

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN040320\
 Data File : BN010407.D
 Acq On : 03 Apr 2020 11:52
 Operator : CG/JU
 Sample : L2095-01DL 4X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBF16DL

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-BN040120MA.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	2.893	14	18	27	rBV3	60767	111277	0.76%	0.092%
2	2.964	27	30	39	rVB	111063	167735	1.14%	0.138%
3	3.246	70	78	92	rBV	572727	911120	6.22%	0.750%
4	3.652	143	147	151	rBV	67889	91264	0.62%	0.075%
5	4.170	230	235	242	rBV	167747	242350	1.65%	0.200%
6	4.334	259	263	271	rBV	230501	343685	2.34%	0.283%
7	4.775	327	338	349	rBV	338330	513401	3.50%	0.423%
8	5.617	475	481	489	rVV	53740	86575	0.59%	0.071%
9	6.793	673	681	688	rBV	82657	144354	0.98%	0.119%
10	6.952	702	708	718	rBV	288602	481307	3.28%	0.396%
11	7.152	736	742	751	rVB	329168	528285	3.60%	0.435%
12	7.611	814	820	827	rVB	1215317	1953257	13.32%	1.608%
13	7.981	876	883	889	rVB	78957	126813	0.87%	0.104%
14	8.140	905	910	916	rVB	39148	59617	0.41%	0.049%
15	8.310	933	939	950	rVB	265623	413089	2.82%	0.340%
16	8.746	1001	1013	1024	rBV	337974	582946	3.98%	0.480%
17	9.457	1128	1134	1144	rBV	240559	404962	2.76%	0.333%
18	9.993	1218	1225	1234	rBV	490059	841252	5.74%	0.693%
19	10.369	1281	1289	1293	rBV	1790772	3043087	20.76%	2.505%
20	10.416	1293	1297	1306	rVV	2153584	3536789	24.13%	2.912%
21	10.505	1306	1312	1325	rVB2	359448	739357	5.04%	0.609%
22	11.904	1545	1550	1558	rBV2	119075	211278	1.44%	0.174%
23	12.034	1565	1572	1581	rBV	1346565	2052816	14.00%	1.690%
24	12.157	1587	1593	1600	rBV4	75835	134828	0.92%	0.111%
25	12.257	1600	1610	1616	rVB	788722	1276071	8.70%	1.051%
26	13.063	1739	1747	1756	rBV	671648	991283	6.76%	0.816%
27	13.257	1775	1780	1790	rVB	199411	357633	2.44%	0.294%
28	13.551	1821	1830	1835	rBV	321573	572925	3.91%	0.472%
29	13.610	1835	1840	1842	rVV	199809	286939	1.96%	0.236%
30	13.646	1842	1846	1852	rVB	1105916	1530917	10.44%	1.260%
31	13.793	1865	1871	1875	rVV	164059	264688	1.81%	0.218%
32	13.922	1886	1893	1898	rBV	1167149	1760317	12.01%	1.449%
33	14.234	1936	1946	1951	rBV	3335775	4952634	33.78%	4.077%
34	14.298	1951	1957	1964	rVV	6292598	9055123	61.77%	7.454%

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 Title : SVOA CALIBRATION

35	14.357	1964	1967	1975	rVV	243611	355143	2.42%	0.292%
36	14.445	1975	1982	1983	rVV3	66770	112260	0.77%	0.092%
37	14.469	1983	1986	1991	rVV	143369	211963	1.45%	0.174%
38	14.551	1995	2000	2005	rVV2	183348	261386	1.78%	0.215%
39	14.634	2005	2014	2023	rVV	3210727	4548969	31.03%	3.745%
40	14.716	2025	2028	2034	rVB2	45227	65092	0.44%	0.054%
41	14.863	2045	2053	2058	rBV5	34219	70429	0.48%	0.058%
42	15.034	2073	2082	2085	rBV5	37158	69038	0.47%	0.057%
43	15.175	2099	2106	2110	rBV8	37657	71481	0.49%	0.059%
44	15.228	2110	2115	2120	rVV	1716613	2546764	17.37%	2.097%
45	15.287	2120	2125	2131	rVV2	4357189	5956750	40.63%	4.904%
46	15.345	2131	2135	2138	rVV2	98288	170585	1.16%	0.140%
47	15.381	2138	2141	2143	rVV	136443	189712	1.29%	0.156%
48	15.410	2143	2146	2149	rVV	216405	318505	2.17%	0.262%
49	15.451	2149	2153	2158	rVV	433706	634040	4.33%	0.522%
50	15.498	2158	2161	2164	rVV2	79934	116245	0.79%	0.096%
51	15.534	2164	2167	2171	rVV	179878	245968	1.68%	0.202%
52	15.587	2171	2176	2182	rVB	358203	498587	3.40%	0.410%
53	15.728	2195	2200	2208	rVV	359012	714799	4.88%	0.588%
54	15.816	2212	2215	2222	rVB2	68804	103618	0.71%	0.085%
55	15.951	2233	2238	2245	rVB2	101779	164794	1.12%	0.136%
56	16.081	2253	2260	2270	rBV	260023	381388	2.60%	0.314%
57	16.251	2285	2289	2291	rBV2	80963	133680	0.91%	0.110%
58	16.292	2291	2296	2301	rVV	208253	352806	2.41%	0.290%
59	16.339	2301	2304	2311	rVV2	79922	105507	0.72%	0.087%
60	16.439	2316	2321	2326	rVB2	101019	158575	1.08%	0.131%
61	16.598	2345	2348	2351	rVV4	39616	55891	0.38%	0.046%
62	16.634	2351	2354	2362	rVB2	59024	89308	0.61%	0.074%
63	16.739	2365	2372	2375	rBV2	71689	146264	1.00%	0.120%
64	16.798	2375	2382	2388	rVB	784082	1113294	7.59%	0.917%
65	16.981	2407	2413	2416	rVV	4346443	6123189	41.77%	5.041%
66	17.022	2416	2420	2425	rVV	10783992	14659875	100.00%	12.069%
67	17.075	2425	2429	2433	rVV	1971464	2956275	20.17%	2.434%
68	17.110	2433	2435	2441	rVV	559833	704072	4.80%	0.580%
69	17.175	2442	2446	2452	rVB4	75569	132484	0.90%	0.109%
70	17.381	2471	2481	2487	rBV	798393	1207137	8.23%	0.994%
71	17.504	2493	2502	2506	rVV5	77346	161165	1.10%	0.133%

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72	17.669	2521	2530	2533	rVV9	34888	86345	0.59%	0.071%
73	17.704	2533	2536	2543	rVB4	32123	61508	0.42%	0.051%
74	17.857	2557	2562	2565	rVV	253302	353389	2.41%	0.291%
75	17.904	2565	2570	2576	rVV	396967	578815	3.95%	0.477%
76	17.981	2578	2583	2587	rVV2	51331	91631	0.63%	0.075%
77	18.051	2588	2595	2599	rVV	822642	1300042	8.87%	1.070%
78	18.086	2599	2601	2605	rVV	114584	129821	0.89%	0.107%
79	18.157	2607	2613	2616	rVV4	35738	67982	0.46%	0.056%
80	18.198	2616	2620	2624	rVV2	65726	97318	0.66%	0.080%
81	18.292	2632	2636	2640	rVV	57841	88089	0.60%	0.073%
82	18.363	2643	2648	2651	rVV	156389	231822	1.58%	0.191%
83	18.392	2651	2653	2659	rVB3	58776	75146	0.51%	0.062%
84	19.045	2758	2764	2770	rVV	2337346	3187019	21.74%	2.624%
85	19.286	2799	2805	2811	rVV4	73547	161183	1.10%	0.133%
86	19.381	2815	2821	2824	rVV	2860794	4415354	30.12%	3.635%
87	19.404	2824	2825	2835	rVV	1453593	1449434	9.89%	1.193%
88	19.480	2836	2838	2843	rVB5	42467	58248	0.40%	0.048%
89	19.786	2887	2890	2895	rVB	91638	117203	0.80%	0.096%
90	19.845	2895	2900	2903	rBV2	71713	127892	0.87%	0.105%
91	19.910	2908	2911	2915	rVB3	70499	86511	0.59%	0.071%
92	20.010	2925	2928	2933	rVB4	93312	119475	0.81%	0.098%
93	21.180	3122	3127	3138	rVB	6712528	8101848	55.27%	6.670%
94	22.251	3306	3309	3314	rVB4	241533	291269	1.99%	0.240%
95	22.745	3390	3393	3402	rVB2	344263	451023	3.08%	0.371%
96	23.216	3467	3473	3480	rBV	2382646	4026541	27.47%	3.315%
97	23.292	3482	3486	3490	rVV	520620	795662	5.43%	0.655%
98	23.351	3490	3496	3505	rVB	4848546	8755404	59.72%	7.208%
99	23.910	3588	3591	3597	rVB2	430125	711471	4.85%	0.586%
100	24.627	3709	3713	3720	rVB	421980	777586	5.30%	0.640%

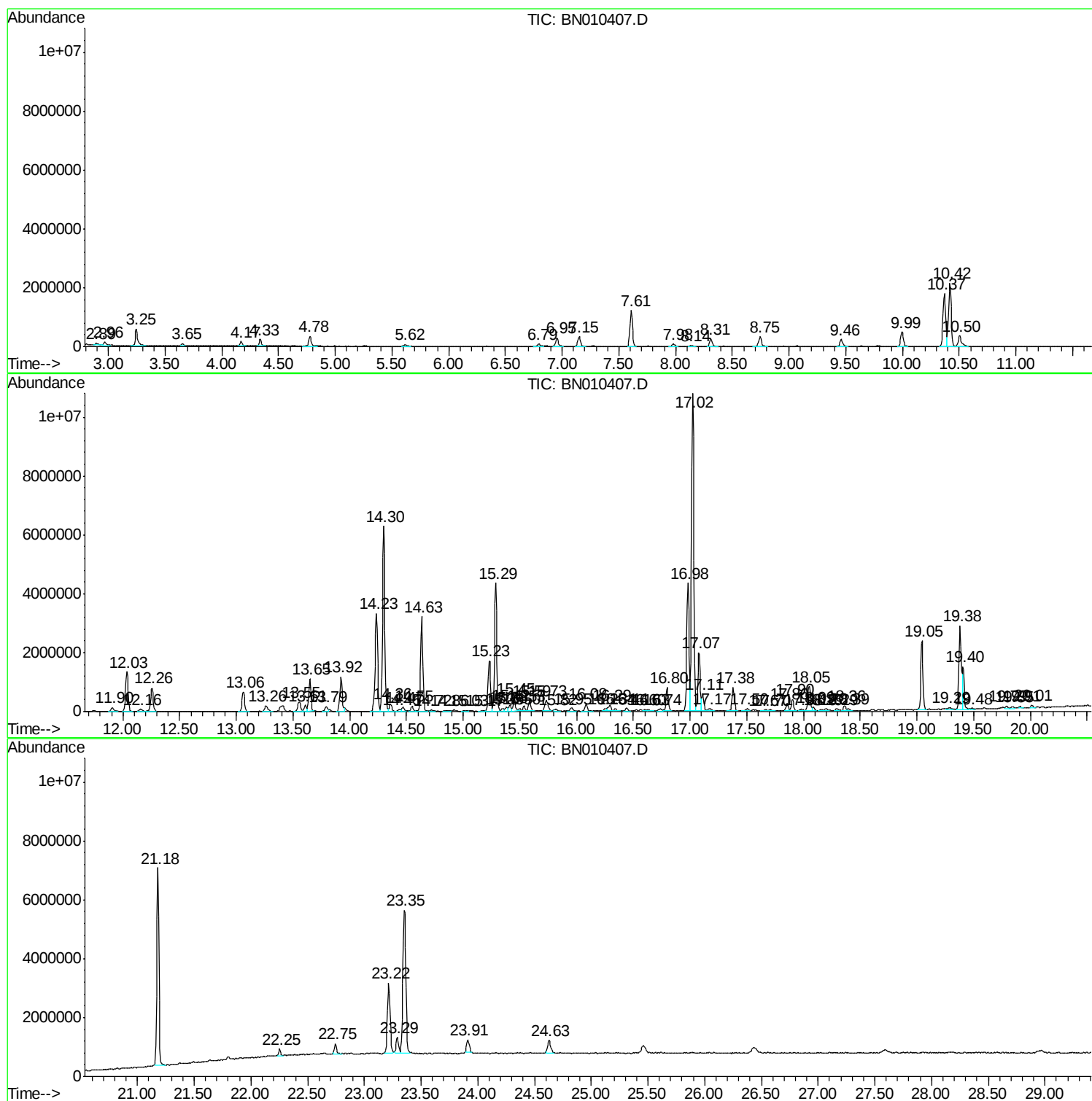
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Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.M
Quant Title : SVOA CALIBRATION

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



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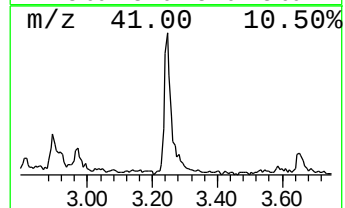
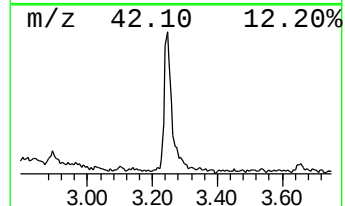
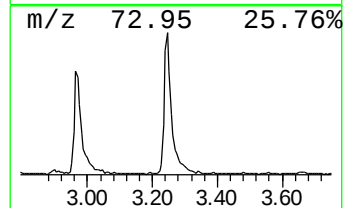
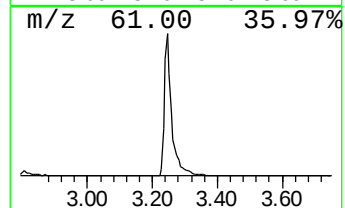
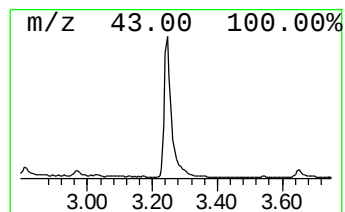
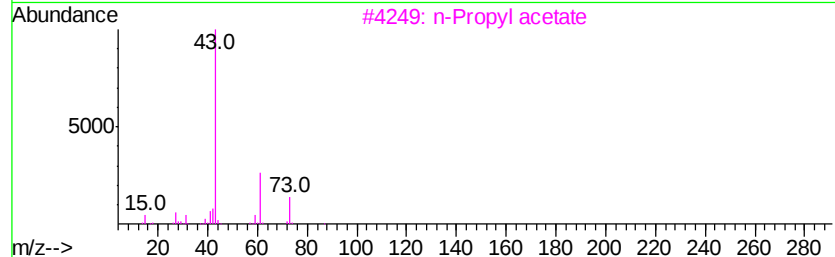
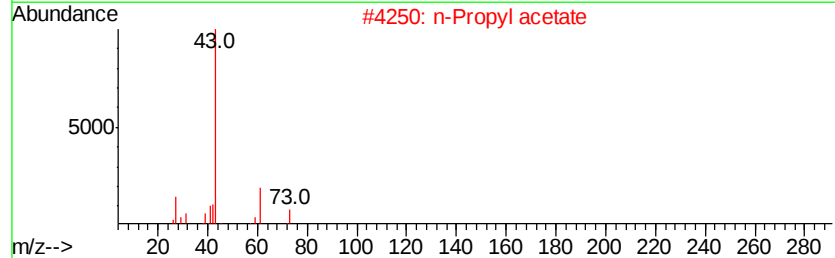
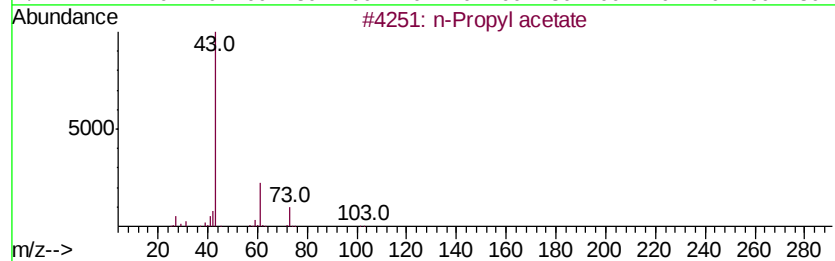
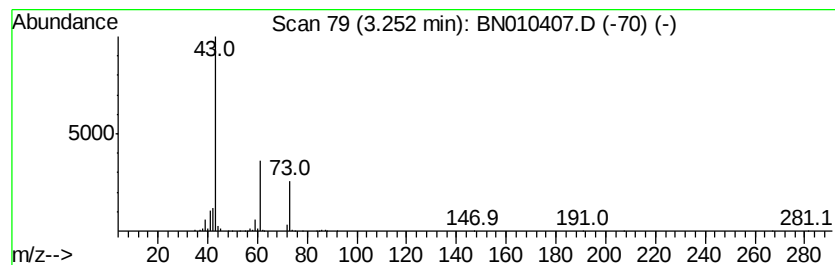
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 n-Propyl acetate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.25	9.33 ng/ul	911120	1,4-Dichlorobenzene-d4	7.61

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Propyl acetate	102	C5H10O2	000109-60-4	78
2			n-Propyl acetate	102	C5H10O2	000109-60-4	59
3			n-Propyl acetate	102	C5H10O2	000109-60-4	38
4			Isopropyl acetate	102	C5H10O2	000108-21-4	9
5			Isopropyl acetate	102	C5H10O2	000108-21-4	9



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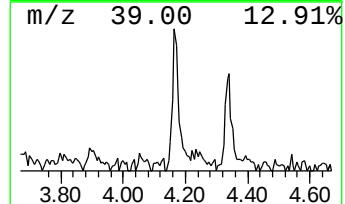
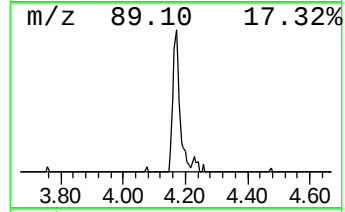
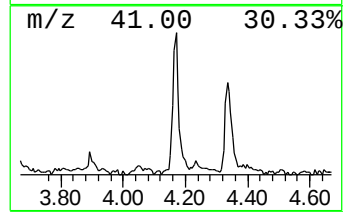
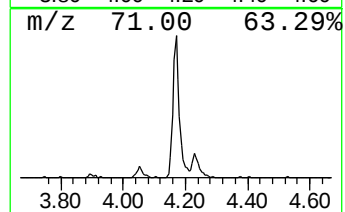
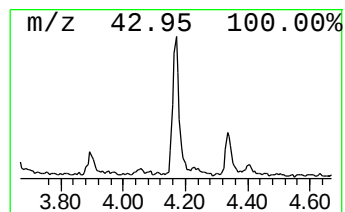
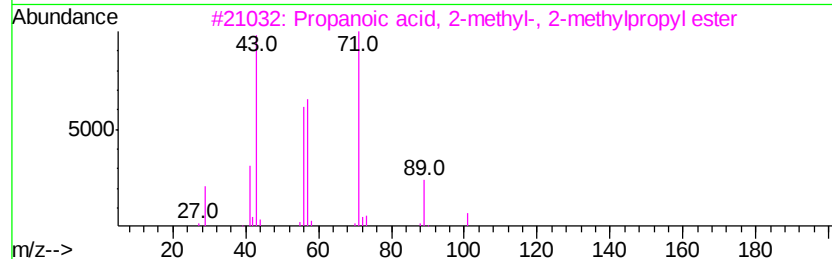
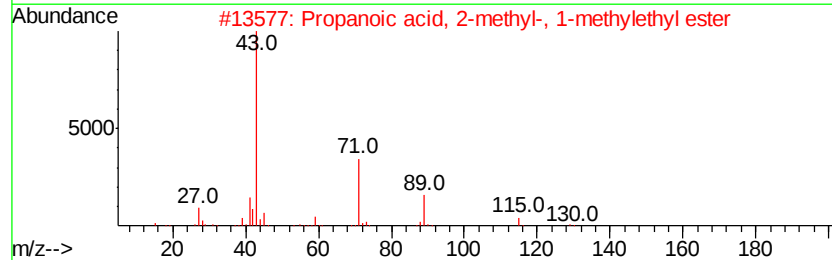
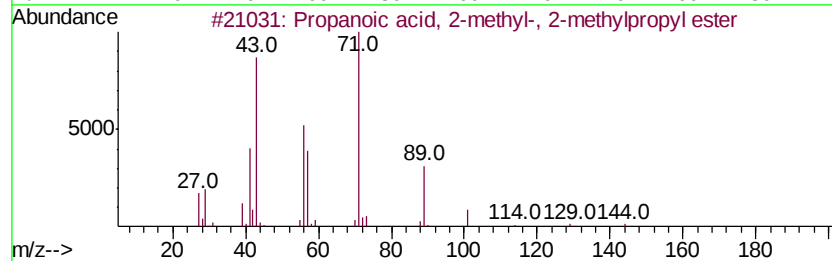
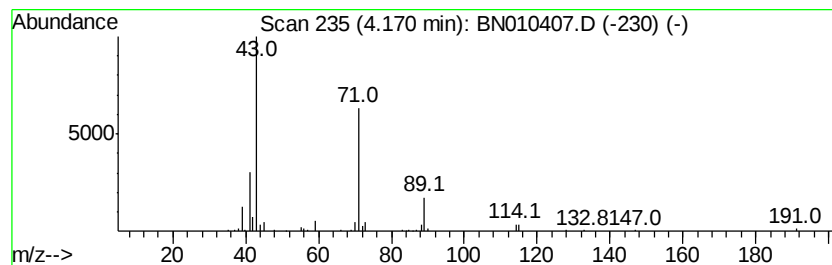
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 Propanoic acid, 2-methyl-, ... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.17	2.48 ng/ul	242350	1,4-Dichlorobenzene-d4	7.61

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propanoic acid, 2-methyl-, 2-met...	144	C8H16O2	000097-85-8	64
2			Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	56
3			Propanoic acid, 2-methyl-, 2-met...	144	C8H16O2	000097-85-8	53
4			Propanoic acid, 2-methyl-, 2-pro...	128	C7H12O2	015727-77-2	47
5			Butanoic acid, 1-methylethyl ester	130	C7H14O2	000638-11-9	45



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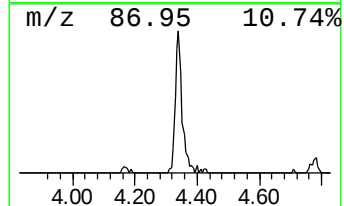
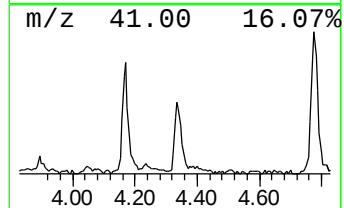
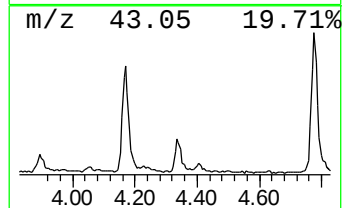
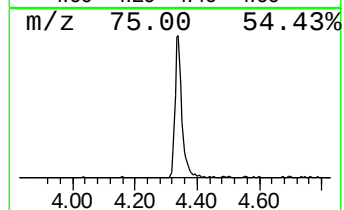
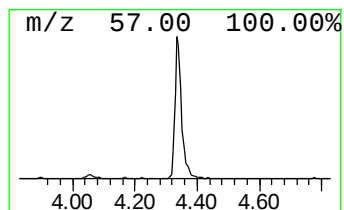
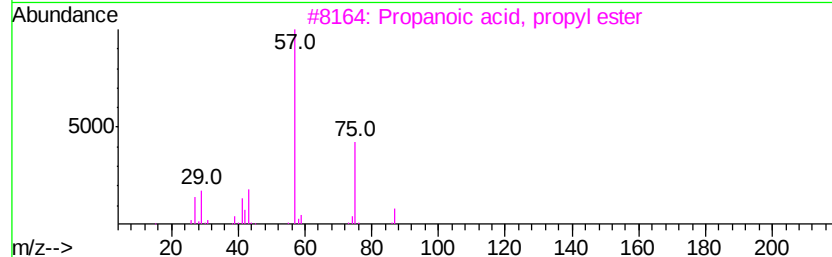
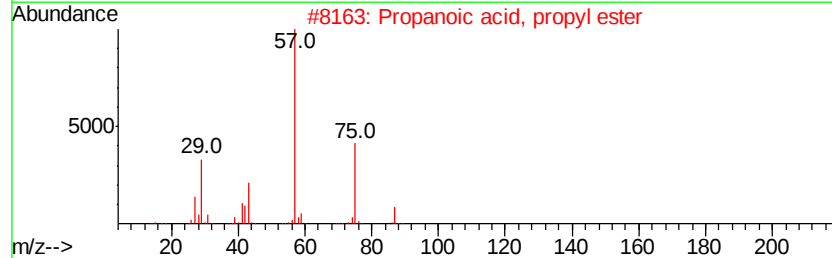
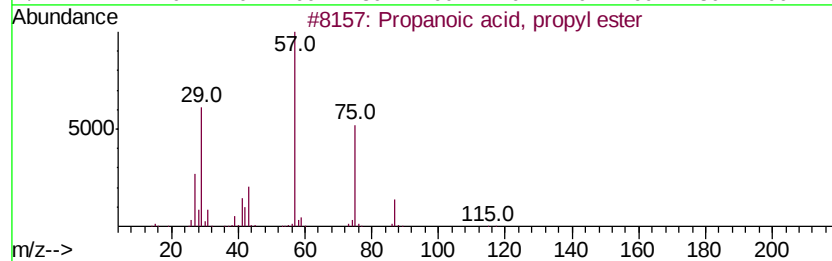
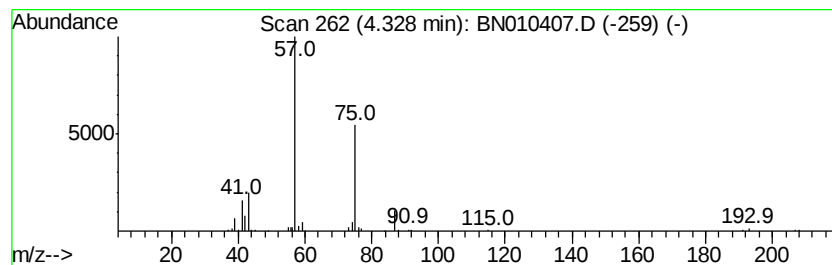
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Peak Number 3 Propanoic acid, propyl ester Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.33	3.52 ng/ul	343685	1,4-Dichlorobenzene-d4	7.61

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	78
3		Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	53
4		Propanoic acid, 1-methylethyl ester	116	C6H12O2	000637-78-5	32
5		Isobutyl ether	130	C8H18O	000628-55-7	12



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Data File : BN010407.D
Acq On : 03 Apr 2020 11:52
Operator : CG/JU
Sample : L2095-01DL 4X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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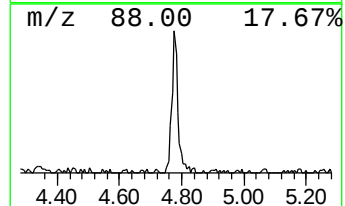
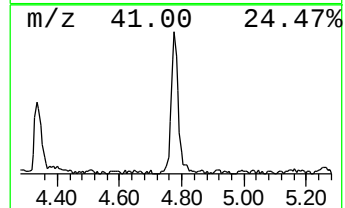
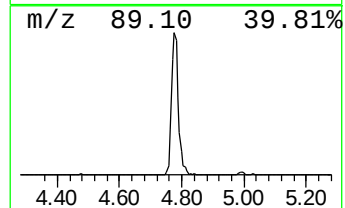
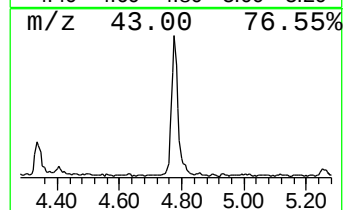
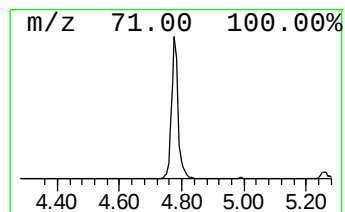
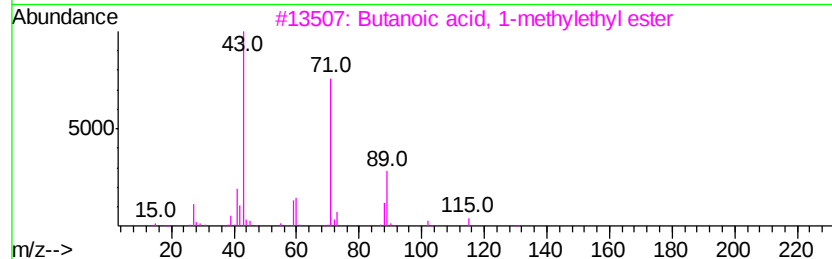
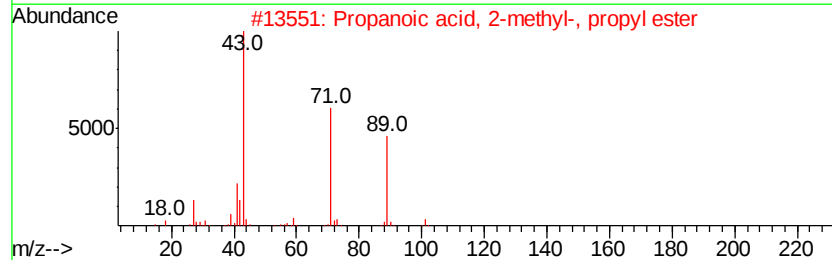
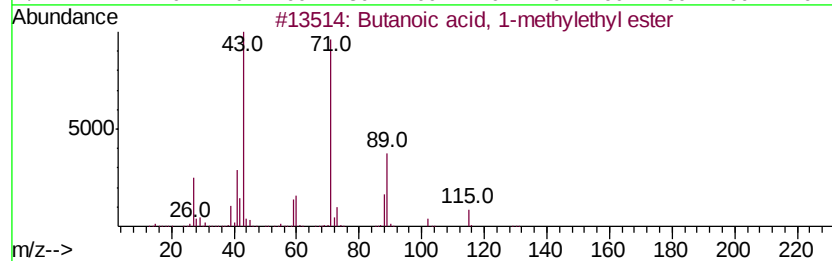
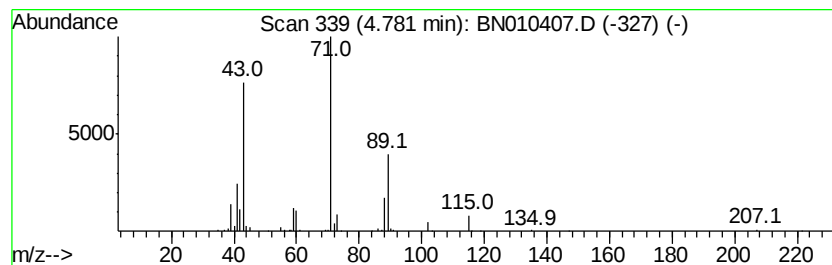
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Butanoic acid, 1-methylethy... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.78	5.26 ng/ul	513401	1,4-Dichlorobenzene-d4	7.61

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butanoic acid, 1-methylethyl ester	130	C7H14O2	000638-11-9	83
2		Propanoic acid, 2-methyl-, propyl ester	130	C7H14O2	000644-49-5	72
3		Butanoic acid, 1-methylethyl ester	130	C7H14O2	000638-11-9	68
4		Butanoic acid, 1-methylpropyl ester	144	C8H16O2	000819-97-6	56
5		Butanoic acid, butyl ester	144	C8H16O2	000109-21-7	45



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Data File : BN010407.D
Acq On : 03 Apr 2020 11:52
Operator : CG/JU
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Misc :
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ClientSampleId :
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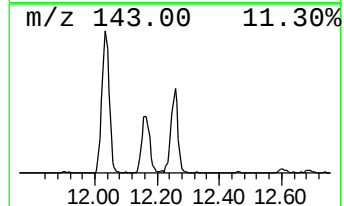
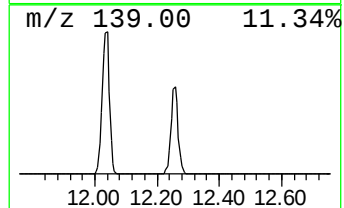
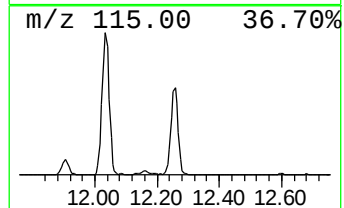
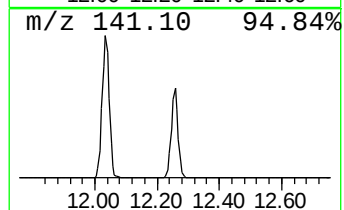
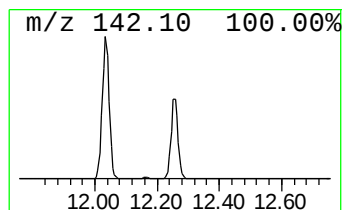
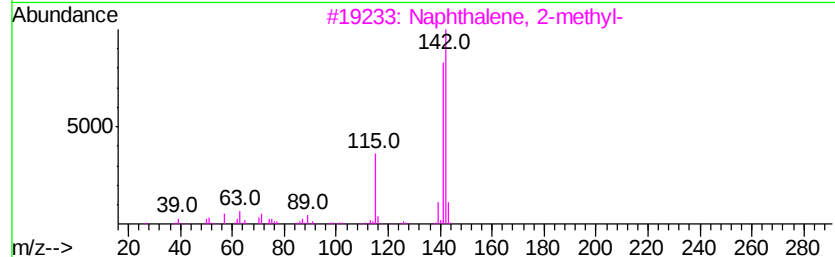
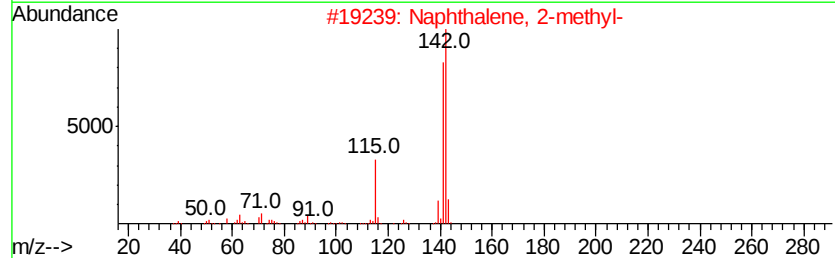
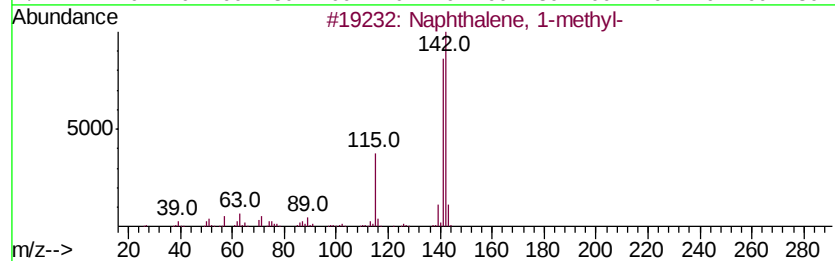
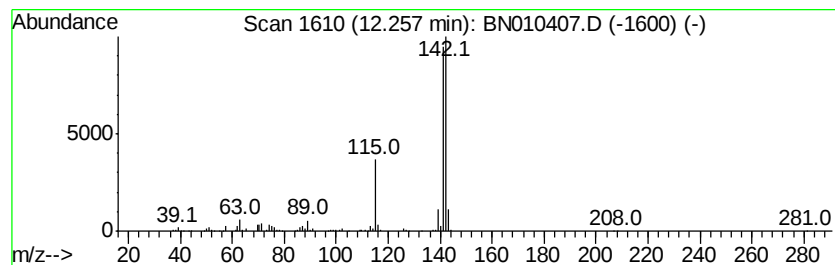
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Naphthalene, 1-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.26	8.39 ng/ul	1276070	Naphthalene-d8	10.37

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		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
5		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN040320\
Data File : BN010407.D
Acq On : 03 Apr 2020 11:52
Operator : CG/JU
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Misc :
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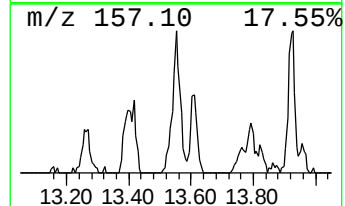
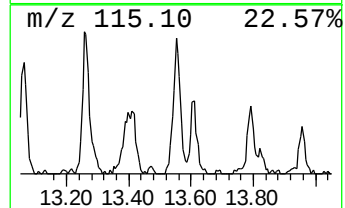
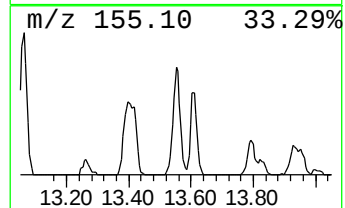
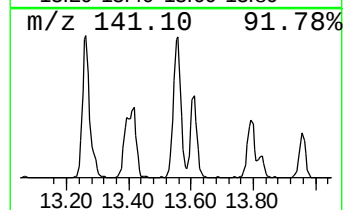
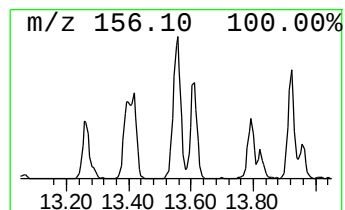
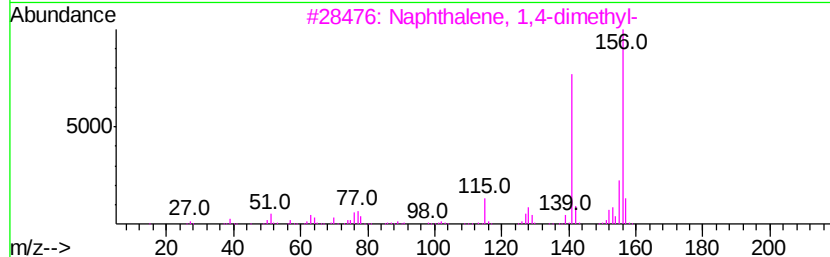
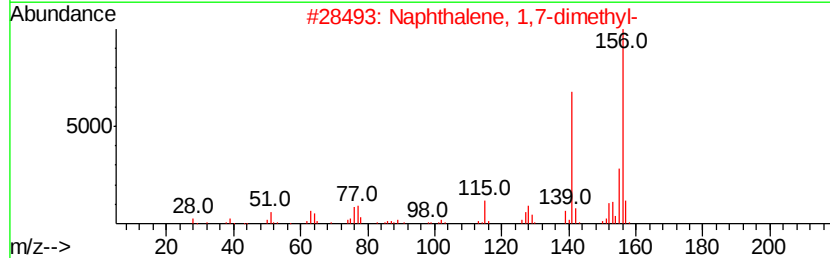
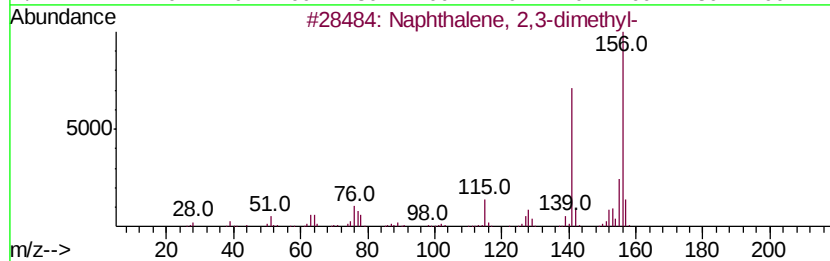
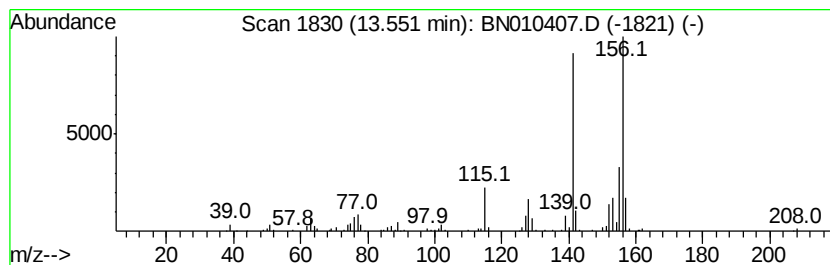
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Naphthalene, 2,3-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.55	2.31 ng/ul	572925	Acenaphthene-d10	14.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
2		Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96
3		Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	96
4		Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	95
5		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	95



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN040320\
Data File : BN010407.D
Acq On : 03 Apr 2020 11:52
Operator : CG/JU
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Misc :
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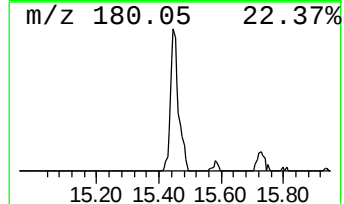
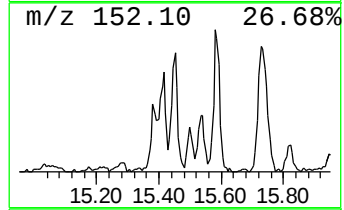
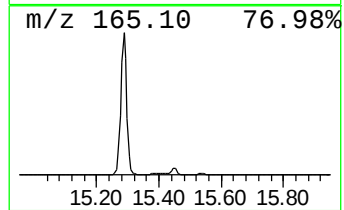
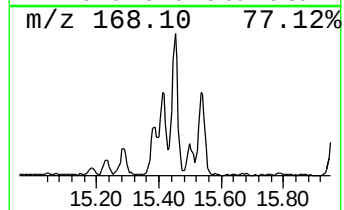
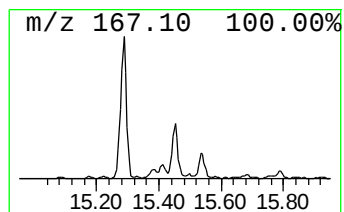
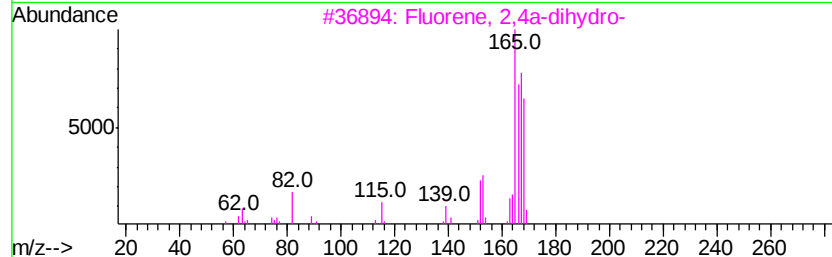
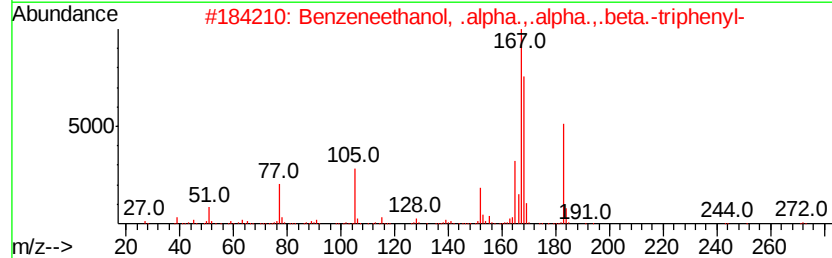
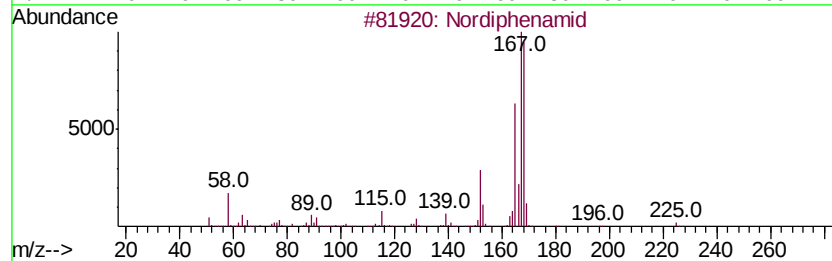
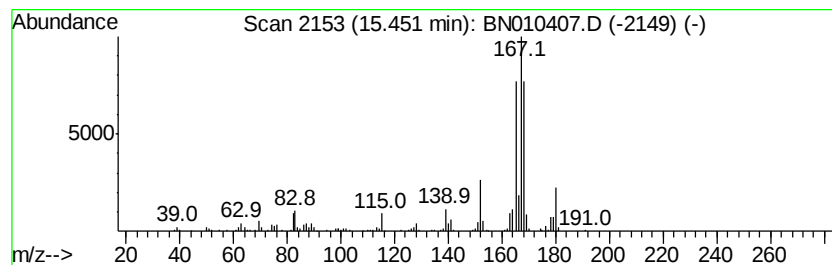
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Nordiphenamid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.45	2.56 ng/ul	634040	Acenaphthene-d10	14.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nordiphenamid	225	C15H15NO	000954-21-2	59
2		Benzeneethanol, .alpha.,.alpha.,...	350	C26H22O	000981-24-8	50
3		Fluorene, 2,4a-dihydro-	168	C13H12	059247-36-8	49
4		Diphenylmethane	168	C13H12	000101-81-5	49
5		Fluorene, 1,4-dihydro-	168	C13H12	041593-21-9	46



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Data File : BN010407.D
Acq On : 03 Apr 2020 11:52
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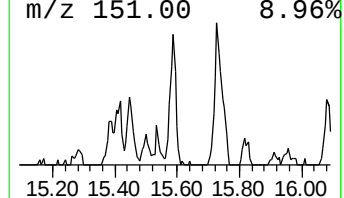
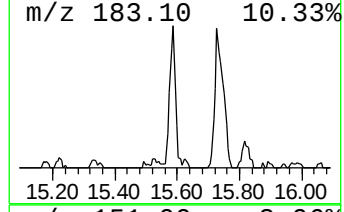
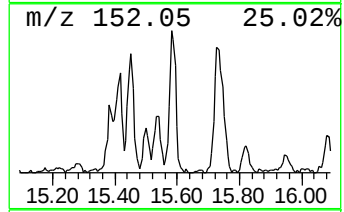
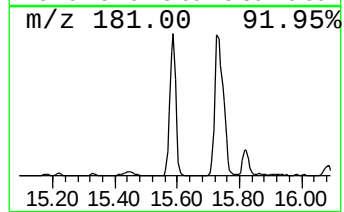
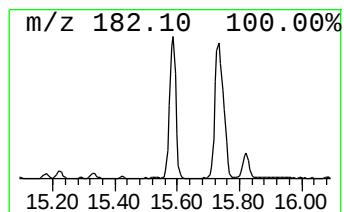
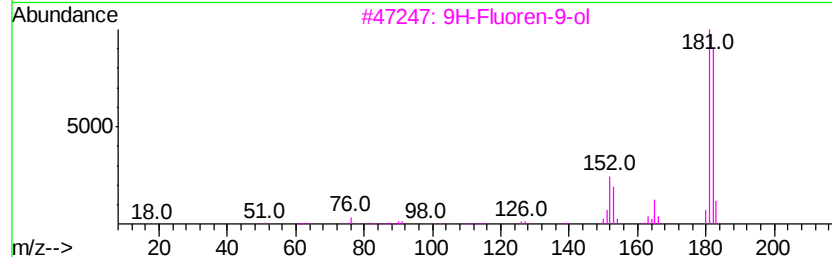
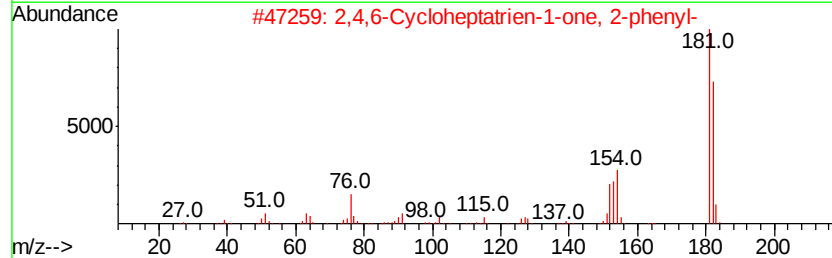
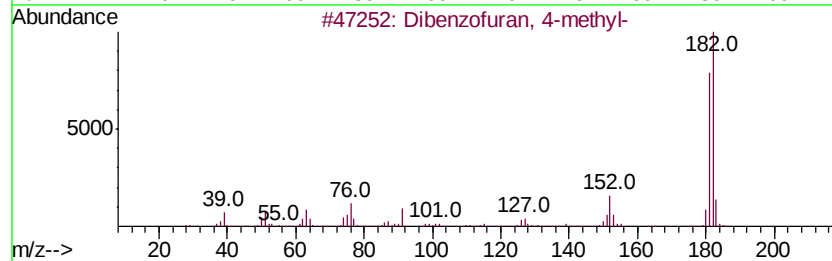
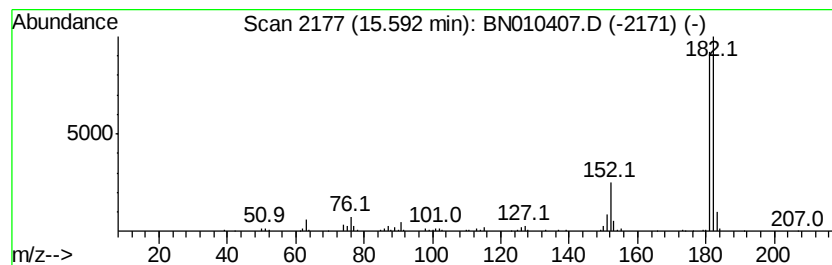
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Dibenzofuran, 4-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.59	2.01 ng/ul	498587	Acenaphthene-d10	14.23

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Dibenzofuran, 4-methyl-	182	C13H10O	007320-53-8	90
2		2,4,6-Cycloheptatrien-1-one, 2-p...	182	C13H10O	014562-09-5	86
3		9H-Fluoren-9-ol	182	C13H10O	001689-64-1	83
4		[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	78
5		Pyrido[2,3-b]indole, 6-methyl-	182	C12H10N2	108349-67-3	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN040320\
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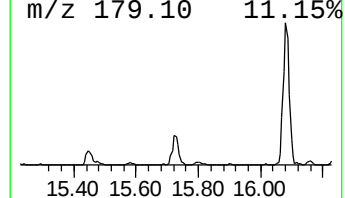
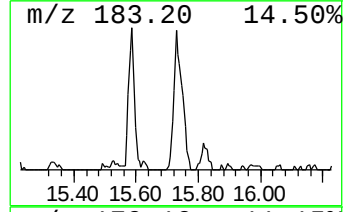
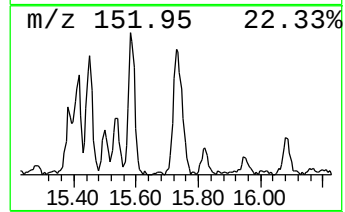
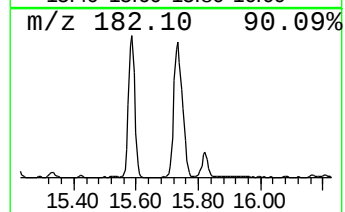
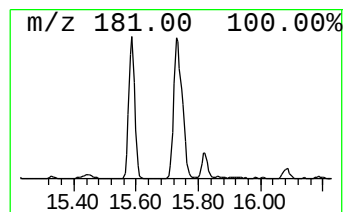
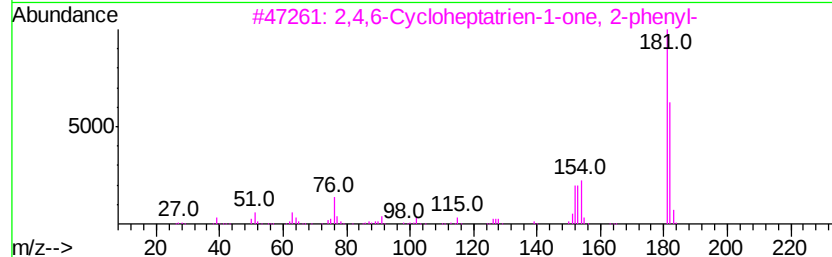
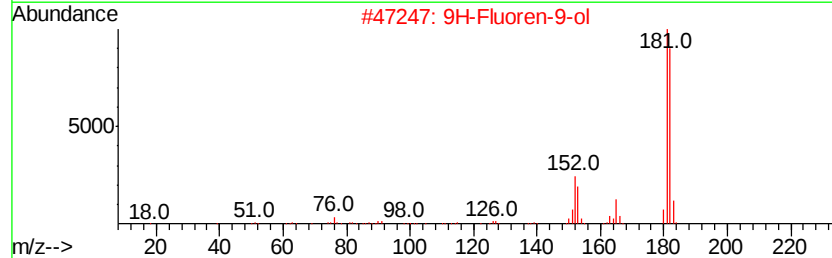
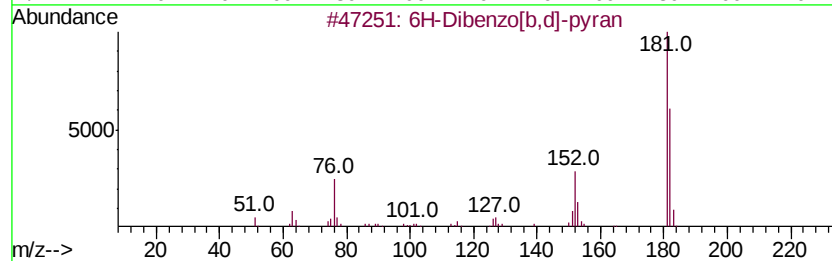
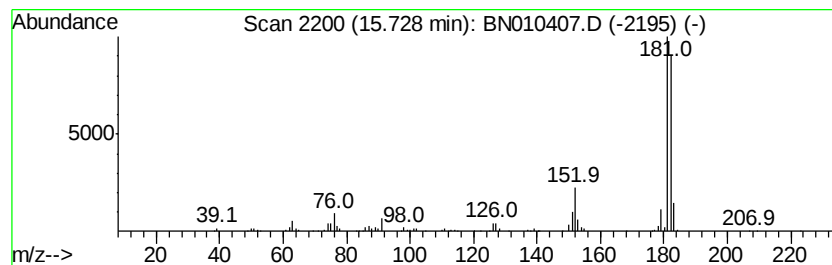
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 6H-Dibenzo[b,d]-pyran Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.		
15.73	2.33 ng/ul	714799	Phenanthrene-d10	16.98		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	6H-Dibenzo[b,d]-pvrn		182	C13H10O	000229-95-8	90
2	9H-Fluoren-9-ol		182	C13H10O	001689-64-1	86
3	2,4,6-Cvcloheptatrien-1-one, 2-p...		182	C13H10O	014562-09-5	83
4	2,4,6-Cvcloheptatrien-1-one, 2-p...		182	C13H10O	014562-09-5	80
5	3-(2-Naphthyl)acrylaldehyde		182	C13H10O	020884-05-3	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN040320\
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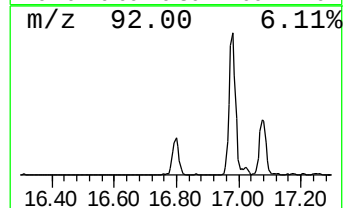
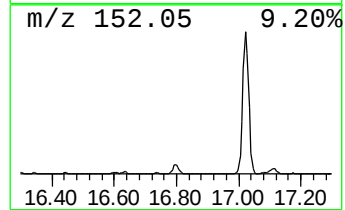
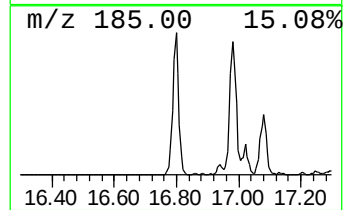
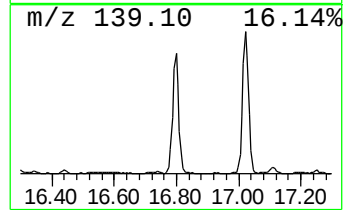
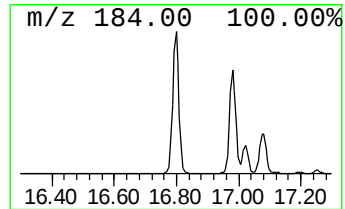
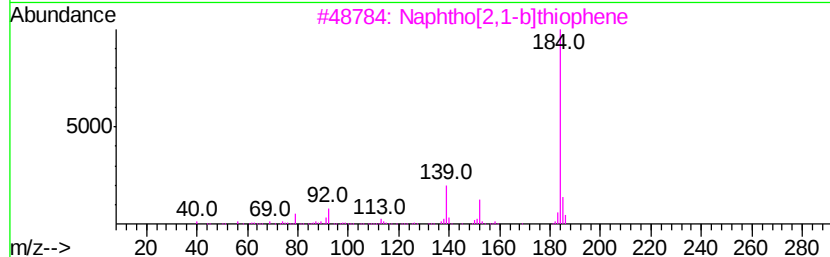
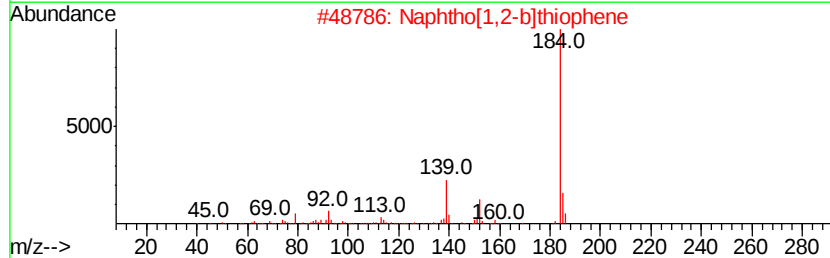
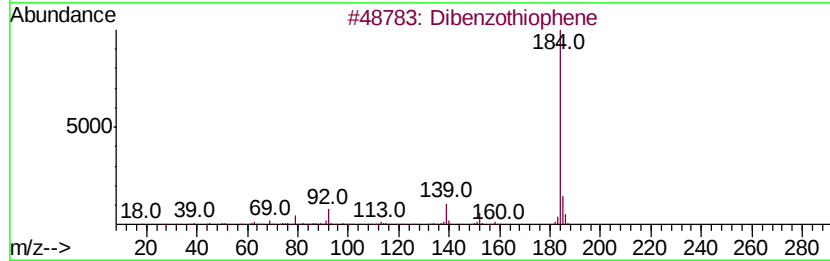
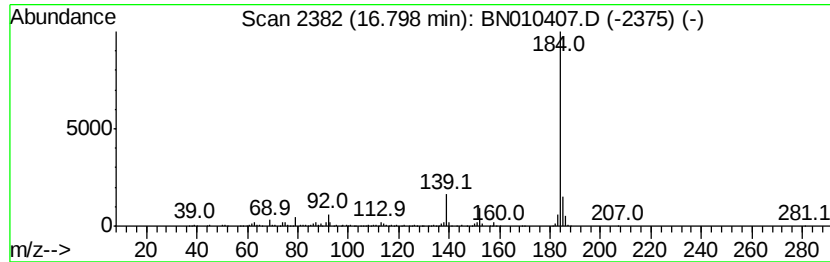
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Dibenzothiophene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.80	3.64 ng/ul	1113290	Phenanthrene-d10	16.98

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Dibenzothiophene	184	C12H8S	000132-65-0	96
2		Naphtho[1,2-b]thiophene	184	C12H8S	000234-41-3	95
3		Naphtho[2,1-b]thiophene	184	C12H8S	000233-02-3	95
4		Naphtho[2,3-b]thiophene	184	C12H8S	000268-77-9	94
5		Dibenzothiophene	184	C12H8S	000132-65-0	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN040320\
Data File : BN010407.D
Acq On : 03 Apr 2020 11:52
Operator : CG/JU
Sample : L2095-01DL 4X
Misc :
ALS Vial : 7 Sample Multiplier: 1

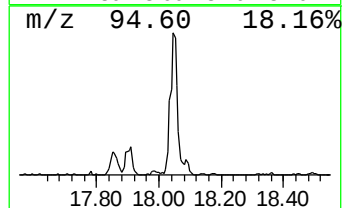
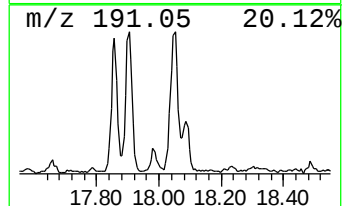
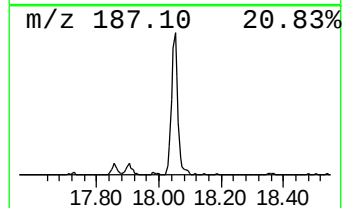
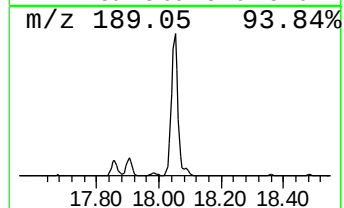
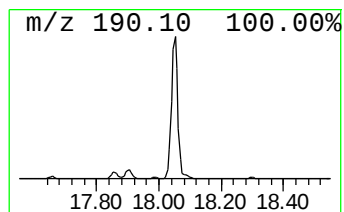
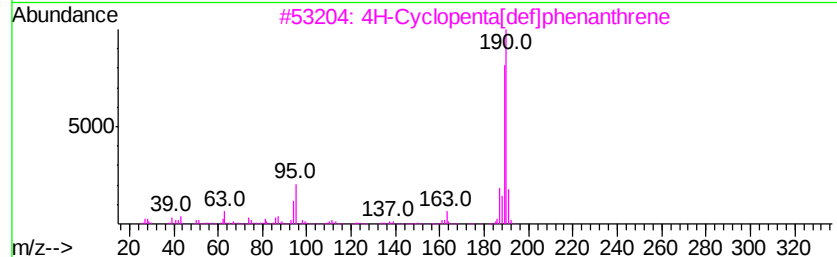
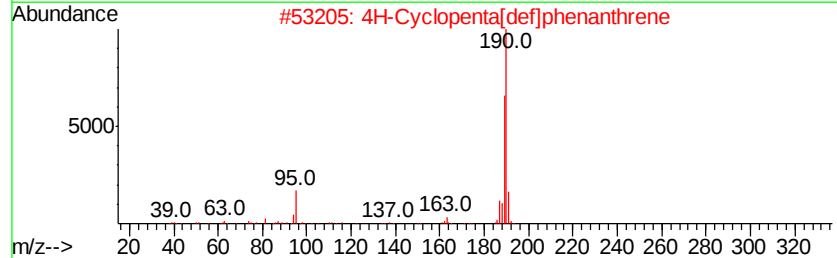
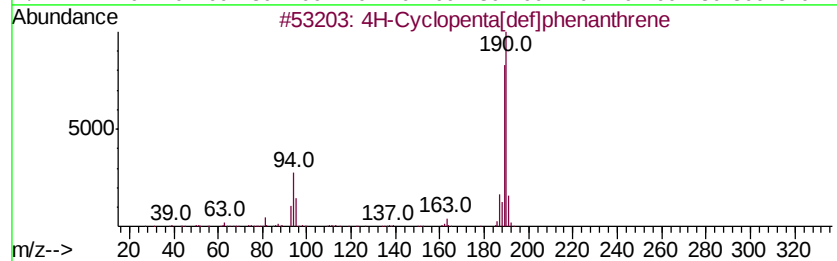
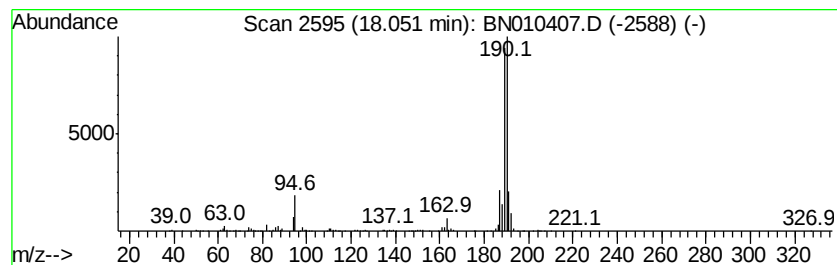
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 11 4H-Cyclopenta[def]phenanthrene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD		R.T.	
18.05	4.25 ng/ul	1300040	Phenanthrene-d10		16.98	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	97
2		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	87
3		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	74
4		Phenol, 4-(2-thienylmethyl)-	190	C11H10OS	091680-55-6	64
5		6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	56



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN040320\
Data File : BN010407.D
Acq On : 03 Apr 2020 11:52
Operator : CG/JU
Sample : L2095-01DL 4X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBF16DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.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
n-Propyl acetate	3.25	9.3	ng/ul	911120	1	7.61	1953260	20.0
Propanoic acid, 2...	4.17	2.5	ng/ul	242350	1	7.61	1953260	20.0
Propanoic acid, p...	4.33	3.5	ng/ul	343685	1	7.61	1953260	20.0
Butanoic acid, 1-...	4.78	5.3	ng/ul	513401	1	7.61	1953260	20.0
Naphthalene, 1-me...	12.26	8.4	ng/ul	1276070	2	10.37	3043090	20.0
Naphthalene, 2,3-...	13.55	2.3	ng/ul	572925	3	14.23	4952630	20.0
Nordiphenamid	15.45	2.6	ng/ul	634040	3	14.23	4952630	20.0
Dibenzofuran, 4-m...	15.59	2.0	ng/ul	498587	3	14.23	4952630	20.0
6H-Dibenzo[b,d]-p...	15.73	2.3	ng/ul	714799	4	16.98	6123190	20.0
Dibenzothiophene	16.80	3.6	ng/ul	1113290	4	16.98	6123190	20.0
4H-Cyclopenta[def...	18.05	4.3	ng/ul	1300040	4	16.98	6123190	20.0