

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
 Data File : BN015968.D
 Acq On : 19 Aug 2021 11:49
 Operator : CG/JU
 Sample : M3399-05 30X
 Misc :
 ALS Vial : 75 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBKA7

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-BN081621MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BN015968.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.405	387	394	401	rBV	530401	861195	1.36%	0.161%
2	7.622	765	771	782	rVB	467900	823781	1.31%	0.154%
3	7.899	809	818	826	rBV	965834	1653991	2.62%	0.309%
4	8.275	874	882	889	rVB	472402	789976	1.25%	0.148%
5	8.440	898	910	923	rBV	4702986	8105156	12.85%	1.515%
6	10.104	1188	1193	1203	rVB3	506750	1118058	1.77%	0.209%
7	10.340	1227	1233	1241	rVB	1481253	2625295	4.16%	0.491%
8	10.710	1286	1296	1299	rVV2	1370588	3130573	4.96%	0.585%
9	10.769	1299	1306	1313	rVV	19782487	35934110	56.96%	6.718%
10	10.881	1319	1325	1331	rVV	965816	1706239	2.70%	0.319%
11	12.140	1534	1539	1544	rVV	746090	1231729	1.95%	0.230%
12	12.204	1544	1550	1556	rVV2	536392	1112971	1.76%	0.208%
13	12.381	1570	1580	1586	rVB	15210476	25768845	40.84%	4.818%
14	12.487	1592	1598	1605	rBV2	350701	719176	1.14%	0.134%
15	12.593	1605	1616	1626	rVB	7889072	13299675	21.08%	2.487%
16	13.381	1742	1750	1757	rBV	5553790	8579927	13.60%	1.604%
17	13.575	1778	1783	1797	rVB	1971463	4085884	6.48%	0.764%
18	13.716	1798	1807	1808	rBV	2620299	4553841	7.22%	0.851%
19	13.869	1826	1833	1838	rVV	3950058	7250509	11.49%	1.356%
20	13.922	1838	1842	1853	rVV	2553741	4527134	7.18%	0.846%
21	14.034	1853	1861	1865	rVV3	397114	876645	1.39%	0.164%
22	14.104	1865	1873	1877	rVV	1546320	2985512	4.73%	0.558%
23	14.140	1877	1879	1885	rVB2	572030	817184	1.30%	0.153%
24	14.269	1893	1901	1907	rVB2	1294536	2158965	3.42%	0.404%
25	14.451	1927	1932	1937	rBV	1970409	2684598	4.26%	0.502%
26	14.528	1939	1945	1946	rVV	1433928	2131366	3.38%	0.398%
27	14.551	1946	1949	1954	rVV	1942263	2936579	4.65%	0.549%
28	14.616	1954	1960	1966	rVV	18874662	31203582	49.46%	5.834%
29	14.675	1966	1970	1976	rVB	1215138	1799837	2.85%	0.337%
30	14.757	1977	1984	1992	rBV2	1017981	2544909	4.03%	0.476%
31	14.857	1992	2001	2005	rVV3	556600	1267187	2.01%	0.237%
32	14.951	2010	2017	2023	rVV	12456605	19731014	31.27%	3.689%
33	15.022	2023	2029	2034	rVV	1320328	2104633	3.34%	0.393%
34	15.092	2034	2041	2046	rVV	5397623	8139722	12.90%	1.522%
35	15.169	2046	2054	2059	rVV3	2131986	3811341	6.04%	0.713%
36	15.222	2059	2063	2067	rVB	974076	1335019	2.12%	0.250%

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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-BN081621MA.M
 Title : SVOA CALIBRATION

37	15.339	2075	2083	2086	rBV2	627473	1455980	2.31%	0.272%
38	15.404	2090	2094	2099	rVB	2126926	2876657	4.56%	0.538%
39	15.510	2108	2112	2120	rVB5	330320	795549	1.26%	0.149%
40	15.604	2120	2128	2133	rBV	13174104	20799281	32.97%	3.889%

41	15.722	2144	2148	2151	rVV3	1320566	2064356	3.27%	0.386%
42	15.757	2151	2154	2159	rVV3	1645322	2546058	4.04%	0.476%
43	15.810	2159	2163	2166	rVV2	1018541	1529639	2.42%	0.286%
44	15.845	2166	2169	2173	rVV	824996	1261169	2.00%	0.236%
45	15.892	2173	2177	2185	rVB	1798923	3267093	5.18%	0.611%

46	15.969	2185	2190	2193	rBV5	361727	735769	1.17%	0.138%
47	16.034	2194	2201	2210	rVB	1681054	4319197	6.85%	0.808%
48	16.116	2210	2215	2225	rVB3	462022	1123838	1.78%	0.210%
49	16.263	2236	2240	2255	rVB2	3534685	6081142	9.64%	1.137%
50	16.392	2256	2262	2269	rBV2	1121375	2108555	3.34%	0.394%

51	16.604	2292	2298	2302	rVV2	1126311	2157650	3.42%	0.403%
52	16.651	2302	2306	2312	rVV4	809136	1455326	2.31%	0.272%
53	16.875	2341	2344	2351	rVB4	753406	1081594	1.71%	0.202%
54	16.951	2351	2357	2364	rVV	10330528	15592138	24.71%	2.915%
55	17.063	2366	2376	2381	rVV3	2281967	4620741	7.32%	0.864%

56	17.116	2381	2385	2395	rVB	3402706	5492052	8.70%	1.027%
57	17.304	2411	2417	2419	rBV	2078151	3455715	5.48%	0.646%
58	17.357	2419	2426	2430	rVV2	38076530	63091546	100.00%	11.796%
59	17.404	2430	2434	2435	rVV2	1161254	1499195	2.38%	0.280%
60	17.433	2435	2439	2444	rVV	4416894	6102387	9.67%	1.141%

61	17.492	2444	2449	2454	rVV	626893	1111694	1.76%	0.208%
62	17.704	2480	2485	2496	rVB	3827791	5953043	9.44%	1.113%
63	17.804	2496	2502	2508	rBV4	2116334	3683871	5.84%	0.689%
64	18.028	2535	2540	2546	rBV	476863	909476	1.44%	0.170%
65	18.180	2561	2566	2569	rBV	2309957	3372600	5.35%	0.631%

66	18.228	2569	2574	2581	rVB2	3636737	5768081	9.14%	1.078%
67	18.304	2581	2587	2593	rBV3	1135094	2234090	3.54%	0.418%
68	18.375	2593	2599	2603	rBV3	4781462	8177669	12.96%	1.529%
69	18.492	2615	2619	2627	rVB2	1232553	1915423	3.04%	0.358%
70	18.675	2645	2650	2655	rBV	1959418	2785987	4.42%	0.521%

71	18.922	2685	2692	2696	rBV2	589887	1020333	1.62%	0.191%
72	18.975	2696	2701	2704	rVV2	563475	897623	1.42%	0.168%
73	19.092	2716	2721	2724	rBV	930568	1536798	2.44%	0.287%
74	19.127	2724	2727	2730	rVB	704653	761367	1.21%	0.142%
75	19.363	2760	2767	2773	rBV	24516162	36297645	57.53%	6.786%

76	19.604	2803	2808	2812	rVB	649755	884000	1.40%	0.165%
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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-BN081621MA.M
 Title : SVOA CALIBRATION

77	19.727	2817	2829	2836	rVV	16406140	33017177	52.33%	6.173%
78	19.792	2836	2840	2846	rVB3	766759	1372832	2.18%	0.257%
79	19.898	2853	2858	2864	rVB	899004	1302533	2.06%	0.244%
80	20.039	2877	2882	2883	rBV2	667632	930603	1.48%	0.174%
81	20.098	2888	2892	2897	rVB	564329	710227	1.13%	0.133%
82	20.180	2898	2906	2907	rBV3	393323	775340	1.23%	0.145%
83	20.216	2907	2912	2923	rVB	2366825	3929814	6.23%	0.735%
84	20.316	2923	2929	2933	rBV	2937161	3930039	6.23%	0.735%
85	20.369	2933	2938	2942	rVV2	757262	1376569	2.18%	0.257%
86	20.410	2942	2945	2949	rVV	590975	757047	1.20%	0.142%
87	20.739	2997	3001	3006	rVV2	656489	833438	1.32%	0.156%
88	21.127	3063	3067	3070	rBV	732898	974703	1.54%	0.182%
89	21.169	3070	3074	3077	rVV	841700	1030654	1.63%	0.193%
90	21.210	3077	3081	3086	rVB2	698839	1170801	1.86%	0.219%
91	21.474	3116	3126	3131	rBV3	6408490	12695362	20.12%	2.374%
92	21.521	3131	3134	3145	rVB	4195443	5683332	9.01%	1.063%
93	21.645	3151	3155	3163	rVB	1518172	2084192	3.30%	0.390%
94	21.810	3178	3183	3189	rBV2	559894	909524	1.44%	0.170%
95	22.316	3265	3269	3274	rVB2	468873	789189	1.25%	0.148%
96	23.168	3407	3414	3420	rBV	1688177	4493940	7.12%	0.840%
97	23.692	3497	3503	3509	rBV	695423	1377569	2.18%	0.258%
98	23.798	3515	3521	3530	rVB2	1224024	2783326	4.41%	0.520%
99	23.910	3530	3540	3545	rBV3	1615032	3761144	5.96%	0.703%
100	26.409	3958	3965	3976	rVB2	311674	951893	1.51%	0.178%

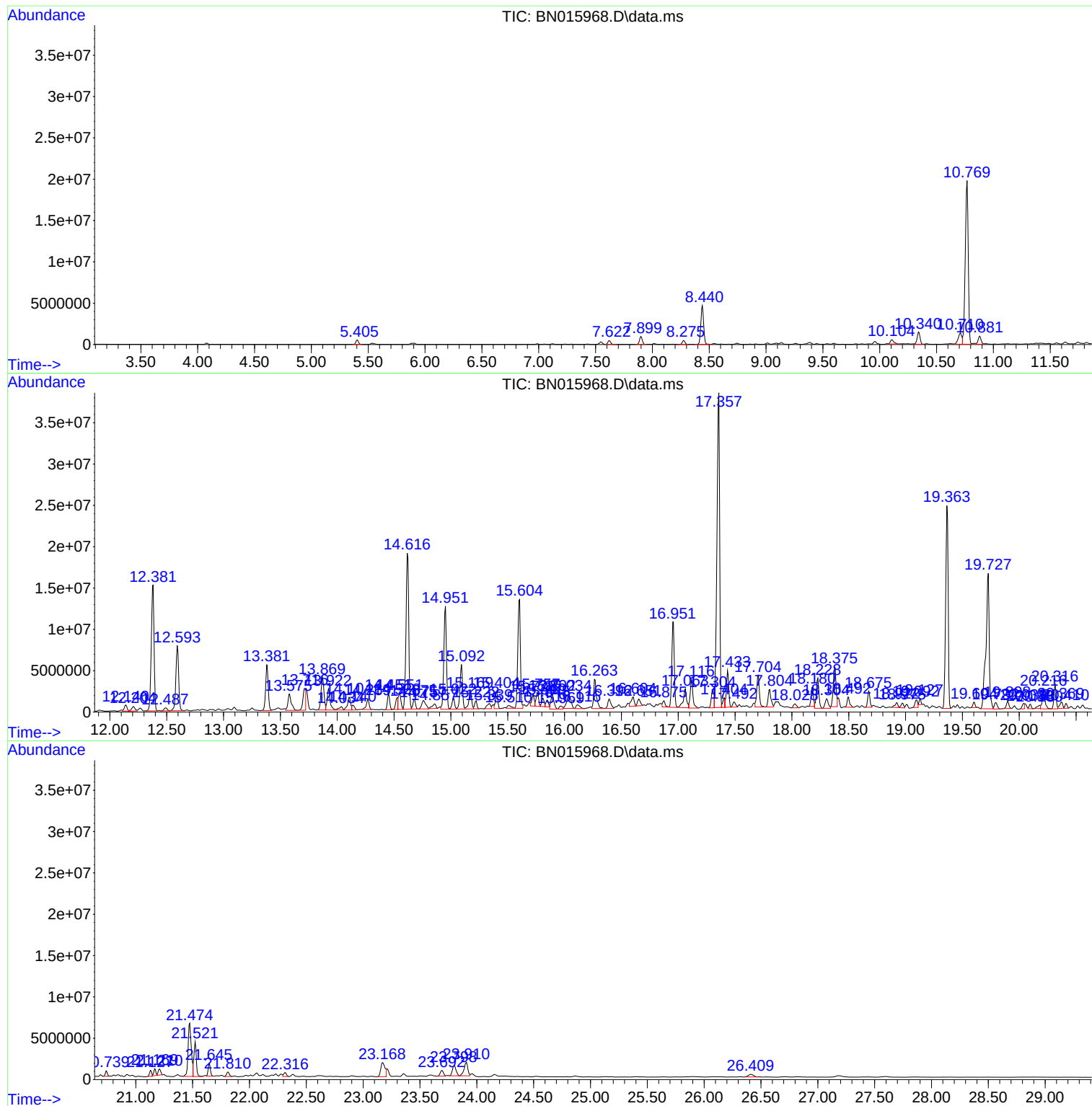
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081621MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
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Instrument :
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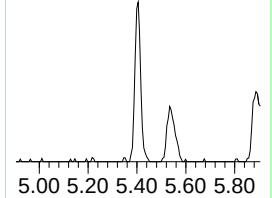
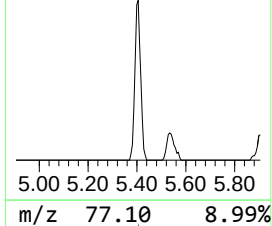
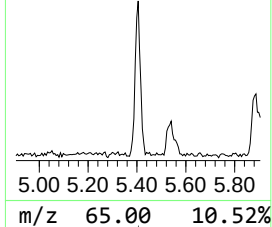
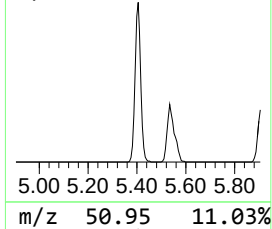
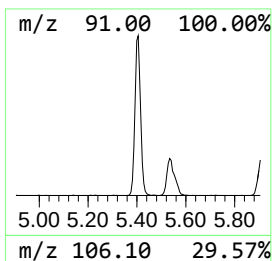
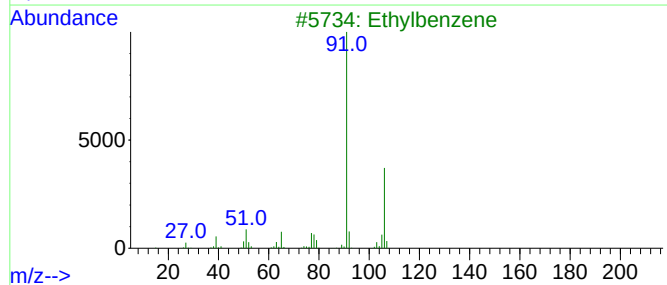
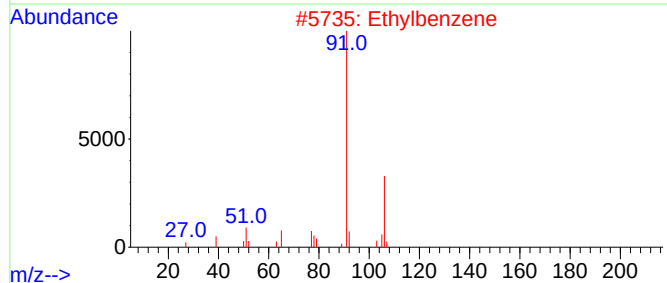
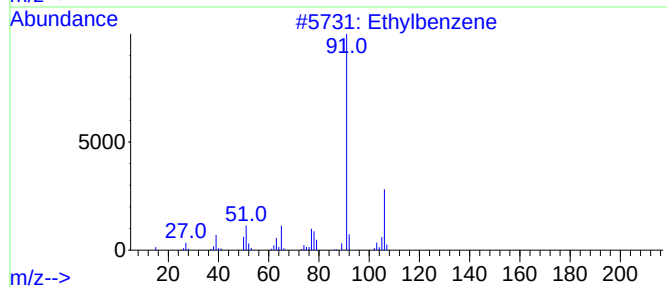
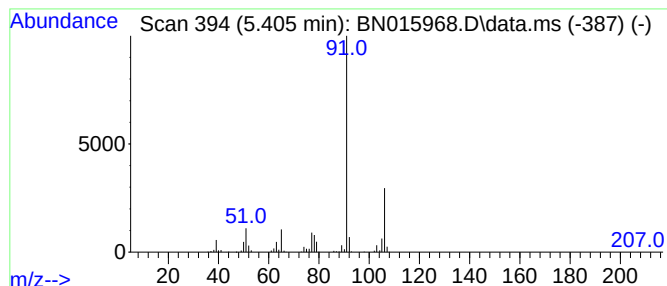
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081621MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Ethylbenzene Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.405	10.41 ng/ul	861195	1,4-Dichlorobenzene-d4	7.899

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethylbenzene	106	C8H10	000100-41-4	91
2			Ethylbenzene	106	C8H10	000100-41-4	91
3			Ethylbenzene	106	C8H10	000100-41-4	91
4			Ethylbenzene	106	C8H10	000100-41-4	91
5			p-Xylene	106	C8H10	000106-42-3	78



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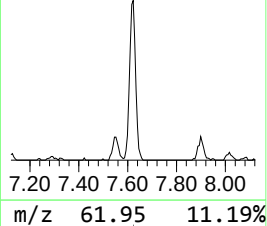
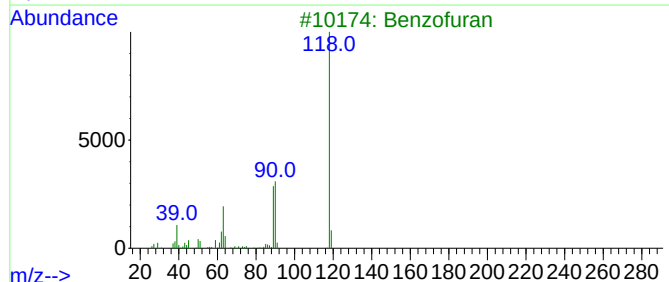
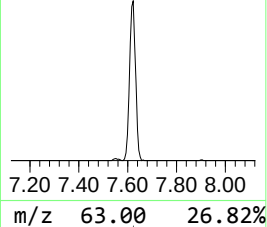
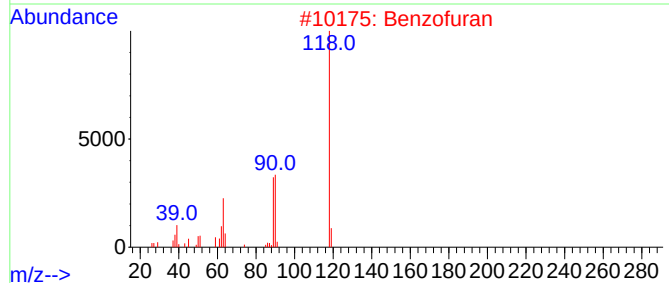
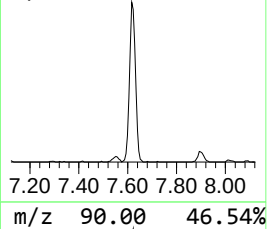
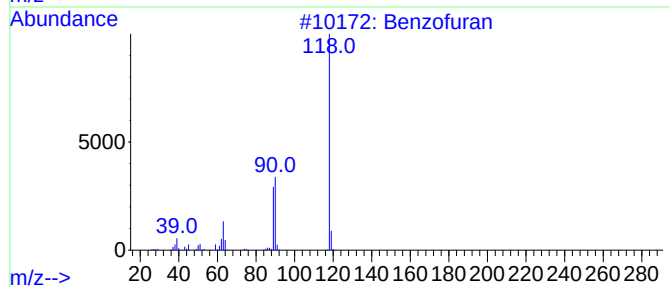
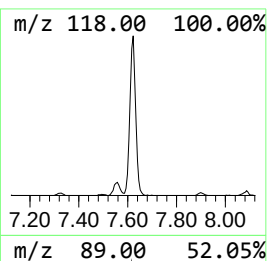
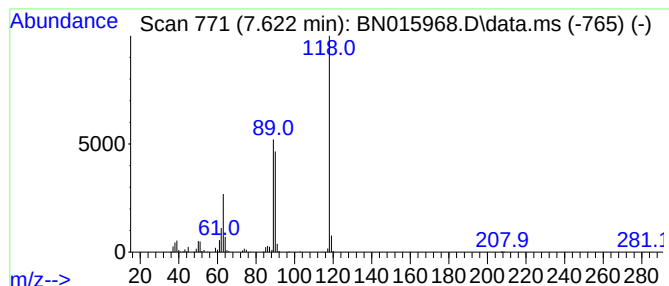
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081621MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzofuran Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.622	9.96 ng/ul	823781	1,4-Dichlorobenzene-d4	7.899

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzofuran	118	C8H6O	000271-89-6	94
2			Benzofuran	118	C8H6O	000271-89-6	94
3			Benzofuran	118	C8H6O	000271-89-6	91
4			Benzofuran	118	C8H6O	000271-89-6	91
5			3-Hydroxyphenylacetylene	118	C8H6O	010401-11-3	87



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Misc :
ALS Vial : 75 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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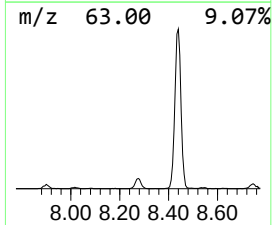
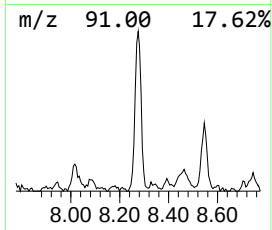
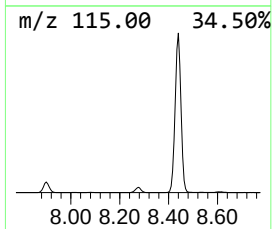
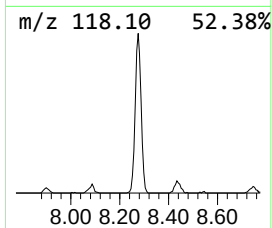
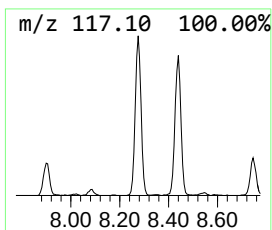
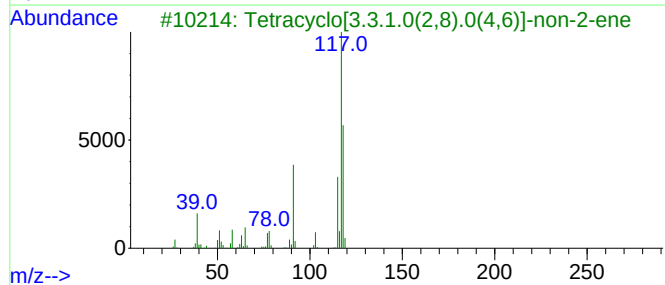
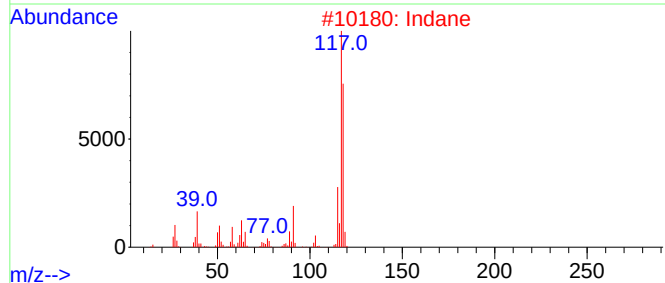
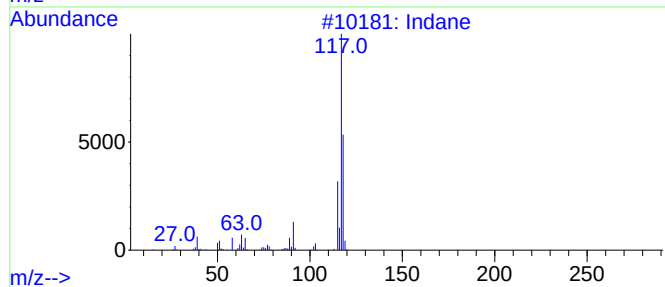
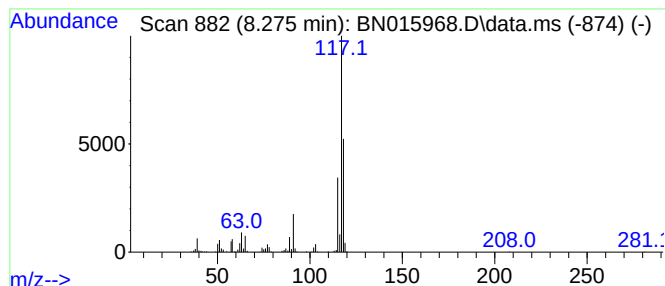
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081621MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Tetracyclo[3.3.1.0(2,8).0(4... Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.275	9.55 ng/ul	789976	1,4-Dichlorobenzene-d4	7.899

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	93
2			Indane	118	C9H10	000496-11-7	90
3			Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	80
4			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	78
5			Benzene, cyclopropyl-	118	C9H10	000873-49-4	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
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Acq On : 19 Aug 2021 11:49
Operator : CG/JU
Sample : M3399-05 30X
Misc :
ALS Vial : 75 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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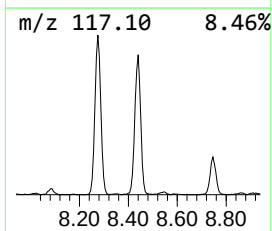
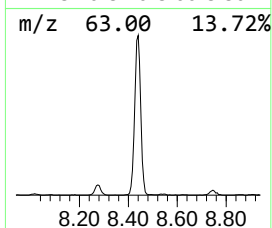
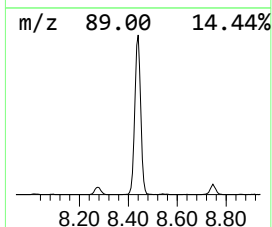
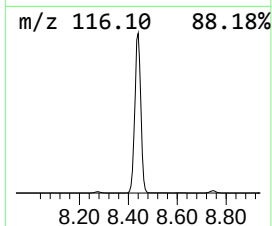
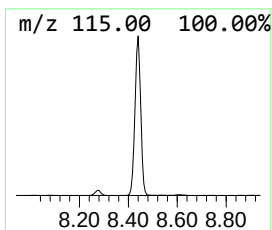
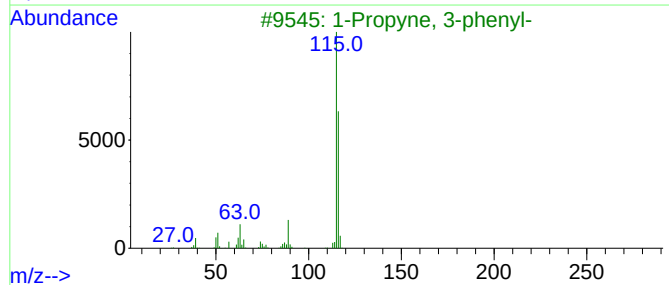
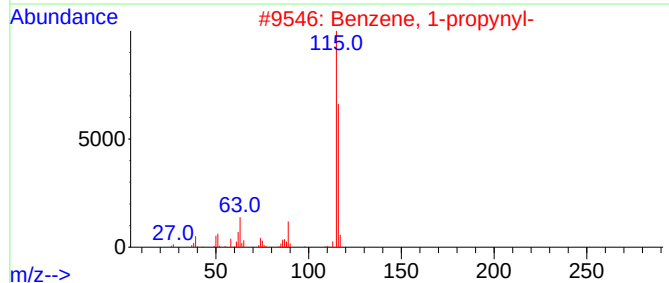
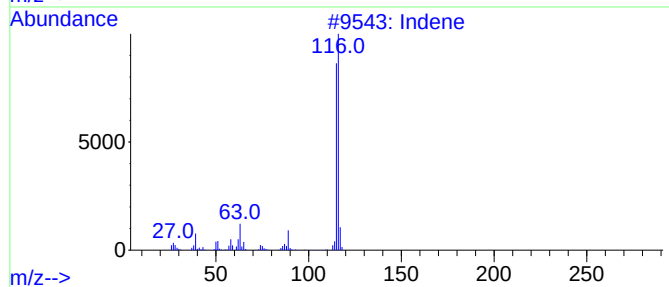
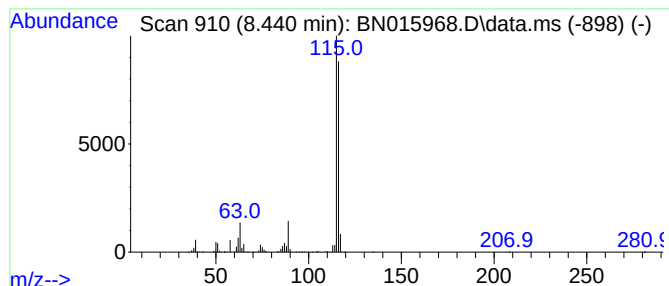
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Indene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.440	98.01 ng/ul	8105160	1,4-Dichlorobenzene-d4	7.899

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indene	116	C9H8	000095-13-6	95
2			Benzene, 1-propynyl-	116	C9H8	000673-32-5	94
3			1-Propyne, 3-phenyl-	116	C9H8	010147-11-2	94
4			Indene	116	C9H8	000095-13-6	91
5			2-Methylphenylacetylene	116	C9H8	000766-47-2	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
Data File : BN015968.D
Acq On : 19 Aug 2021 11:49
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Sample : M3399-05 30X
Misc :
ALS Vial : 75 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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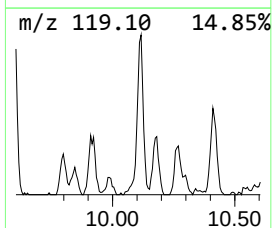
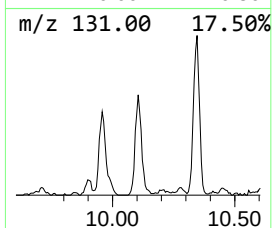
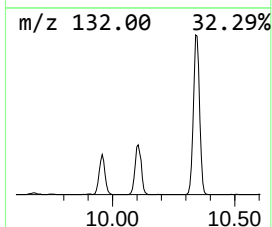
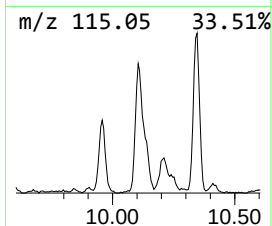
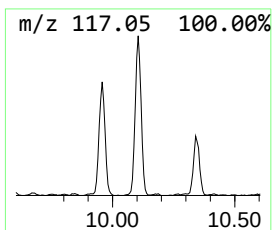
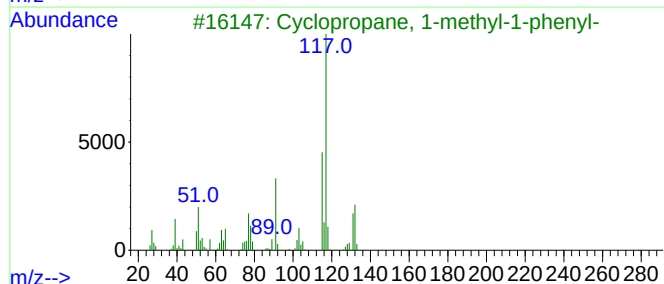
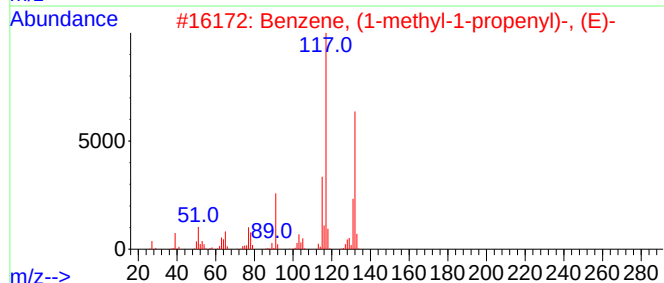
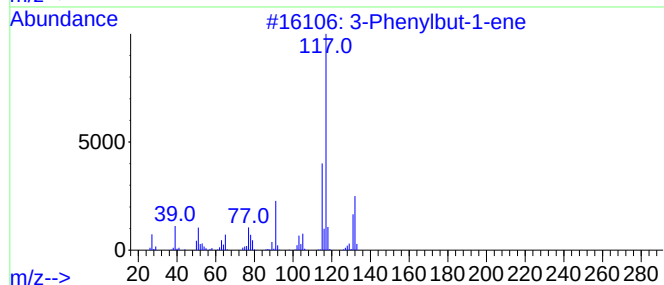
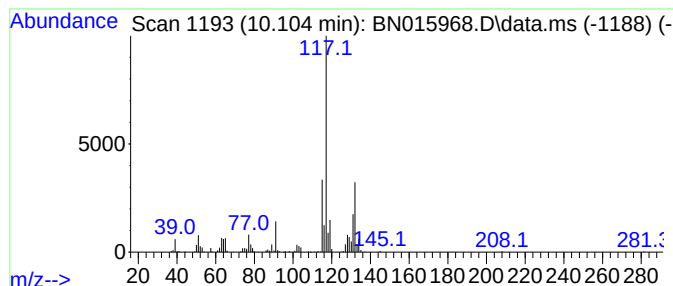
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 3-Phenylbut-1-ene Concentration Rank 40

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.104	7.14 ng/ul	1118060	Naphthalene-d8	10.710

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Phenylbut-1-ene	132	C10H12	000934-10-1	87
2			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	83
3			Cyclopropane, 1-methyl-1-phenyl-	132	C10H12	002214-14-4	81
4			Indan, 1-methyl-	132	C10H12	000767-58-8	81
5			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
Data File : BN015968.D
Acq On : 19 Aug 2021 11:49
Operator : CG/JU
Sample : M3399-05 30X
Misc :
ALS Vial : 75 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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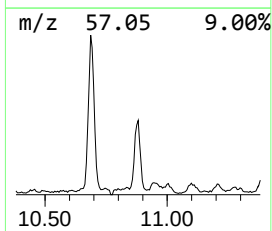
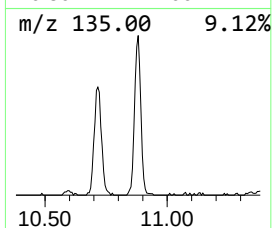
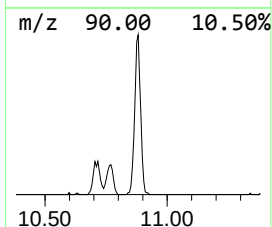
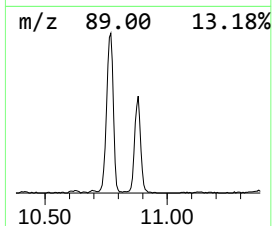
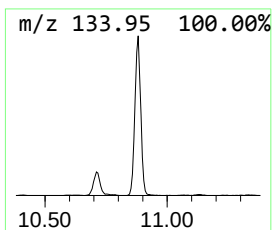
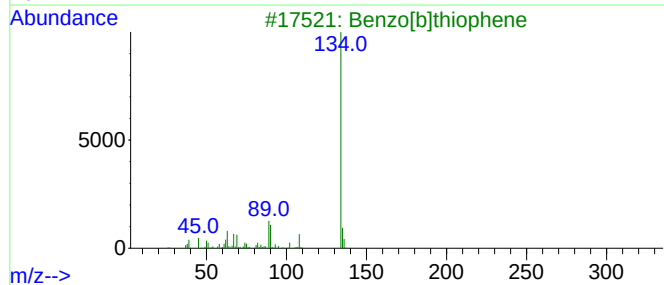
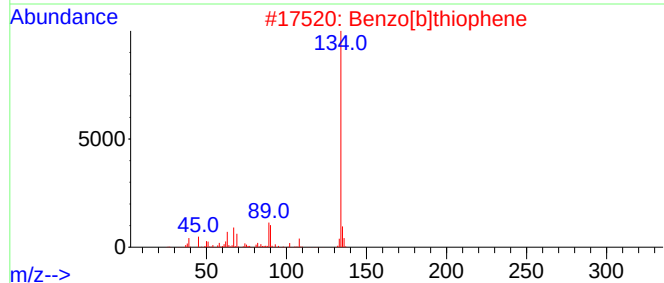
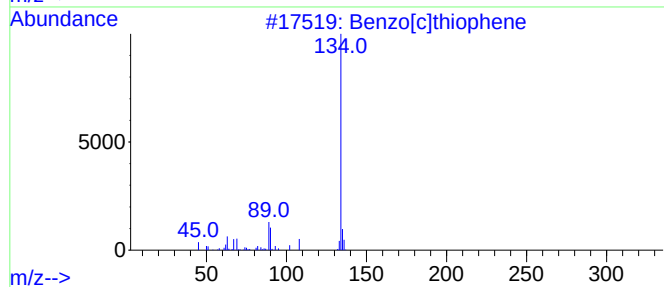
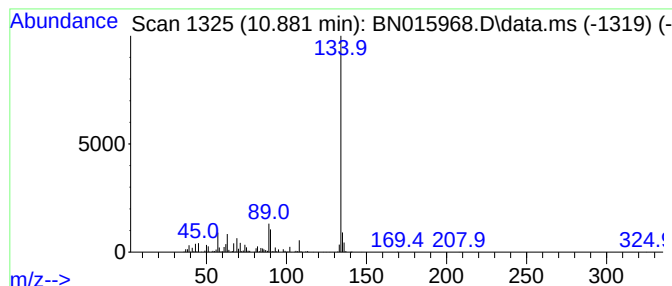
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzo[c]thiophene Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.881	10.90 ng/ul	1706240	Naphthalene-d8	10.710

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[c]thiophene	134	C8H6S	000270-82-6	95
2			Benzo[b]thiophene	134	C8H6S	000095-15-8	94
3			Benzo[b]thiophene	134	C8H6S	000095-15-8	94
4			Cyclopenta[c]thiapyran	134	C8H6S	000270-63-3	90
5			Benzo[b]thiophene	134	C8H6S	000095-15-8	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
Data File : BN015968.D
Acq On : 19 Aug 2021 11:49
Operator : CG/JU
Sample : M3399-05 30X
Misc :
ALS Vial : 75 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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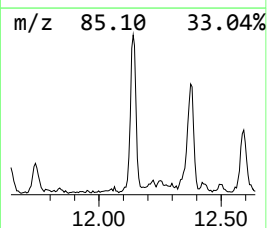
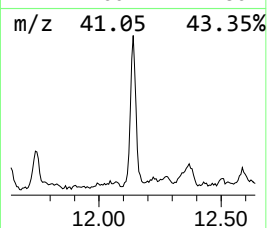
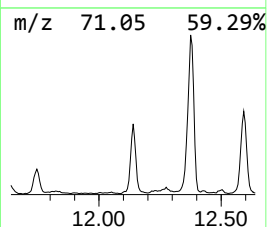
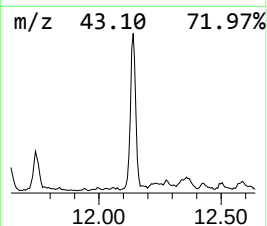
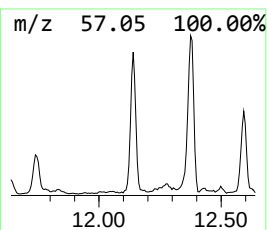
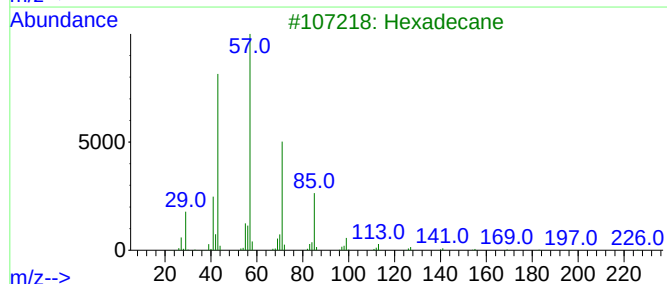
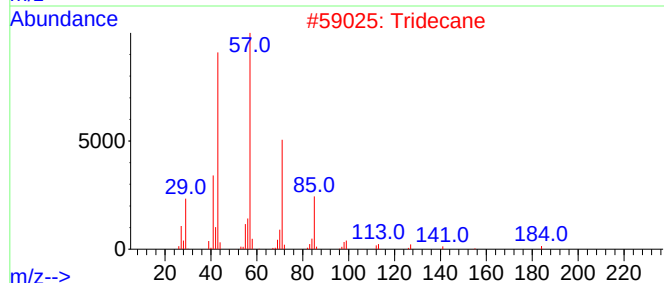
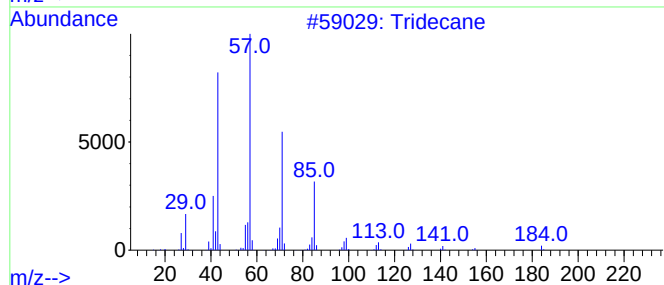
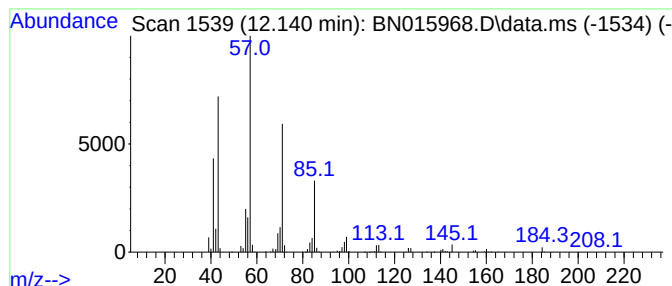
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.140	7.87 ng/ul	1231730	Naphthalene-d8	10.710

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane	184	C13H28	000629-50-5	94
2			Tridecane	184	C13H28	000629-50-5	91
3			Hexadecane	226	C16H34	000544-76-3	86
4			Carbonic acid, octadecyl prop-1-...	354	C22H42O3	1000383-11-5	86
5			Tetradecane	198	C14H30	000629-59-4	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
Data File : BN015968.D
Acq On : 19 Aug 2021 11:49
Operator : CG/JU
Sample : M3399-05 30X
Misc :
ALS Vial : 75 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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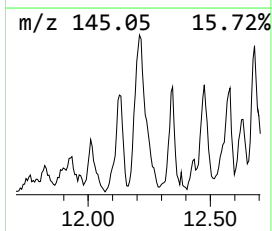
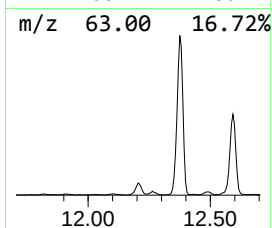
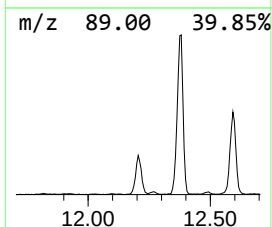
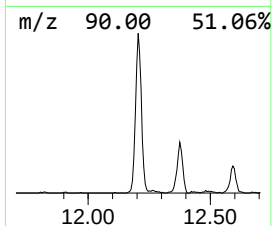
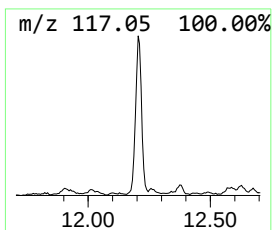
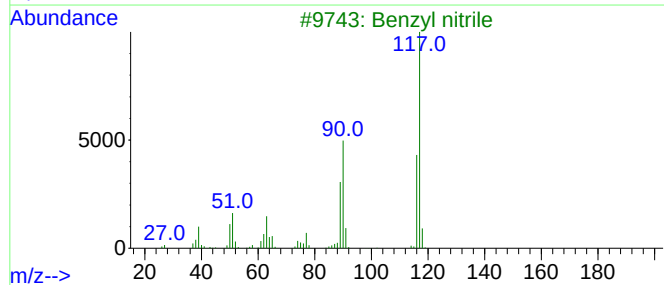
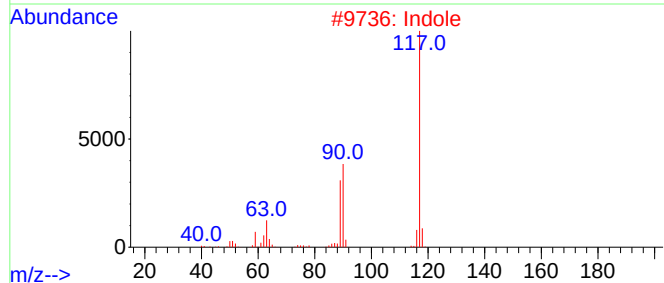
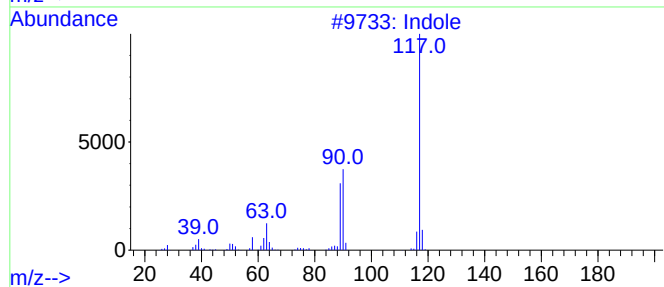
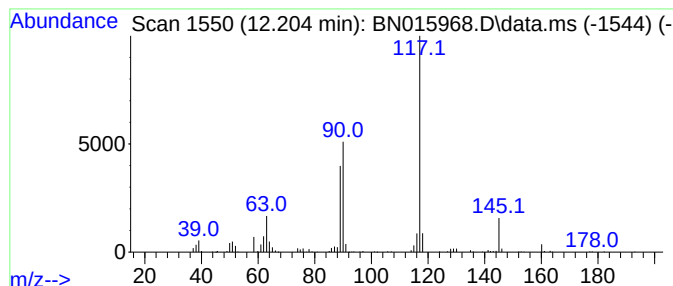
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Indole Concentration Rank 41

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.204	7.11 ng/ul	1112970	Naphthalene-d8	10.710

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indole	117	C8H7N	000120-72-9	96
2			Indole	117	C8H7N	000120-72-9	95
3			Benzyl nitrile	117	C8H7N	000140-29-4	90
4			Indolizine	117	C8H7N	000274-40-8	90
5			Benzyl nitrile	117	C8H7N	000140-29-4	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
Data File : BN015968.D
Acq On : 19 Aug 2021 11:49
Operator : CG/JU
Sample : M3399-05 30X
Misc :
ALS Vial : 75 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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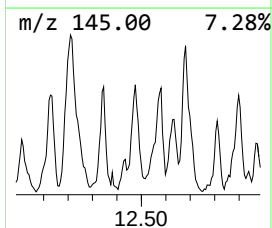
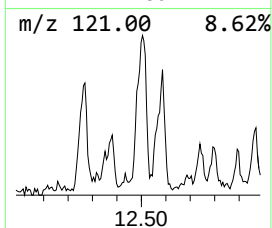
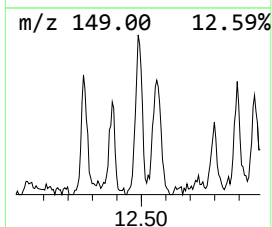
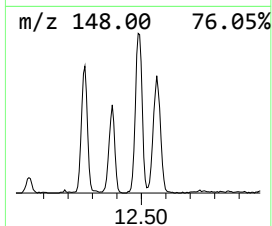
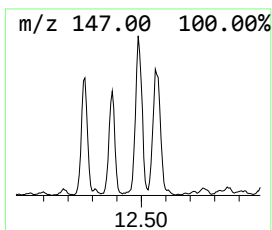
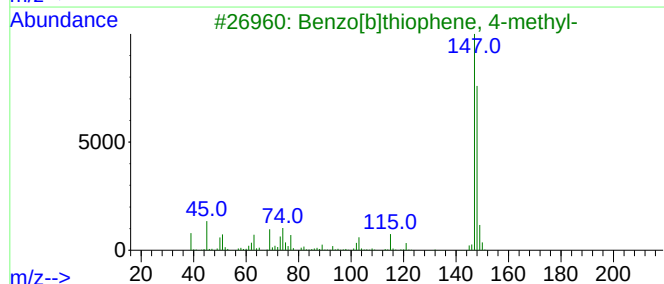
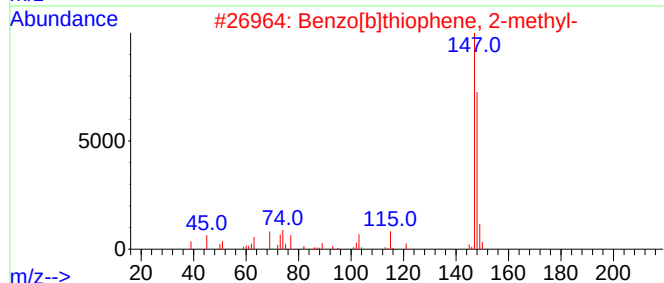
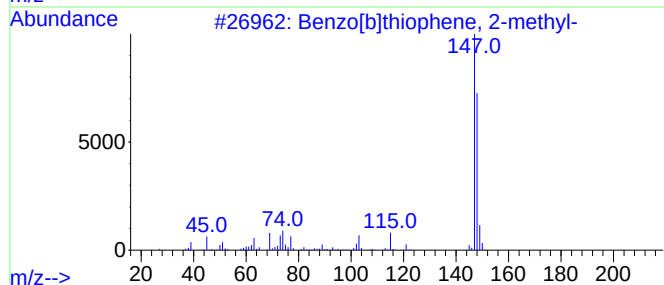
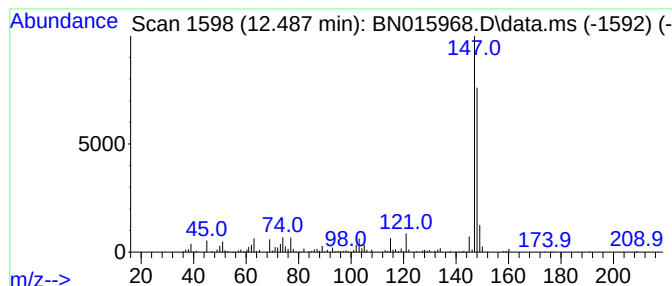
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Benzo[b]thiophene, 2-methyl- Concentration Rank 52

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.487	4.59 ng/ul	719176	Naphthalene-d8	10.710

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[b]thiophene, 2-methyl-	148	C9H8S	001195-14-8	91
2			Benzo[b]thiophene, 2-methyl-	148	C9H8S	001195-14-8	91
3			Benzo[b]thiophene, 4-methyl-	148	C9H8S	014315-11-8	91
4			3-Methylbenzothiophene	148	C9H8S	001455-18-1	91
5			Benzo[b]thiophene, 5-methyl-	148	C9H8S	014315-14-1	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
Data File : BN015968.D
Acq On : 19 Aug 2021 11:49
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Misc :
ALS Vial : 75 Sample Multiplier: 1

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ClientSampleId :
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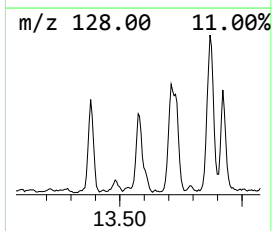
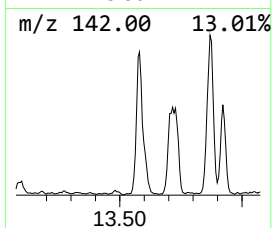
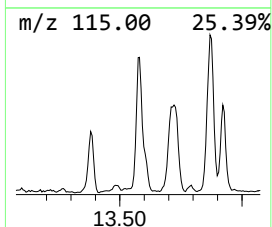
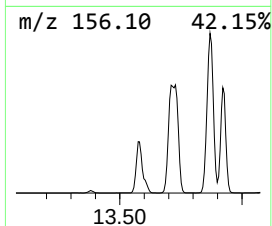
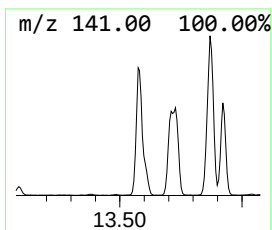
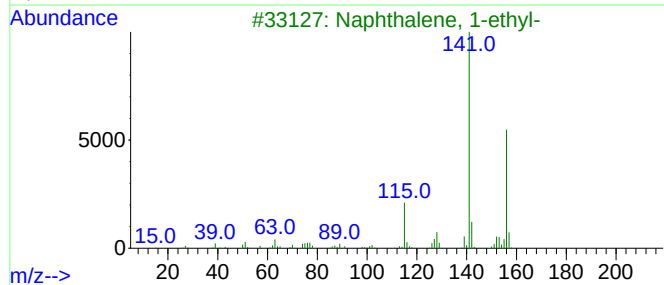
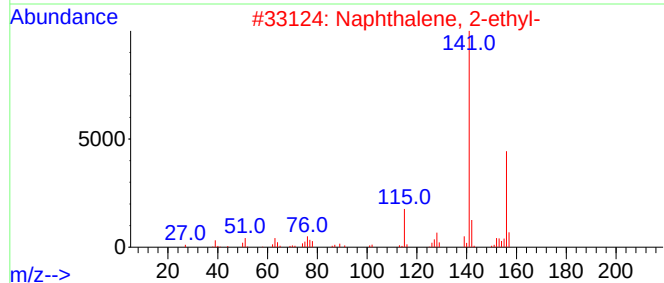
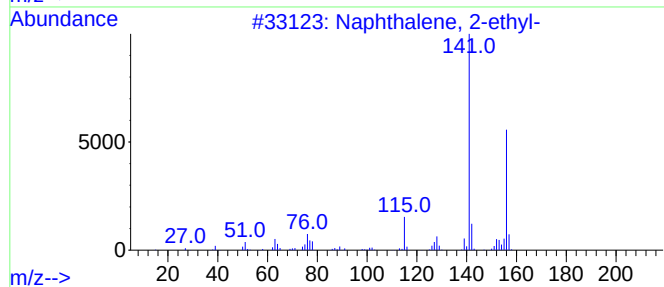
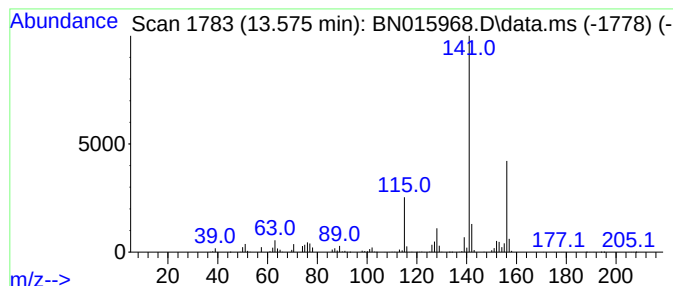
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Naphthalene, 2-ethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.575	27.83 ng/ul	4085880	Acenaphthene-d10	14.551

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	96
2		Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	95
3		Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	95
4		Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	94
5		Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
Data File : BN015968.D
Acq On : 19 Aug 2021 11:49
Operator : CG/JU
Sample : M3399-05 30X
Misc :
ALS Vial : 75 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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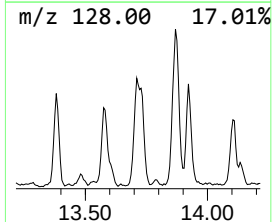
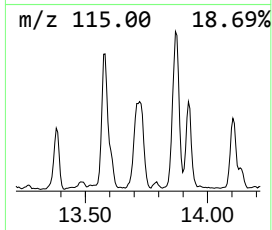
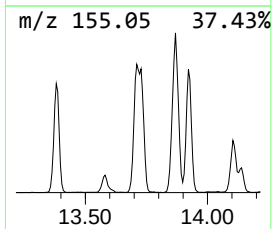
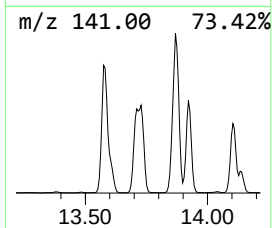
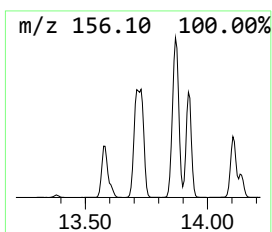
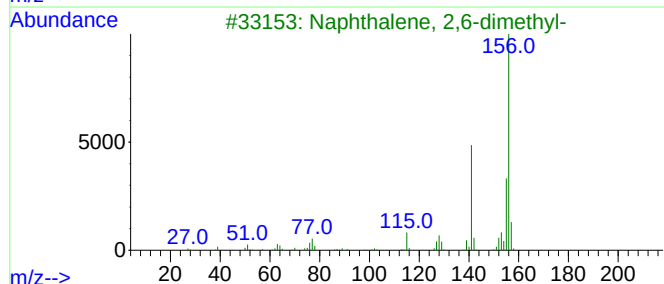
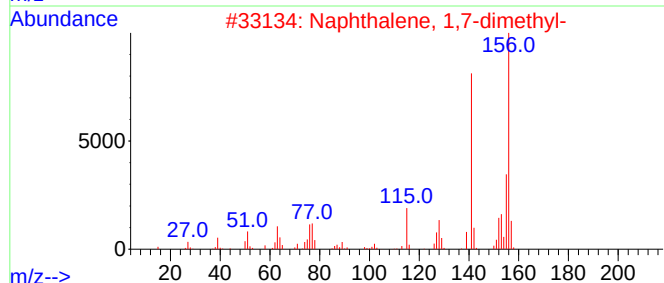
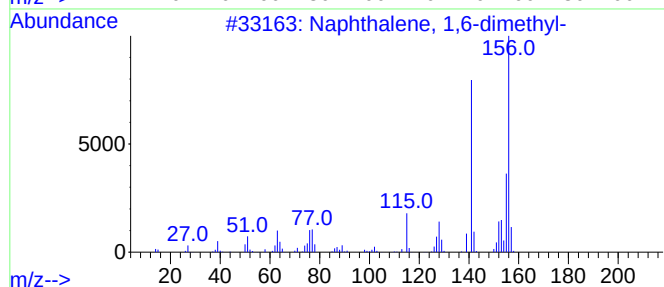
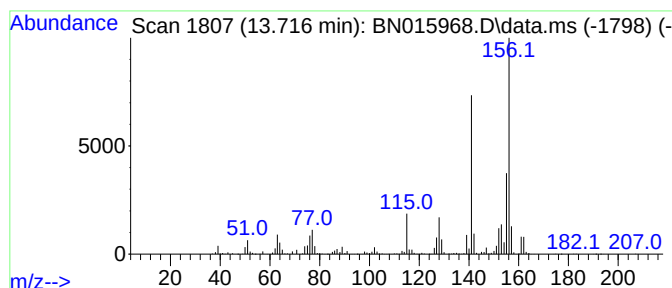
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Naphthalene, 1,6-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.716	31.01 ng/ul	4553840	Acenaphthene-d10	14.551

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	96
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\BN081721\
Data File : BN015968.D
Acq On : 19 Aug 2021 11:49
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Sample : M3399-05 30X
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ALS Vial : 75 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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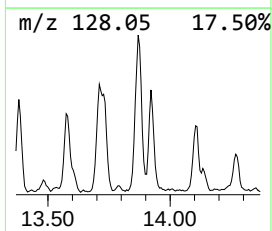
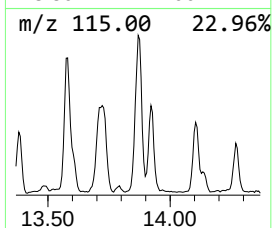
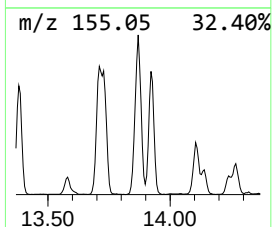
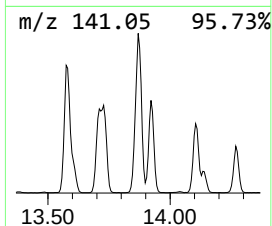
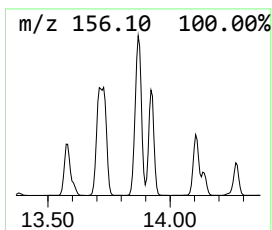
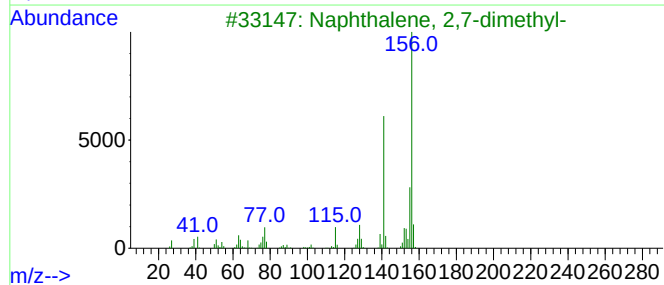
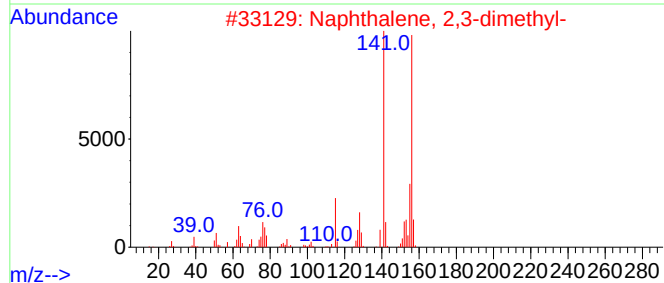
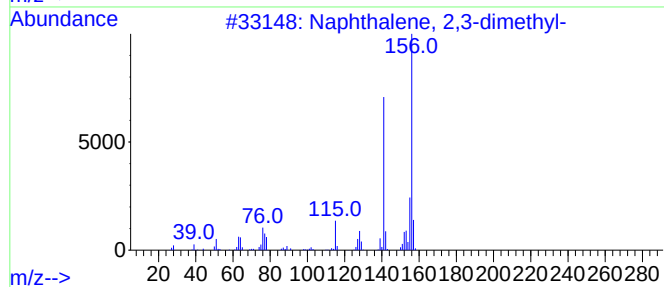
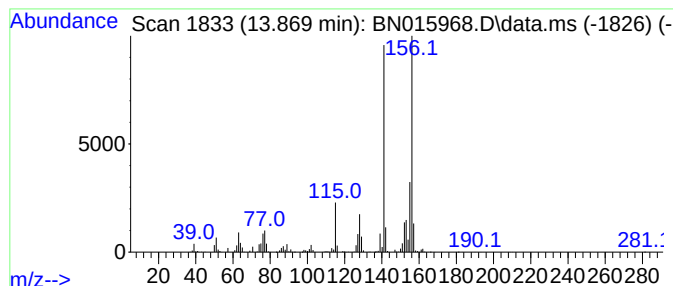
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.869	49.38 ng/ul	7250510	Acenaphthene-d10	14.551

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
Data File : BN015968.D
Acq On : 19 Aug 2021 11:49
Operator : CG/JU
Sample : M3399-05 30X
Misc :
ALS Vial : 75 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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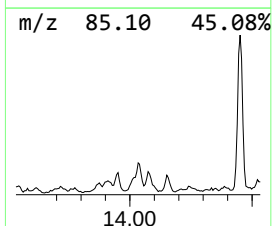
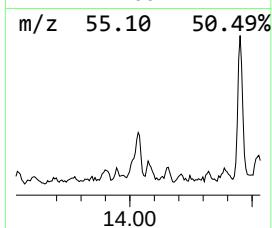
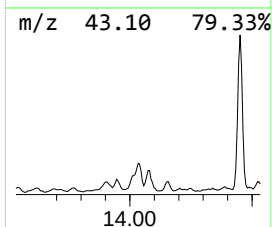
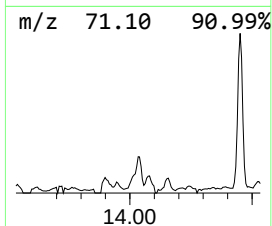
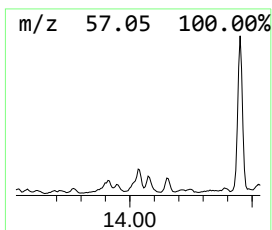
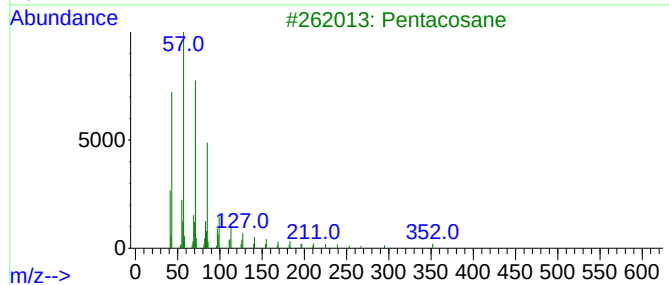
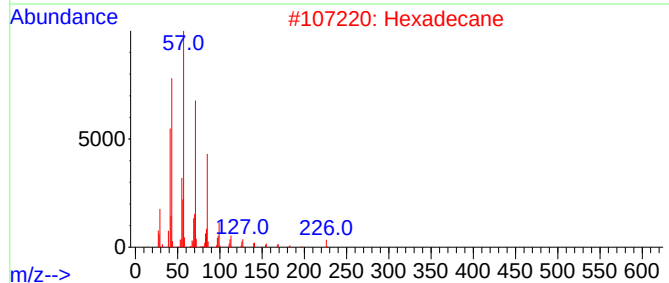
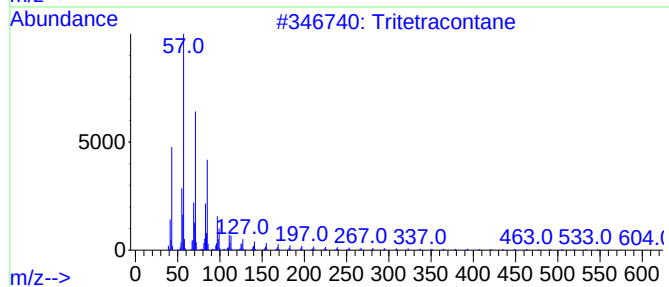
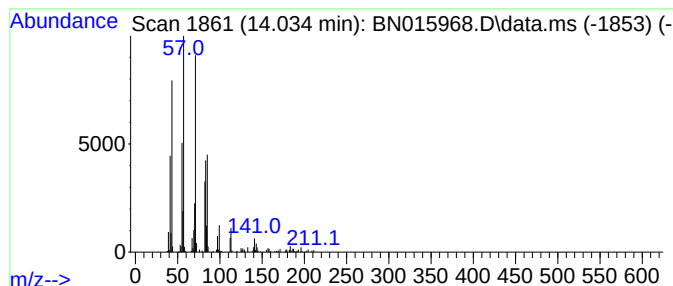
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 47

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.034	5.97 ng/ul	876645	Acenaphthene-d10	14.551

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tritetracontane	605	C43H88	007098-21-7	64
2		Hexadecane	226	C16H34	000544-76-3	59
3		Pentacosane	352	C25H52	000629-99-2	59
4		3-Ethyl-2,6,10-trimethylundecane	226	C16H34	1000432-25-9	58
5		Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
Data File : BN015968.D
Acq On : 19 Aug 2021 11:49
Operator : CG/JU
Sample : M3399-05 30X
Misc :
ALS Vial : 75 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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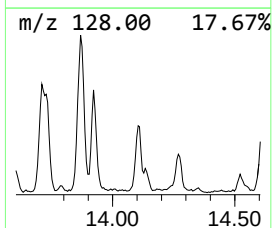
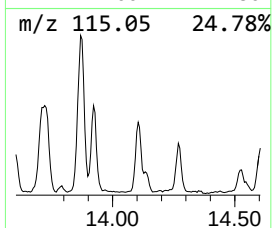
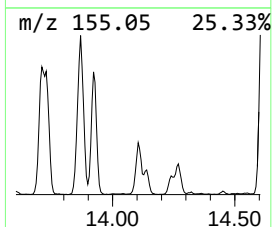
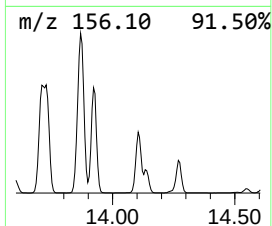
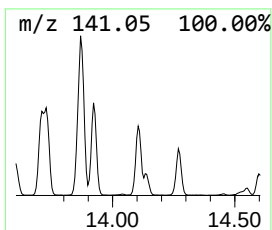
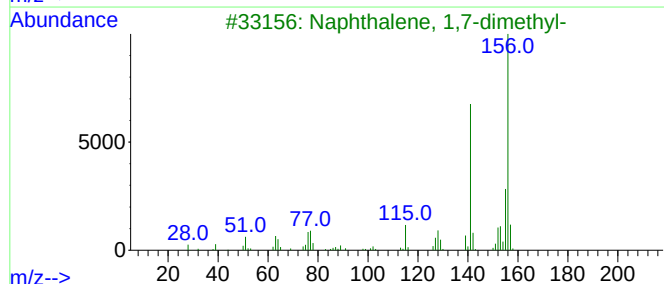
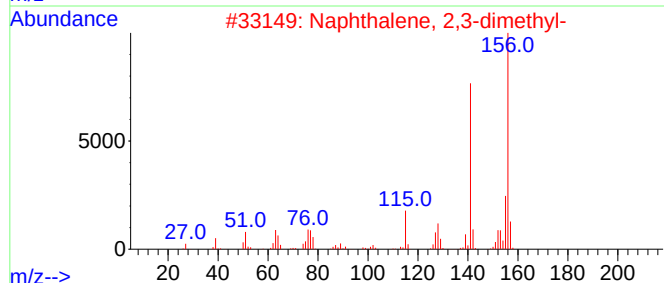
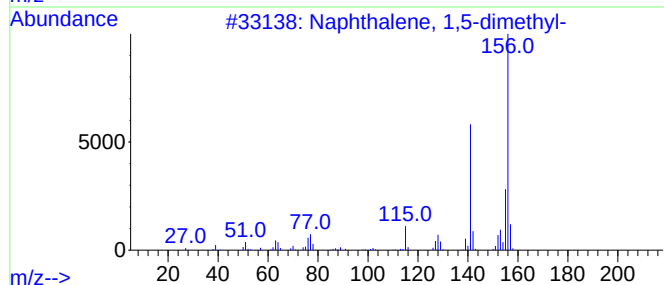
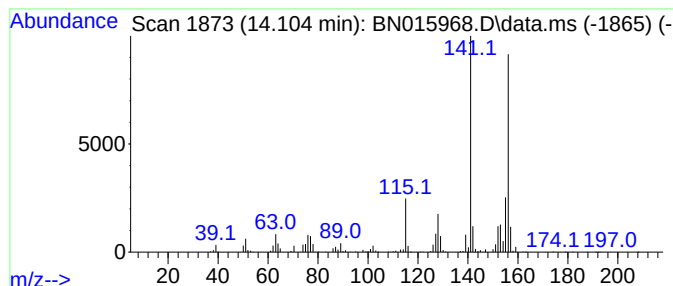
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 Naphthalene, 1,5-dimethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.104	20.33 ng/ul	2985510	Acenaphthene-d10	14.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97
2			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
3			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
4			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97
5			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
Data File : BN015968.D
Acq On : 19 Aug 2021 11:49
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Sample : M3399-05 30X
Misc :
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Instrument :
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ClientSampleId :
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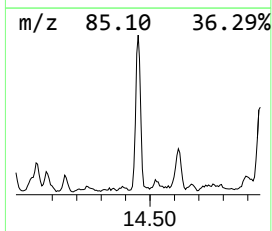
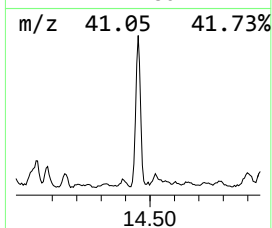
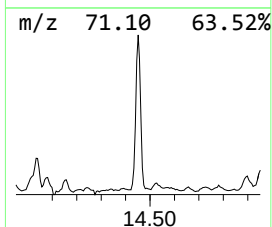
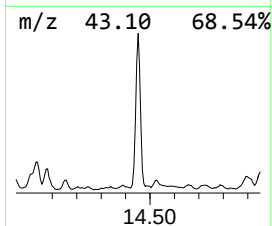
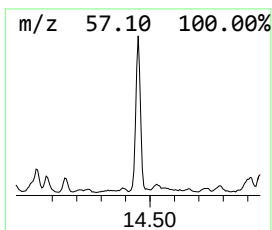
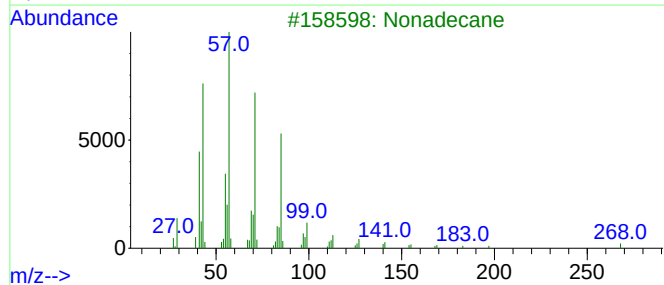
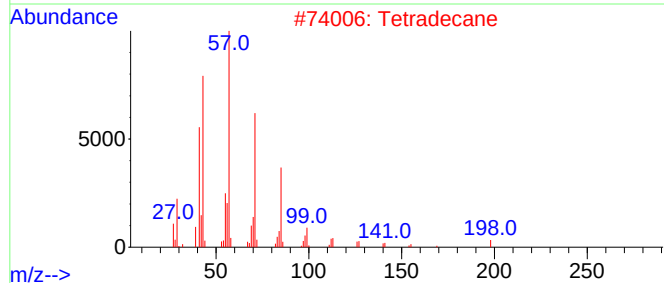
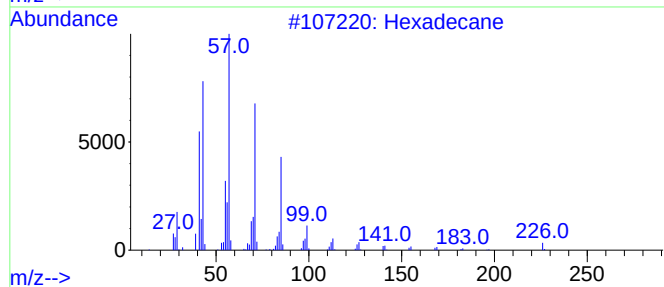
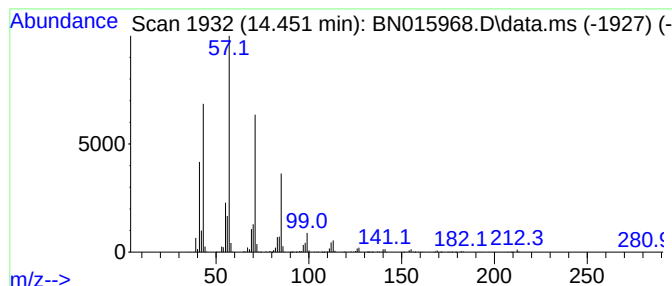
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.451	18.28 ng/ul	2684600	Acenaphthene-d10	14.551

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexadecane	226	C16H34	000544-76-3	91
2		Tetradecane	198	C14H30	000629-59-4	91
3		Nonadecane	268	C19H40	000629-92-5	90
4		Hexadecane	226	C16H34	000544-76-3	90
5		Hexadecane	226	C16H34	000544-76-3	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
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Acq On : 19 Aug 2021 11:49
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Sample : M3399-05 30X
Misc :
ALS Vial : 75 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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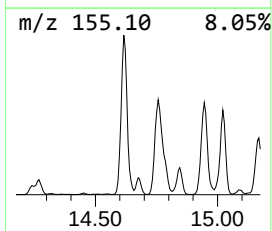
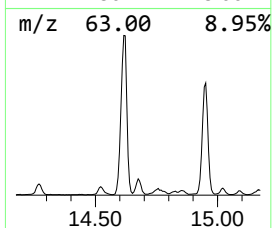
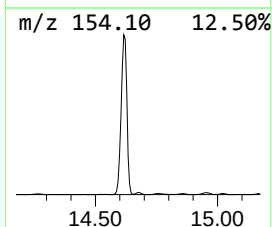
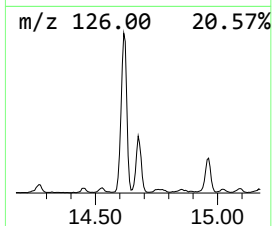
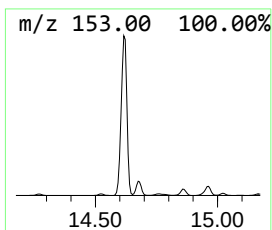
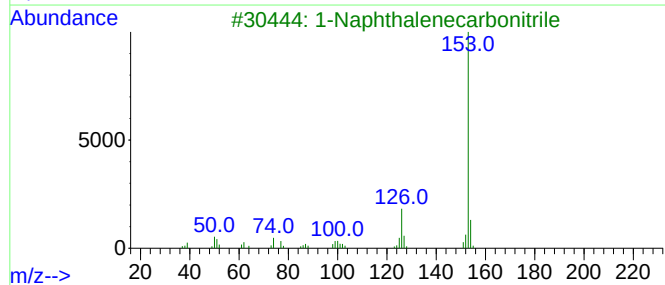
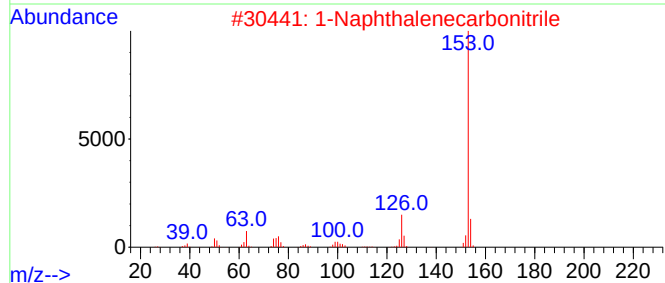
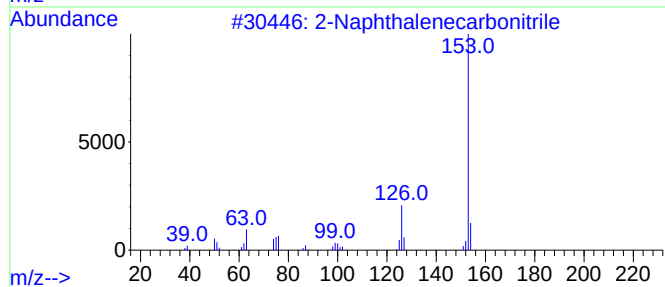
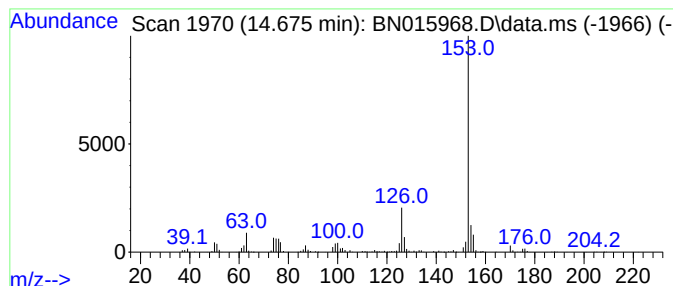
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 2-Naphthalenecarbonitrile Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.675	12.26 ng/ul	1799840	Acenaphthene-d10	14.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Naphthalenecarbonitrile		153	C11H7N	000613-46-7	97	
2	1-Naphthalenecarbonitrile		153	C11H7N	000086-53-3	91	
3	1-Naphthalenecarbonitrile		153	C11H7N	000086-53-3	90	
4	Naphthalene, 1-isocyano-		153	C11H7N	001984-04-9	90	
5	2-Naphthalenecarbonitrile		153	C11H7N	000613-46-7	87	



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
Data File : BN015968.D
Acq On : 19 Aug 2021 11:49
Operator : CG/JU
Sample : M3399-05 30X
Misc :
ALS Vial : 75 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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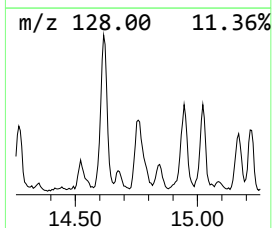
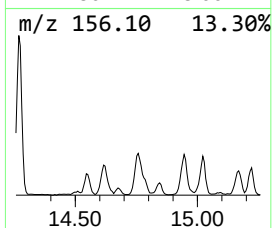
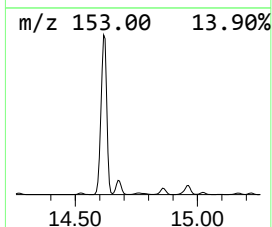
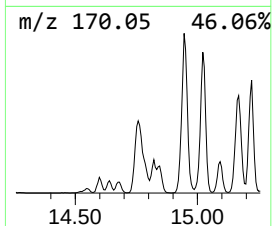
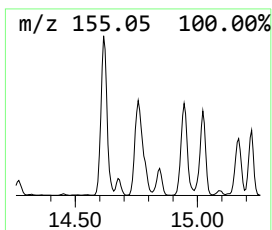
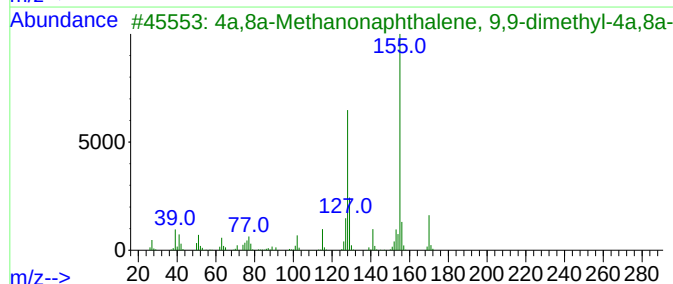
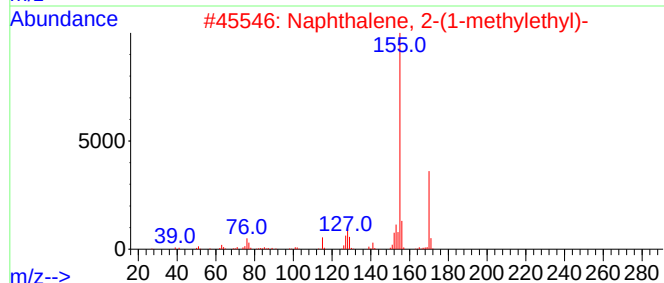
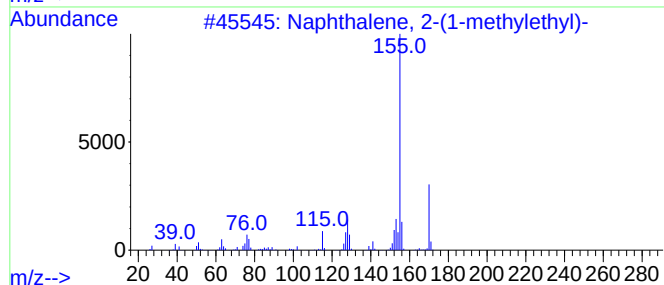
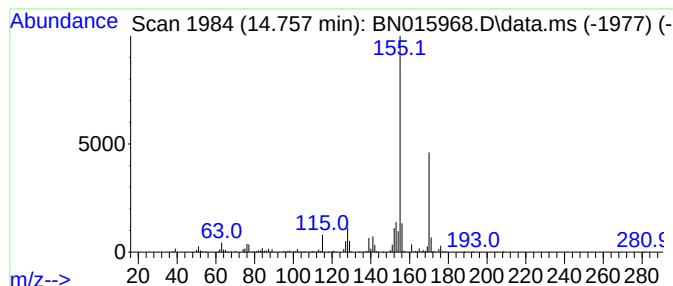
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 Naphthalene, 2-(1-methyleth... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.757	17.33 ng/ul	2544910	Acenaphthene-d10	14.551

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 2-(1-methylethyl)-	170	C13H14	002027-17-0	91
2		Naphthalene, 2-(1-methylethyl)-	170	C13H14	002027-17-0	91
3		4a,8a-Methanonaphthalene, 9,9-di...	170	C13H14	038963-97-2	86
4		Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	76
5		Ethanone, 1-(5-chloro-2-hydroxyp...	170	C8H7ClO2	001450-74-4	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
Data File : BN015968.D
Acq On : 19 Aug 2021 11:49
Operator : CG/JU
Sample : M3399-05 30X
Misc :
ALS Vial : 75 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBKA7

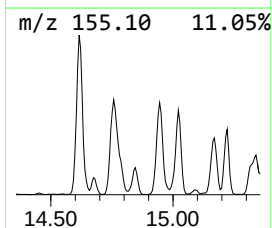
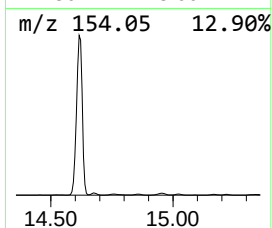
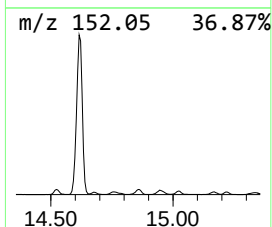
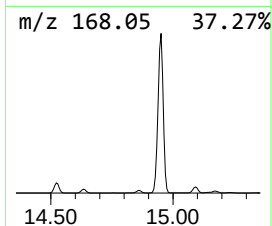
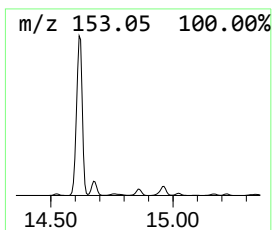
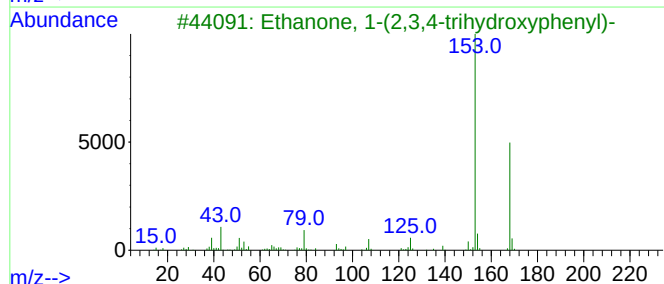
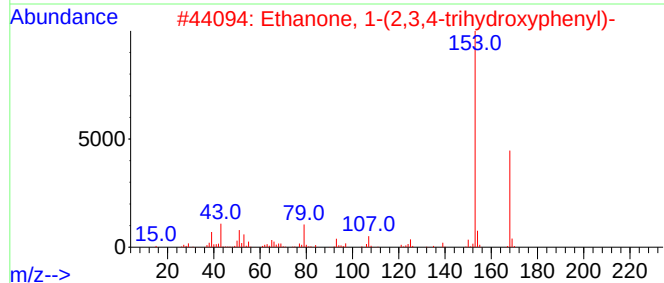
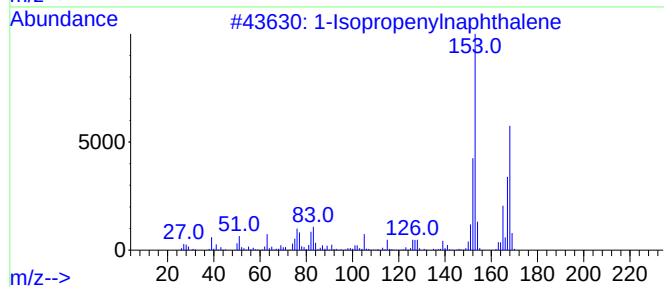
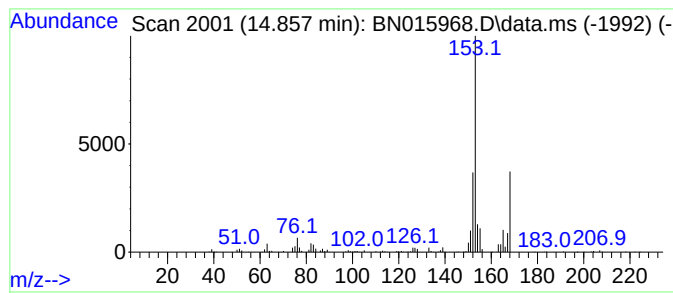
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 1-Isopropenylnaphthalene Concentration Rank 35

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.857	8.63 ng/ul	1267190	Acenaphthene-d10	14.551

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Isopropenylnaphthalene	168	C13H12	001855-47-6	59
2		Ethanone, 1-(2,3,4-trihydroxyphe...	168	C8H8O4	000528-21-2	52
3		Ethanone, 1-(2,3,4-trihydroxyphe...	168	C8H8O4	000528-21-2	52
4		Ethanone, 1-(2,4,6-trihydroxyphe...	168	C8H8O4	000480-66-0	50
5		Ethanone, 1-(2,4,6-trihydroxyphe...	168	C8H8O4	000480-66-0	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
Data File : BN015968.D
Acq On : 19 Aug 2021 11:49
Operator : CG/JU
Sample : M3399-05 30X
Misc :
ALS Vial : 75 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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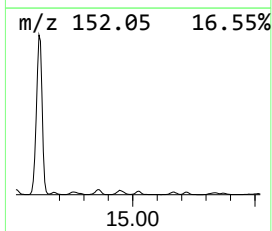
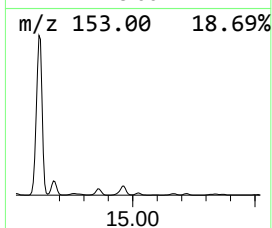
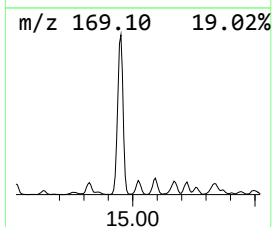
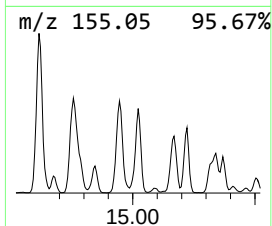
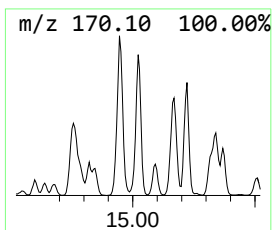
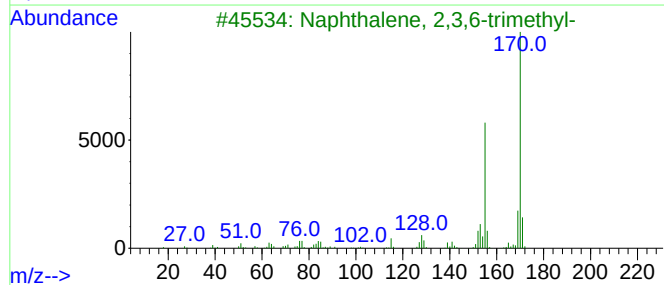
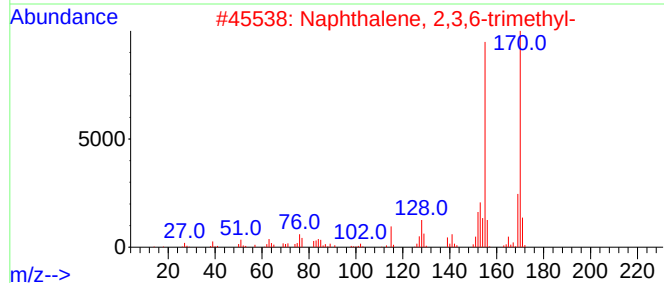
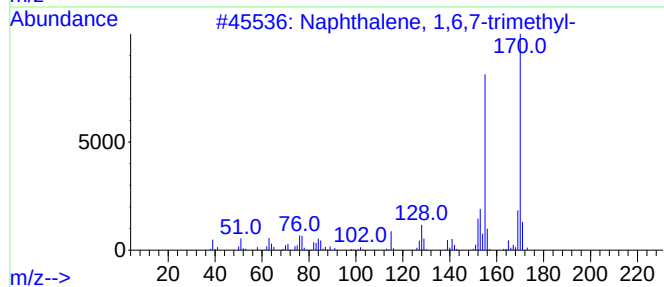
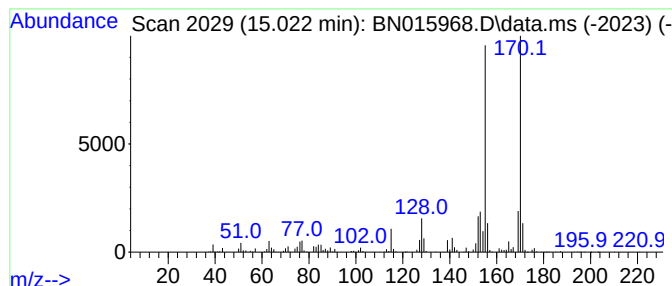
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 Naphthalene, 1,6,7-trimethyl- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.022	14.33 ng/ul	2104630	Acenaphthene-d10	14.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	97
2			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	97
3			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	97
4			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	96
5			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
Data File : BN015968.D
Acq On : 19 Aug 2021 11:49
Operator : CG/JU
Sample : M3399-05 30X
Misc :
ALS Vial : 75 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBKA7

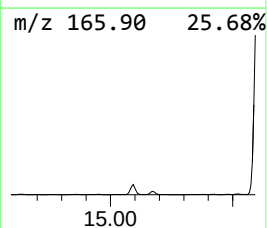
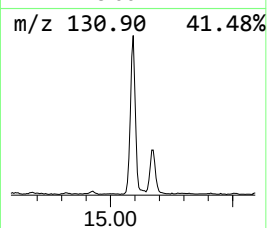
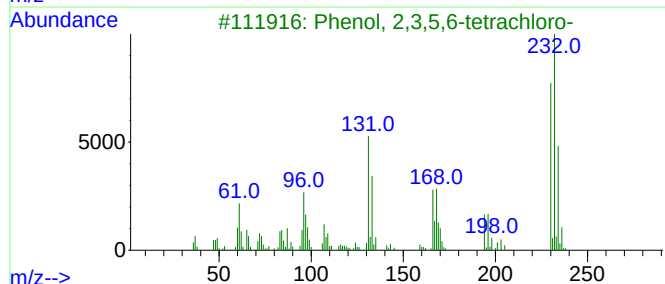
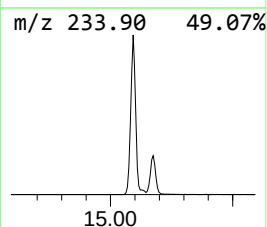
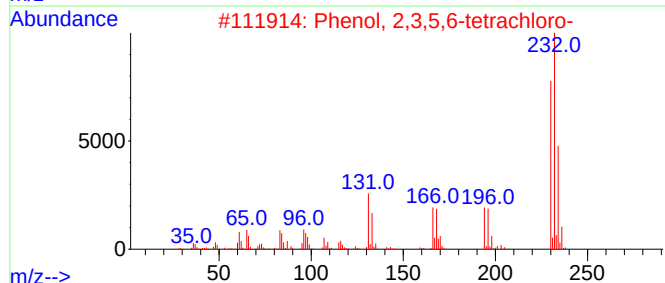
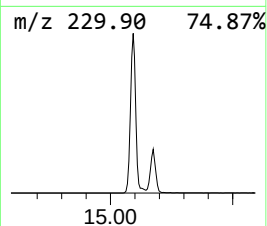
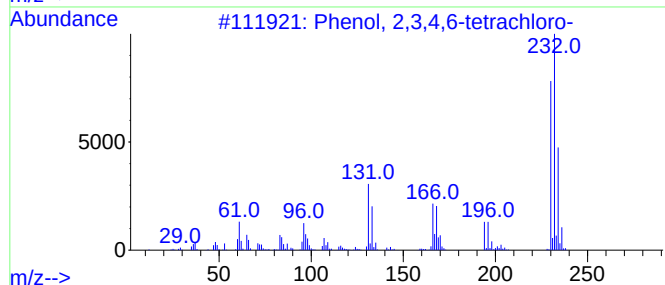
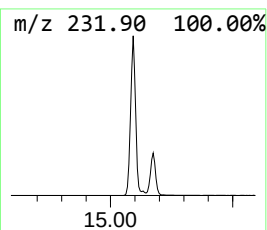
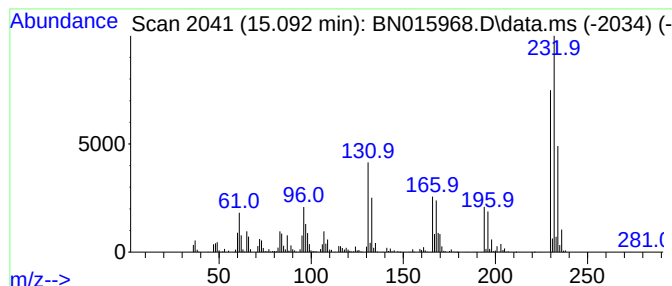
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 Phenol, 2,3,4,6-tetrachloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.092	55.44 ng/ul	8139720	Acenaphthene-d10	14.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 2,3,4,6-tetrachloro-	230	C6H2Cl4O	000058-90-2	99
2			Phenol, 2,3,5,6-tetrachloro-	230	C6H2Cl4O	000935-95-5	99
3			Phenol, 2,3,5,6-tetrachloro-	230	C6H2Cl4O	000935-95-5	99
4			Phenol, 2,3,4,6-tetrachloro-	230	C6H2Cl4O	000058-90-2	99
5			Phenol, 2,3,4,5-tetrachloro-	230	C6H2Cl4O	004901-51-3	98



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
Data File : BN015968.D
Acq On : 19 Aug 2021 11:49
Operator : CG/JU
Sample : M3399-05 30X
Misc :
ALS Vial : 75 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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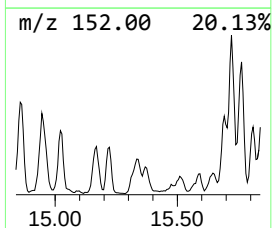
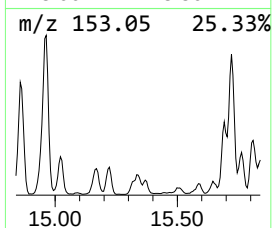
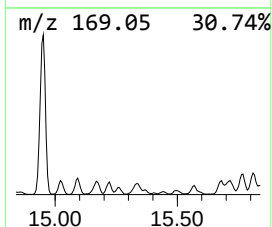
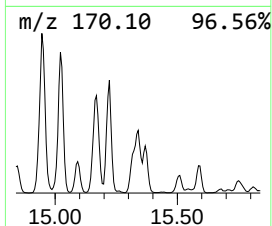
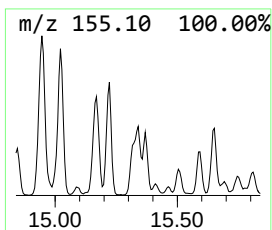
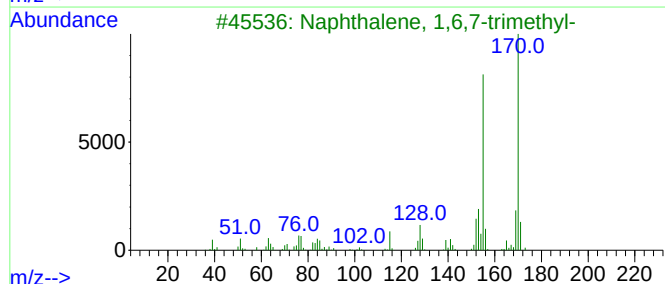
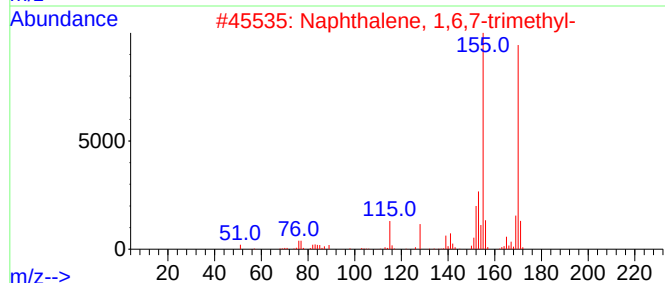
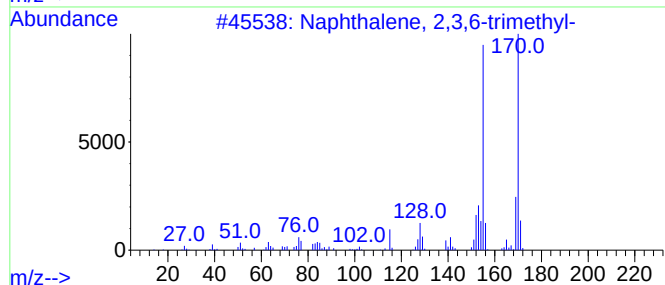
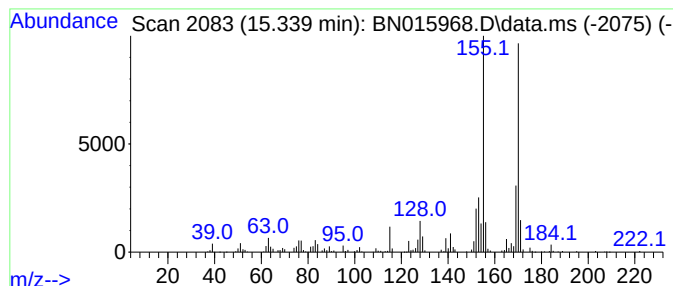
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 Naphthalene, 2,3,6-trimethyl- Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.339	9.92 ng/ul	1455980	Acenaphthene-d10	14.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	98
2			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	98
3			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	97
4			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	97
5			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
Data File : BN015968.D
Acq On : 19 Aug 2021 11:49
Operator : CG/JU
Sample : M3399-05 30X
Misc :
ALS Vial : 75 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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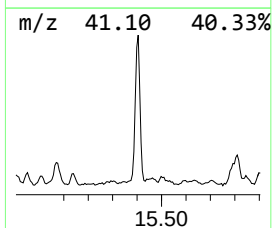
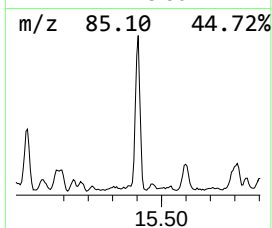
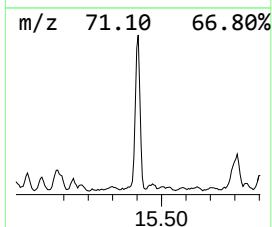
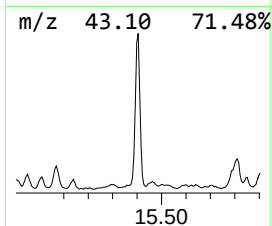
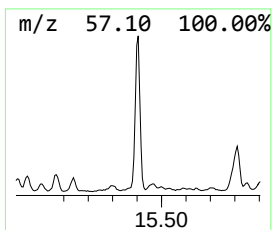
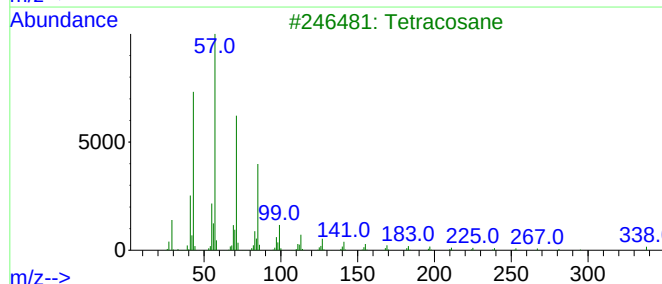
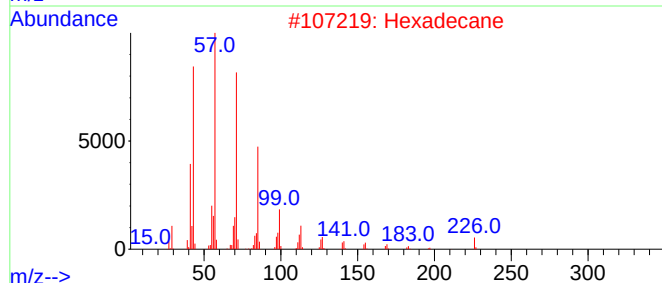
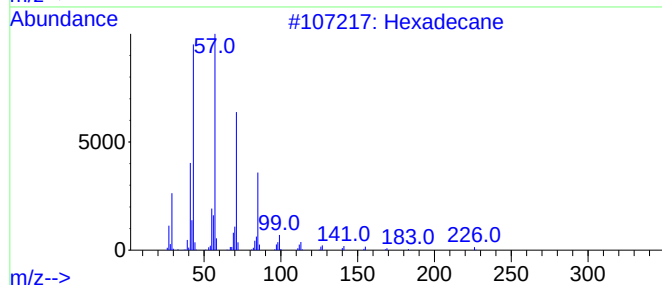
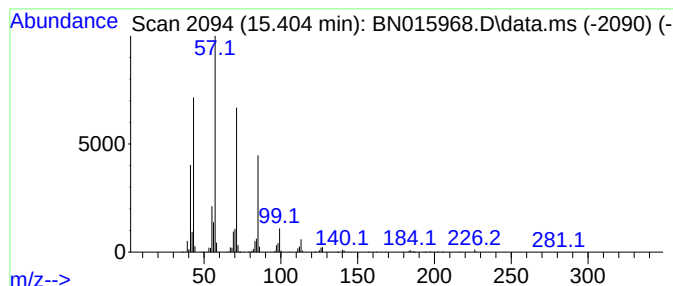
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Quant Title : SVOA CALIBRATION

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

Peak Number 23 (DEL) Alkane: Straight-Chai... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.404	19.59 ng/ul	2876660	Acenaphthene-d10	14.551

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexadecane	226	C16H34	000544-76-3	93
2		Hexadecane	226	C16H34	000544-76-3	91
3		Tetracosane	338	C24H50	000646-31-1	90
4		Nonane, 5-butyl-	184	C13H28	017312-63-9	89
5		Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
Data File : BN015968.D
Acq On : 19 Aug 2021 11:49
Operator : CG/JU
Sample : M3399-05 30X
Misc :
ALS Vial : 75 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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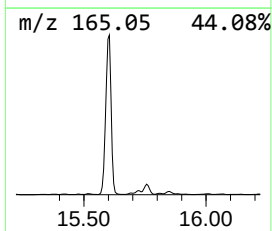
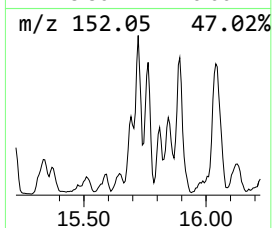
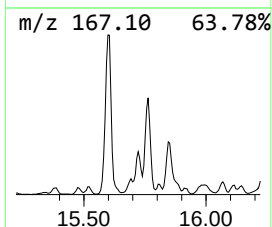
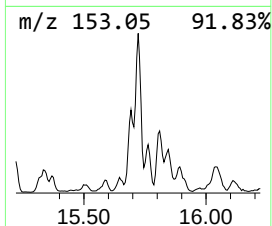
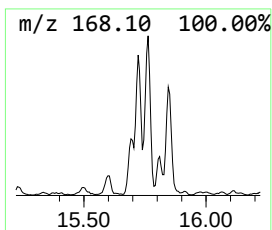
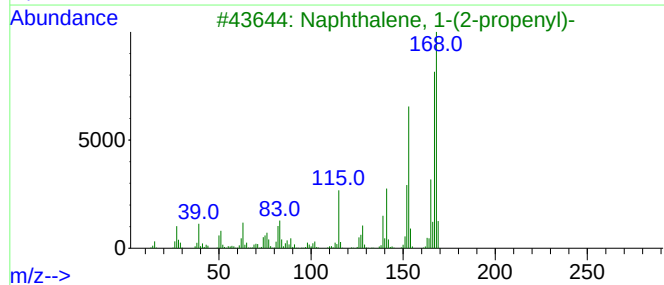
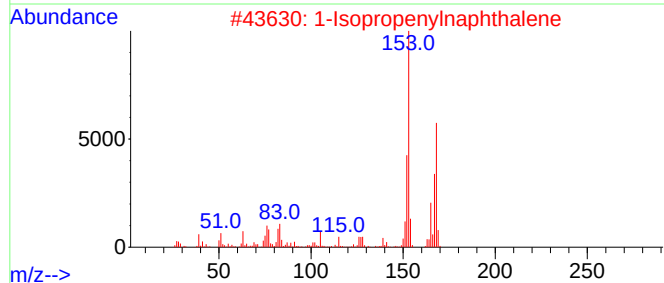
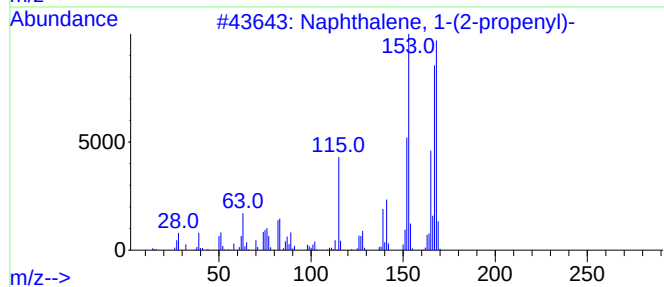
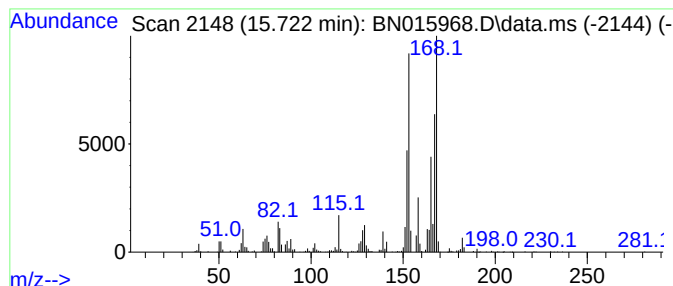
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Quant Title : SVOA CALIBRATION

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

Peak Number 24 Naphthalene, 1-(2-propenyl)- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.722	14.06 ng/ul	2064360	Acenaphthene-d10	14.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1-(2-propenyl)-	168	C13H12	002489-86-3	74
2			1-Isopropenylnaphthalene	168	C13H12	001855-47-6	64
3			Naphthalene, 1-(2-propenyl)-	168	C13H12	002489-86-3	58
4			Benzene, [1-(2,4-cyclopentadien-...	168	C13H12	002320-32-3	53
5			1,1'-Biphenyl, 2-methyl-	168	C13H12	000643-58-3	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
Data File : BN015968.D
Acq On : 19 Aug 2021 11:49
Operator : CG/JU
Sample : M3399-05 30X
Misc :
ALS Vial : 75 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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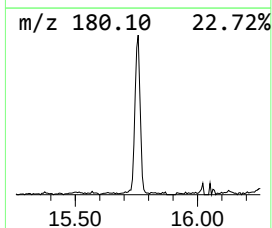
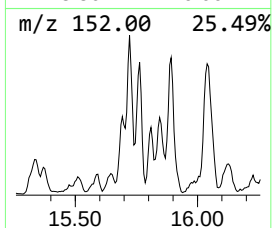
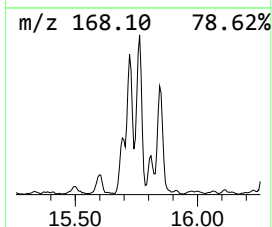
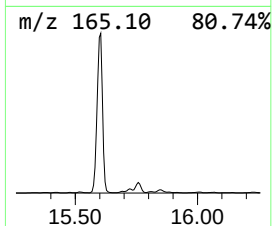
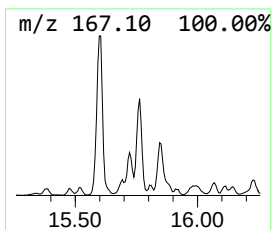
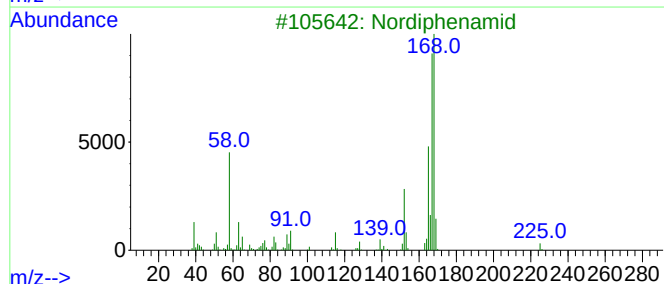
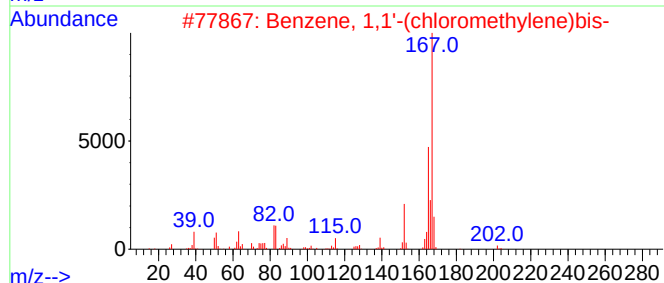
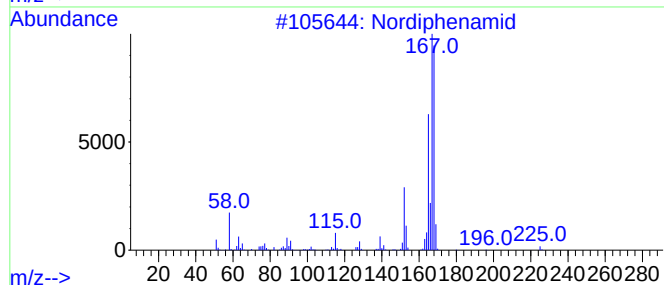
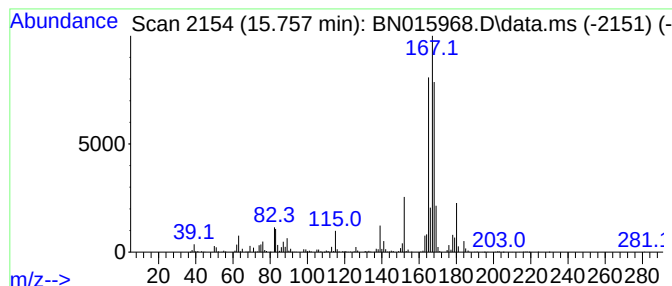
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Quant Title : SVOA CALIBRATION

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

Peak Number 25 Nordiphenamid Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.757	17.34 ng/ul	2546060	Acenaphthene-d10	14.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nordiphenamid	225	C15H15NO	000954-21-2	59
2			Benzene, 1,1'-(chloromethylene)bis-	202	C13H11Cl	000090-99-3	50
3			Nordiphenamid	225	C15H15NO	000954-21-2	50
4			L-Cysteine, S-(diphenylmethyl)-	287	C16H17NO2S	005191-80-0	50
5			Diphenylmethane	168	C13H12	000101-81-5	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
Data File : BN015968.D
Acq On : 19 Aug 2021 11:49
Operator : CG/JU
Sample : M3399-05 30X
Misc :
ALS Vial : 75 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBKA7

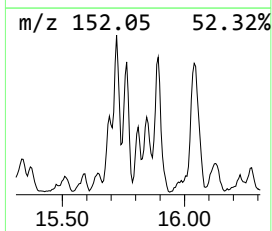
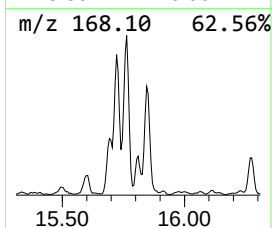
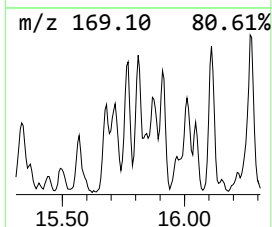
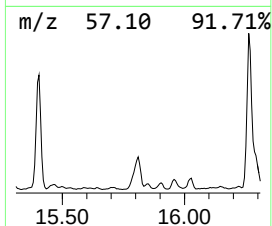
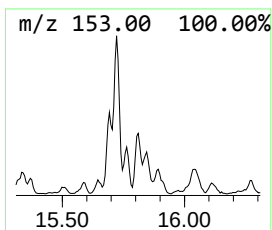
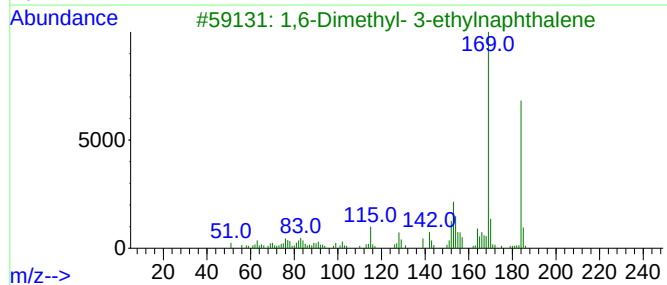
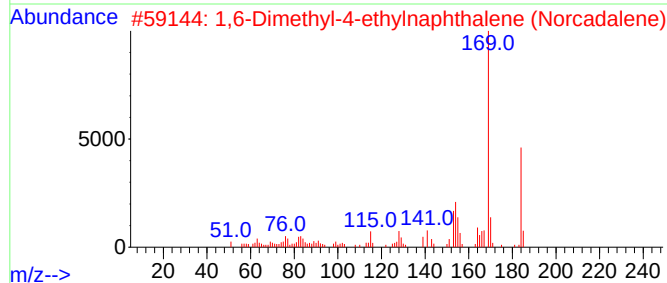
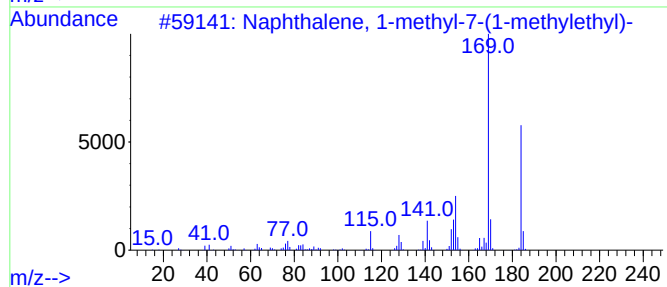
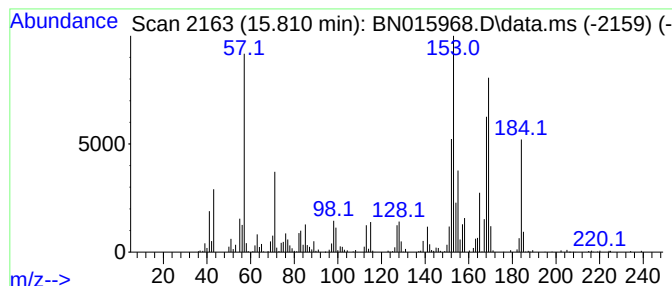
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Quant Title : SVOA CALIBRATION

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

Peak Number 26 Naphthalene, 1-methyl-7-(1-... Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.810	10.42 ng/ul	1529640	Acenaphthene-d10	14.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1-methyl-7-(1-methy...	184	C14H16	000490-65-3	50
2			1,6-Dimethyl-4-ethylnaphthalene ...	184	C14H16	1000383-71-7	50
3			1,6-Dimethyl- 3-ethylnaphthalene	184	C14H16	1000383-71-8	49
4			Naphthalene, 1,2,3,4-tetramethyl-	184	C14H16	003031-15-0	46
5			2-Isopropyl-3-methylnaphthalene	184	C14H16	1000383-71-2	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
Data File : BN015968.D
Acq On : 19 Aug 2021 11:49
Operator : CG/JU
Sample : M3399-05 30X
Misc :
ALS Vial : 75 Sample Multiplier: 1

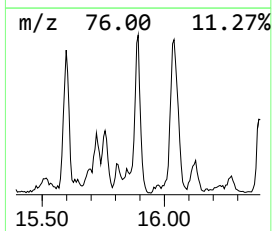
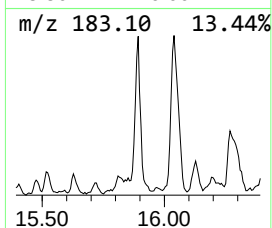
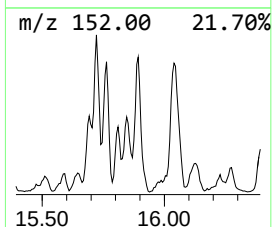
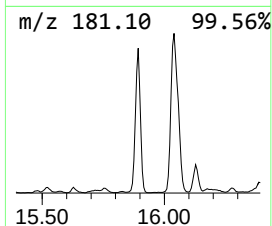
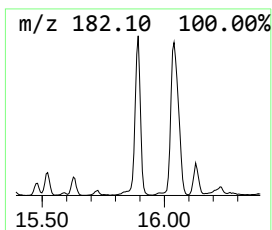
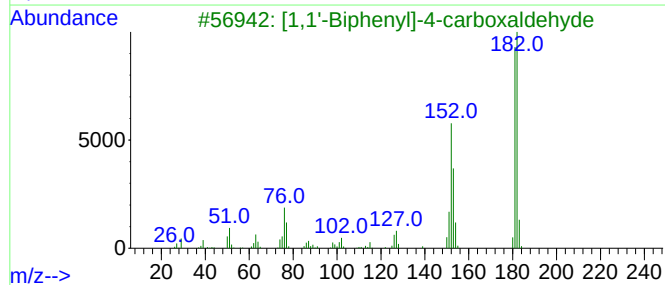
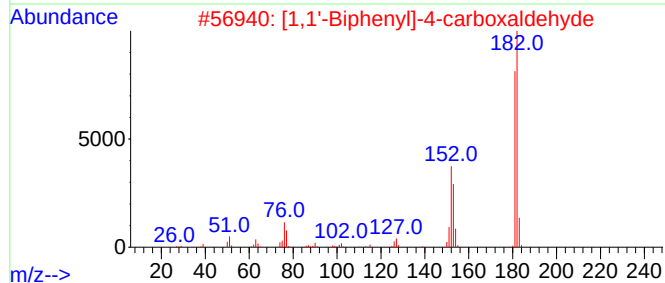
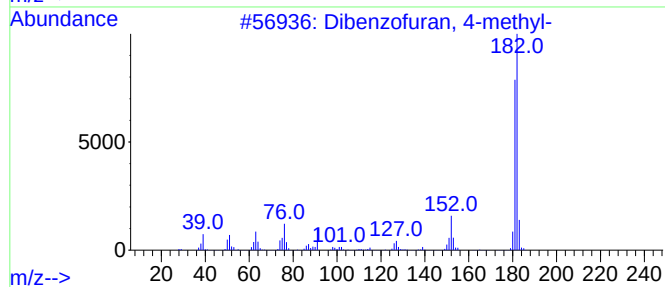
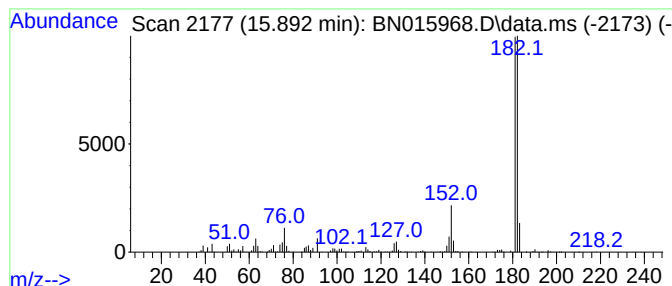
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081621MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 28 Dibenzofuran, 4-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.	
15.892	22.25 ng/ul	3267090	Acenaphthene-d10		14.551	
Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Dibenzofuran, 4-methyl-	182	C13H10O	007320-53-8	91
2		[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	83
3		[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	78
4		[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	78
5		9H-Fluoren-9-ol	182	C13H10O	001689-64-1	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
Data File : BN015968.D
Acq On : 19 Aug 2021 11:49
Operator : CG/JU
Sample : M3399-05 30X
Misc :
ALS Vial : 75 Sample Multiplier: 1

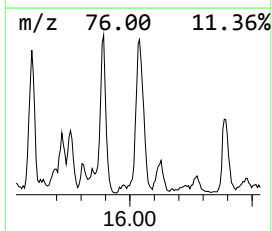
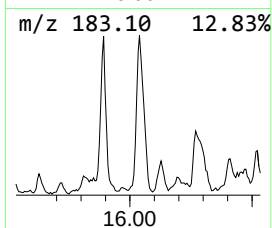
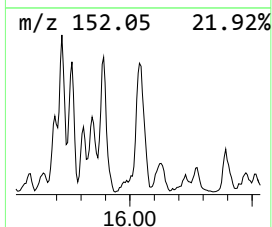
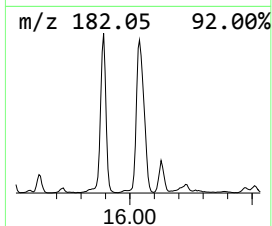
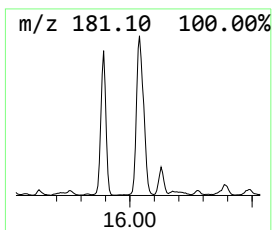
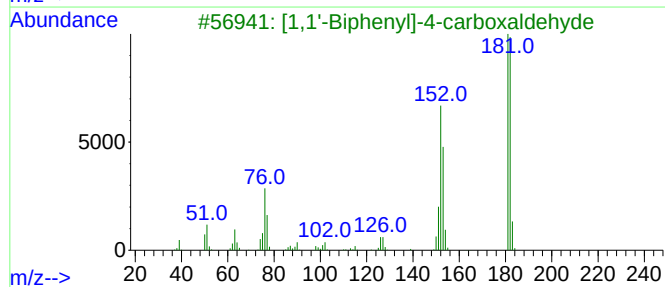
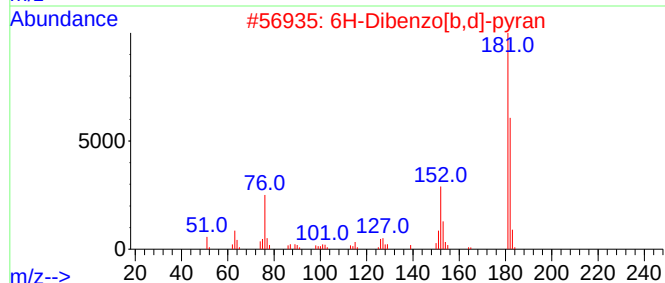
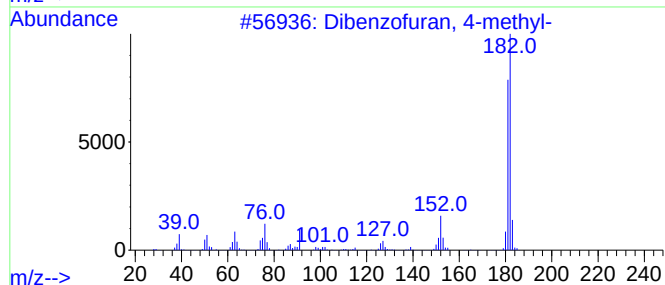
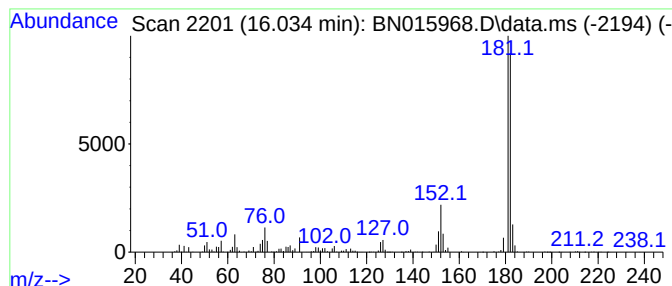
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 30 6H-Dibenzo[b,d]-pyran Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.034	25.00 ng/ul	4319200	Phenanthrene-d10	17.304	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dibenzofuran, 4-methyl-	182	C13H10O	007320-53-8	91
2	6H-Dibenzo[b,d]-pyran	182	C13H10O	000229-95-8	91
3	[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	91
4	[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	87
5	9H-Fluoren-9-ol	182	C13H10O	001689-64-1	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
Data File : BN015968.D
Acq On : 19 Aug 2021 11:49
Operator : CG/JU
Sample : M3399-05 30X
Misc :
ALS Vial : 75 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBKA7

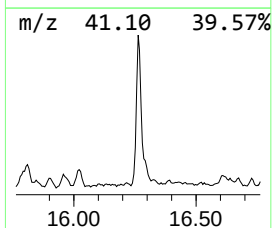
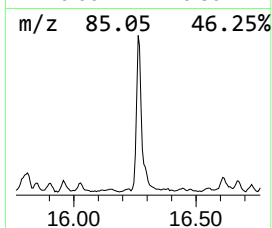
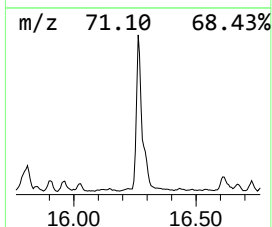
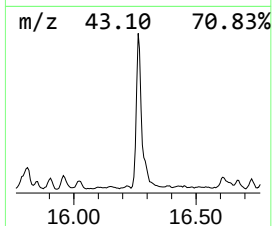
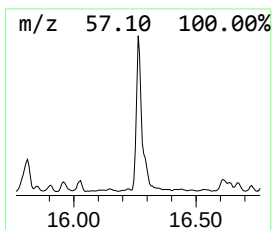
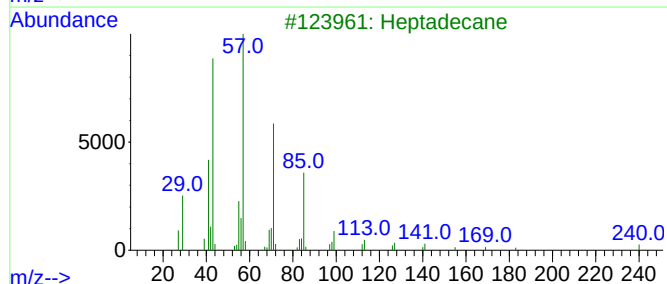
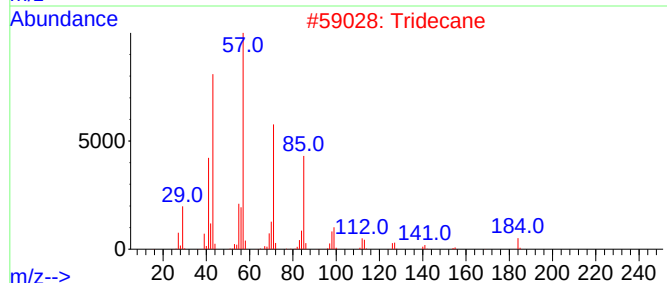
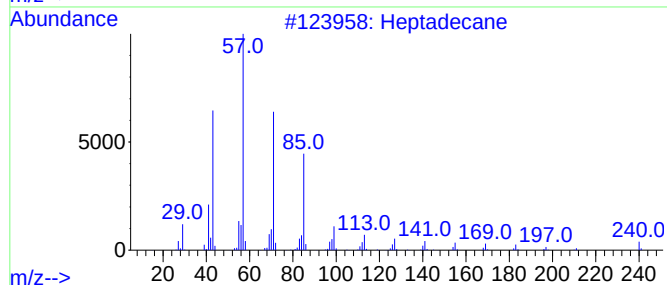
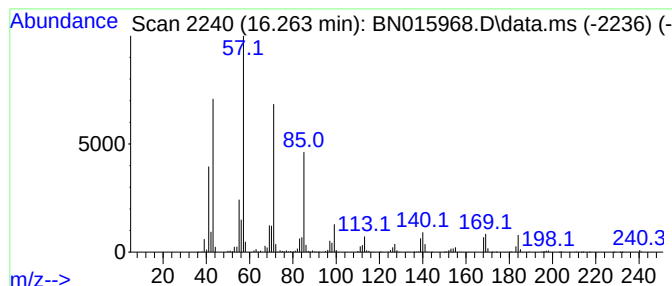
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Quant Title : SVOA CALIBRATION

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

Peak Number 32 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.263	35.19 ng/ul	6081140	Phenanthrene-d10	17.304

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane	240	C17H36	000629-78-7	94
2		Tridecane	184	C13H28	000629-50-5	90
3		Heptadecane	240	C17H36	000629-78-7	81
4		Pentadecane	212	C15H32	000629-62-9	80
5		Tetradecane	198	C14H30	000629-59-4	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
Data File : BN015968.D
Acq On : 19 Aug 2021 11:49
Operator : CG/JU
Sample : M3399-05 30X
Misc :
ALS Vial : 75 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBKA7

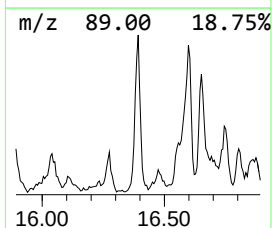
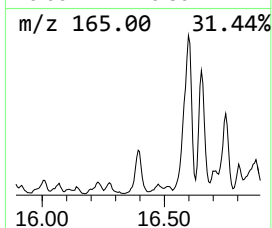
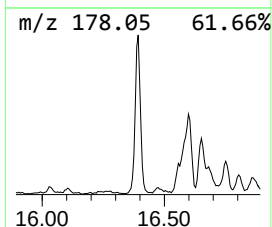
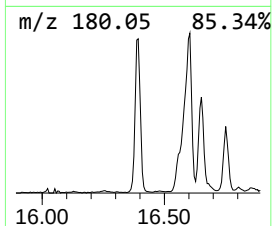
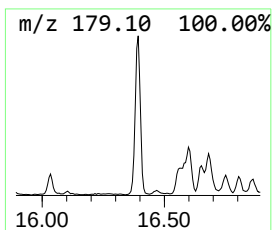
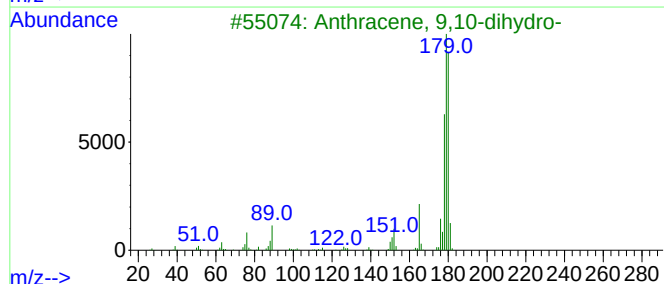
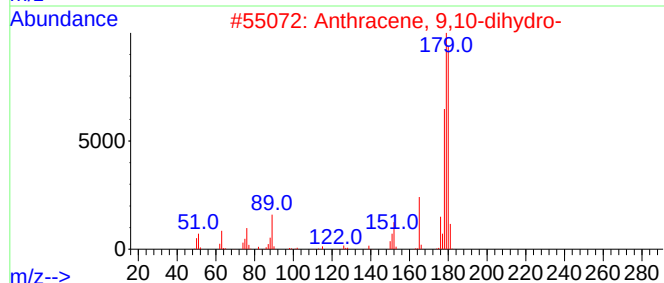
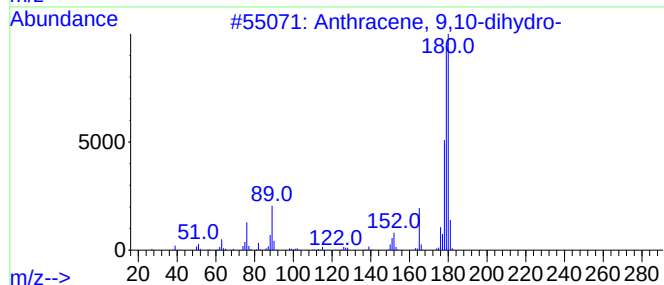
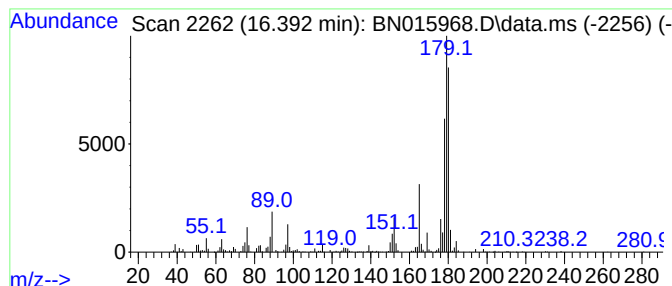
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Quant Title : SVOA CALIBRATION

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

Peak Number 33 Anthracene, 9,10-dihydro- Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.392	12.20 ng/ul	2108560	Phenanthrene-d10	17.304

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 9,10-dihydro-	180	C14H12	000613-31-0	96
2			Anthracene, 9,10-dihydro-	180	C14H12	000613-31-0	96
3			Anthracene, 9,10-dihydro-	180	C14H12	000613-31-0	93
4			Phenanthrene, 9,10-dihydro-	180	C14H12	000776-35-2	91
5			Phenanthrene, 9,10-dihydro-	180	C14H12	000776-35-2	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
Data File : BN015968.D
Acq On : 19 Aug 2021 11:49
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Sample : M3399-05 30X
Misc :
ALS Vial : 75 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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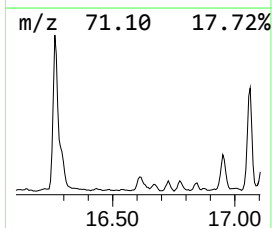
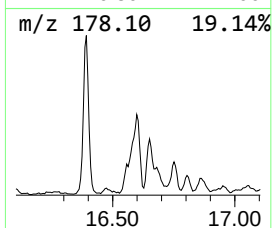
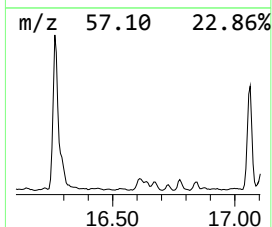
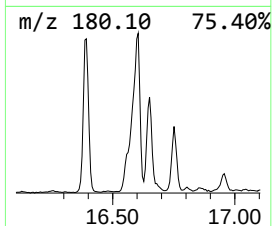
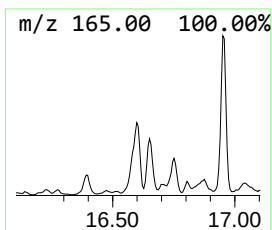
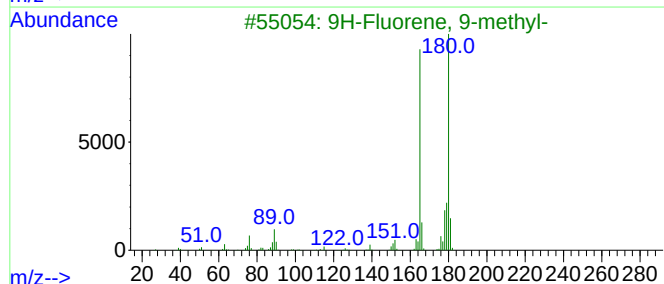
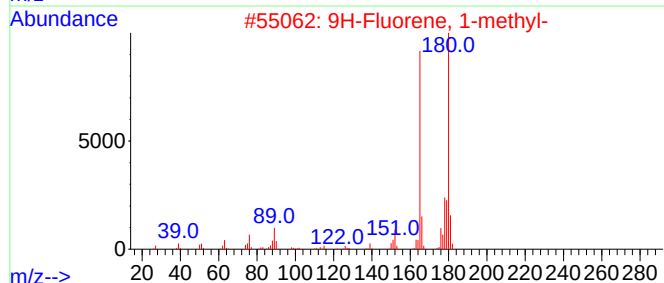
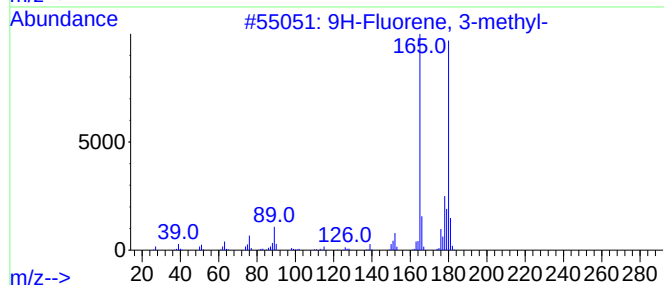
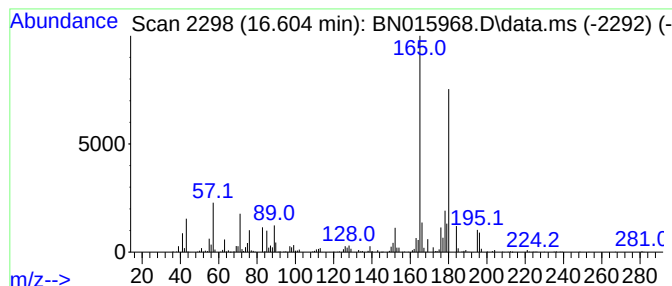
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Quant Title : SVOA CALIBRATION

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

Peak Number 34 9H-Fluorene, 3-methyl- Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.604	12.49 ng/ul	2157650	Phenanthrene-d10	17.304

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9H-Fluorene, 3-methyl-	180	C14H12	002523-39-9	93
2			9H-Fluorene, 1-methyl-	180	C14H12	001730-37-6	89
3			9H-Fluorene, 9-methyl-	180	C14H12	002523-37-7	76
4			9H-Fluorene, 1-methyl-	180	C14H12	001730-37-6	76
5			9H-Fluorene, 9-methyl-	180	C14H12	002523-37-7	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
Data File : BN015968.D
Acq On : 19 Aug 2021 11:49
Operator : CG/JU
Sample : M3399-05 30X
Misc :
ALS Vial : 75 Sample Multiplier: 1

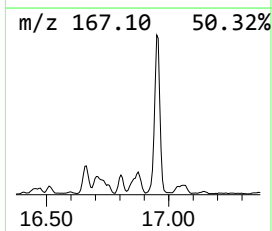
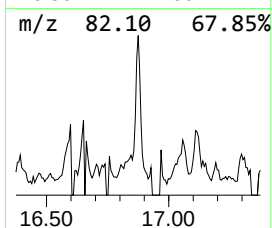
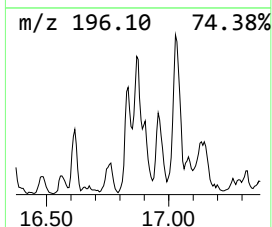
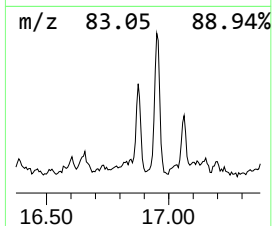
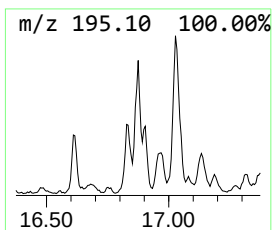
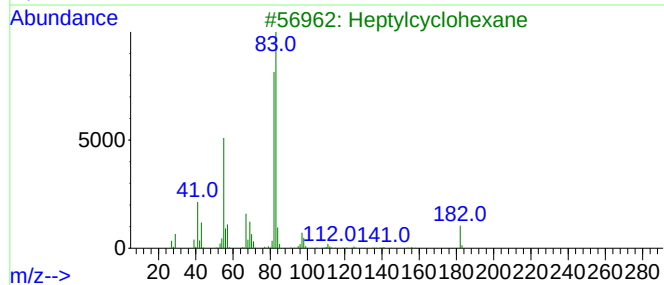
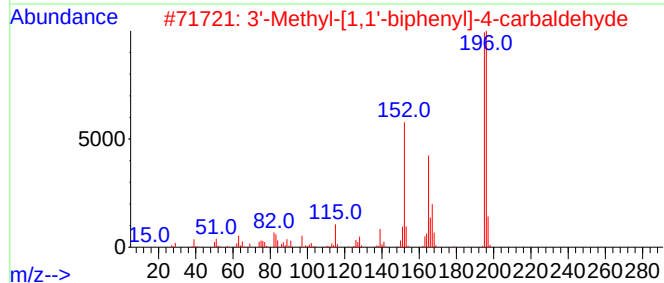
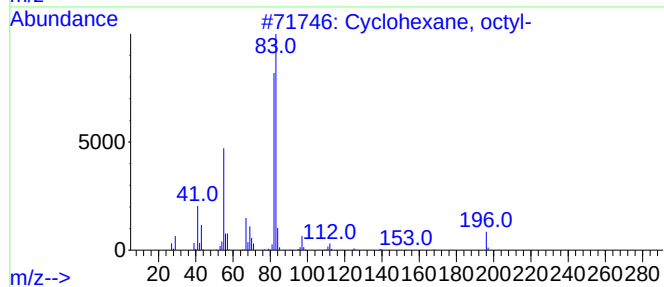
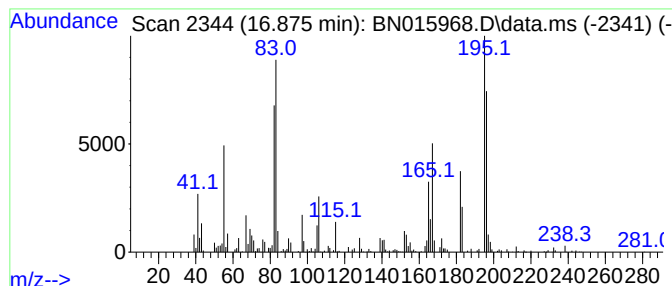
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081621MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 36 (DEL) Alkane: Cyclic16.875 Concentration Rank 44

R.T.	EstConc	Area	Relative to ISTD			R.T.
16.875	6.26 ng/ul	1081590	Phenanthrene-d10			17.304
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Cyclohexane, octyl-		196	C14H28	001795-15-9	50
2	3'-Methyl-[1,1'-biphenyl]-4-carb...		196	C14H12O	400744-83-4	41
3	Heptylcyclohexane		182	C13H26	005617-41-4	38
4	4-(4-Methylphenyl)benzaldehyde		196	C14H12O	036393-42-7	38
5	Cyclohexane, undecyl-		238	C17H34	054105-66-7	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
Data File : BN015968.D
Acq On : 19 Aug 2021 11:49
Operator : CG/JU
Sample : M3399-05 30X
Misc :
ALS Vial : 75 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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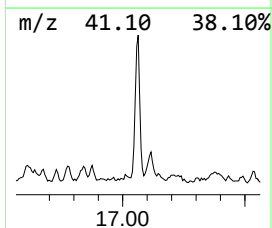
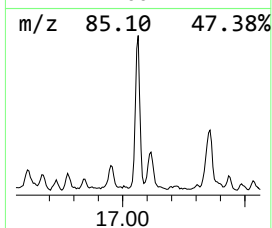
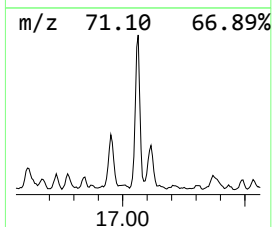
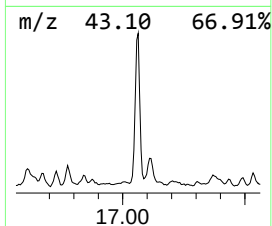
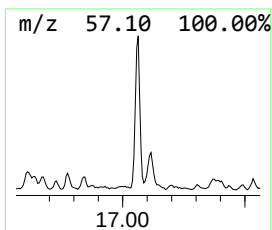
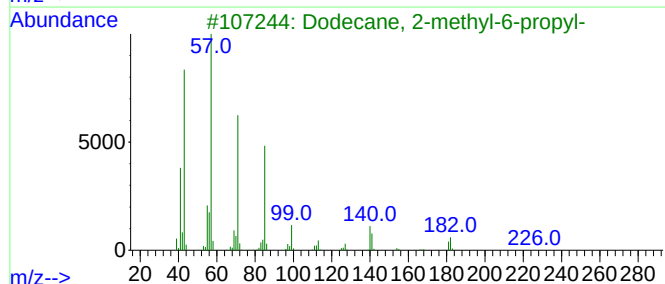
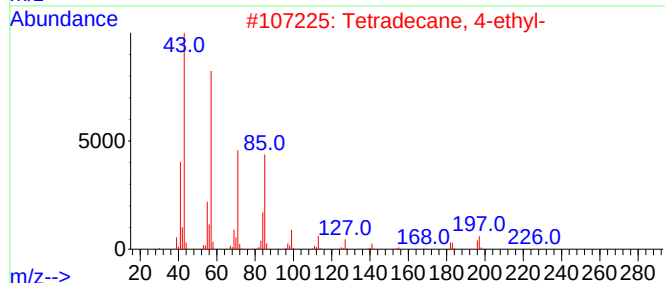
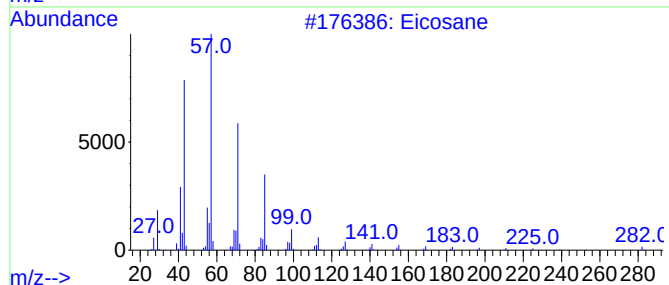
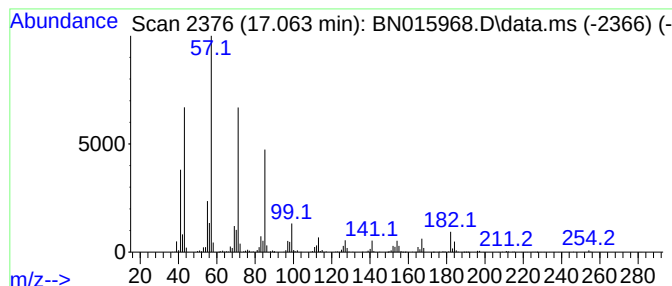
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Quant Title : SVOA CALIBRATION

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

Peak Number 37 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.063	26.74 ng/ul	4620740	Phenanthrene-d10	17.304

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	95
2			Tetradecane, 4-ethyl-	226	C16H34	055045-14-2	87
3			Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	87
4			Heneicosane	296	C21H44	000629-94-7	81
5			Pentadecane	212	C15H32	000629-62-9	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
Data File : BN015968.D
Acq On : 19 Aug 2021 11:49
Operator : CG/JU
Sample : M3399-05 30X
Misc :
ALS Vial : 75 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBKA7

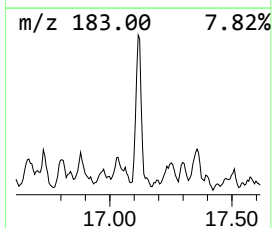
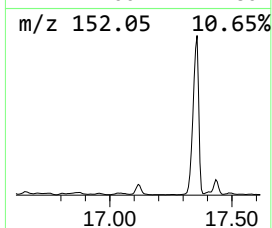
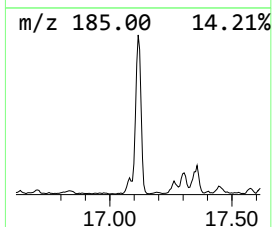
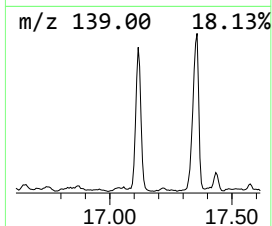
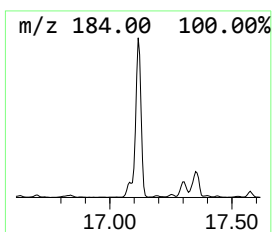
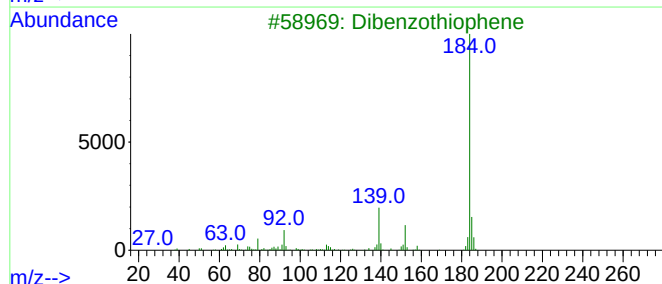
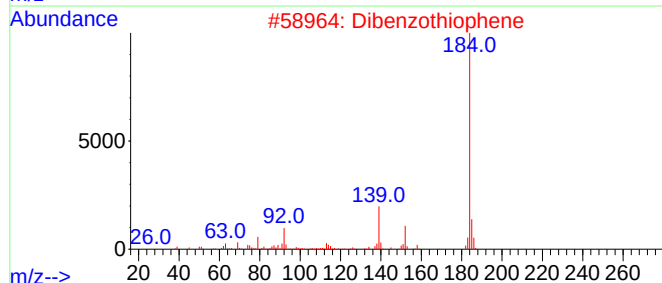
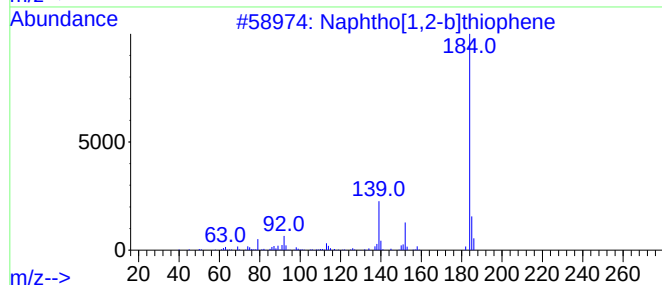
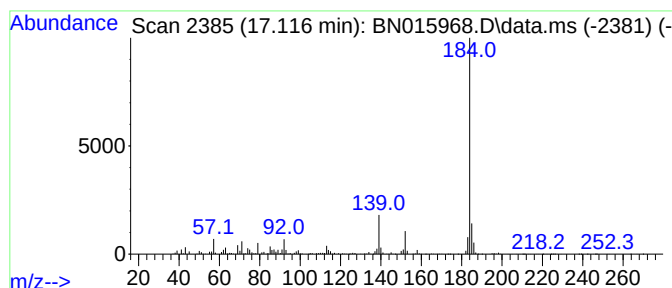
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Quant Title : SVOA CALIBRATION

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

Peak Number 38 Naphtho[1,2-b]thiophene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.116	31.79 ng/ul	5492050	Phenanthrene-d10	17.304

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphtho[1,2-b]thiophene	184	C12H8S	000234-41-3	97
2			Dibenzothiophene	184	C12H8S	000132-65-0	97
3			Dibenzothiophene	184	C12H8S	000132-65-0	97
4			Naphtho[2,1-b]thiophene	184	C12H8S	000233-02-3	97
5			Dibenzothiophene	184	C12H8S	000132-65-0	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
Data File : BN015968.D
Acq On : 19 Aug 2021 11:49
Operator : CG/JU
Sample : M3399-05 30X
Misc :
ALS Vial : 75 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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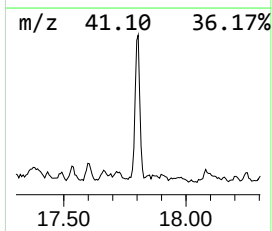
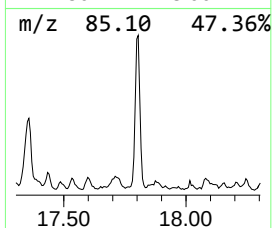
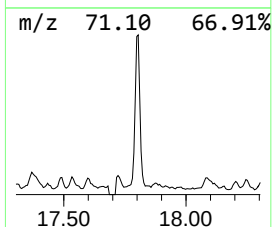
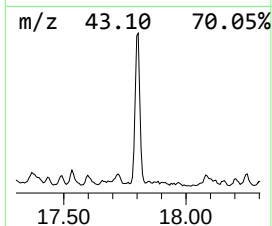
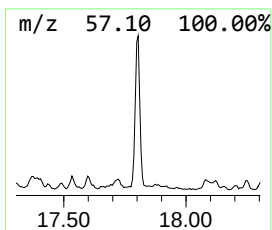
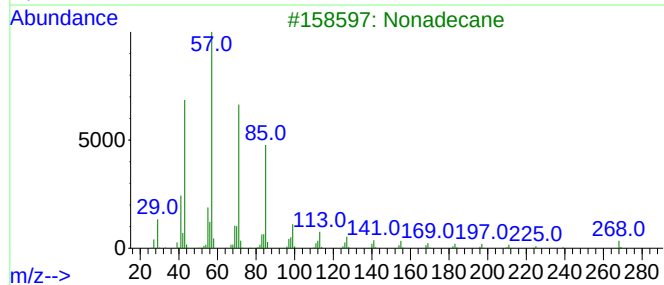
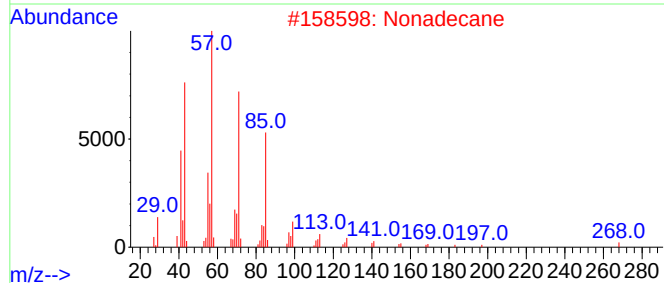
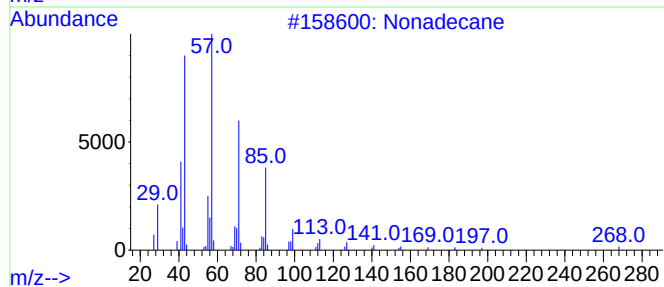
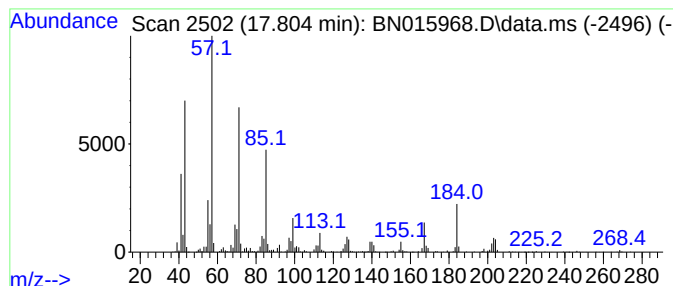
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Quant Title : SVOA CALIBRATION

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

Peak Number 40 (DEL) Alkane: Straight-Chai... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.804	21.32 ng/ul	3683870	Phenanthrene-d10	17.304

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	97
2			Nonadecane	268	C19H40	000629-92-5	91
3			Nonadecane	268	C19H40	000629-92-5	86
4			Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	70
5			Decane, 1-iodo-	268	C10H21I	002050-77-3	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
Data File : BN015968.D
Acq On : 19 Aug 2021 11:49
Operator : CG/JU
Sample : M3399-05 30X
Misc :
ALS Vial : 75 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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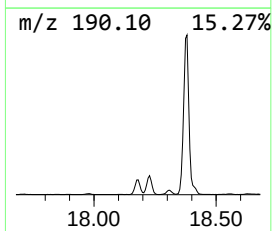
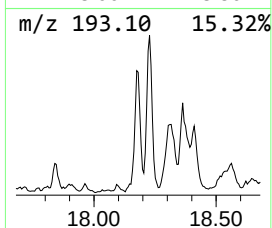
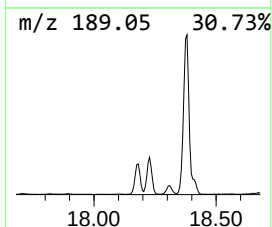
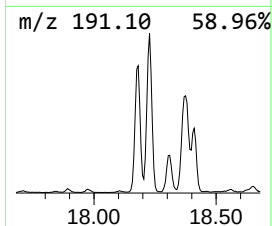
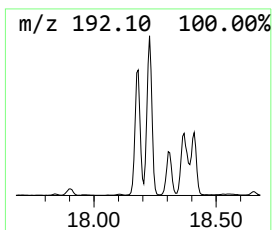
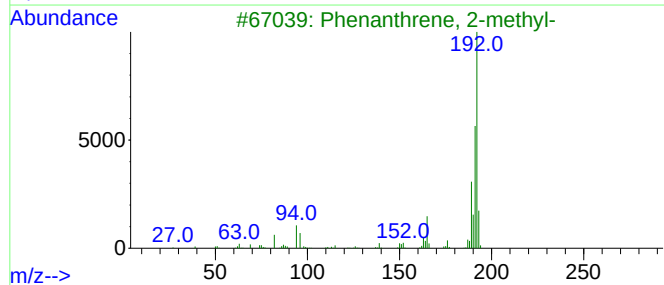
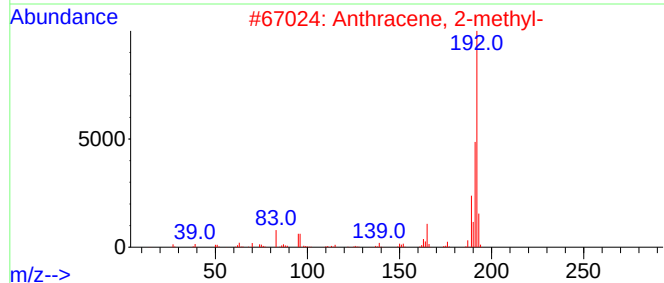
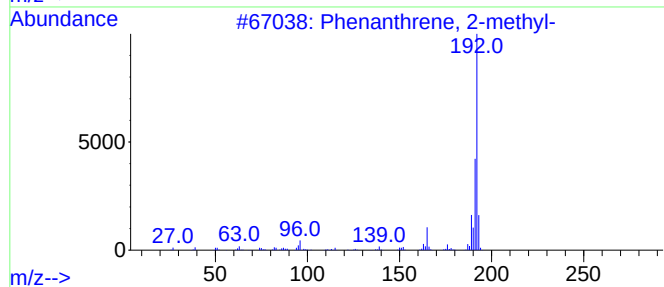
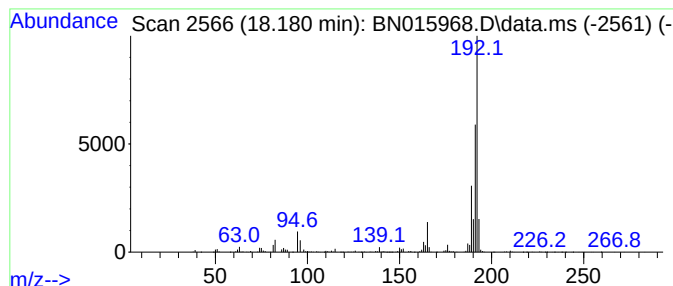
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Quant Title : SVOA CALIBRATION

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

Peak Number 42 Anthracene, 2-methyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.180	19.52 ng/ul	3372600	Phenanthrene-d10	17.304

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
2			Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
3			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
4			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
5			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
Data File : BN015968.D
Acq On : 19 Aug 2021 11:49
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Sample : M3399-05 30X
Misc :
ALS Vial : 75 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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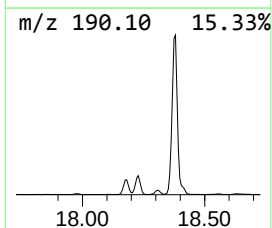
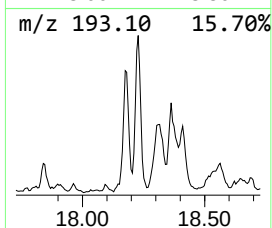
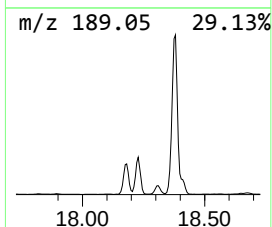
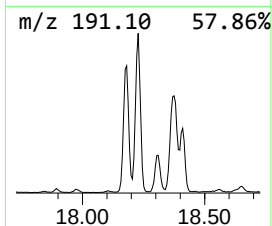
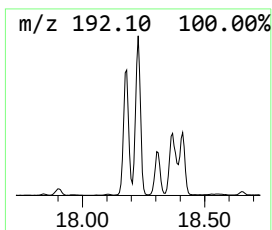
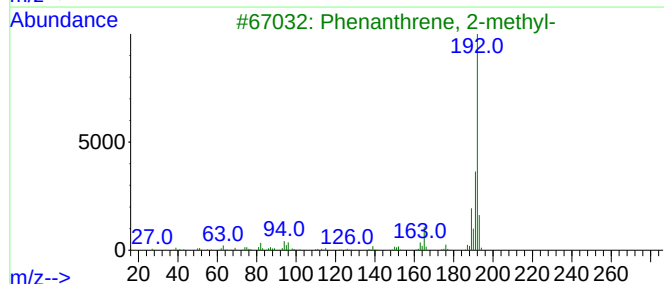
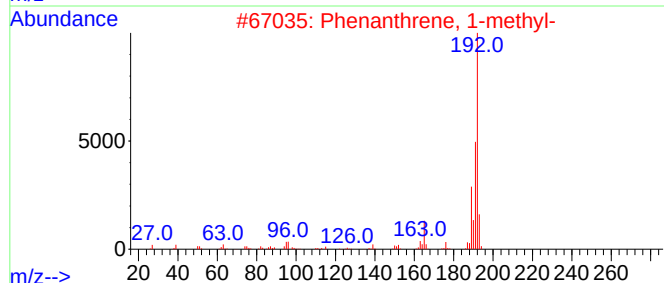
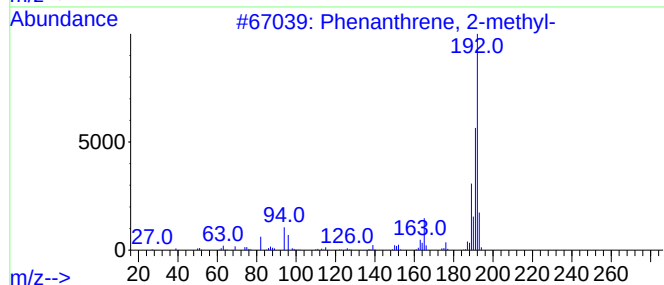
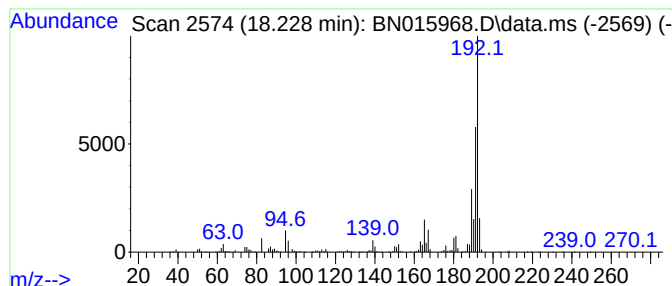
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Quant Title : SVOA CALIBRATION

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

Peak Number 43 Phenanthrene, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.227	33.38 ng/ul	5768080	Phenanthrene-d10	17.304

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
2			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
3			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	94
4			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
5			Phenanthrene, 3-methyl-	192	C15H12	000832-71-3	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
Data File : BN015968.D
Acq On : 19 Aug 2021 11:49
Operator : CG/JU
Sample : M3399-05 30X
Misc :
ALS Vial : 75 Sample Multiplier: 1

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BNA_N
ClientSampleId :
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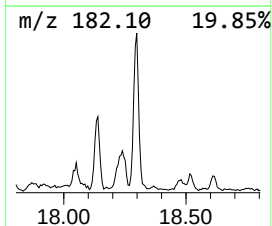
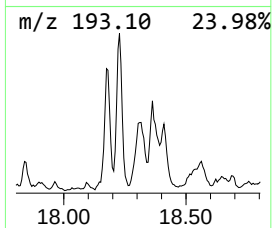
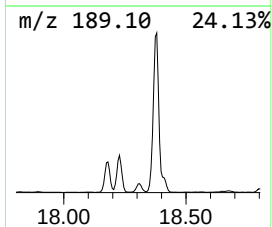
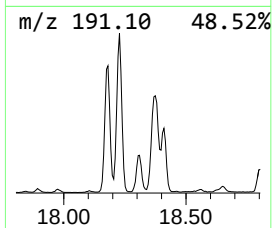
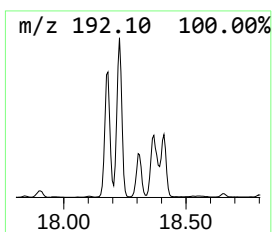
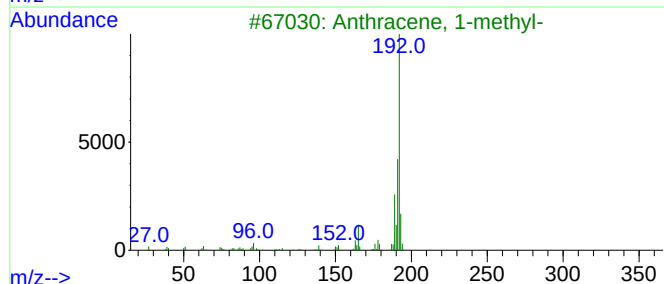
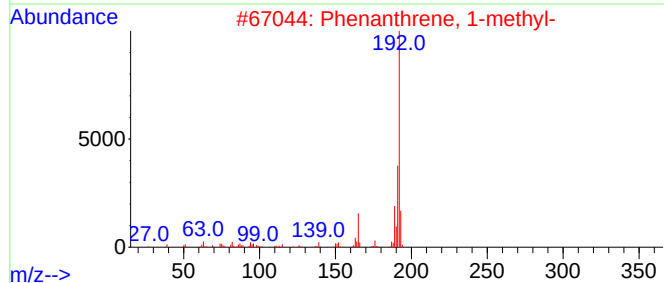
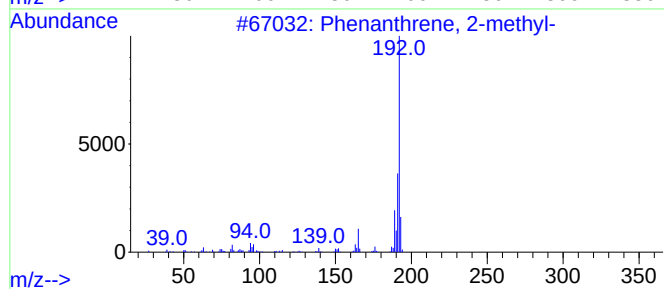
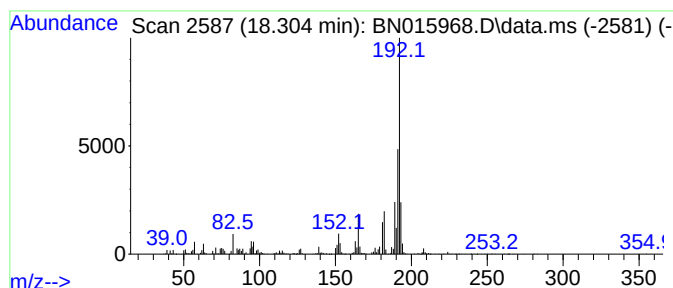
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Quant Title : SVOA CALIBRATION

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

Peak Number 44 Phenanthrene, 1-methyl- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.304	12.93 ng/ul	2234090	Phenanthrene-d10	17.304

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
2			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
3			Anthracene, 1-methyl-	192	C15H12	000610-48-0	92
4			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	91
5			Anthracene, 2-methyl-	192	C15H12	000613-12-7	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
Data File : BN015968.D
Acq On : 19 Aug 2021 11:49
Operator : CG/JU
Sample : M3399-05 30X
Misc :
ALS Vial : 75 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBKA7

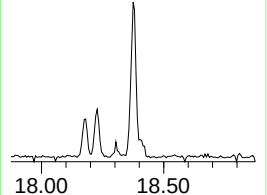
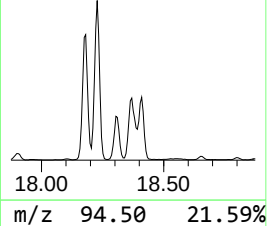
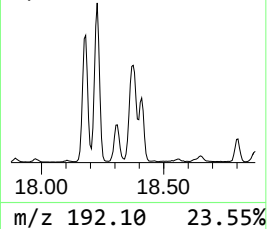
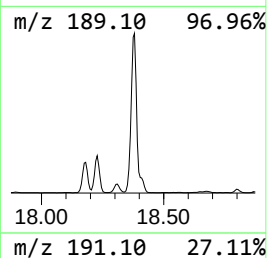
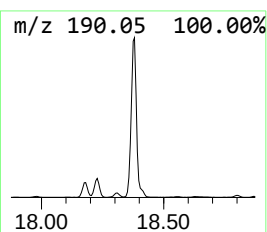
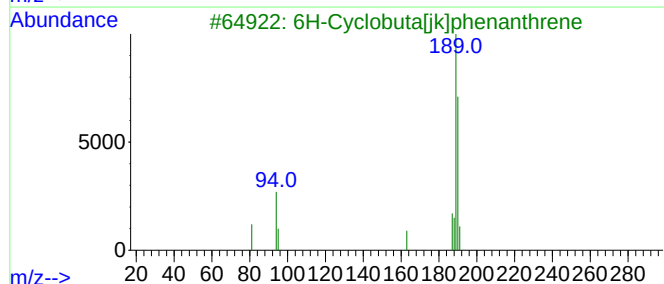
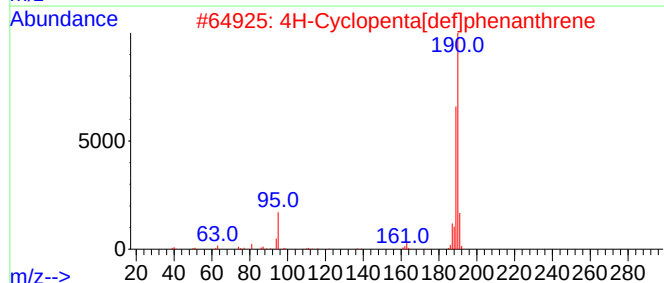
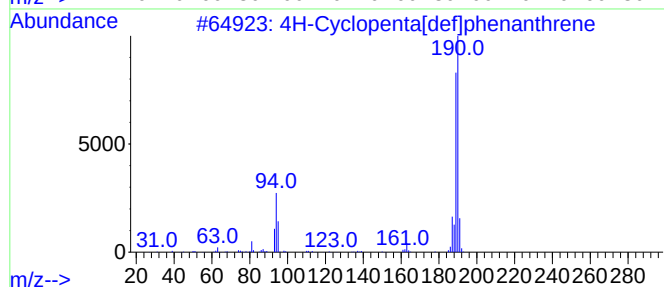
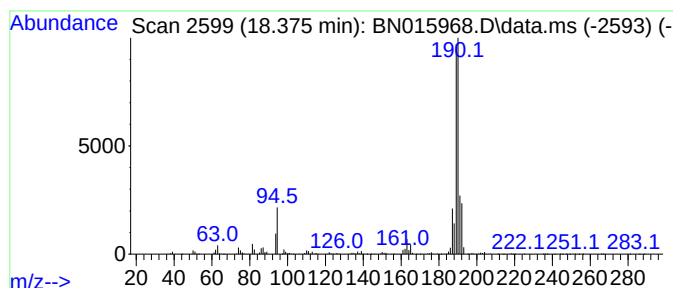
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Quant Title : SVOA CALIBRATION

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

Peak Number 45 4H-Cyclopenta[def]phenanthrene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.375	47.33 ng/ul	8177670	Phenanthrene-d10	17.304

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	93
2			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	64
3			6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	64
4			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	60
5			2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	45



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
Data File : BN015968.D
Acq On : 19 Aug 2021 11:49
Operator : CG/JU
Sample : M3399-05 30X
Misc :
ALS Vial : 75 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBKA7

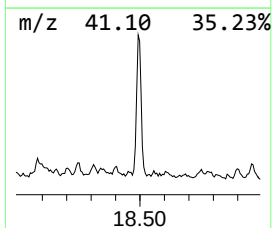
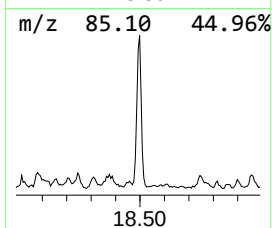
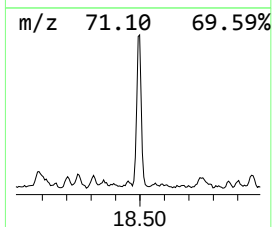
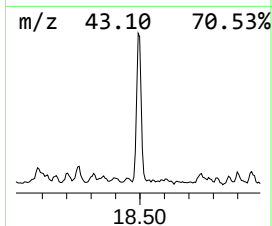
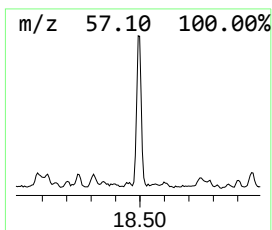
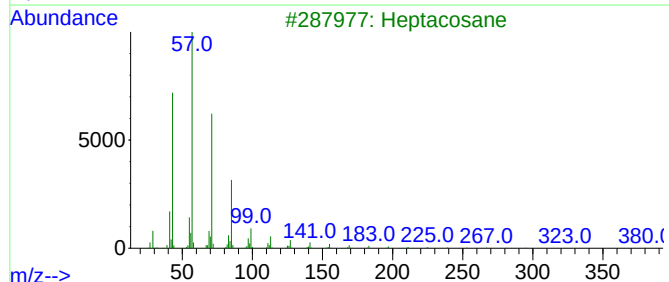
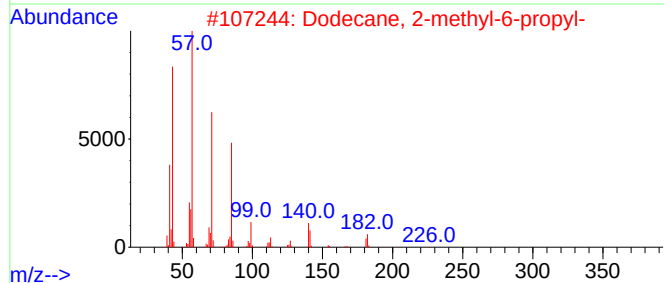
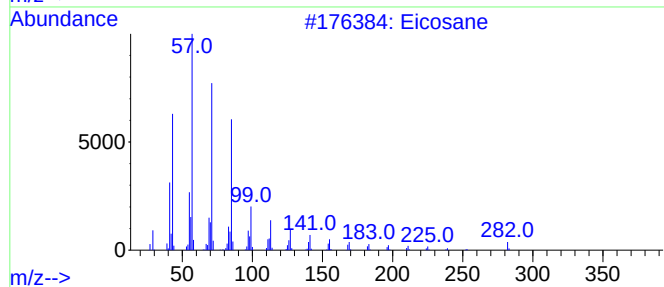
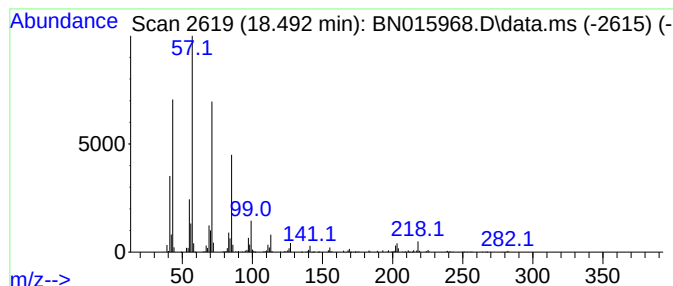
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Quant Title : SVOA CALIBRATION

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

Peak Number 46 (DEL) Alkane: Straight-Chai... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.492	11.09 ng/ul	1915420	Phenanthrene-d10	17.304

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Eicosane	282	C20H42	000112-95-8	94
2		Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	93
3		Heptacosane	380	C27H56	000593-49-7	87
4		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	86
5		2-Bromo dodecane	248	C12H25Br	013187-99-0	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
Data File : BN015968.D
Acq On : 19 Aug 2021 11:49
Operator : CG/JU
Sample : M3399-05 30X
Misc :
ALS Vial : 75 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBKA7

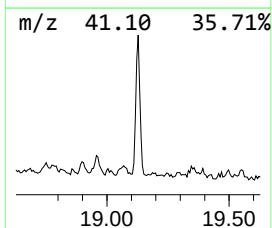
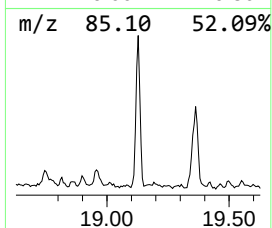
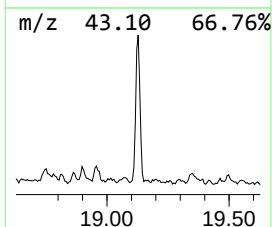
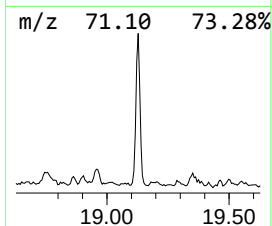
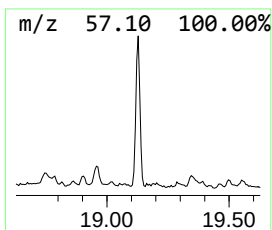
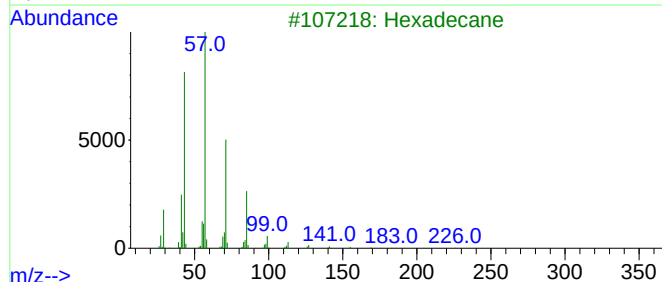
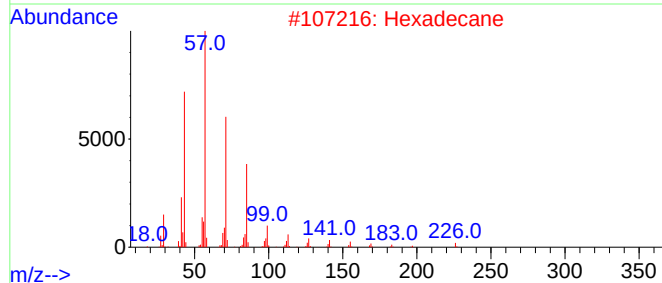
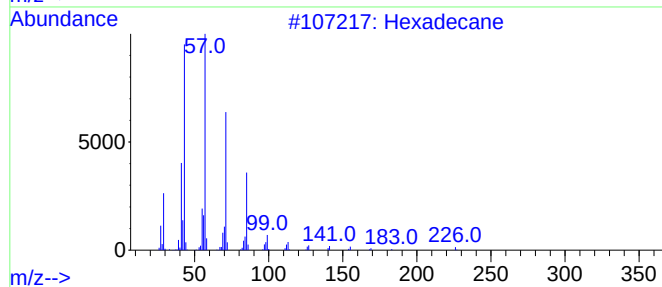
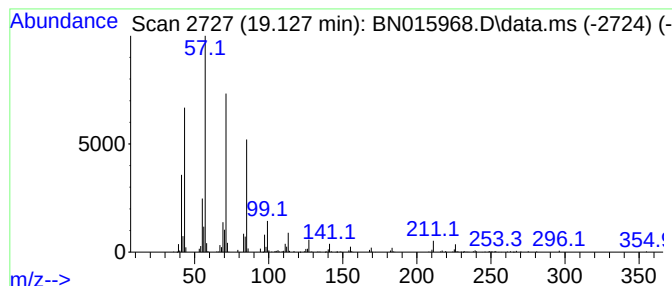
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081621MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 51 (DEL) Alkane: Straight-Chai... Concentration Rank 53

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.128	4.41 ng/ul	761367	Phenanthrene-d10	17.304

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexadecane	226	C16H34	000544-76-3	94
2		Hexadecane	226	C16H34	000544-76-3	94
3		Hexadecane	226	C16H34	000544-76-3	93
4		Hexadecane	226	C16H34	000544-76-3	91
5		Tetracosane	338	C24H50	000646-31-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
 Data File : BN015968.D
 Acq On : 19 Aug 2021 11:49
 Operator : CG/JU
 Sample : M3399-05 30X
 Misc :
 ALS Vial : 75 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBKA7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081621MA.M
 Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Ethylbenzene	5.405	10.4	ng/ul	861195	1	7.899	1653990	20.0
Benzo[<i>a</i>]pyrene	7.622	10.0	ng/ul	823781	1	7.899	1653990	20.0
Tetracyclo[3.3.1.0 ^{2,6} .0 ^{3,4}]	8.275	9.6	ng/ul	789976	1	7.899	1653990	20.0
Indene	8.440	98.0	ng/ul	8105160	1	7.899	1653990	20.0
3-Phenylbut-1-ene	10.104	7.1	ng/ul	1118060	2	10.710	3130570	20.0
Benzo[<i>c</i>]thiophene	10.881	10.9	ng/ul	1706240	2	10.710	3130570	20.0
(DEL) Alkane: S...	12.140	7.9	ng/ul	1231730	2	10.710	3130570	20.0
Indole	12.204	7.1	ng/ul	1112970	2	10.710	3130570	20.0
Benzo[<i>b</i>]thiophe...	12.487	4.6	ng/ul	719176	2	10.710	3130570	20.0
Naphthalene, 2-...	13.575	27.8	ng/ul	4085880	3	14.551	2936580	20.0
Naphthalene, 1-...	13.716	31.0	ng/ul	4553840	3	14.551	2936580	20.0
Naphthalene, 2-...	13.869	49.4	ng/ul	7250510	3	14.551	2936580	20.0
(DEL) Alkane: S...	14.034	6.0	ng/ul	876645	3	14.551	2936580	20.0
Naphthalene, 1-...	14.104	20.3	ng/ul	2985510	3	14.551	2936580	20.0
(DEL) Alkane: S...	14.451	18.3	ng/ul	2684600	3	14.551	2936580	20.0
2-Naphthaleneca...	14.675	12.3	ng/ul	1799840	3	14.551	2936580	20.0
Naphthalene, 2-...	14.757	17.3	ng/ul	2544910	3	14.551	2936580	20.0
1-Isopropenylna...	14.857	8.6	ng/ul	1267190	3	14.551	2936580	20.0
Naphthalene, 1-...	15.022	14.3	ng/ul	2104630	3	14.551	2936580	20.0
Phenol, 2,3,4,6...	15.092	55.4	ng/ul	8139720	3	14.551	2936580	20.0
Naphthalene, 2-...	15.339	9.9	ng/ul	1455980	3	14.551	2936580	20.0
(DEL) Alkane: S...	15.404	19.6	ng/ul	2876660	3	14.551	2936580	20.0
Naphthalene, 1-...	15.722	14.1	ng/ul	2064360	3	14.551	2936580	20.0
Nordiphenamid	15.757	17.3	ng/ul	2546060	3	14.551	2936580	20.0
Naphthalene, 1-...	15.810	10.4	ng/ul	1529640	3	14.551	2936580	20.0
Dibenzofuran, 4...	15.892	22.3	ng/ul	3267090	3	14.551	2936580	20.0
6H-Dibenzo[<i>b,d</i>]...	16.034	25.0	ng/ul	4319200	4	17.304	3455720	20.0
(DEL) Alkane: S...	16.263	35.2	ng/ul	6081140	4	17.304	3455720	20.0
Anthracene, 9,1...	16.392	12.2	ng/ul	2108560	4	17.304	3455720	20.0
9H-Fluorene, 3-...	16.604	12.5	ng/ul	2157650	4	17.304	3455720	20.0
(DEL) Alkane: C...	16.875	6.3	ng/ul	1081590	4	17.304	3455720	20.0
(DEL) Alkane: S...	17.063	26.7	ng/ul	4620740	4	17.304	3455720	20.0
Naphtho[1,2- <i>b</i>]t...	17.116	31.8	ng/ul	5492050	4	17.304	3455720	20.0
(DEL) Alkane: S...	17.804	21.3	ng/ul	3683870	4	17.304	3455720	20.0
Anthracene, 2-m...	18.180	19.5	ng/ul	3372600	4	17.304	3455720	20.0
Phenanthrene, 2...	18.227	33.4	ng/ul	5768080	4	17.304	3455720	20.0
Phenanthrene, 1...	18.304	12.9	ng/ul	2234090	4	17.304	3455720	20.0
4H-Cyclopenta[<i>d</i> ...	18.375	47.3	ng/ul	8177670	4	17.304	3455720	20.0
(DEL) Alkane: S...	18.492	11.1	ng/ul	1915420	4	17.304	3455720	20.0
(DEL) Alkane: S...	19.128	4.4	ng/ul	761367	4	17.304	3455720	20.0