

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060722\
 Data File : BN020042.D
 Acq On : 07 Jun 2022 23:31
 Operator : CG/JU
 Sample : N3171-09
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBTZ4

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN060622.M
 Title : SVOA CALIBRATION

Signal : TIC: BN020042.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.281	29	33	47	rVB	26310	50615	1.11%	0.096%
2	3.693	99	103	119	rBV	101332	180040	3.93%	0.340%
3	5.346	379	384	391	rVB	43790	67820	1.48%	0.128%
4	6.987	657	663	672	rBV	111383	183111	4.00%	0.346%
5	7.152	683	691	699	rBV	466094	765747	16.72%	1.447%
6	7.222	699	703	714	rVV2	22636	38196	0.83%	0.072%
7	7.352	718	725	735	rBV	424269	707434	15.45%	1.337%
8	7.822	797	805	814	rBV	461291	753582	16.46%	1.424%
9	8.199	861	869	877	rBV	649830	1086297	23.72%	2.053%
10	8.358	891	896	906	rVB	21812	36613	0.80%	0.069%
11	8.522	918	924	936	rBV	361937	600047	13.10%	1.134%
12	8.975	988	1001	1013	rBV	599932	1098286	23.98%	2.076%
13	9.175	1030	1035	1041	rVV2	20059	32628	0.71%	0.062%
14	9.299	1049	1056	1062	rVV	34239	77973	1.70%	0.147%
15	9.693	1116	1123	1135	rBV	455905	791720	17.29%	1.496%
16	9.869	1147	1153	1160	rVB	50295	85635	1.87%	0.162%
17	10.016	1168	1178	1188	rBV2	91348	178126	3.89%	0.337%
18	10.234	1207	1215	1225	rBV	676483	1199578	26.20%	2.267%
19	10.610	1272	1279	1283	rVV	705451	1205822	26.33%	2.279%
20	10.663	1283	1288	1295	rVV	788272	1388344	30.32%	2.624%
21	10.746	1295	1302	1319	rVB2	373949	795466	17.37%	1.504%
22	11.999	1511	1515	1526	rVB9	17485	44583	0.97%	0.084%
23	12.169	1538	1544	1548	rBV	50779	92193	2.01%	0.174%
24	12.275	1555	1562	1569	rBV	293420	487387	10.64%	0.921%
25	12.387	1575	1581	1588	rVV	60814	113896	2.49%	0.215%
26	12.493	1588	1599	1609	rVB	640008	1087126	23.74%	2.055%
27	12.704	1629	1635	1642	rVB2	20750	33093	0.72%	0.063%
28	12.869	1658	1663	1667	rVV2	26623	42029	0.92%	0.079%
29	13.287	1730	1734	1742	rVV3	37129	69419	1.52%	0.131%
30	13.416	1749	1756	1761	rVV4	26237	50662	1.11%	0.096%
31	13.487	1761	1768	1778	rVB5	41972	129013	2.82%	0.244%
32	13.628	1778	1792	1800	rBV4	53879	177234	3.87%	0.335%
33	13.775	1811	1817	1822	rVV	115222	214741	4.69%	0.406%
34	13.857	1822	1831	1841	rVV	1553368	2324184	50.75%	4.393%
35	14.010	1850	1857	1860	rVV	50646	82316	1.80%	0.156%
36	14.040	1860	1862	1868	rVB2	24446	29520	0.64%	0.056%

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Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN060622.M
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37	14.140	1870	1879	1890	rBV	1610544	2596996	56.71%	4.909%
38	14.451	1920	1932	1937	rBV	1131475	1704102	37.21%	3.221%
39	14.516	1937	1943	1949	rVV	1317782	2026232	44.25%	3.830%
40	14.575	1949	1953	1959	rVV	21885	34564	0.75%	0.065%
41	14.640	1959	1964	1972	rVB2	137098	215629	4.71%	0.408%
42	14.722	1972	1978	1982	rBV	76630	118101	2.58%	0.223%
43	14.845	1993	1999	2006	rVV	226732	359314	7.85%	0.679%
44	14.922	2006	2012	2020	rVB6	19313	38726	0.85%	0.073%
45	15.075	2027	2038	2042	rBV8	17581	37123	0.81%	0.070%
46	15.151	2042	2051	2056	rBV5	17457	36869	0.81%	0.070%
47	15.304	2072	2077	2084	rVB3	14878	32658	0.71%	0.062%
48	15.440	2090	2100	2105	rBV	2208613	3272837	71.47%	6.186%
49	15.492	2105	2109	2115	rVB	509036	762820	16.66%	1.442%
50	15.557	2115	2120	2127	rBV	842385	1164101	25.42%	2.200%
51	15.622	2127	2131	2134	rBV3	44234	60824	1.33%	0.115%
52	15.787	2156	2159	2168	rVB4	36525	66131	1.44%	0.125%
53	15.875	2168	2174	2177	rBV4	15494	30913	0.68%	0.058%
54	15.934	2177	2184	2191	rVB4	44273	103670	2.26%	0.196%
55	15.998	2191	2195	2204	rVB3	15729	29373	0.64%	0.056%
56	16.163	2204	2223	2230	rBV	160472	331010	7.23%	0.626%
57	16.222	2230	2233	2238	rVV6	19647	31517	0.69%	0.060%
58	16.287	2238	2244	2249	rVV2	24693	43159	0.94%	0.082%
59	16.345	2249	2254	2257	rVV2	26489	44177	0.96%	0.084%
60	16.381	2257	2260	2267	rVV6	18734	35564	0.78%	0.067%
61	16.451	2267	2272	2276	rVV	26167	42508	0.93%	0.080%
62	16.551	2284	2289	2295	rVB8	16554	33397	0.73%	0.063%
63	16.651	2302	2306	2313	rBV5	49107	96149	2.10%	0.182%
64	16.975	2348	2361	2364	rBV2	71522	151892	3.32%	0.287%
65	17.010	2364	2367	2372	rVV	41681	72680	1.59%	0.137%
66	17.104	2378	2383	2387	rVV5	23454	59673	1.30%	0.113%
67	17.145	2387	2390	2392	rVV	34861	46938	1.02%	0.089%
68	17.187	2392	2397	2402	rVV	1338925	2071320	45.23%	3.915%
69	17.234	2402	2405	2409	rVV	293114	449005	9.80%	0.849%
70	17.287	2409	2414	2425	rVV2	2437459	3866111	84.42%	7.307%
71	17.387	2425	2431	2437	rVB9	21087	49402	1.08%	0.093%
72	17.586	2454	2465	2471	rBV	485617	786443	17.17%	1.486%
73	17.686	2478	2482	2487	rBV4	24684	47300	1.03%	0.089%
74	17.834	2501	2507	2508	rBV6	19259	32154	0.70%	0.061%
75	17.875	2511	2514	2521	rVB2	33137	57981	1.27%	0.110%
76	17.939	2521	2525	2529	rBV4	20121	30440	0.66%	0.058%

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77	18.028	2536	2540	2546	rVB	43791	64296	1.40%	0.122%
78	18.098	2546	2552	2563	rVV3	144229	334222	7.30%	0.632%
79	18.186	2563	2567	2573	rVV	65223	99379	2.17%	0.188%
80	18.263	2573	2580	2585	rVB4	27267	61414	1.34%	0.116%
81	18.369	2591	2598	2601	rBV4	28829	67826	1.48%	0.128%
82	18.410	2601	2605	2609	rVB	43971	66906	1.46%	0.126%
83	18.504	2616	2621	2627	rBV2	38585	67652	1.48%	0.128%
84	18.563	2627	2631	2635	rVB4	20240	30376	0.66%	0.057%
85	18.957	2691	2698	2701	rBV8	15329	32768	0.72%	0.062%
86	19.145	2726	2730	2738	rBV2	39275	80549	1.76%	0.152%
87	19.216	2738	2742	2745	rVV2	36956	59187	1.29%	0.112%
88	19.245	2745	2747	2752	rVB	31252	40913	0.89%	0.077%
89	19.375	2765	2769	2778	rBV2	70573	108093	2.36%	0.204%
90	19.581	2798	2804	2821	rVB	2961231	4446475	97.10%	8.404%
91	19.980	2866	2872	2878	rBV3	36181	88797	1.94%	0.168%
92	20.045	2878	2883	2893	rVV	150014	222151	4.85%	0.420%
93	20.145	2896	2900	2907	rVB	64828	86841	1.90%	0.164%
94	20.692	2988	2993	3005	rVB2	82145	139652	3.05%	0.264%
95	21.104	3058	3063	3069	rBV	95636	170332	3.72%	0.322%
96	21.375	3103	3109	3121	rBV2	1753187	2266798	49.50%	4.285%
97	23.010	3383	3387	3394	rVB	46202	66505	1.45%	0.126%
98	23.557	3471	3480	3488	rBV2	2177505	4579409	100.00%	8.656%
99	23.704	3497	3505	3515	rVB2	1135874	2236652	48.84%	4.228%
100	24.304	3602	3607	3614	rVB2	58466	118880	2.60%	0.225%

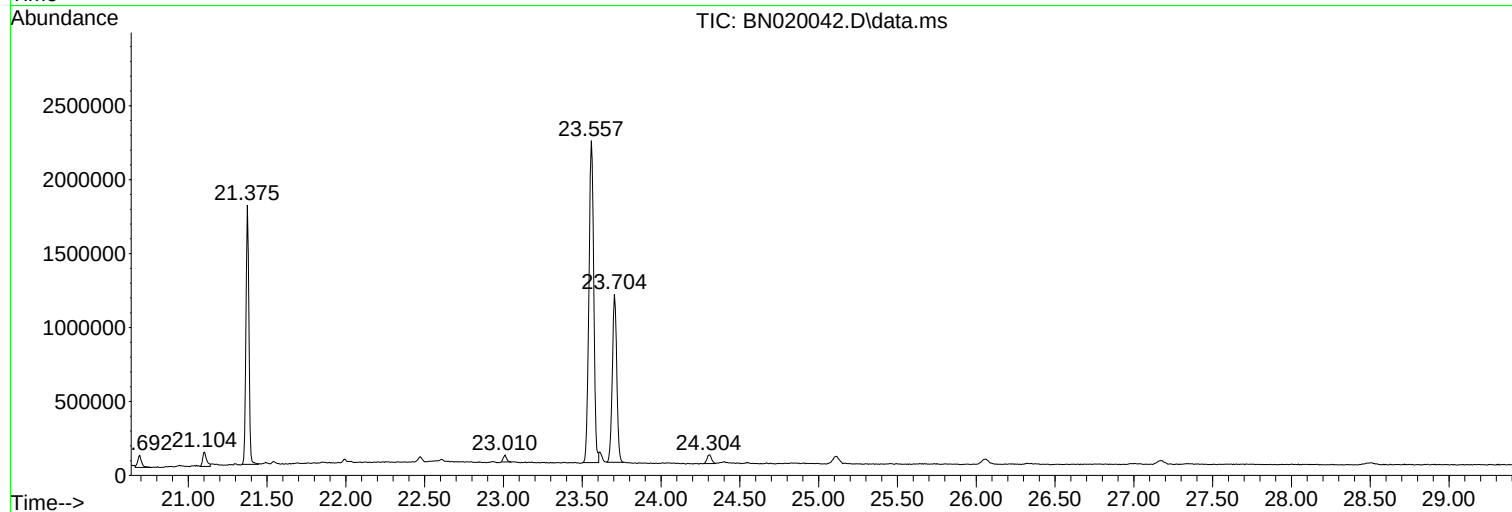
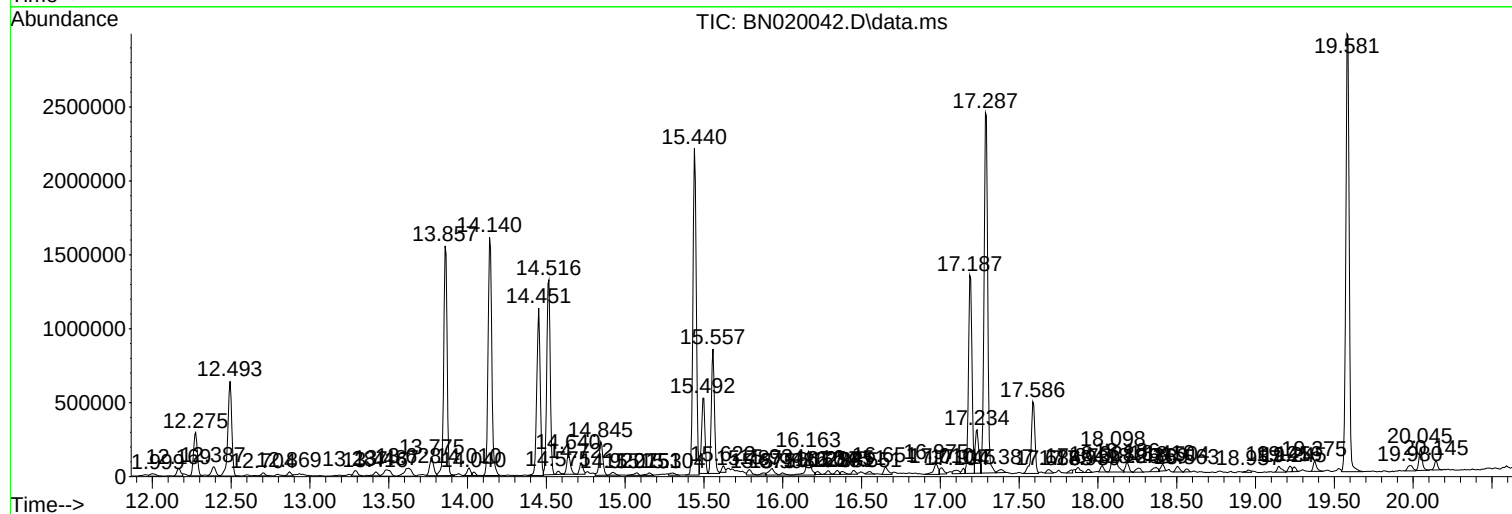
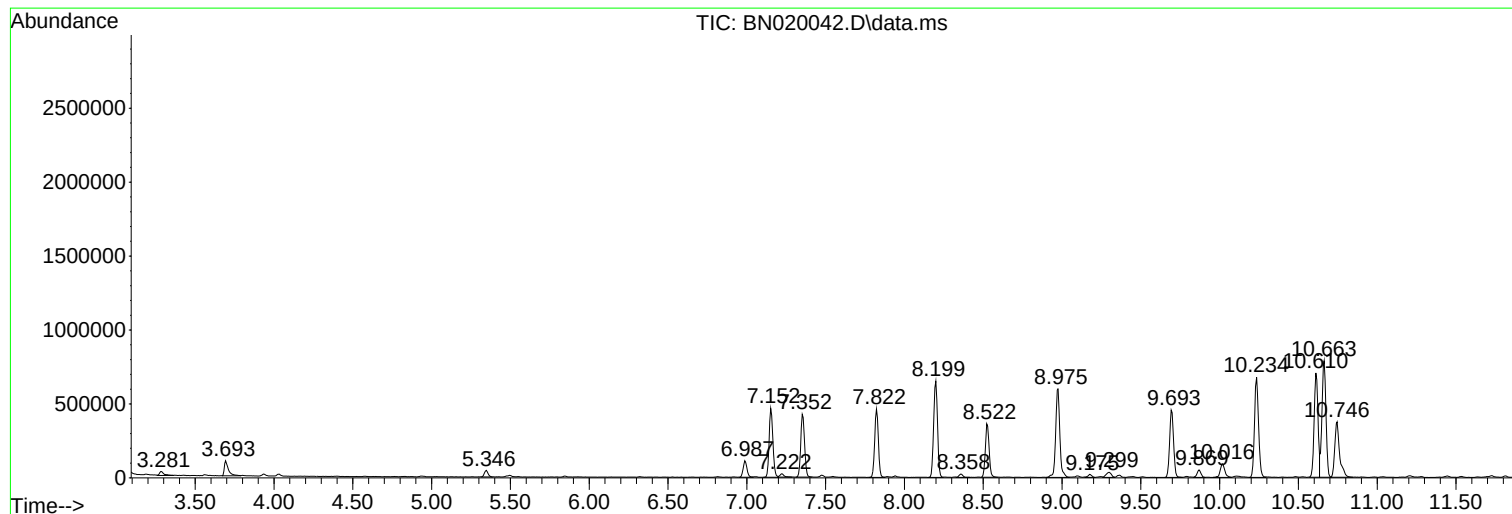
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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN060622.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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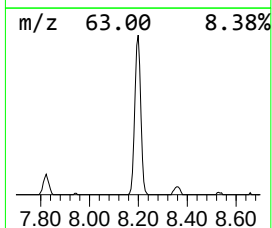
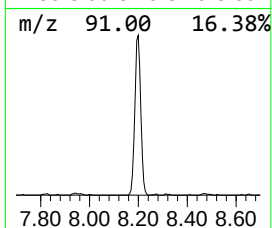
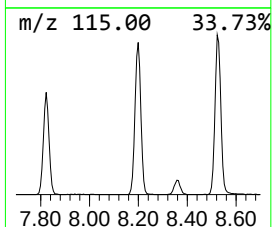
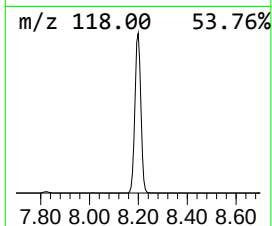
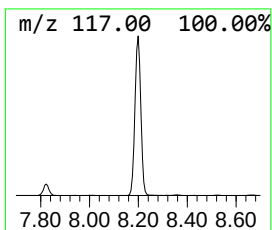
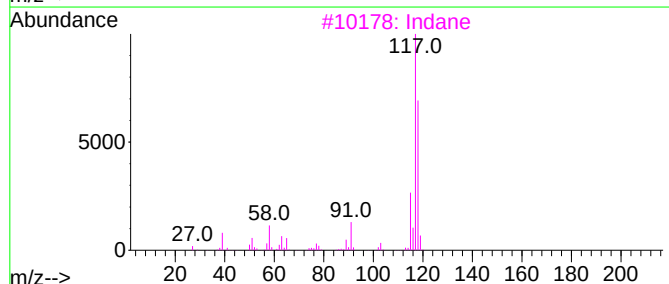
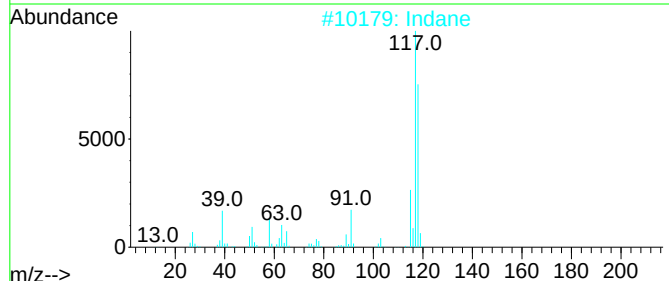
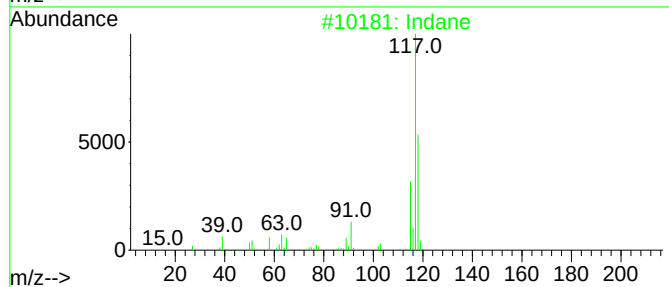
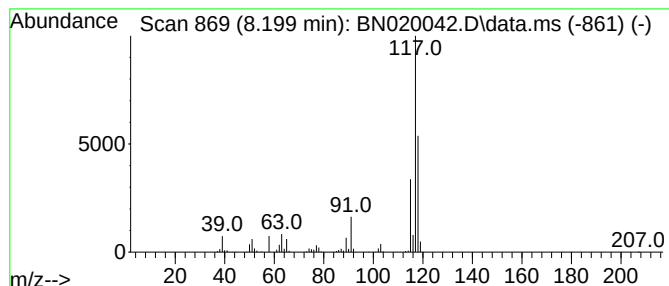
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN060622.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.199	28.83 ng/ul	1086300	1,4-Dichlorobenzene-d4	7.822

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane			118	C9H10	000496-11-7	93
2	Indane			118	C9H10	000496-11-7	91
3	Indane			118	C9H10	000496-11-7	83
4	Indane			118	C9H10	000496-11-7	72
5	Benzene, 1-ethenyl-2-methyl-			118	C9H10	000611-15-4	60



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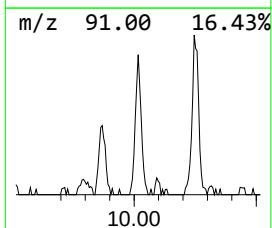
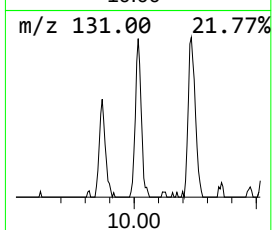
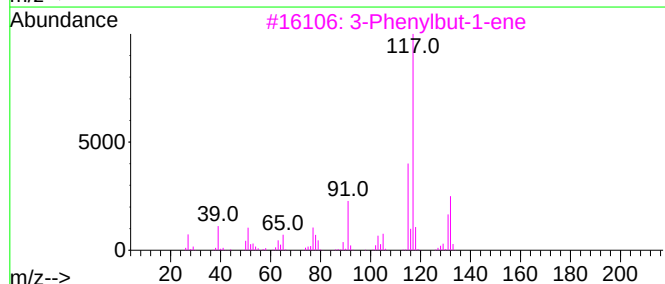
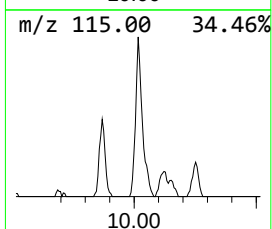
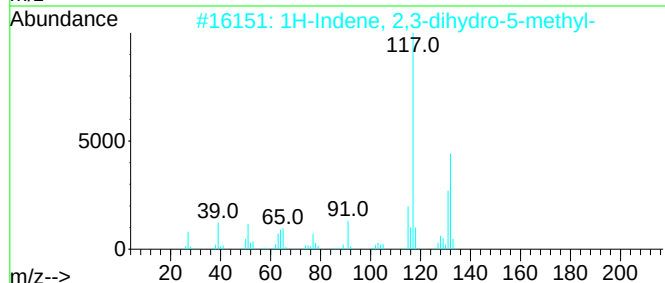
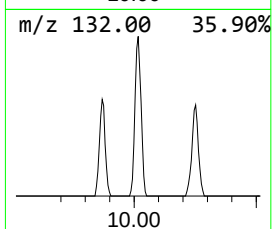
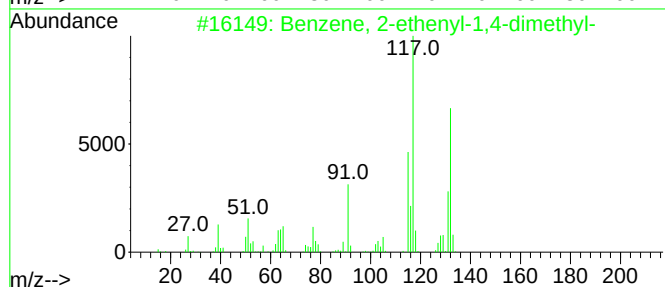
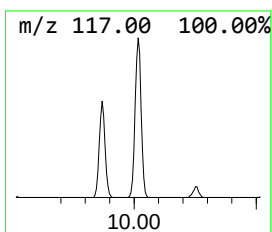
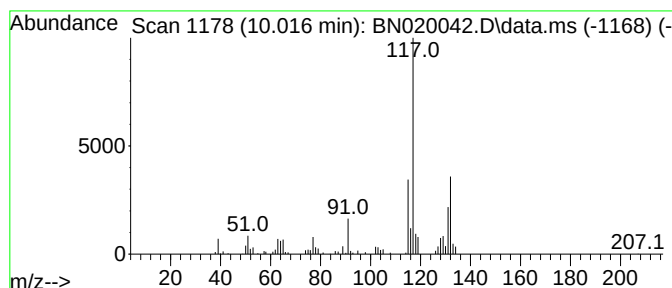
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN060622.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.016	2.95 ng/ul	178126	Naphthalene-d8	10.610

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
3		3-Phenylbut-1-ene	132	C10H12	000934-10-1	91
4		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90
5		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87



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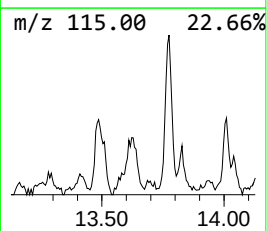
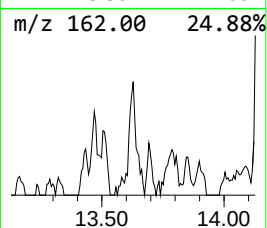
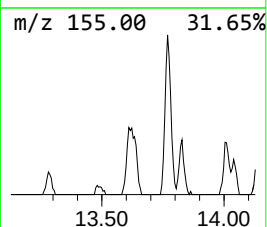
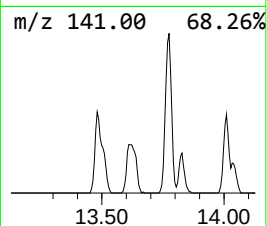
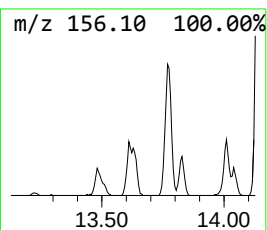
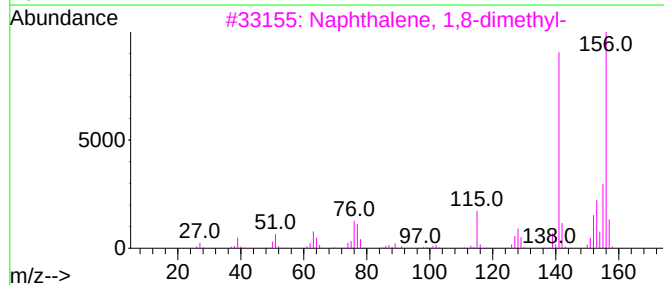
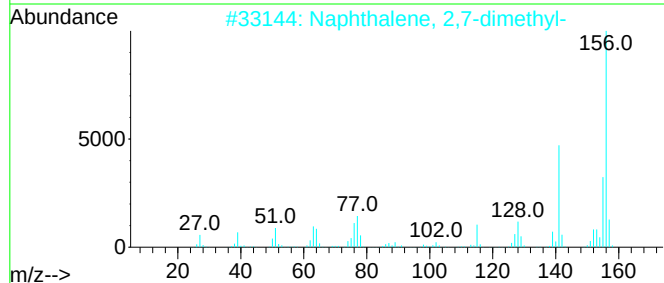
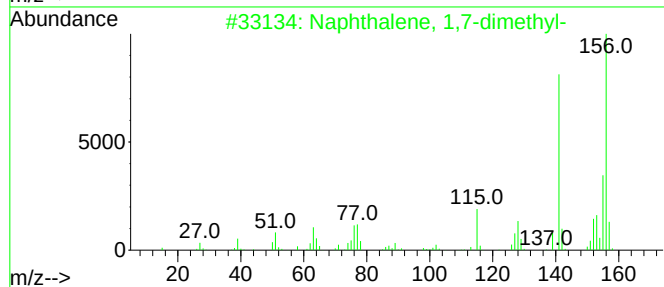
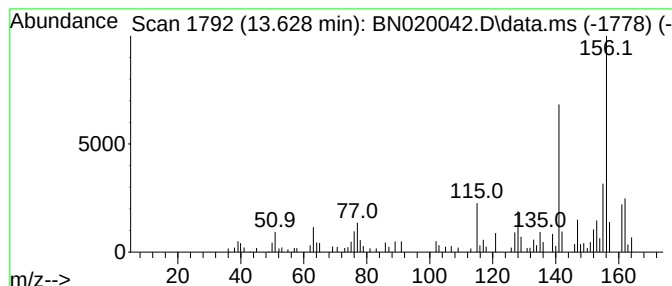
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN060622.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Naphthalene, 1,7-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.628	2.08 ng/ul	177234	Acenaphthene-d10	14.451

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96
2			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96
3			Naphthalene, 1,8-dimethyl-	156	C12H12	000569-41-5	96
4			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	96
5			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96



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Data File : BN020042.D
Acq On : 07 Jun 2022 23:31
Operator : CG/JU
Sample : N3171-09
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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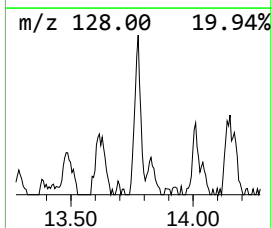
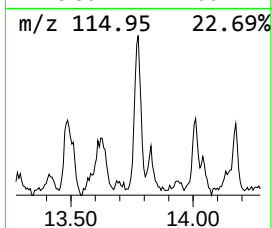
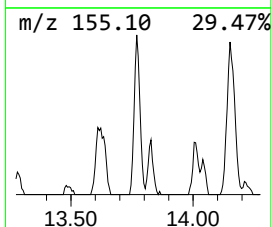
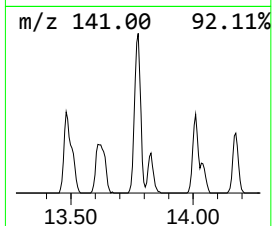
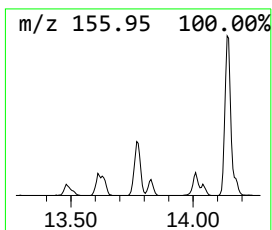
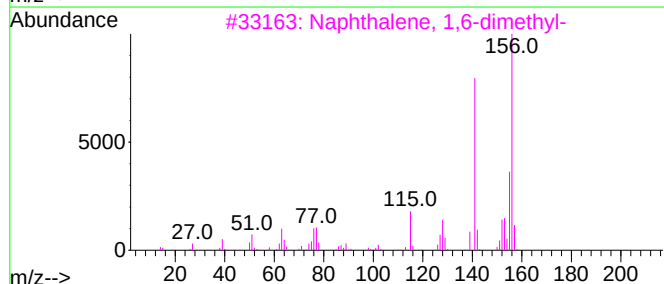
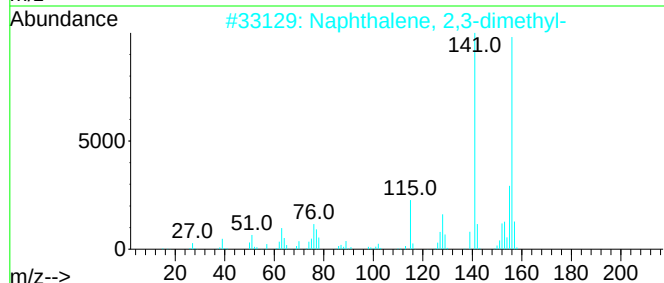
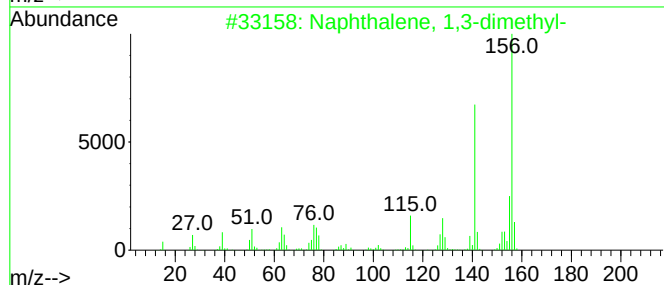
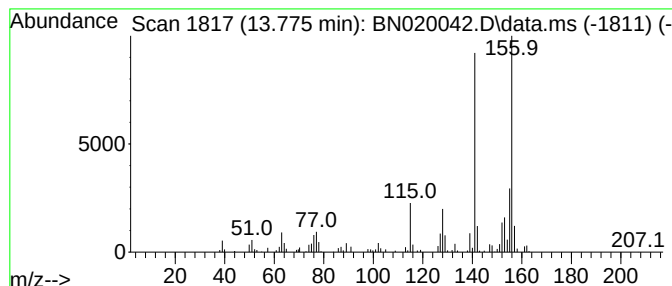
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Naphthalene, 1,3-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.775	2.52 ng/ul	214741	Acenaphthene-d10	14.451

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	97
2			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96
3			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
4			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	96
5			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060722\
Data File : BN020042.D
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Sample : N3171-09
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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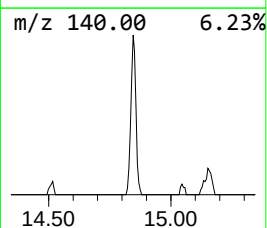
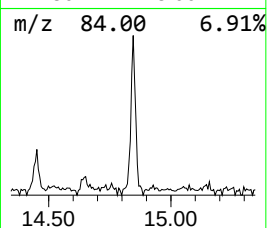
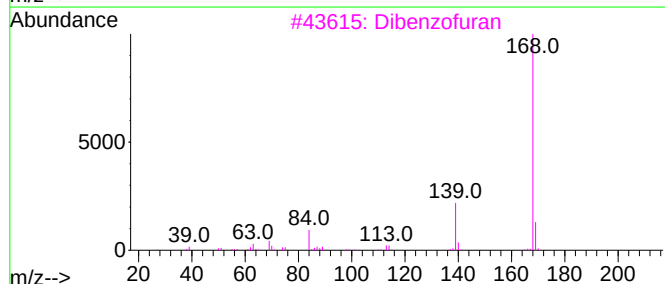
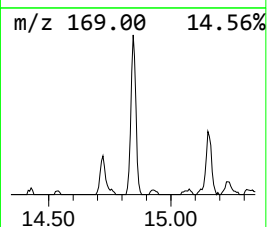
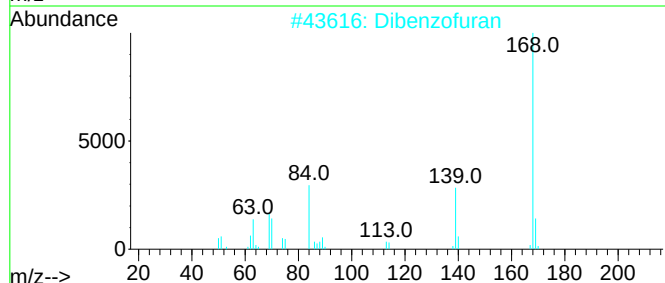
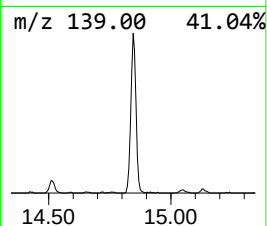
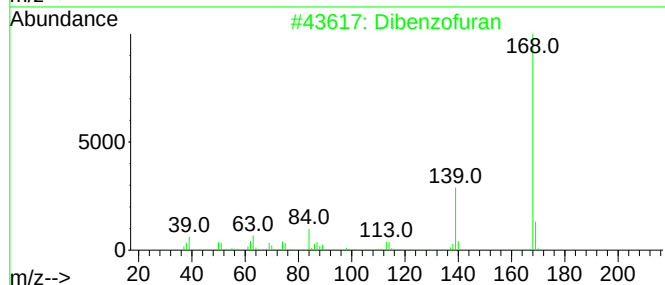
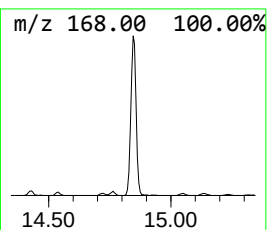
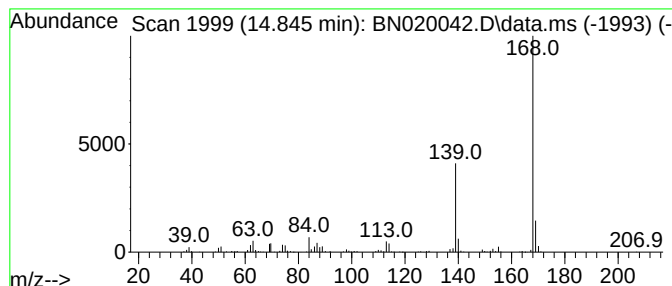
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Dibenzo furan Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.845	4.22 ng/ul	359314	Acenaphthene-d10	14.451

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dibenzofuran	168	C12H8O	000132-64-9	91
2			Dibenzofuran	168	C12H8O	000132-64-9	86
3			Dibenzofuran	168	C12H8O	000132-64-9	78
4			Benzene, 1,3,5-trimethoxy-	168	C9H12O3	000621-23-8	72
5			Pyrimidine, 2-(dimethylamino)-5-...	168	C6H8N4O2	014233-44-4	56



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060722\
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Operator : CG/JU
Sample : N3171-09
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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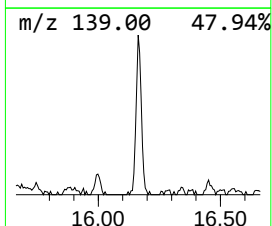
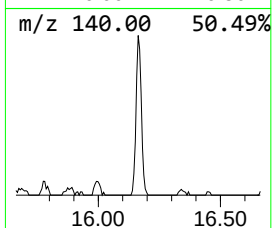
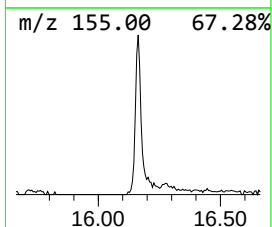
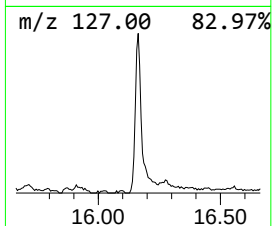
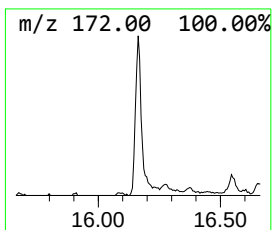
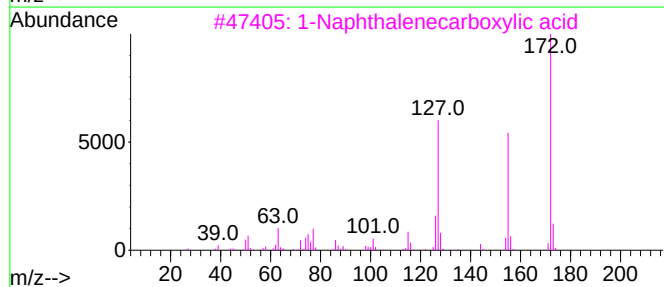
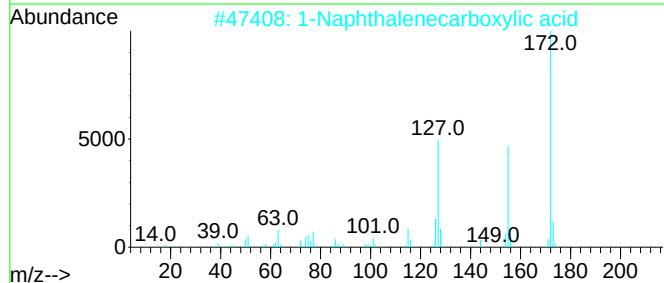
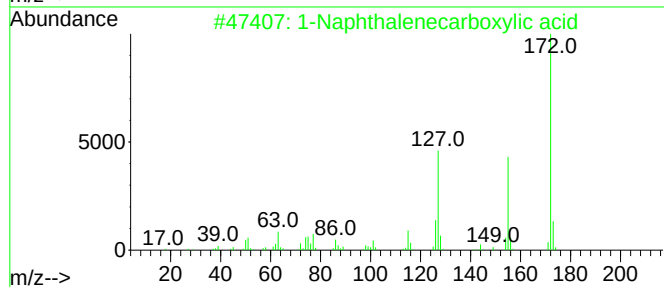
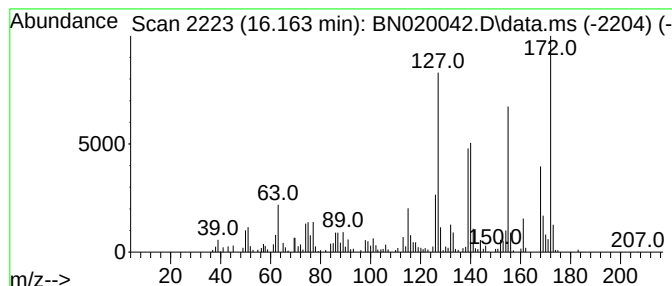
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 1-Naphthalenecarboxylic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.163	3.20 ng/ul	331010	Phenanthrene-d10	17.186

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Naphthalenecarboxylic acid	172	C11H8O2	000086-55-5	96		
2	1-Naphthalenecarboxylic acid	172	C11H8O2	000086-55-5	93		
3	1-Naphthalenecarboxylic acid	172	C11H8O2	000086-55-5	83		
4	2-Naphthalenecarboxylic acid	172	C11H8O2	000093-09-4	83		
5	2-Naphthalenecarboxylic acid	172	C11H8O2	000093-09-4	60		



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060722\
Data File : BN020042.D
Acq On : 07 Jun 2022 23:31
Operator : CG/JU
Sample : N3171-09
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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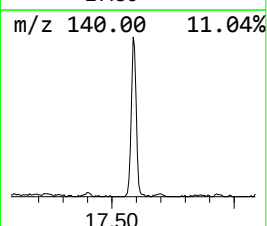
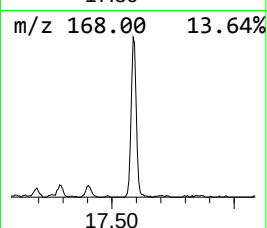
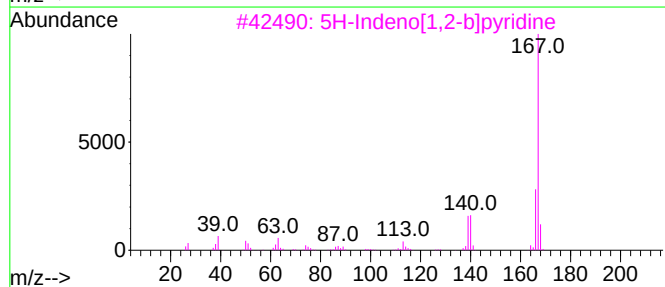
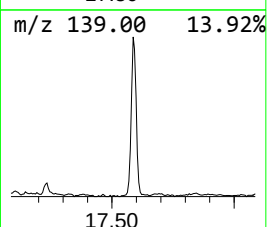
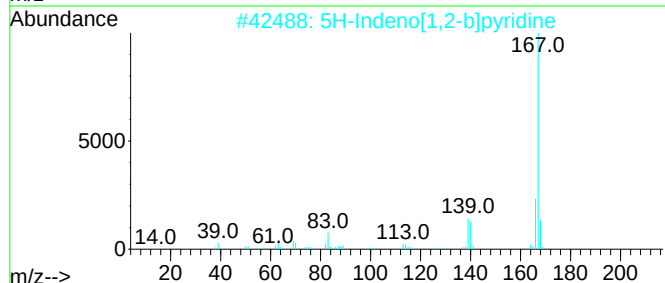
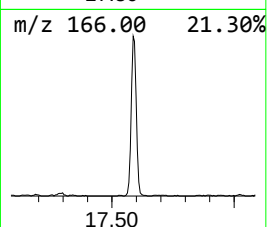
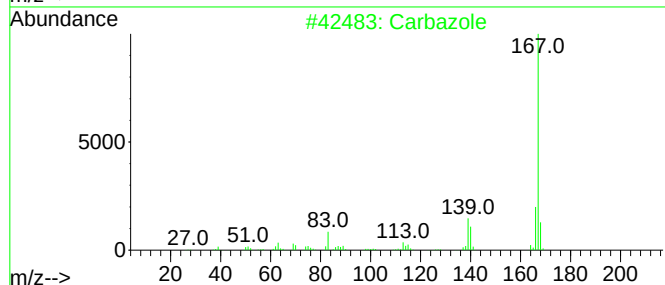
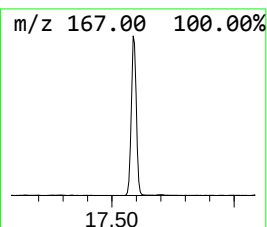
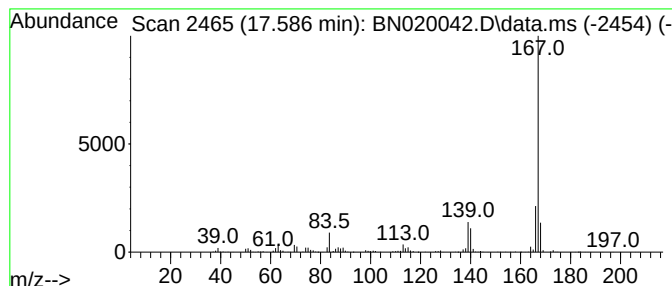
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Carba zole Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.587	7.59 ng/ul	786443	Phenanthrene-d10	17.186

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbazole	167	C12H9N	000086-74-8	96
2			5H-Indeno[1,2-b]pyridine	167	C12H9N	000244-99-5	95
3			5H-Indeno[1,2-b]pyridine	167	C12H9N	000244-99-5	90
4			Carbazole	167	C12H9N	000086-74-8	90
5			9H-Carbazole-9-methanol	197	C13H11NO	002409-36-1	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060722\
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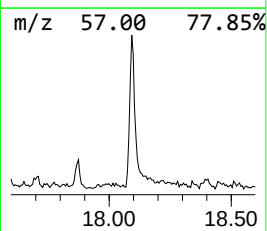
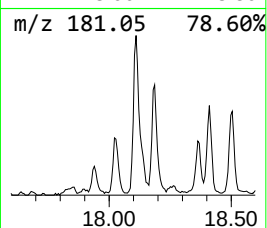
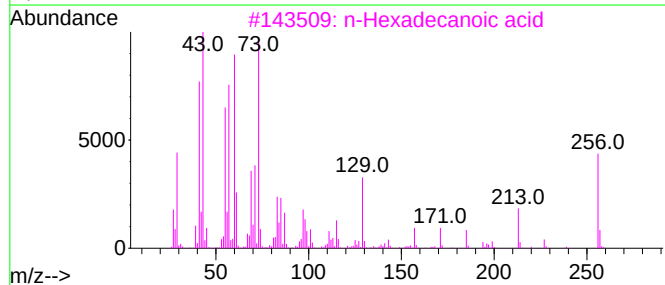
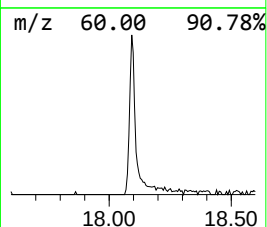
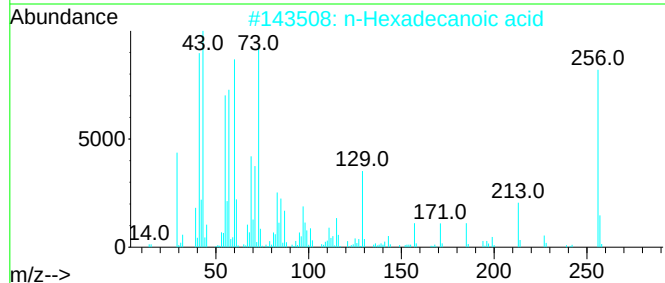
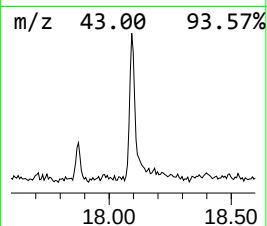
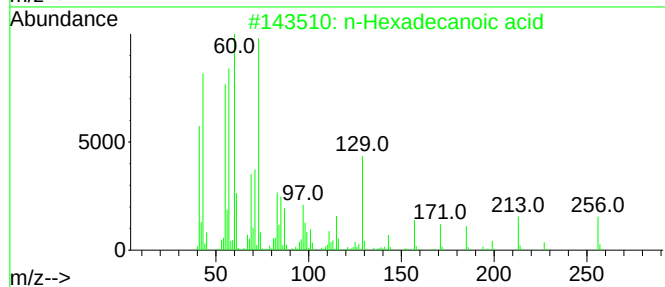
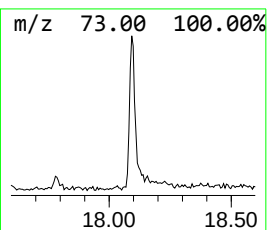
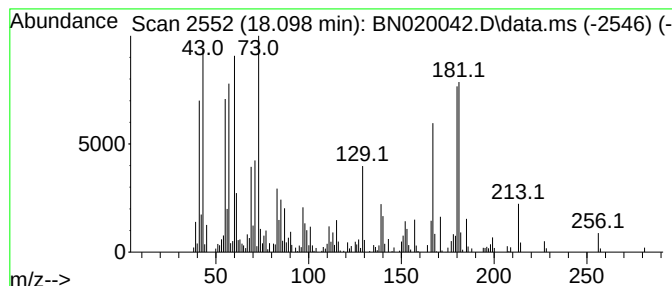
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.098	3.23 ng/ul	334222	Phenanthrene-d10	17.186

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
5			5H-Indeno[1,2-b]pyridine, 4-methyl-	181	C13H11N	064292-01-9	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060722\
Data File : BN020042.D
Acq On : 07 Jun 2022 23:31
Operator : CG/JU
Sample : N3171-09
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

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					#	RT	Resp	Conc
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Benzene, 2-ethe...	10.016	3.0	ng/ul	178126	2	10.610	1205820	20.0
Naphthalene, 1,...	13.628	2.1	ng/ul	177234	3	14.451	1704100	20.0
Naphthalene, 1,...	13.775	2.5	ng/ul	214741	3	14.451	1704100	20.0
Dibenzo furan	14.845	4.2	ng/ul	359314	3	14.451	1704100	20.0
1-Naphthaleneca...	16.163	3.2	ng/ul	331010	4	17.186	2071320	20.0
Carba zole	17.587	7.6	ng/ul	786443	4	17.186	2071320	20.0
n-Hexadecanoic ...	18.098	3.2	ng/ul	334222	4	17.186	2071320	20.0