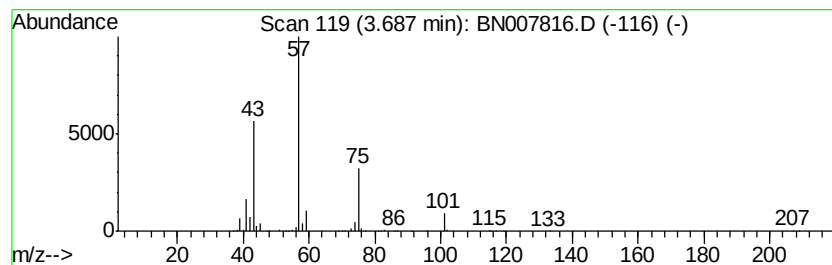
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN091619MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.69	4.73 ng/ul	793921	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	83
3		Propanoic acid, 1-methylethyl ester	116	C6H12O2	000637-78-5	64
4		Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	53
5		Propanoic acid, 1-methylethyl ester	116	C6H12O2	000637-78-5	53



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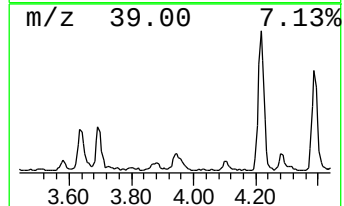
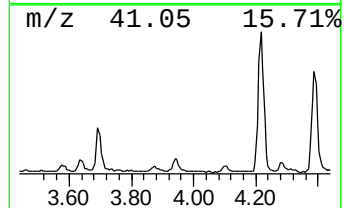
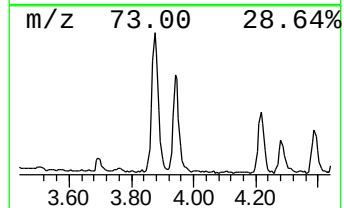
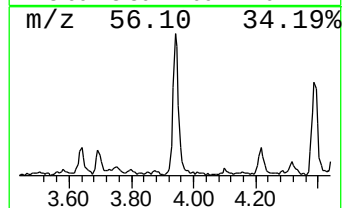
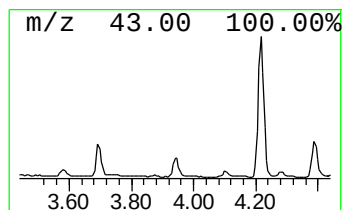
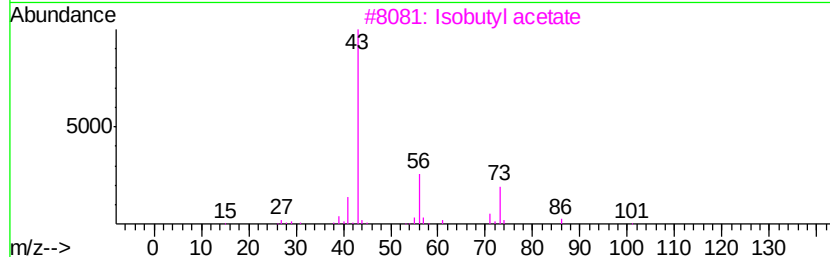
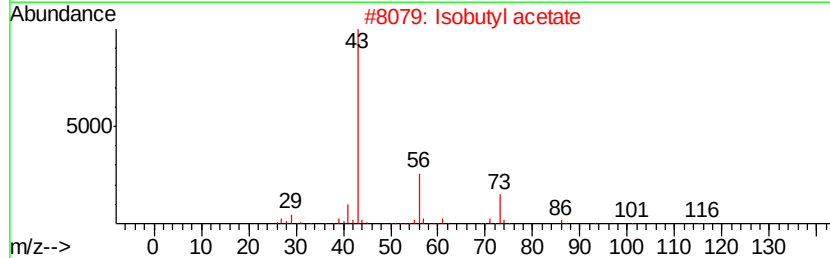
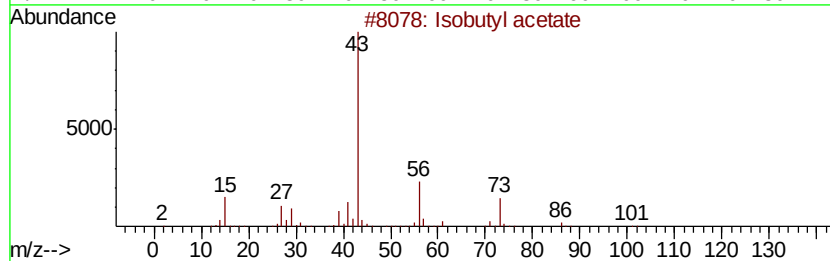
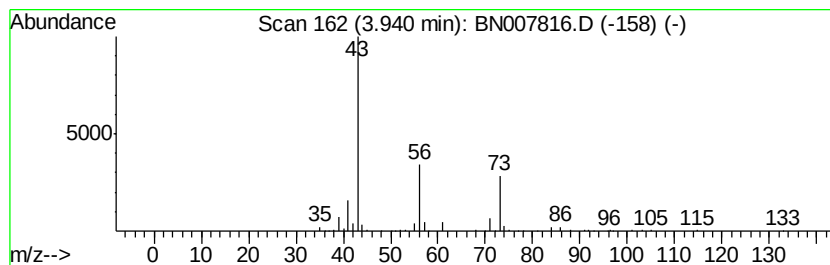
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 Isobutyl acetate Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.94	2.31 ng/ul	387824	1,4-Dichlorobenzene-d4	7.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Isobutyl acetate	116	C6H12O2	000110-19-0	78
2		Isobutyl acetate	116	C6H12O2	000110-19-0	78
3		Isobutyl acetate	116	C6H12O2	000110-19-0	72
4		Isobutyl acetate	116	C6H12O2	000110-19-0	64
5		Isobutyl acetate	116	C6H12O2	000110-19-0	50



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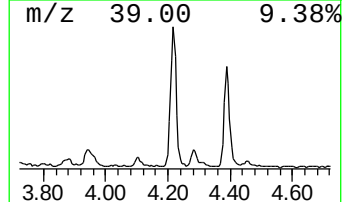
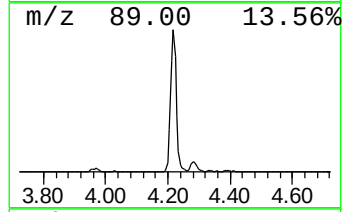
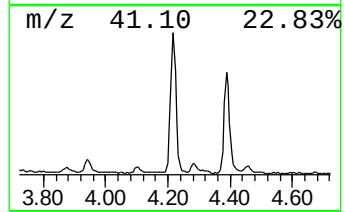
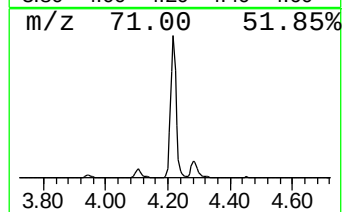
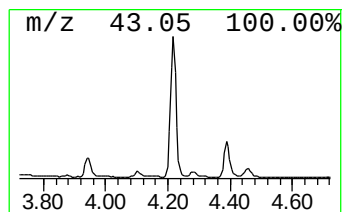
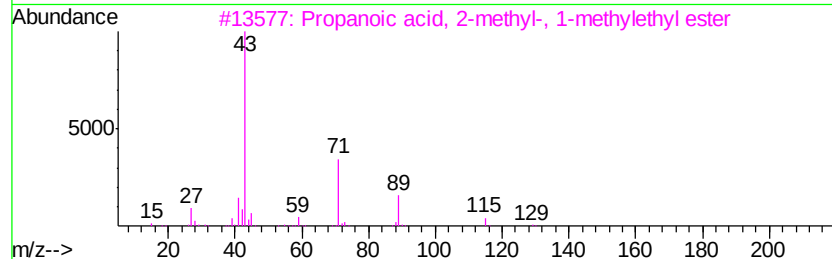
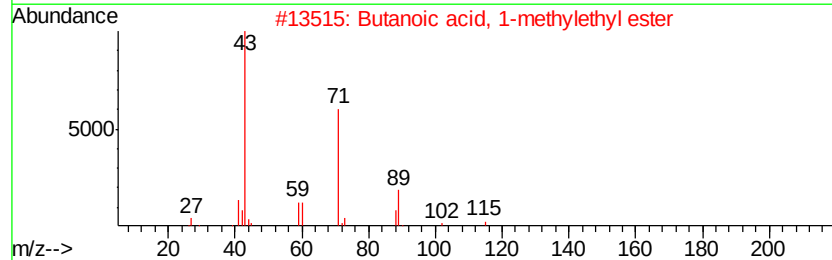
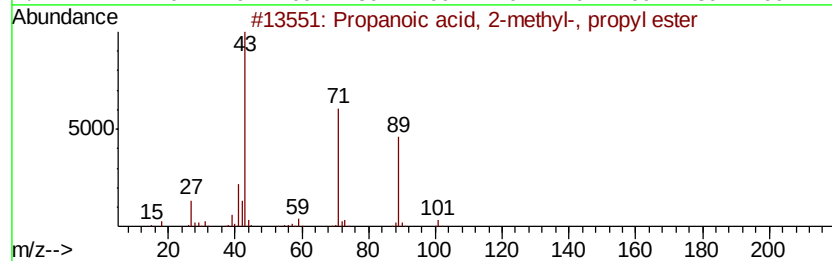
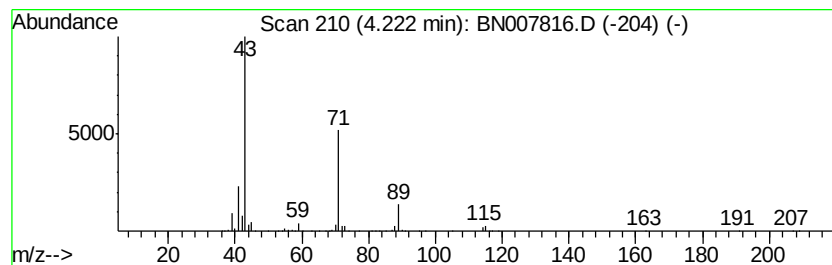
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TIC Integration Parameters: LSCINT.P

Peak Number 3 Propanoic acid, 2-methyl-, ... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.22	11.00 ng/ul	1846080	1,4-Dichlorobenzene-d4	7.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propanoic acid, 2-methyl-, propy...	130	C7H14O2	000644-49-5	64
2		Butanoic acid, 1-methylethyl ester	130	C7H14O2	000638-11-9	59
3		Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	53
4		Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	53
5		Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	50



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Data File : BN007816.D
Acq On : 23 Sep 2019 23:52
Operator : HP/JU
Sample : K4895-08
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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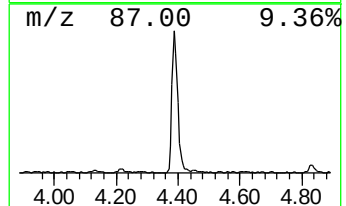
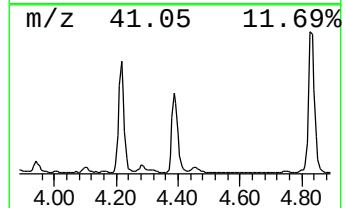
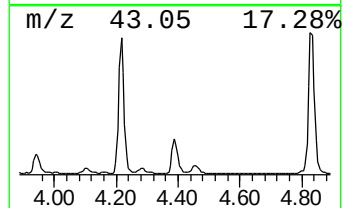
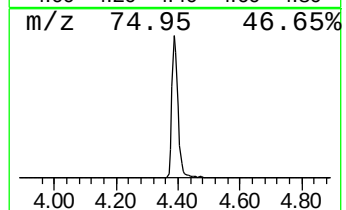
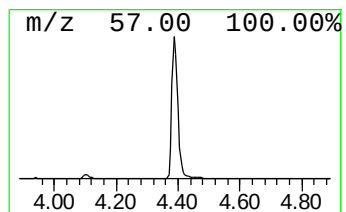
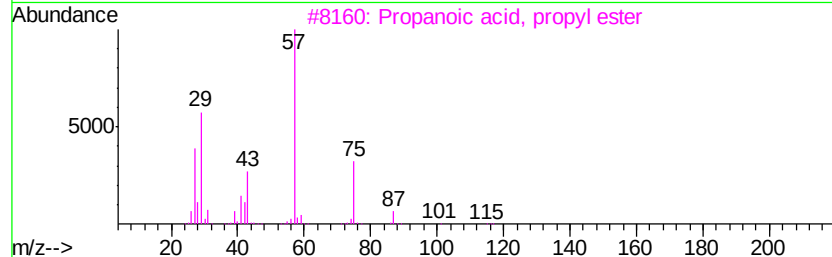
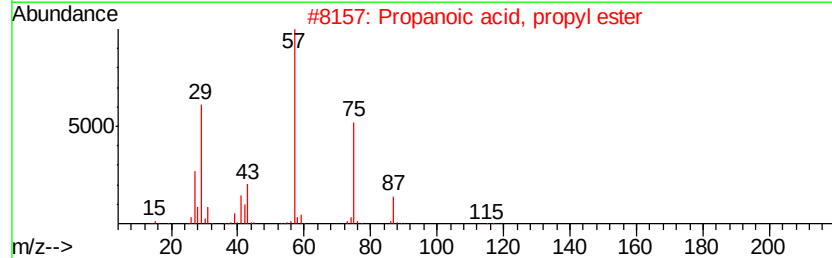
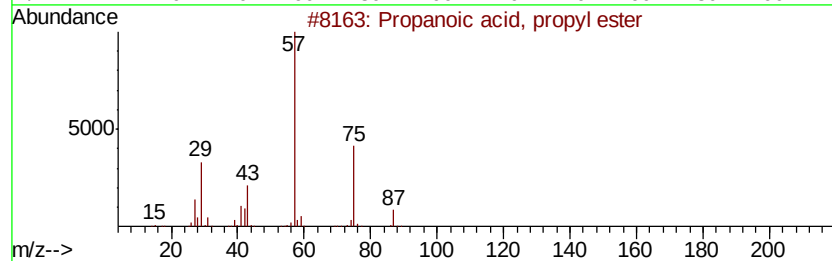
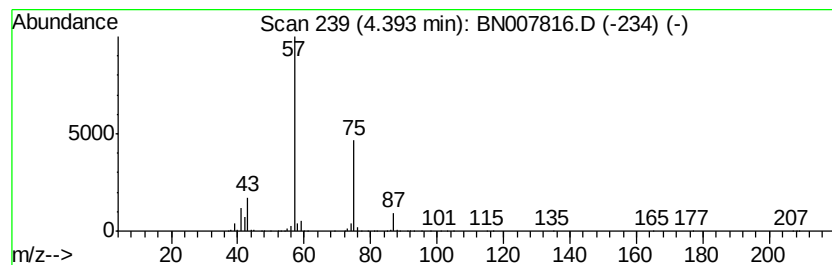
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Propanoic acid, propyl ester Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.39	15.39 ng/ul	2582690	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	83
2		Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	74
3		Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	59
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\
Data File : BN007816.D
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Sample : K4895-08
Misc :
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Instrument :
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ClientSampleId :
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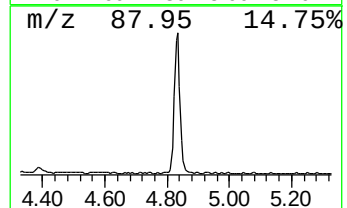
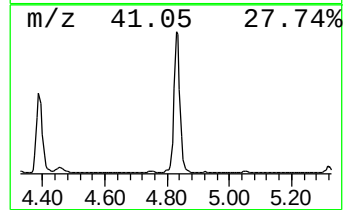
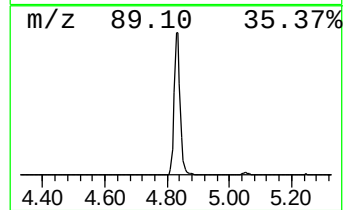
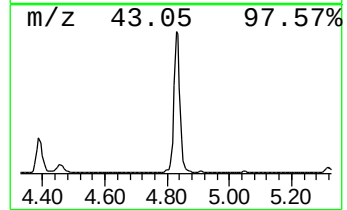
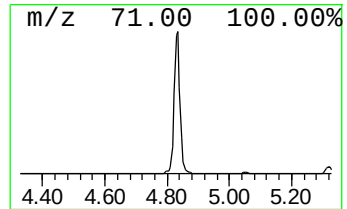
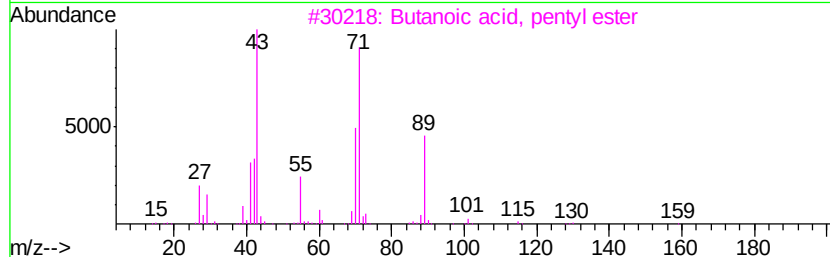
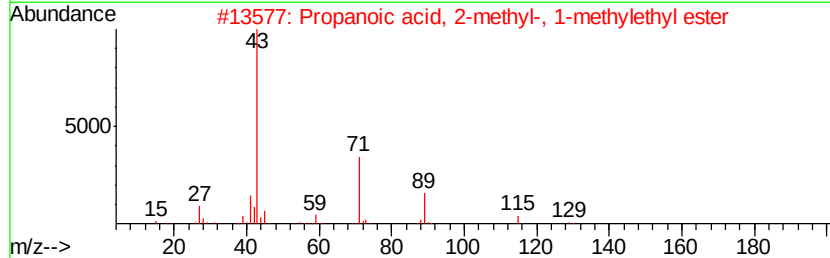
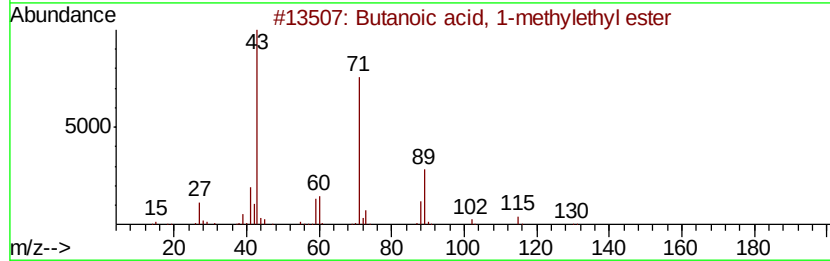
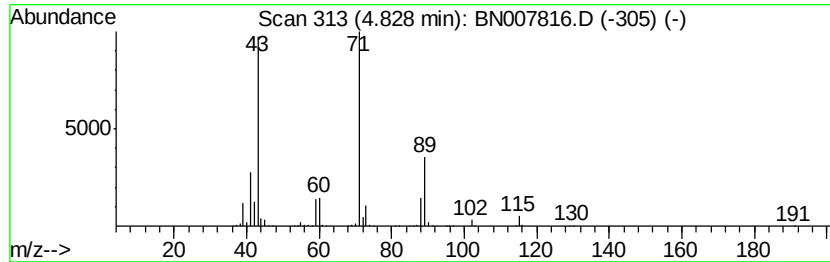
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Butanoic acid, 1-methylethy... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.83	20.65 ng/ul	3466590	1,4-Dichlorobenzene-d4	7.69

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Butanoic acid, 1-methylethyl ester	130	C7H14O2	000638-11-9	83
2		Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	72
3		Butanoic acid, pentyl ester	158	C9H18O2	000540-18-1	64
4		Butanoic acid, 1-methylethyl ester	130	C7H14O2	000638-11-9	64
5		Butanoic acid, 1-methylpropyl ester	144	C8H16O2	000819-97-6	56



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN092319\
Data File : BN007816.D
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Operator : HP/JU
Sample : K4895-08
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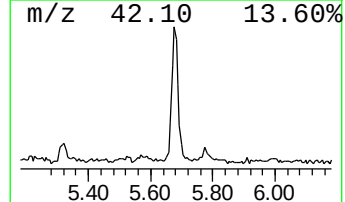
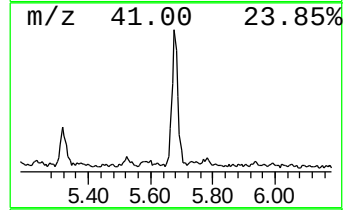
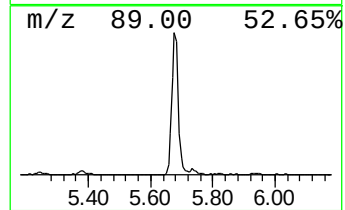
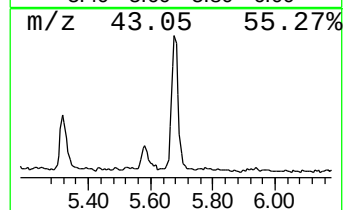
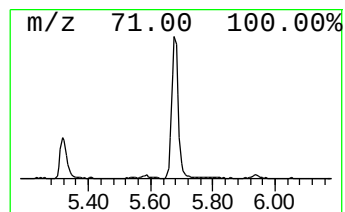
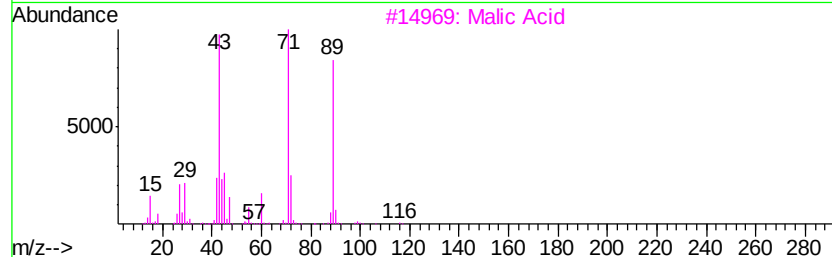
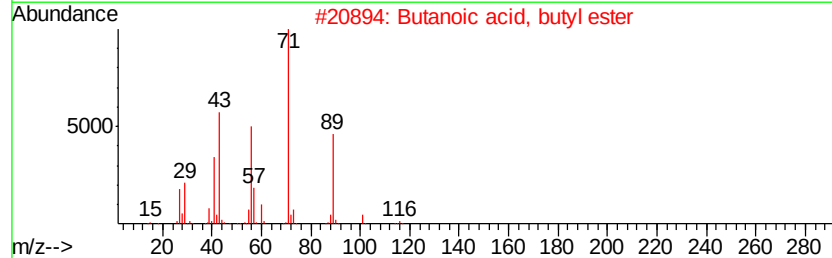
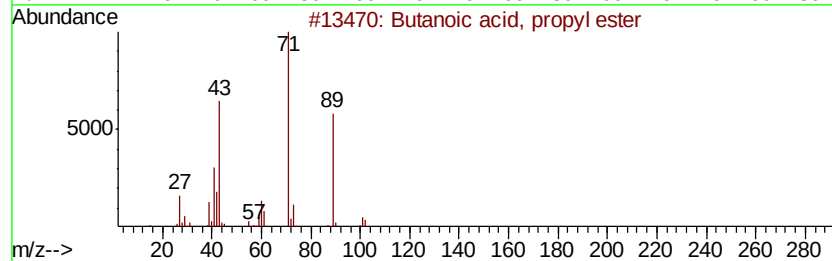
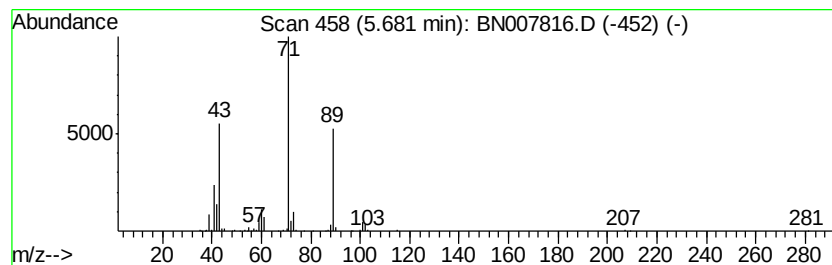
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Butanoic acid, propyl ester Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.	
5.68	3.12 ng/ul	524446	1,4-Dichlorobenzene-d4	7.69	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanoic acid, propyl ester	130	C7H14O2	000105-66-8	90
2	Butanoic acid, butyl ester	144	C8H16O2	000109-21-7	64
3	Malic Acid	134	C4H6O5	006915-15-7	64
4	Butanoic acid, hexyl ester	172	C10H20O2	002639-63-6	64
5	Butanoic acid, octyl ester	200	C12H24O2	000110-39-4	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN092319\
Data File : BN007816.D
Acq On : 23 Sep 2019 23:52
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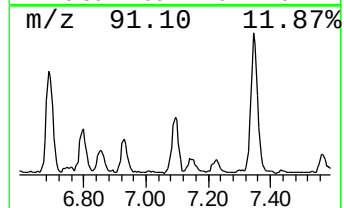
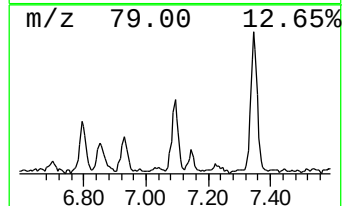
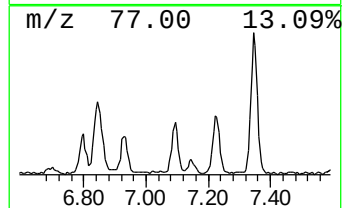
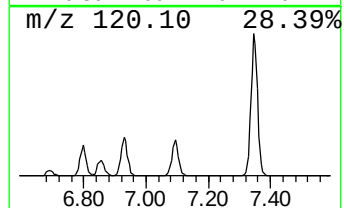
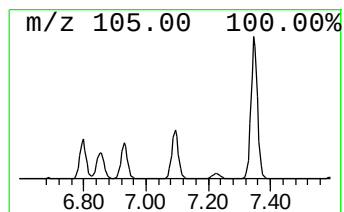
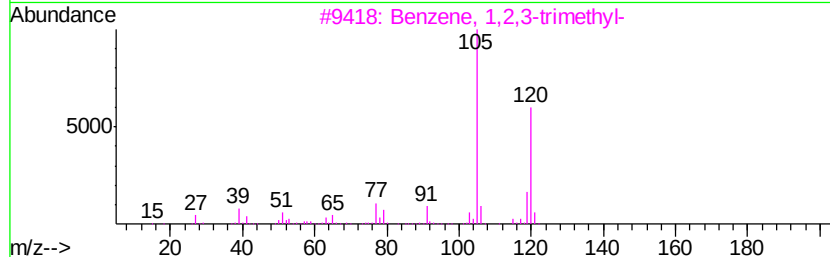
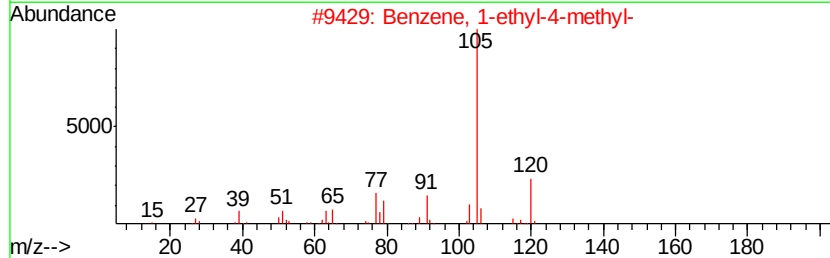
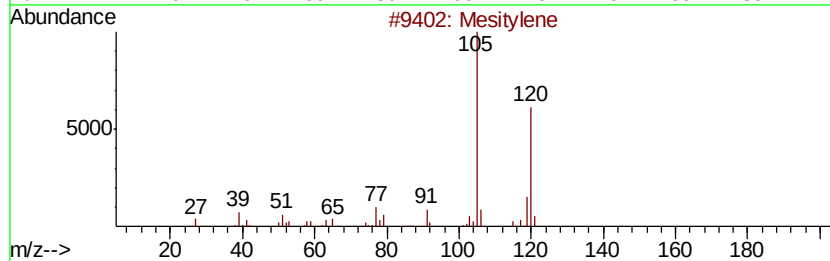
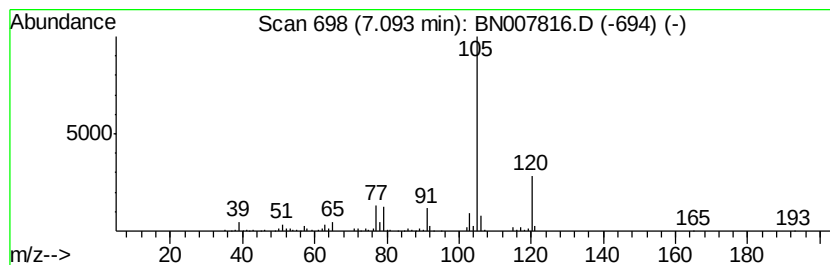
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Mesitylene Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.09	2.69 ng/ul	451703	1,4-Dichlorobenzene-d4	7.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Mesitylene	120	C9H12	000108-67-8	91
2		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
3		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
4		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	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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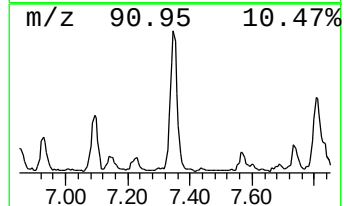
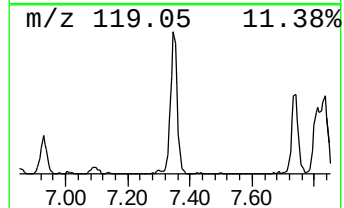
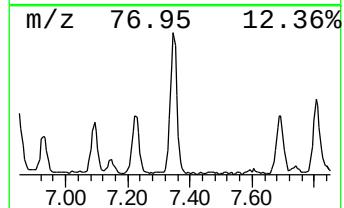
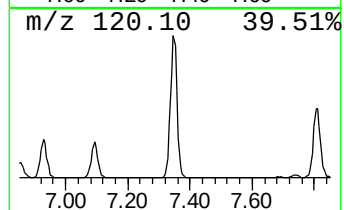
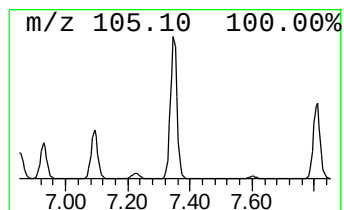
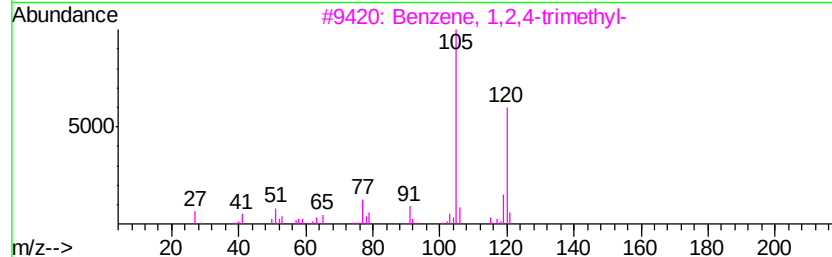
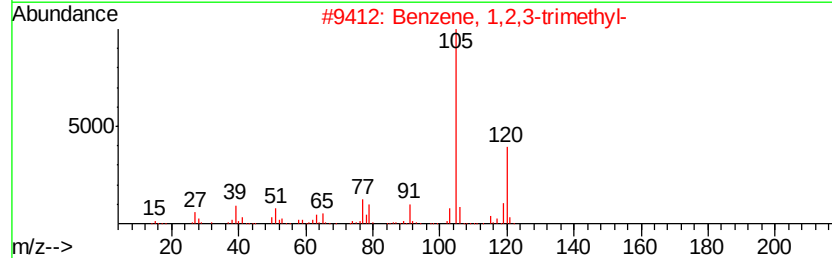
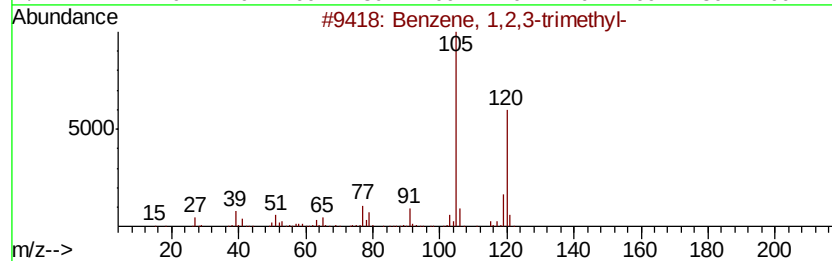
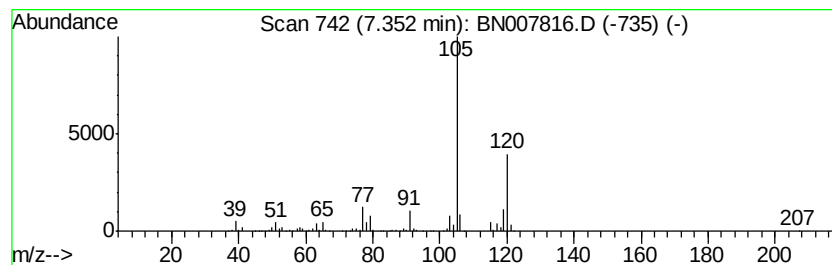
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.35	6.68 ng/ul	1121790	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	94
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	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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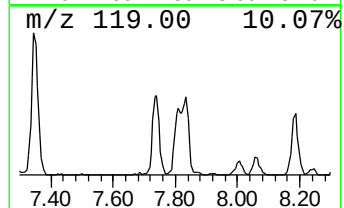
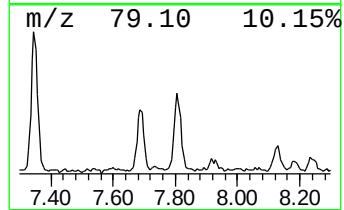
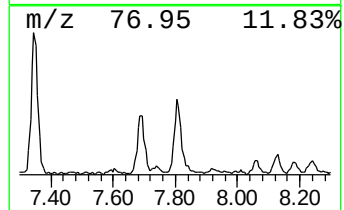
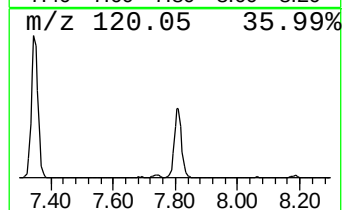
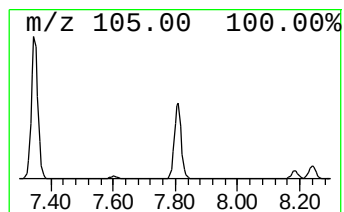
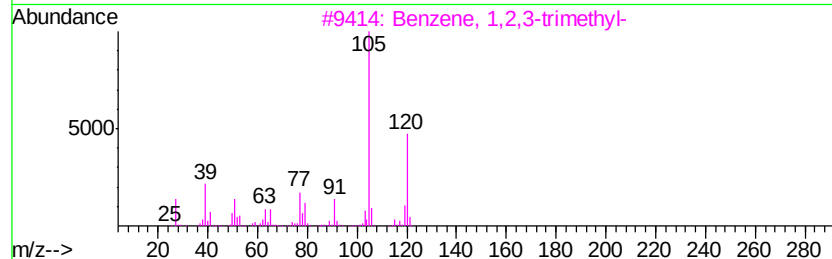
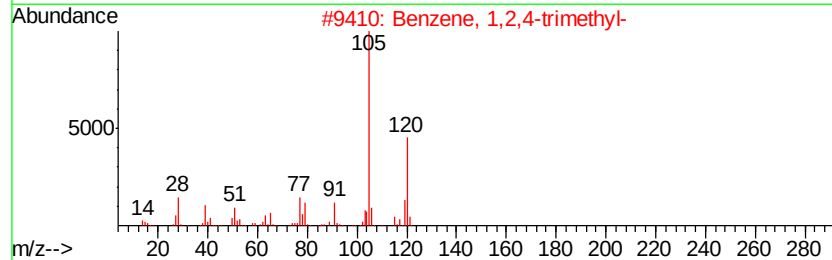
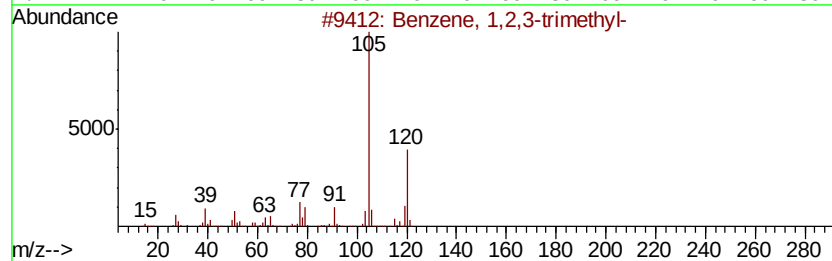
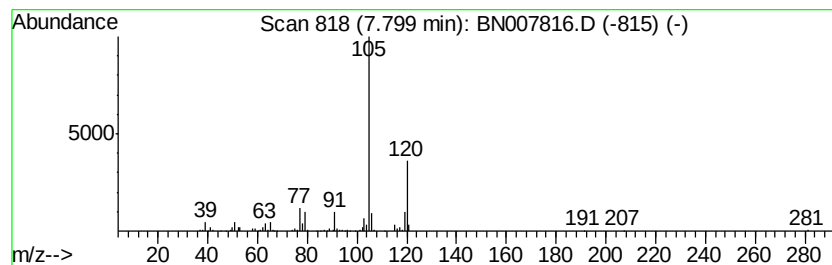
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Benzene, 1,2,4-trimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.80	3.81 ng/ul	640006	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, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
3		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
4		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5		Mesitylene	120	C9H12	000108-67-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN092319\
Data File : BN007816.D
Acq On : 23 Sep 2019 23:52
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ALS Vial : 18 Sample Multiplier: 1

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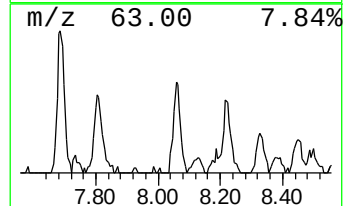
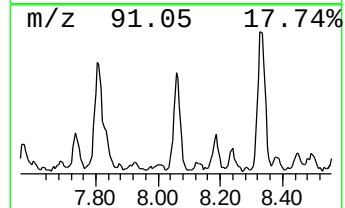
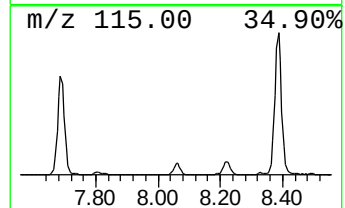
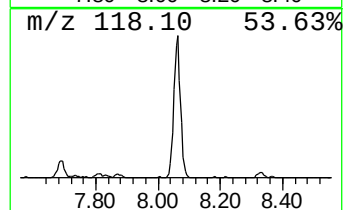
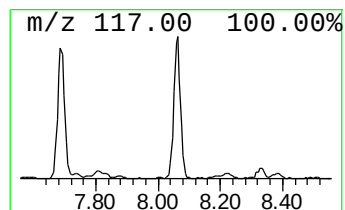
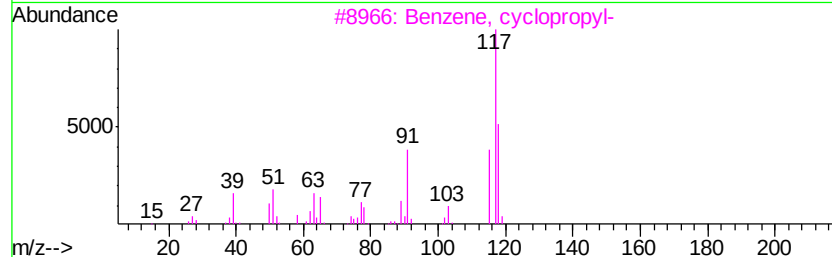
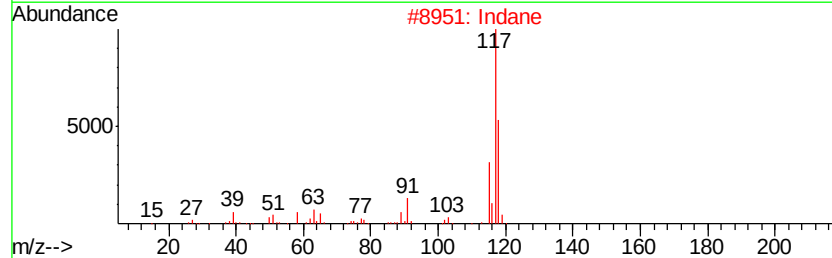
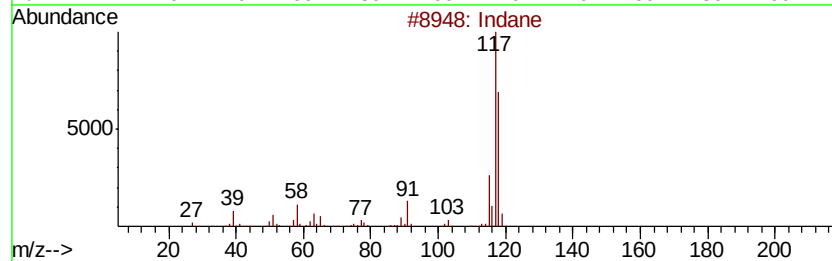
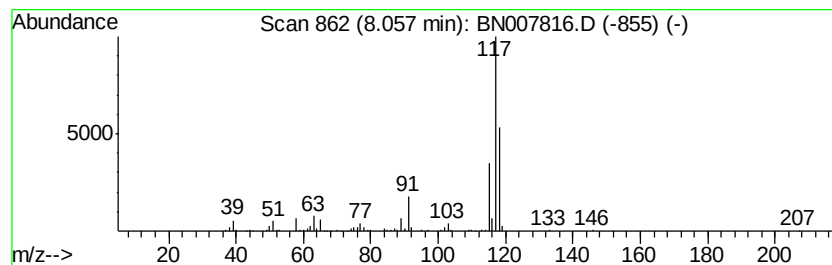
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Indane Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.06	2.26 ng/ul	378820	1,4-Dichlorobenzene-d4	7.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indane	118	C9H10	000496-11-7	87
2		Indane	118	C9H10	000496-11-7	81
3		Benzene, cyclopropyl-	118	C9H10	000873-49-4	74
4		Benzene, cyclopropyl-	118	C9H10	000873-49-4	72
5		Benzene, cyclopropyl-	118	C9H10	000873-49-4	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN092319\
Data File : BN007816.D
Acq On : 23 Sep 2019 23:52
Operator : HP/JU
Sample : K4895-08
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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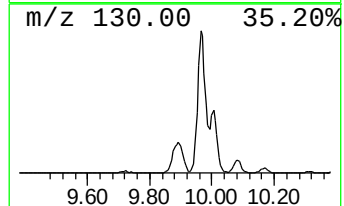
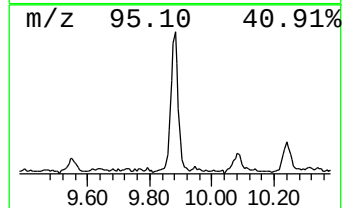
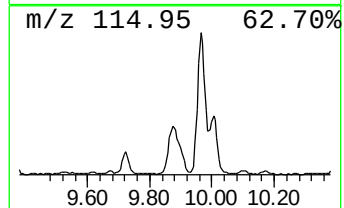
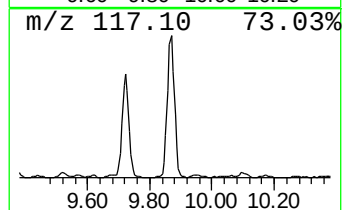
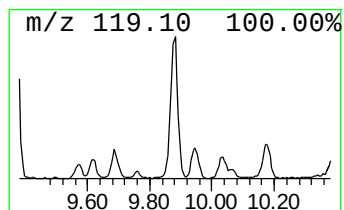
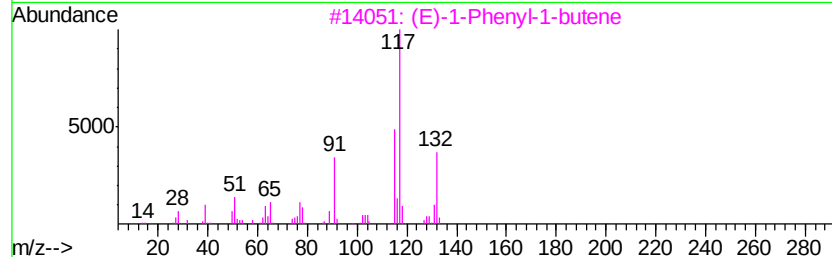
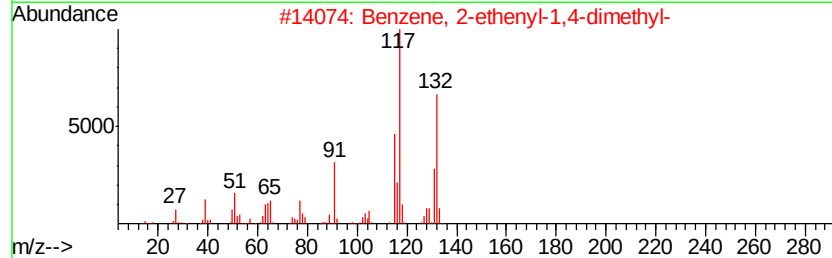
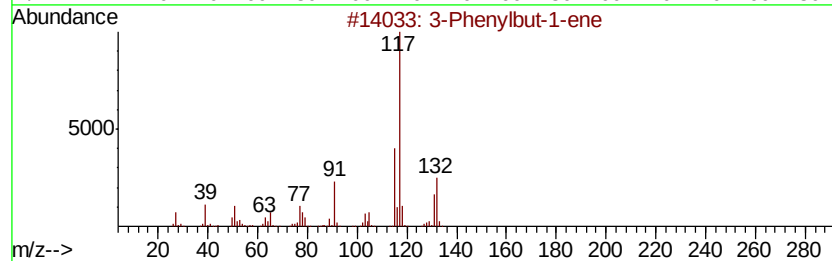
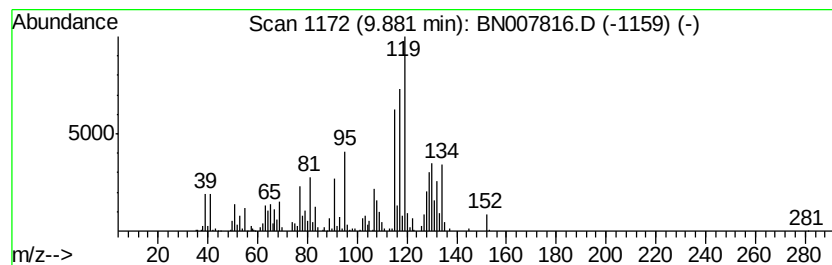
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 3-Phenylbut-1-ene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.88	5.40 ng/ul	1419880	Napthalene-d8	10.46

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Phenylbut-1-ene	132	C10H12	000934-10-1	50
2			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	50
3			(E)-1-Phenyl-1-butene	132	C10H12	001005-64-7	46
4			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	46
5			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	46



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN092319\
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Sample : K4895-08
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Instrument :
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ClientSampleId :
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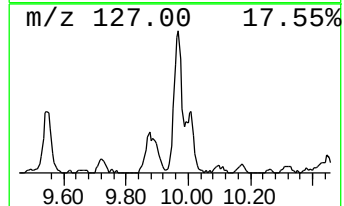
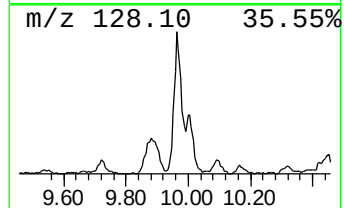
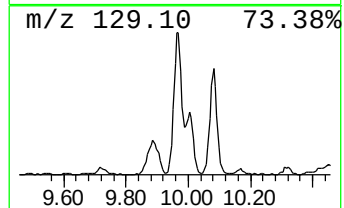
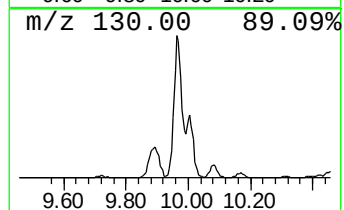
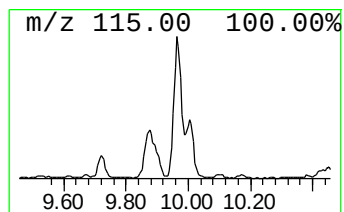
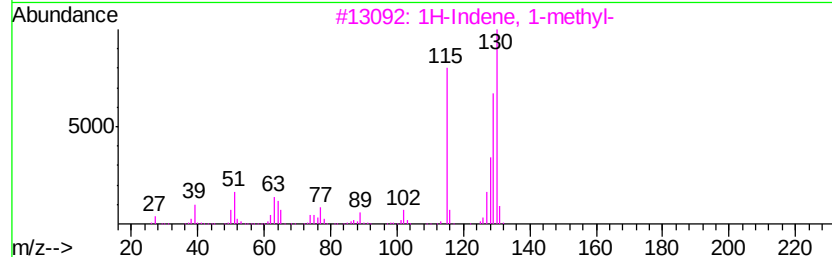
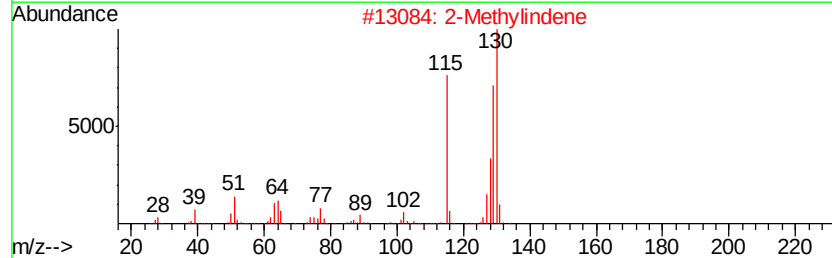
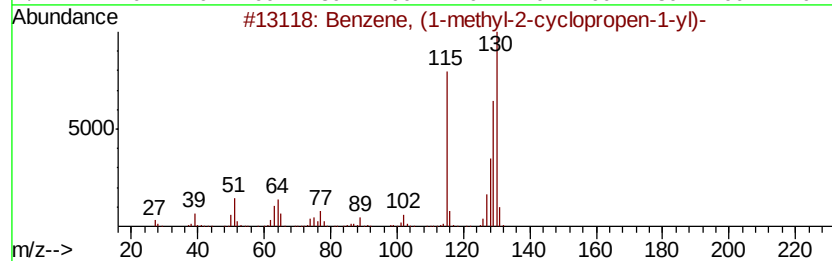
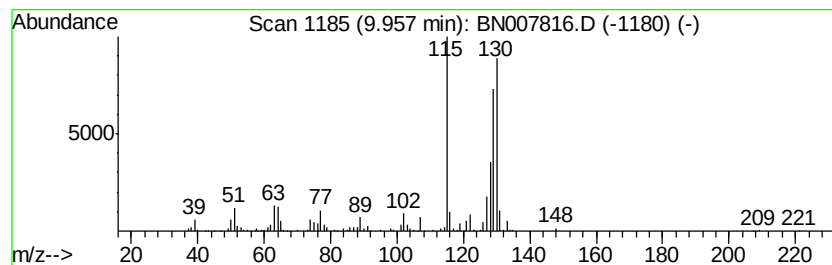
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Benzene, (1-methyl-2-cyclop... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.96	4.83 ng/ul	1271070	Naphthalene-d8	10.46

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (1-methyl-2-cyclopropen...	130	C10H10	065051-83-4	96
2		2-Methylindene	130	C10H10	002177-47-1	96
3		1H-Indene, 1-methyl-	130	C10H10	000767-59-9	95
4		Cycloprop[alindene, 1,1a,6,6a-te...	130	C10H10	015677-15-3	93
5		1H-Indene, 3-methyl-	130	C10H10	000767-60-2	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN092319\
Data File : BN007816.D
Acq On : 23 Sep 2019 23:52
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Sample : K4895-08
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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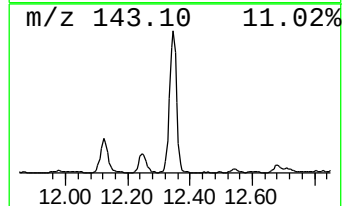
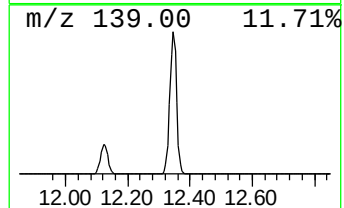
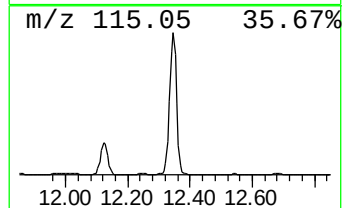
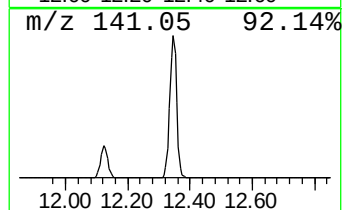
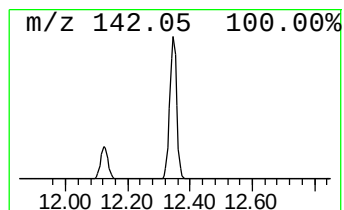
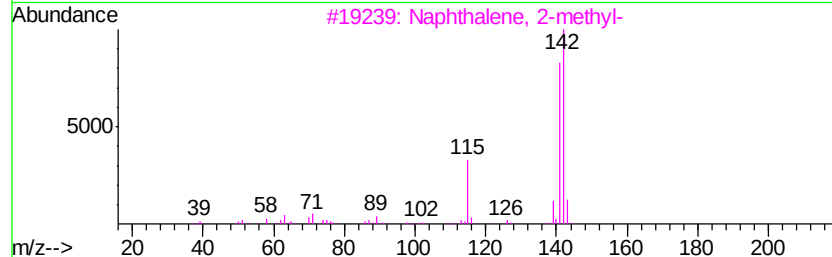
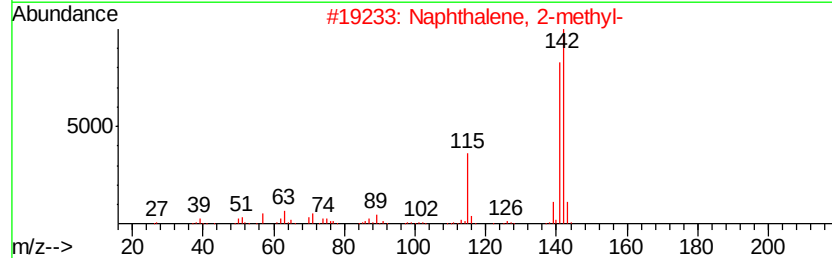
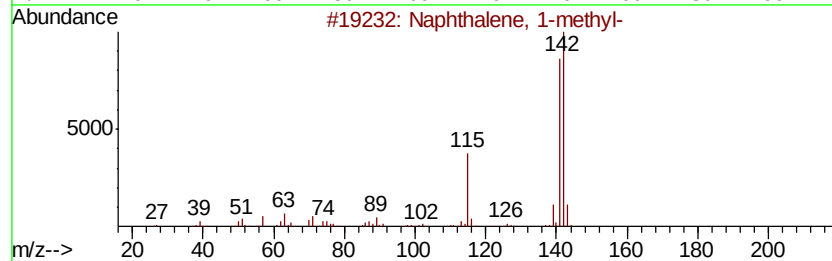
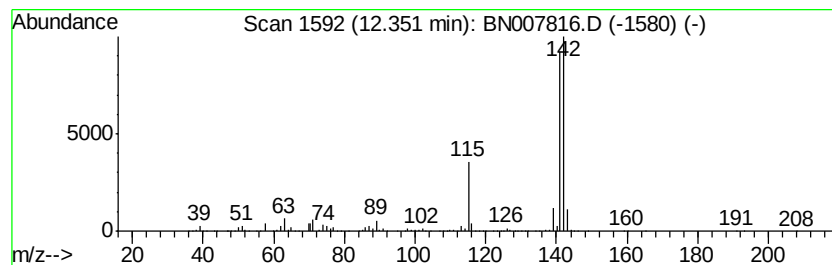
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 Naphthalene, 1-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.35	32.84 ng/ul	8640740	Naphthalene-d8	10.46

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
4		1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	94
5		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94



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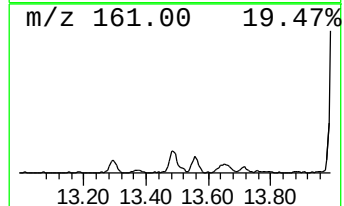
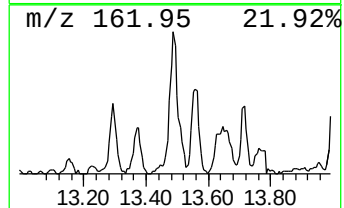
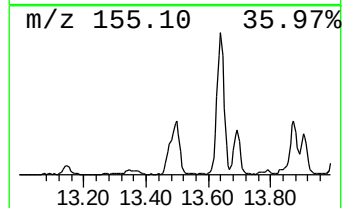
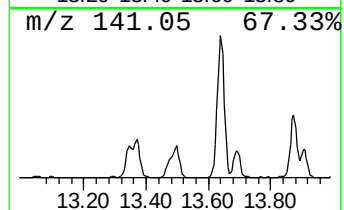
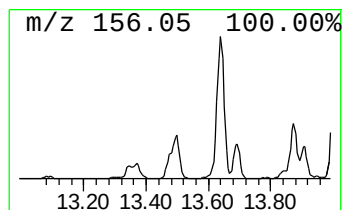
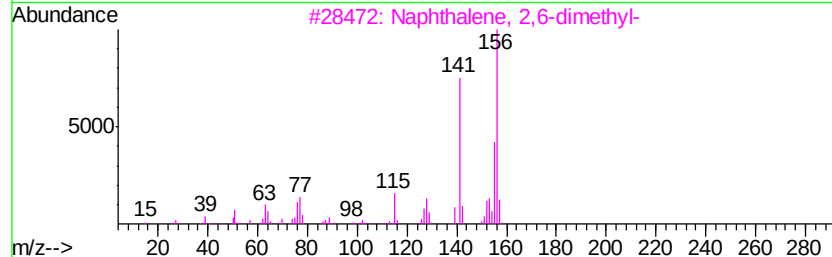
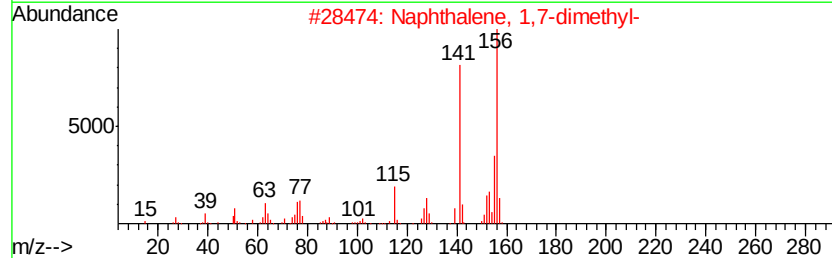
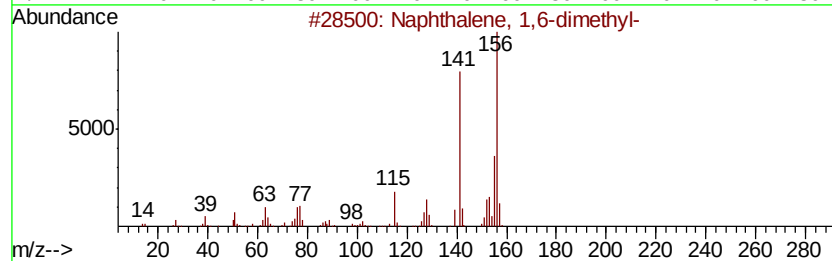
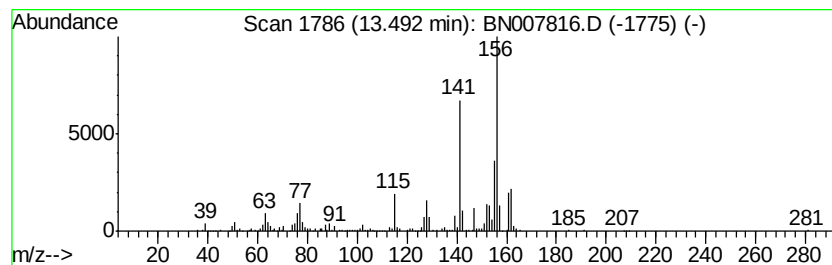
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 Naphthalene, 1,6-dimethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.49	2.84 ng/ul	998475	Acenaphthene-d10	14.32
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,6-dimethyl-	156 C12H12	000575-43-9 98
2		Naphthalene, 1,7-dimethyl-	156 C12H12	000575-37-1 98
3		Naphthalene, 2,6-dimethyl-	156 C12H12	000581-42-0 97
4		Naphthalene, 1,7-dimethyl-	156 C12H12	000575-37-1 96
5		Naphthalene, 2,7-dimethyl-	156 C12H12	000582-16-1 96



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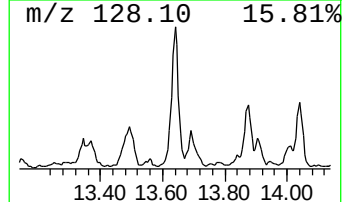
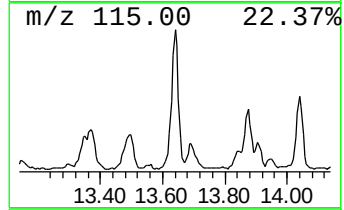
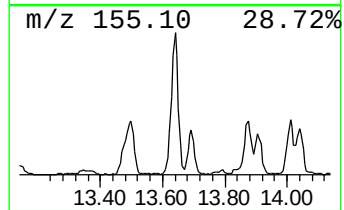
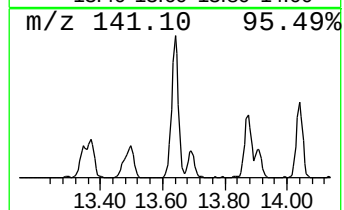
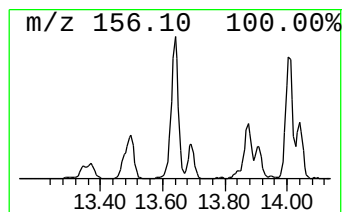
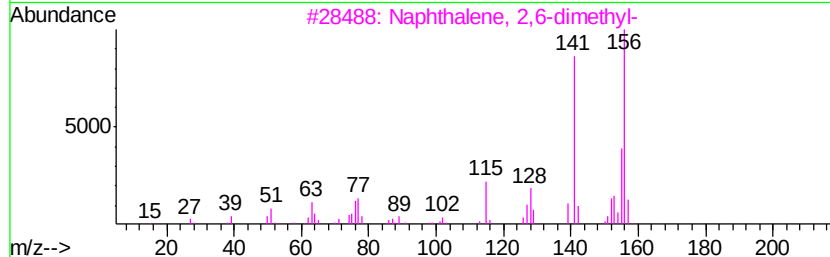
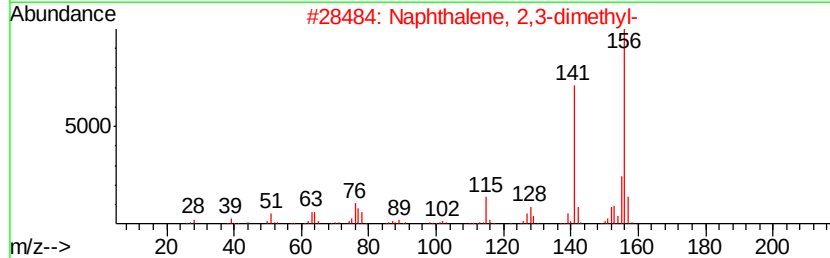
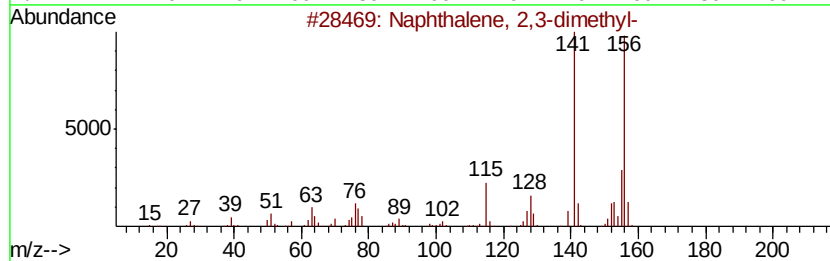
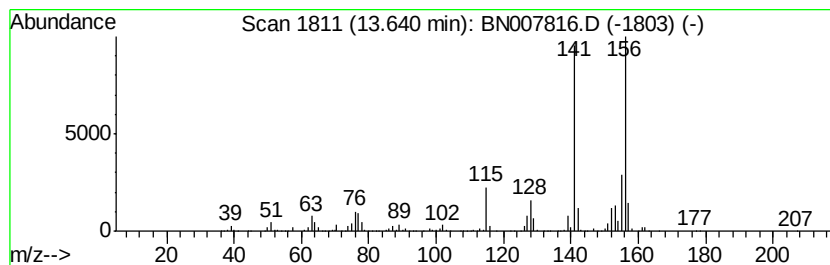
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 Naphthalene, 2,3-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.64	5.90 ng/ul	2078640	Acenaphthene-d10	14.32
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 2,3-dimethyl-	156 C12H12	000581-40-8 98
2		Naphthalene, 2,3-dimethyl-	156 C12H12	000581-40-8 97
3		Naphthalene, 2,6-dimethyl-	156 C12H12	000581-42-0 96
4		Naphthalene, 1,3-dimethyl-	156 C12H12	000575-41-7 96
5		Naphthalene, 2,3-dimethyl-	156 C12H12	000581-40-8 96



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN092319\
Data File : BN007816.D
Acq On : 23 Sep 2019 23:52
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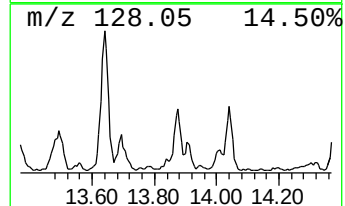
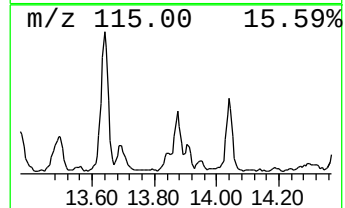
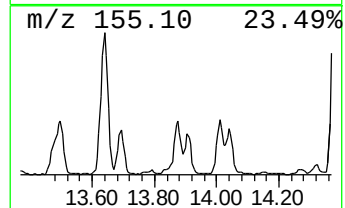
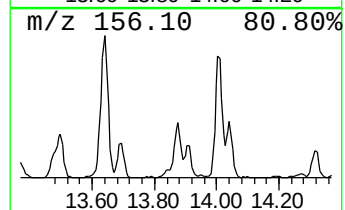
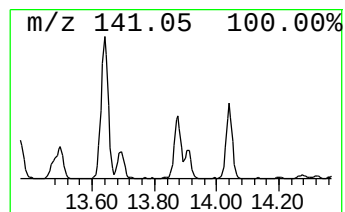
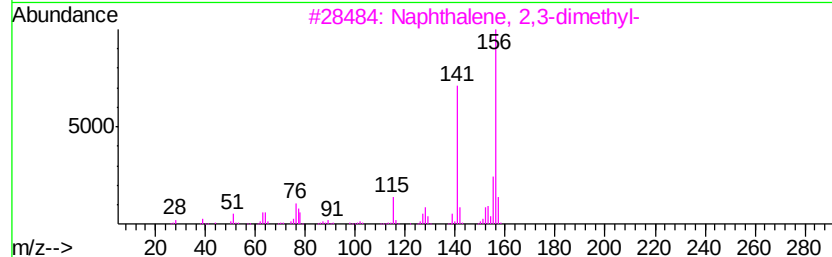
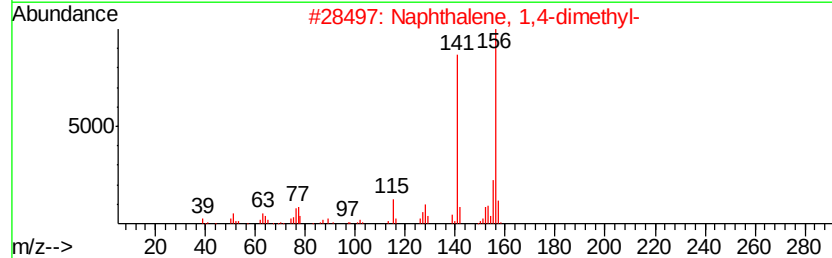
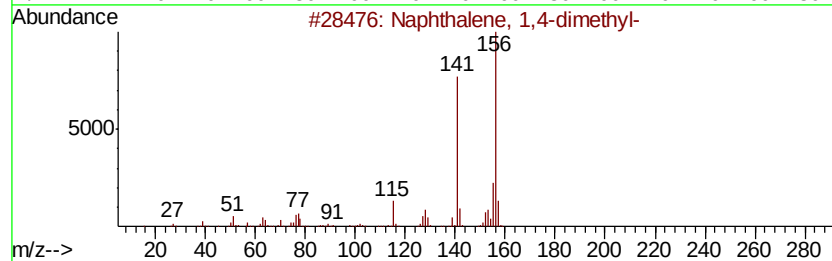
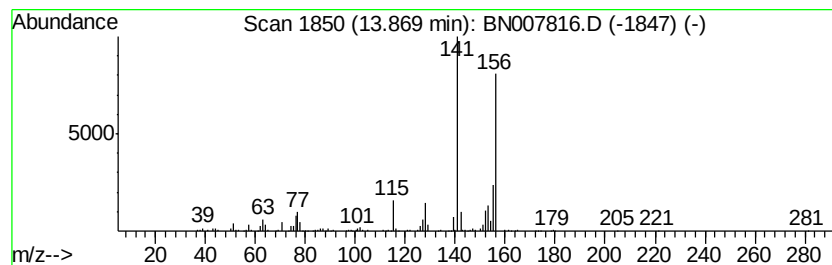
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 Naphthalene, 1,4-dimethyl- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.87	2.01 ng/ul	707753	Acenaphthene-d10	14.32

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97
2			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97
3			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	95
4			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	94
5			Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN092319\
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Sample : K4895-08
Misc :
ALS Vial : 18 Sample Multiplier: 1

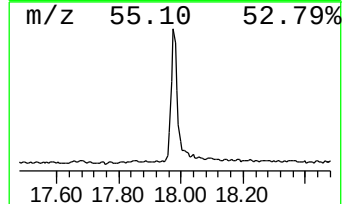
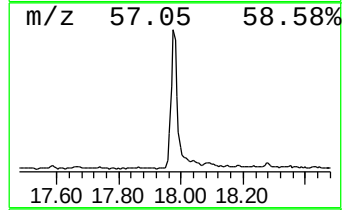
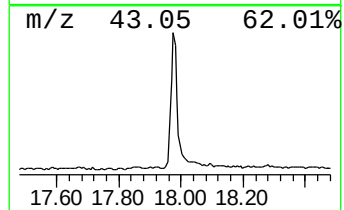
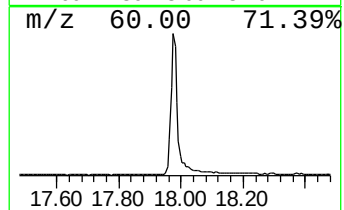
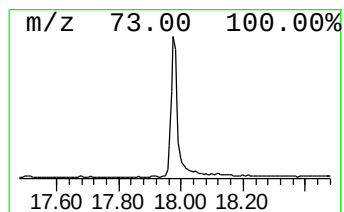
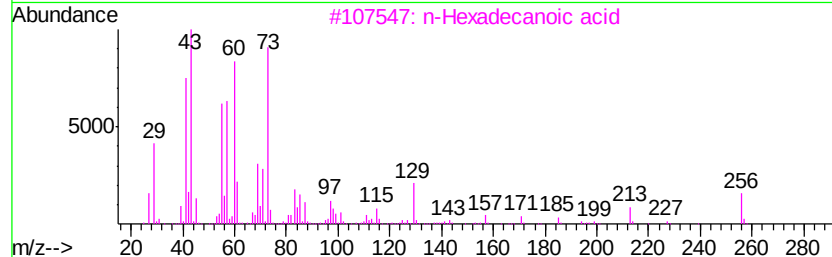
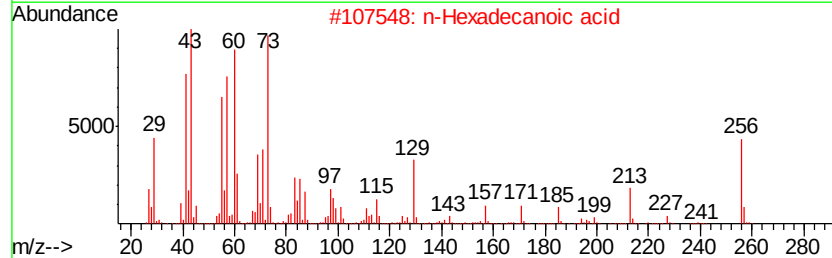
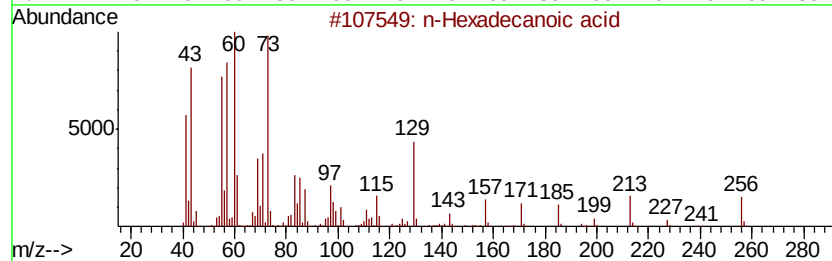
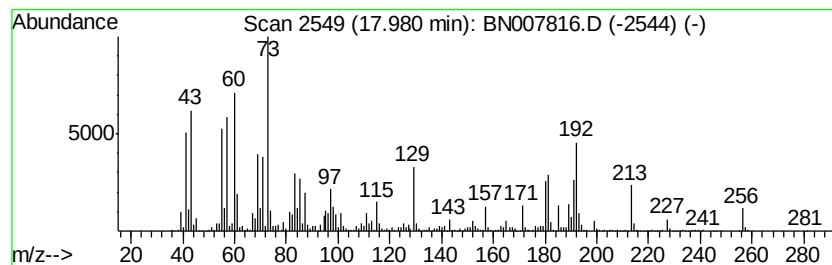
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN091619MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 18 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.98	8.03 ng/ul	3222810	Phenanthrene-d10	17.06	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
4	Tridecanoic acid	214	C13H26O2	000638-53-9	55
5	Tridecanoic acid	214	C13H26O2	000638-53-9	46



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN092319\
Data File : BN007816.D
Acq On : 23 Sep 2019 23:52
Operator : HP/JU
Sample : K4895-08
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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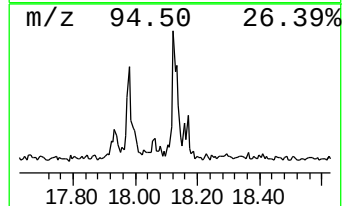
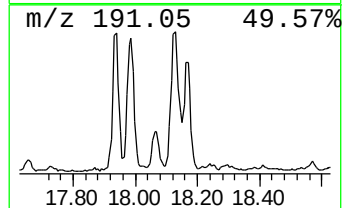
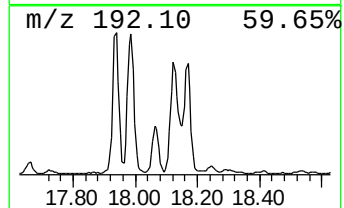
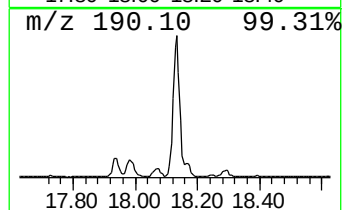
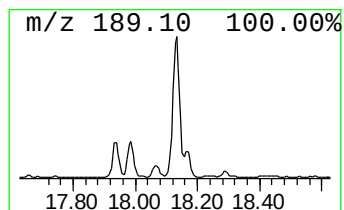
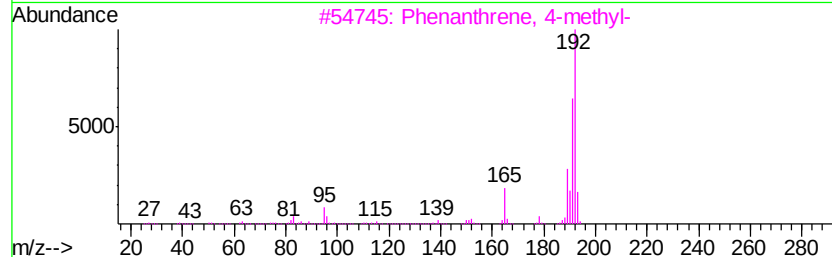
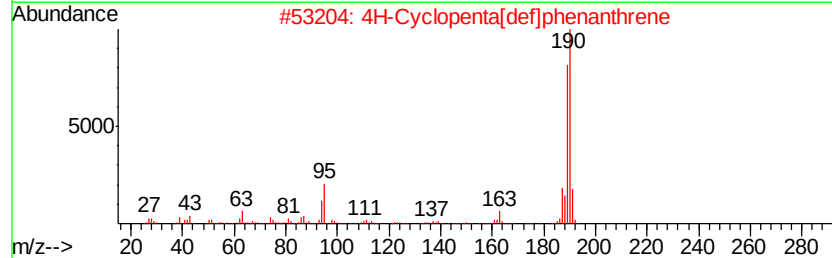
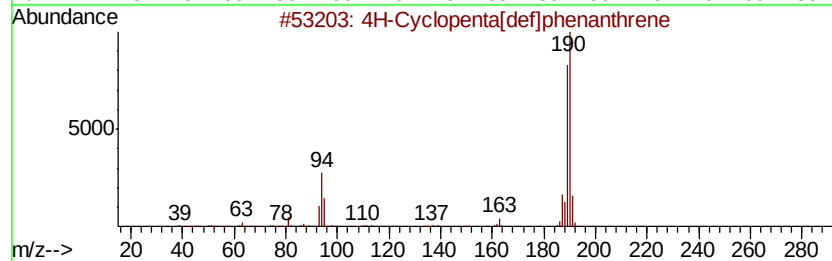
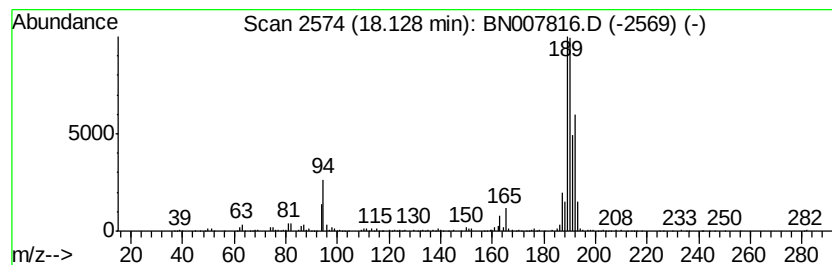
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 4H-Cyclopenta[def]phenanthrene Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.13	2.91 ng/ul	1168930	Phenanthrene-d10	17.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	70
2		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	55
3		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	35
4		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	30
5		Benzene, 1,1'-(1,2-propadienyld...	192	C15H12	014251-57-1	25



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN092319\
Data File : BN007816.D
Acq On : 23 Sep 2019 23:52
Operator : HP/JU
Sample : K4895-08
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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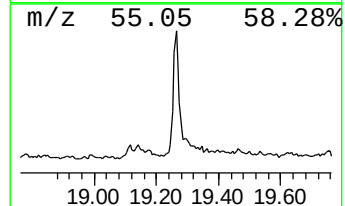
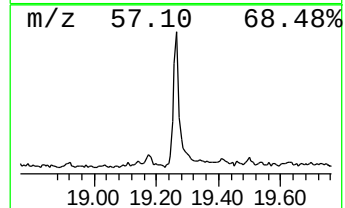
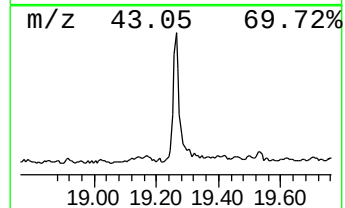
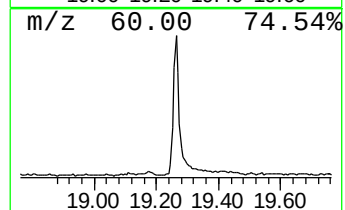
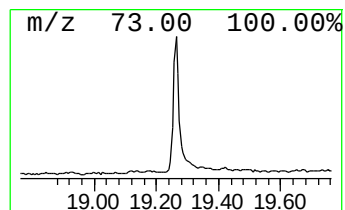
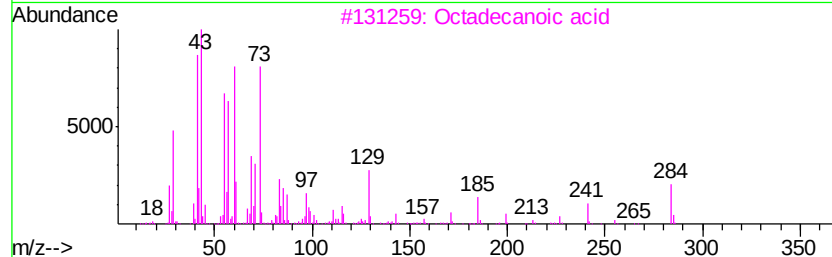
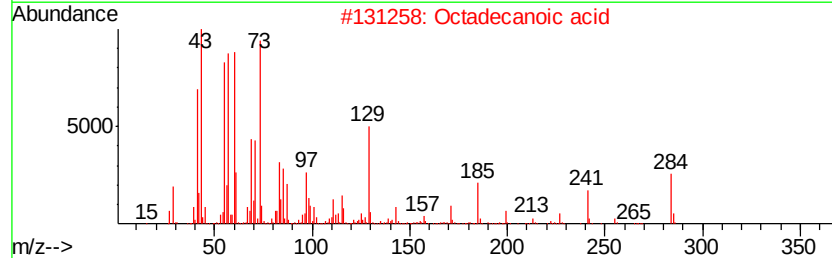
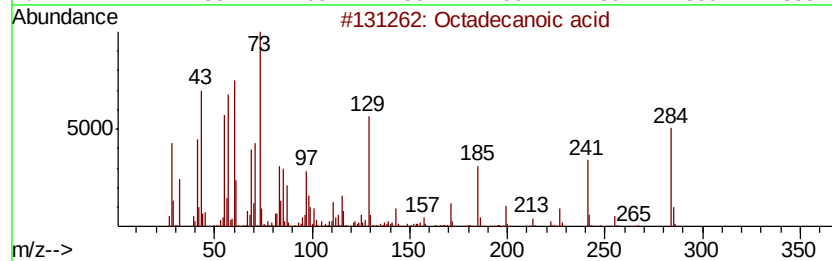
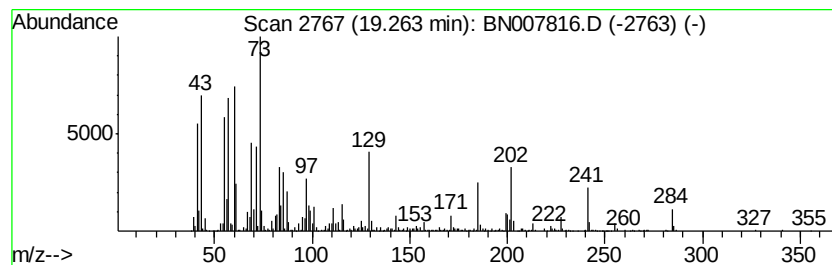
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 Octadecanoic acid Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.26	3.26 ng/ul	1689330	Chrysene-d12	21.26

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Octadecanoic acid	284	C18H36O2	000057-11-4	99
3			Octadecanoic acid	284	C18H36O2	000057-11-4	96
4			Octadecanoic acid	284	C18H36O2	000057-11-4	95
5			Octadecanoic acid	284	C18H36O2	000057-11-4	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN092319\
Data File : BN007816.D
Acq On : 23 Sep 2019 23:52
Operator : HP/JU
Sample : K4895-08
Misc :
ALS Vial : 18 Sample Multiplier: 1

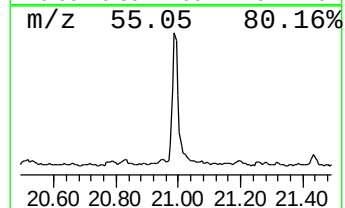
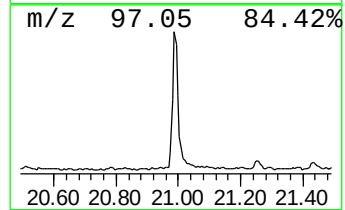
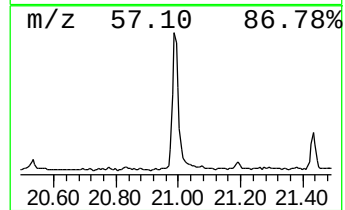
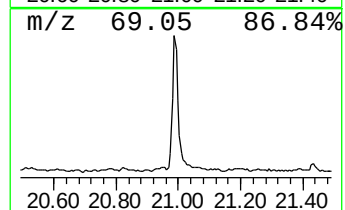
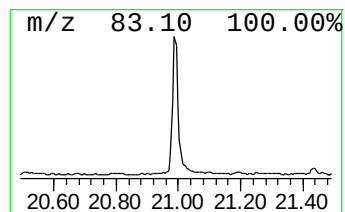
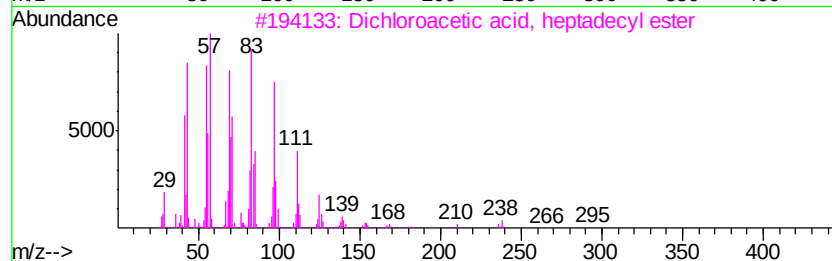
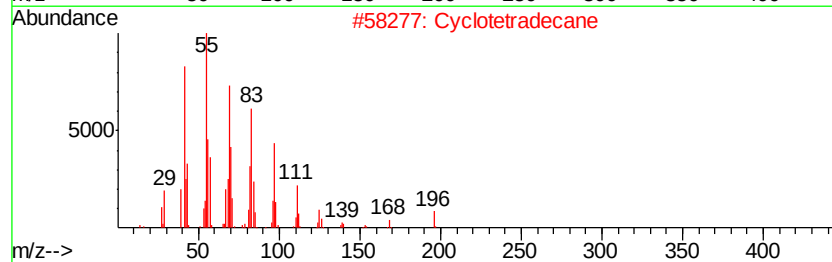
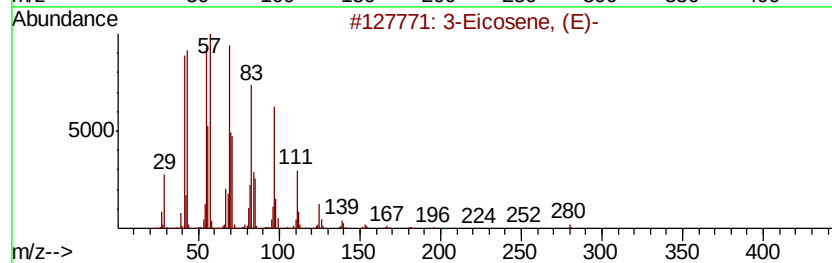
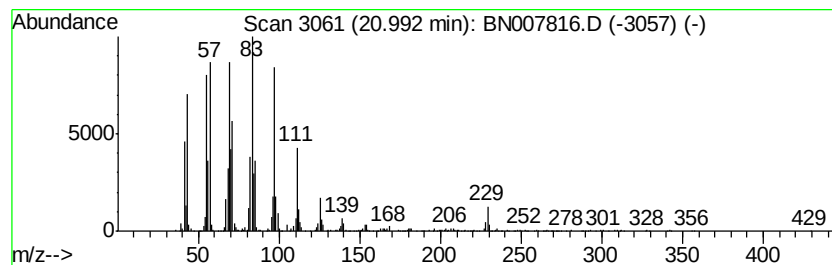
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN091619MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 21 (DEL) Alkane: Cyclic20.99 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.99	3.86 ng/ul	2001620	Chrysene-d12	21.26
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		3-Eicosene, (E)-	280 C20H40	074685-33-9 99
2		Cyclotetradecane	196 C14H28	000295-17-0 93
3		Dichloroacetic acid, heptadecyl ...	366 C19H36Cl2O2	1000282-98-2 91
4		Bromoacetic acid, octadecyl ester	390 C20H39BrO2	018992-03-5 91
5		Heptadecyl heptafluorobutyrate	452 C21H35F7O2	959085-66-6 91



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN092319\
Data File : BN007816.D
Acq On : 23 Sep 2019 23:52
Operator : HP/JU
Sample : K4895-08
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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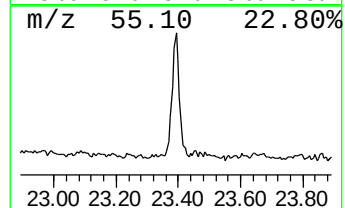
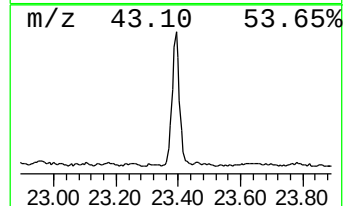
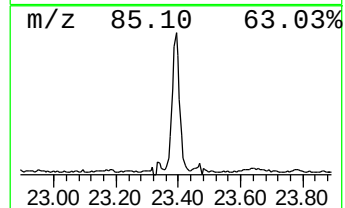
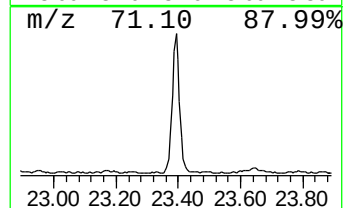
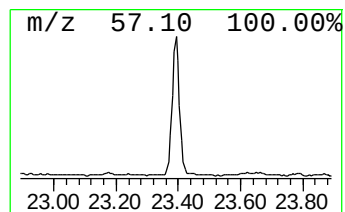
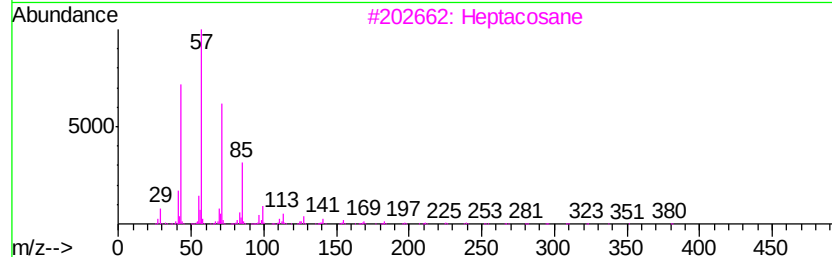
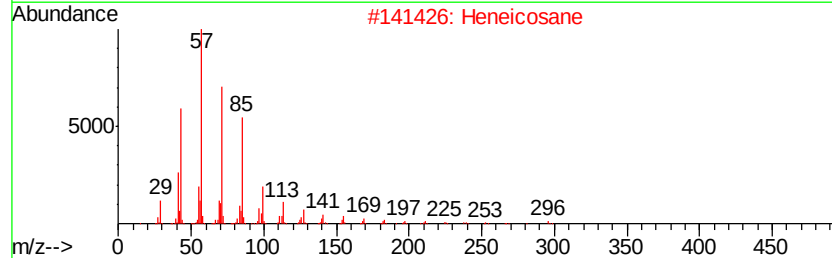
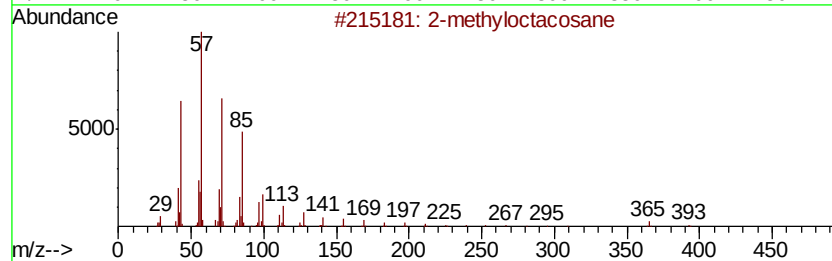
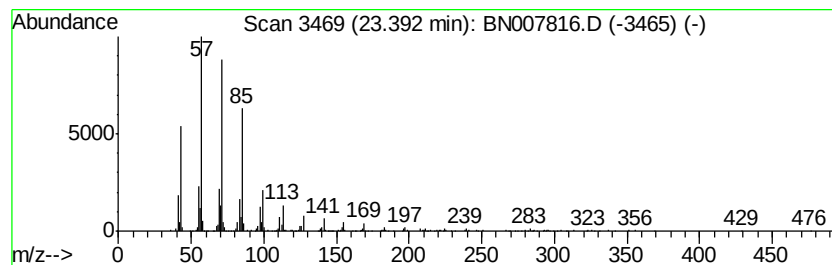
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 (DEL) Alkane: Straight-Chai... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.39	2.38 ng/ul	1271960	Perylene-d12	23.47

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-methyloctacosane	408	C29H60	1000376-72-8	91
2			Heneicosane	296	C21H44	000629-94-7	91
3			Heptacosane	380	C27H56	000593-49-7	91
4			triacontane	422	C30H62	000638-68-6	90
5			Heptacosane	380	C27H56	000593-49-7	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN092319\
Data File : BN007816.D
Acq On : 23 Sep 2019 23:52
Operator : HP/JU
Sample : K4895-08
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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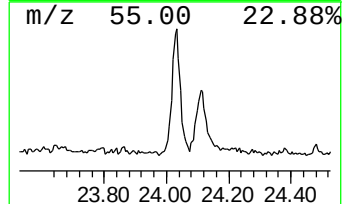
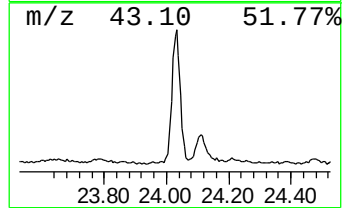
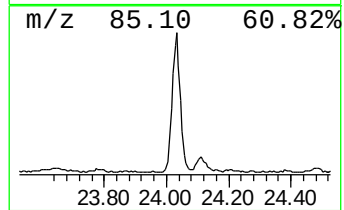
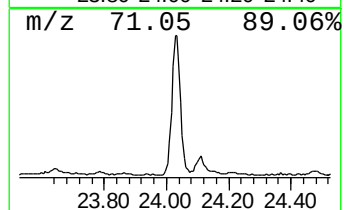
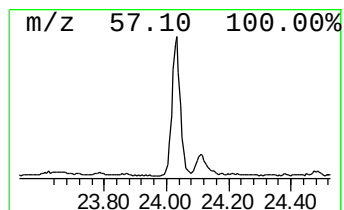
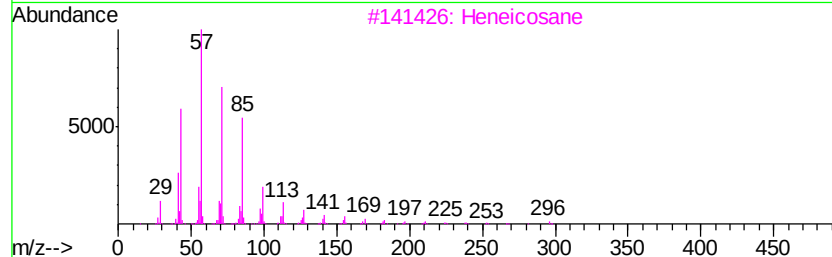
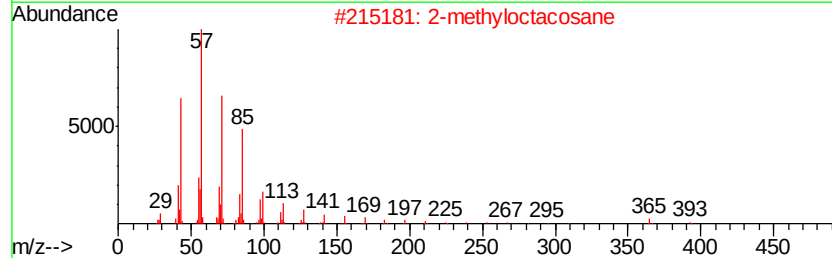
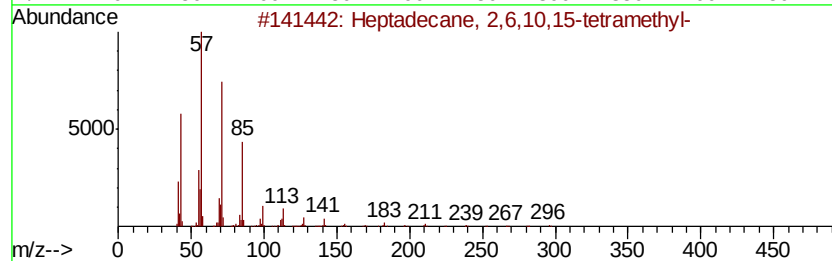
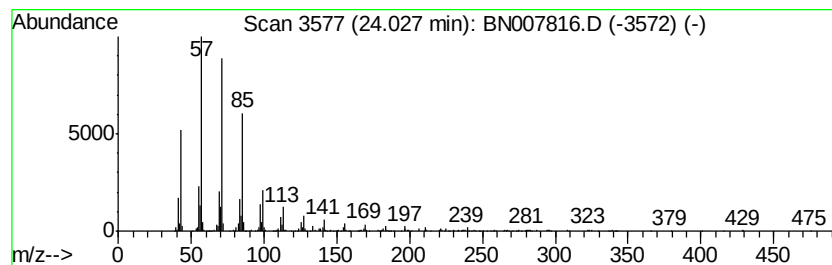
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Quant Title : SVOA CALIBRATION

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

Peak Number 23 (DEL) Alkane: Straight-Chai... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.03	2.29 ng/ul	1227010	Perylene-d12	23.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	91
2		2-methyloctacosane	408	C29H60	1000376-72-8	91
3		Heneicosane	296	C21H44	000629-94-7	91
4		Nonadecane	268	C19H40	000629-92-5	91
5		Hexacosane	366	C26H54	000630-01-3	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN092319\
 Data File : BN007816.D
 Acq On : 23 Sep 2019 23:52
 Operator : HP/JU
 Sample : K4895-08
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BFDH8

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	793921	1	7.69	3356900	20.0
Isobutyl acetate	3.94	2.3	ng/ul	387824	1	7.69	3356900	20.0
Propanoic acid, 2...	4.22	11.0	ng/ul	1846080	1	7.69	3356900	20.0
Propanoic acid, p...	4.39	15.4	ng/ul	2582690	1	7.69	3356900	20.0
Butanoic acid, 1-...	4.83	20.6	ng/ul	3466590	1	7.69	3356900	20.0
Butanoic acid, pr...	5.68	3.1	ng/ul	524446	1	7.69	3356900	20.0
Mesitylene	7.09	2.7	ng/ul	451703	1	7.69	3356900	20.0
Benzene, 1,2,3-tr...	7.35	6.7	ng/ul	1121790	1	7.69	3356900	20.0
Benzene, 1,2,4-tr...	7.80	3.8	ng/ul	640006	1	7.69	3356900	20.0
Indane	8.06	2.3	ng/ul	378820	1	7.69	3356900	20.0
3-Phenylbut-1-ene	9.88	5.4	ng/ul	1419880	2	10.46	5262120	20.0
Benzene, (1-methy...	9.96	4.8	ng/ul	1271070	2	10.46	5262120	20.0
Naphthalene, 1-me...	12.35	32.8	ng/ul	8640740	2	10.46	5262120	20.0
Naphthalene, 1,6-...	13.49	2.8	ng/ul	998475	3	14.32	7040340	20.0
Naphthalene, 2,3-...	13.64	5.9	ng/ul	2078640	3	14.32	7040340	20.0
Naphthalene, 1,4-...	13.87	2.0	ng/ul	707753	3	14.32	7040340	20.0
n-Hexadecanoic acid	17.98	8.0	ng/ul	3222810	4	17.06	8026960	20.0
4H-Cyclopenta[def...	18.13	2.9	ng/ul	1168930	4	17.06	8026960	20.0
Octadecanoic acid	19.26	3.3	ng/ul	1689330	5	21.26	10368500	20.0
(DEL) Alkane: Cyc...	20.99	3.9	ng/ul	2001620	5	21.26	10368500	20.0
(DEL) Alkane: Str...	23.39	2.4	ng/ul	1271960	6	23.47	10693800	20.0
(DEL) Alkane: Str...	24.03	2.3	ng/ul	1227010	6	23.47	10693800	20.0