

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060722\
 Data File : BN020038.D
 Acq On : 07 Jun 2022 21:06
 Operator : CG/JU
 Sample : N3171-08
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBTZ3

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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

Peak separation: 5

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

Signal : TIC: BN020038.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.281	30	33	39	rBV	19824	23932	0.59%	0.056%
2	3.687	99	102	121	rBV	132669	223052	5.51%	0.520%
3	5.346	379	384	391	rVV	33718	52716	1.30%	0.123%
4	6.987	657	663	671	rBV	52670	91752	2.27%	0.214%
5	7.152	685	691	699	rBV	343442	550000	13.59%	1.283%
6	7.222	699	703	712	rVB	15375	26926	0.67%	0.063%
7	7.358	718	726	736	rBV	200625	339302	8.38%	0.791%
8	7.822	797	805	813	rBV	408281	674868	16.68%	1.574%
9	8.199	860	869	878	rBV	502362	821501	20.30%	1.916%
10	8.357	891	896	902	rVB	14639	24134	0.60%	0.056%
11	8.528	918	925	935	rVB	197341	331888	8.20%	0.774%
12	8.975	989	1001	1012	rBV	421368	786635	19.44%	1.835%
13	9.181	1026	1036	1042	rVV	13105	24761	0.61%	0.058%
14	9.299	1049	1056	1063	rVV	25794	58402	1.44%	0.136%
15	9.693	1117	1123	1132	rBV	197988	348574	8.61%	0.813%
16	9.869	1147	1153	1161	rVB	35805	62396	1.54%	0.146%
17	10.016	1167	1178	1188	rBV2	66657	132126	3.26%	0.308%
18	10.234	1208	1215	1225	rBV2	332441	599979	14.82%	1.399%
19	10.610	1270	1279	1284	rBV	600921	1085615	26.82%	2.532%
20	10.663	1284	1288	1295	rVV	583788	975190	24.10%	2.274%
21	10.746	1295	1302	1317	rVB2	375025	765264	18.91%	1.785%
22	12.004	1511	1516	1525	rVB7	11246	29020	0.72%	0.068%
23	12.134	1528	1538	1542	rBV	118635	209730	5.18%	0.489%
24	12.275	1555	1562	1568	rVB	223676	359983	8.89%	0.840%
25	12.393	1575	1582	1588	rVB	49139	85496	2.11%	0.199%
26	12.493	1588	1599	1606	rVB	470091	808120	19.97%	1.885%
27	12.704	1630	1635	1642	rVB2	16849	28271	0.70%	0.066%
28	12.869	1659	1663	1668	rBV	21170	32087	0.79%	0.075%
29	13.287	1730	1734	1744	rVV2	32278	58526	1.45%	0.136%
30	13.416	1751	1756	1761	rVV5	19205	35942	0.89%	0.084%
31	13.487	1761	1768	1779	rVB3	32637	96461	2.38%	0.225%
32	13.628	1779	1792	1800	rBV7	41356	137479	3.40%	0.321%
33	13.775	1811	1817	1822	rVV	91472	166822	4.12%	0.389%
34	13.857	1822	1831	1838	rVV	1220445	1840527	45.48%	4.292%
35	14.010	1850	1857	1860	rVV	40860	65676	1.62%	0.153%
36	14.045	1860	1863	1867	rVV3	17457	23395	0.58%	0.055%

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060722\
 Data File : BN020038.D
 Acq On : 07 Jun 2022 21:06
 Operator : CG/JU
 Sample : N3171-08
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBTZ3

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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

Peak separation: 5

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

37	14.140	1871	1879	1894	rVV	1214453	1947102	48.11%	4.541%
38	14.451	1921	1932	1937	rVV	983628	1479718	36.56%	3.451%
39	14.516	1937	1943	1949	rVV	1032324	1585180	39.17%	3.697%
40	14.575	1949	1953	1957	rVV	19378	29019	0.72%	0.068%
41	14.640	1959	1964	1973	rVB2	50739	89473	2.21%	0.209%
42	14.722	1973	1978	1982	rBV2	38001	58971	1.46%	0.138%
43	14.845	1994	1999	2006	rVV	172489	280255	6.92%	0.654%
44	15.439	2091	2100	2105	rBV	1778242	2588832	63.97%	6.037%
45	15.498	2105	2110	2115	rVB	421608	635058	15.69%	1.481%
46	15.557	2115	2120	2127	rBV	357557	514924	12.72%	1.201%
47	15.622	2127	2131	2134	rBV4	37428	52886	1.31%	0.123%
48	15.792	2156	2160	2169	rVB3	29248	50930	1.26%	0.119%
49	15.934	2177	2184	2191	rVB6	32381	73140	1.81%	0.171%
50	15.992	2191	2194	2205	rVB3	11373	26197	0.65%	0.061%
51	16.163	2214	2223	2229	rVV3	82592	154076	3.81%	0.359%
52	16.222	2230	2233	2238	rVB2	18247	26053	0.64%	0.061%
53	16.286	2238	2244	2250	rBV	19599	33755	0.83%	0.079%
54	16.345	2250	2254	2257	rVV4	20781	28765	0.71%	0.067%
55	16.451	2266	2272	2276	rBV2	19081	31145	0.77%	0.073%
56	16.492	2276	2279	2285	rVV6	13317	26525	0.66%	0.062%
57	16.551	2285	2289	2295	rVB8	15235	26932	0.67%	0.063%
58	16.651	2301	2306	2313	rBV2	45057	78737	1.95%	0.184%
59	16.975	2351	2361	2364	rVV2	49705	102724	2.54%	0.240%
60	17.010	2364	2367	2373	rVV	34109	61487	1.52%	0.143%
61	17.110	2377	2384	2387	rVV6	19710	54927	1.36%	0.128%
62	17.145	2387	2390	2392	rVV2	29021	39821	0.98%	0.093%
63	17.192	2392	2398	2402	rVV	1175820	1818117	44.92%	4.240%
64	17.233	2402	2405	2409	rVV	251162	379970	9.39%	0.886%
65	17.286	2409	2414	2425	rVV2	2073951	3159763	78.07%	7.369%
66	17.381	2425	2430	2443	rVB2	27096	87540	2.16%	0.204%
67	17.592	2455	2466	2472	rBV	392224	638761	15.78%	1.490%
68	17.686	2478	2482	2488	rBV4	18037	32121	0.79%	0.075%
69	17.875	2501	2514	2521	rBV3	67305	141607	3.50%	0.330%
70	18.028	2536	2540	2545	rVB	38775	52105	1.29%	0.122%
71	18.098	2546	2552	2561	rVV4	143175	323708	8.00%	0.755%
72	18.186	2563	2567	2573	rVV	50678	84145	2.08%	0.196%
73	18.263	2573	2580	2585	rVB5	22058	47045	1.16%	0.110%
74	18.369	2590	2598	2601	rVV4	26242	70147	1.73%	0.164%
75	18.410	2601	2605	2609	rVV	38356	66696	1.65%	0.156%
76	18.504	2617	2621	2627	rVV2	36469	62632	1.55%	0.146%

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060722\
 Data File : BN020038.D
 Acq On : 07 Jun 2022 21:06
 Operator : CG/JU
 Sample : N3171-08
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBTZ3

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF Filtering: 5
 Sampling : 1 Min Area: 0.1 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

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

77	18.569	2627	2632	2636	rVV3	18951	33423	0.83%	0.078%
78	18.957	2691	2698	2702	rBV8	11959	24259	0.60%	0.057%
79	19.045	2709	2713	2717	rBV2	25619	33744	0.83%	0.079%
80	19.145	2726	2730	2738	rBV2	32311	68843	1.70%	0.161%
81	19.222	2738	2743	2745	rVV	29160	46058	1.14%	0.107%
82	19.251	2745	2748	2752	rVV	30131	40528	1.00%	0.095%
83	19.375	2765	2769	2778	rBV2	58858	92500	2.29%	0.216%
84	19.586	2798	2805	2823	rVB	2547387	3672580	90.75%	8.565%
85	19.975	2867	2871	2872	rBV2	23199	31358	0.77%	0.073%
86	20.045	2879	2883	2890	rBV	116608	168710	4.17%	0.393%
87	20.145	2896	2900	2906	rVB	52936	70172	1.73%	0.164%
88	20.692	2989	2993	3000	rVB	68232	105023	2.59%	0.245%
89	21.098	3058	3062	3078	rBV	249091	475187	11.74%	1.108%
90	21.304	3093	3097	3101	rVB	33203	36761	0.91%	0.086%
91	21.380	3104	3110	3121	rVV	1492260	2068990	51.12%	4.825%
92	21.545	3135	3138	3144	rVB3	19393	25692	0.63%	0.060%
93	21.992	3212	3214	3227	rVB5	24511	66661	1.65%	0.155%
94	22.474	3294	3296	3301	rVB2	39828	54528	1.35%	0.127%
95	23.016	3383	3388	3395	rBV	58764	96991	2.40%	0.226%
96	23.563	3472	3481	3487	rBV	2024699	4047132	100.00%	9.438%
97	23.710	3498	3506	3519	rVB2	1029164	2195258	54.24%	5.120%
98	24.310	3603	3608	3615	rVB2	64836	137273	3.39%	0.320%
99	25.110	3739	3744	3755	rVB2	66918	163691	4.04%	0.382%
100	26.068	3900	3907	3914	rVB4	39418	106764	2.64%	0.249%

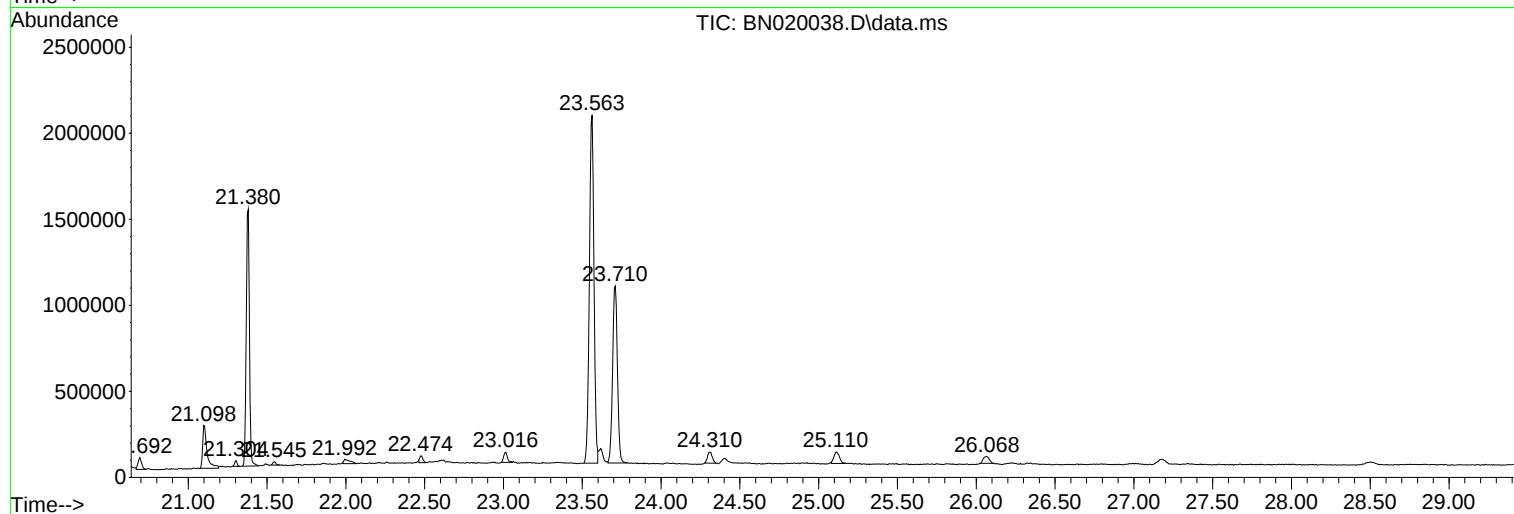
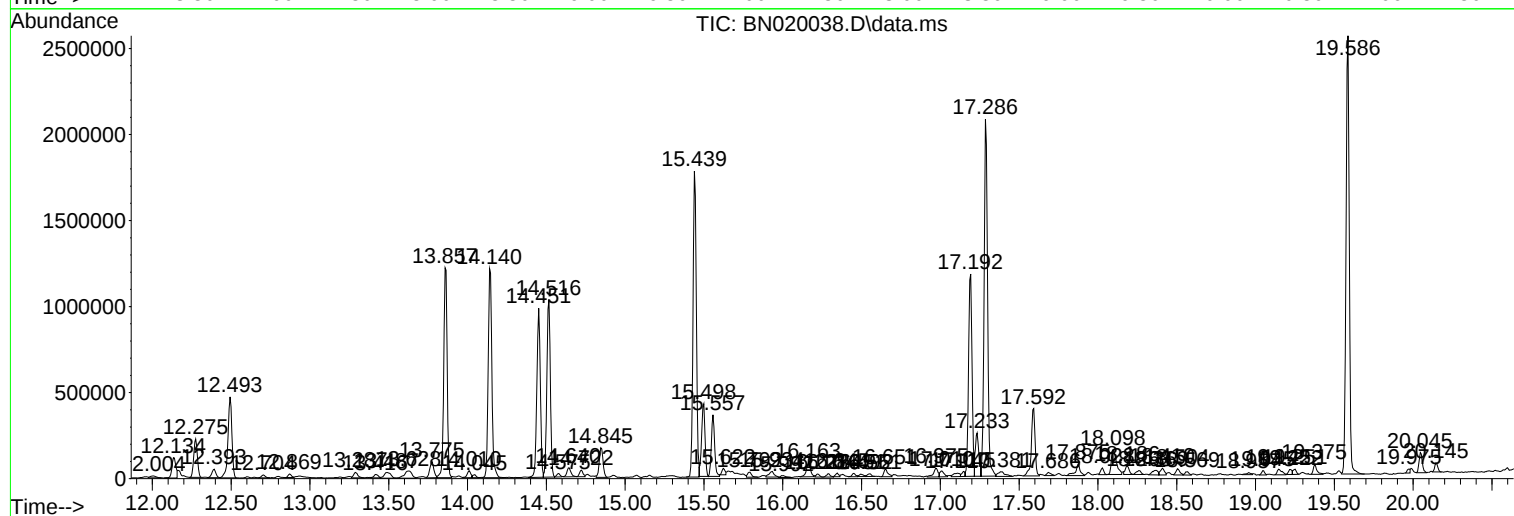
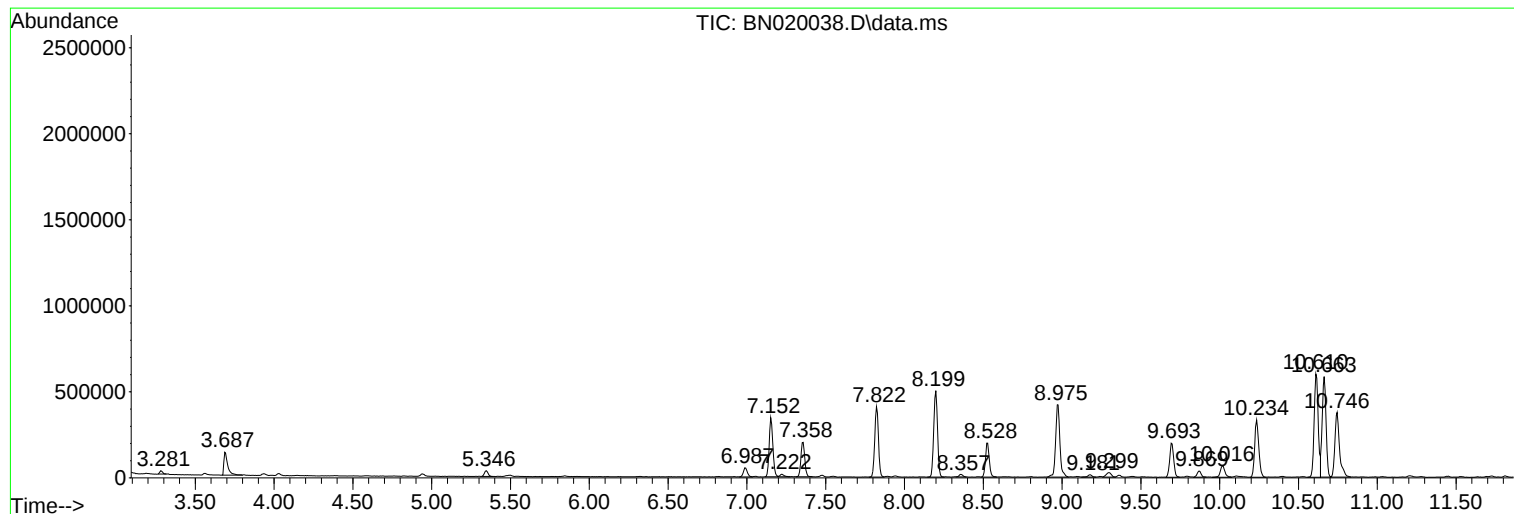
Sum of corrected areas: 42879663

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060722\
Data File : BN020038.D
Acq On : 07 Jun 2022 21:06
Operator : CG/JU
Sample : N3171-08
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBTZ3

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

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060722\
Data File : BN020038.D
Acq On : 07 Jun 2022 21:06
Operator : CG/JU
Sample : N3171-08
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBTZ3

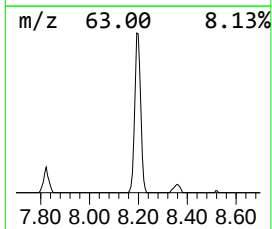
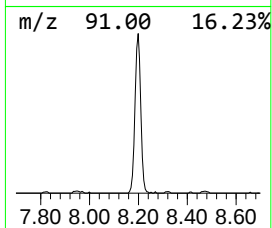
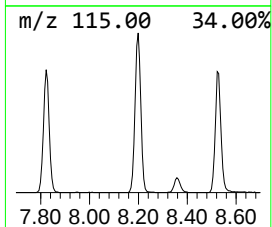
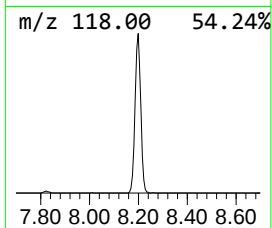
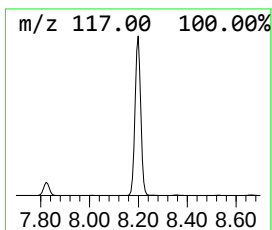
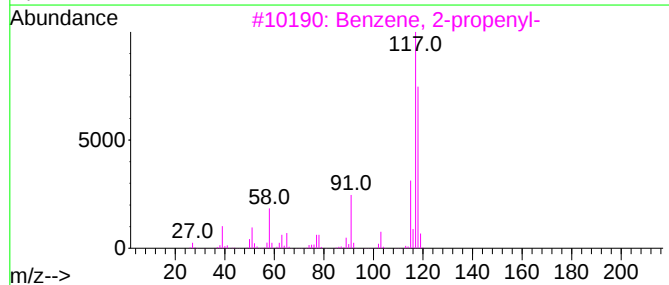
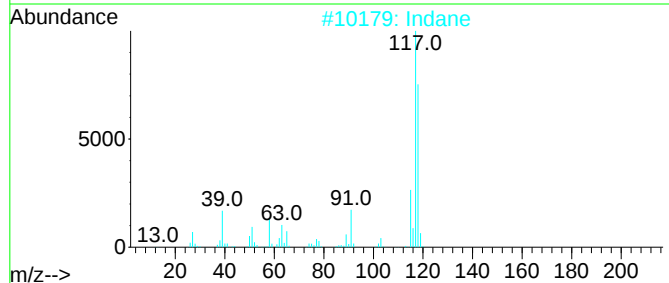
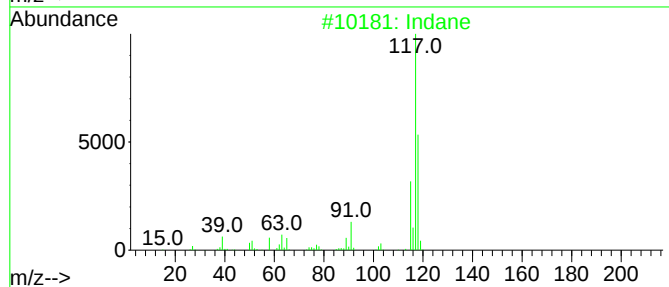
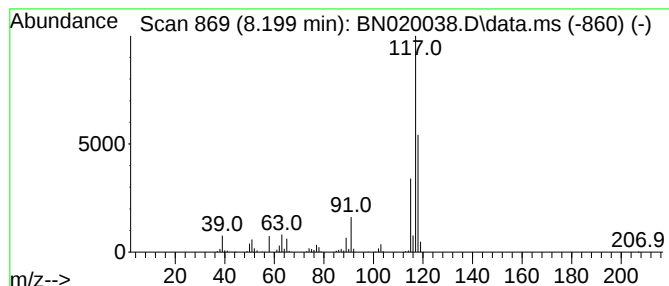
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN060622.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.199	24.35 ng/ul	821501	1,4-Dichlorobenzene-d4	7.822

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane			118	C9H10	000496-11-7	93
2	Indane			118	C9H10	000496-11-7	91
3	Benzene, 2-propenyl-			118	C9H10	000300-57-2	74
4	Deltacyclene			118	C9H10	007785-10-6	72
5	Indane			118	C9H10	000496-11-7	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060722\
Data File : BN020038.D
Acq On : 07 Jun 2022 21:06
Operator : CG/JU
Sample : N3171-08
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBTZ3

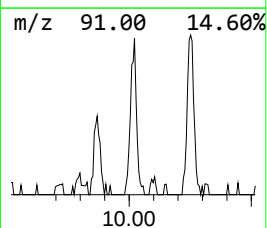
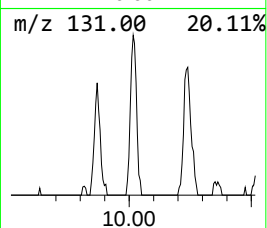
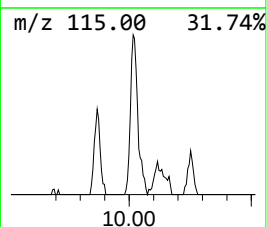
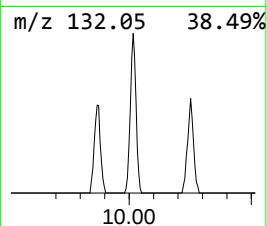
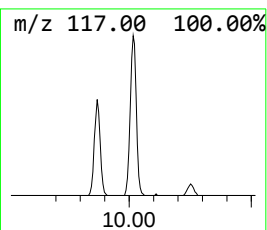
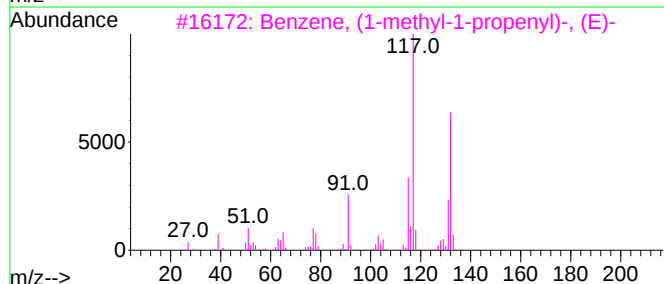
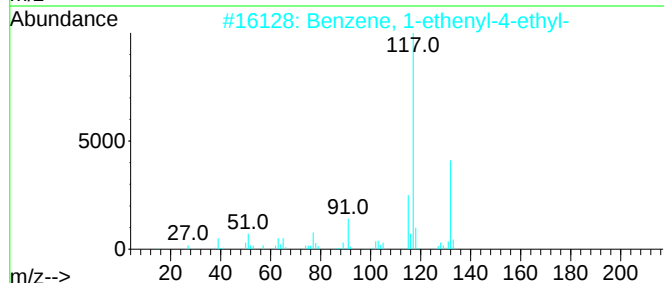
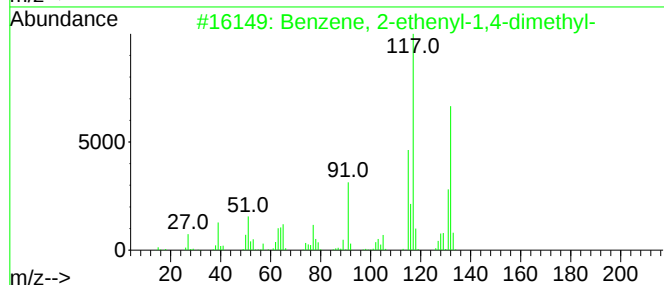
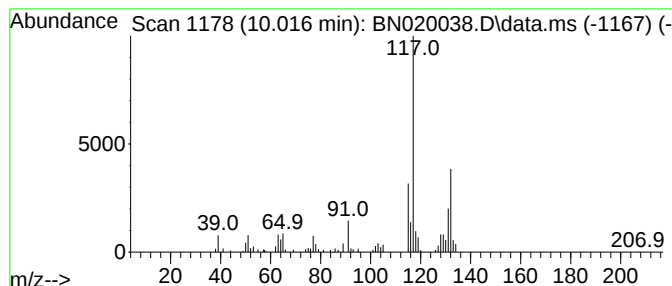
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN060622.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.016	2.43 ng/ul	132126	Naphthalene-d8	10.610

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	91
3			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	91
4			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	90
5			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060722\
Data File : BN020038.D
Acq On : 07 Jun 2022 21:06
Operator : CG/JU
Sample : N3171-08
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBTZ3

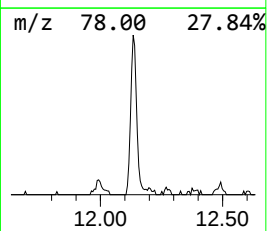
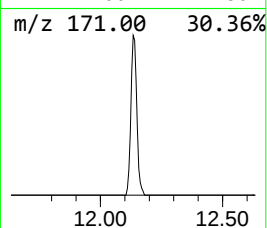
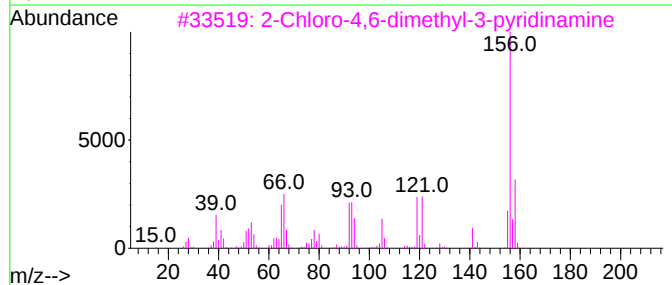
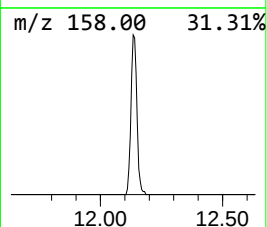
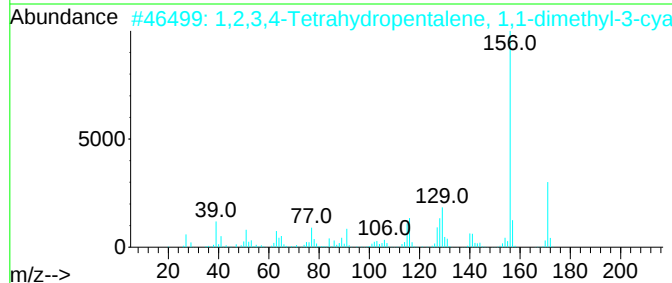
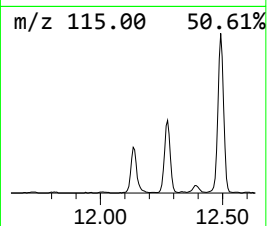
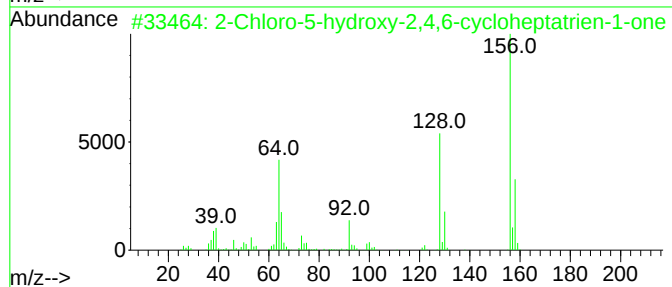
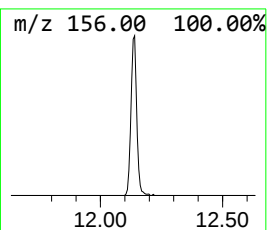
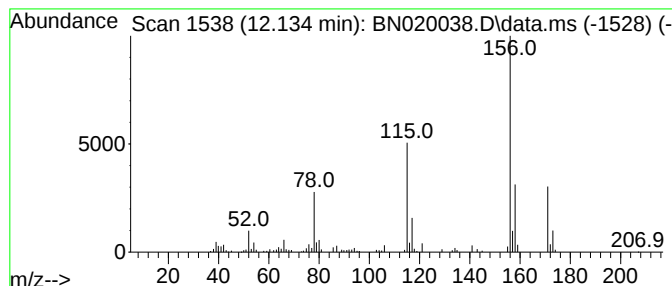
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN060622.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 unknown-01 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.134	3.86 ng/ul	209730	Naphthalene-d8	10.610

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Chloro-5-hydroxy-2,4,6-cyclohe...	156	C7H5ClO2	007009-16-7	43
2			1,2,3,4-Tetrahydropentalene, 1,1...	171	C12H13N	1000217-18-6	43
3			2-Chloro-4,6-dimethyl-3-pyridina...	156	C7H9ClN2	140413-40-7	32
4			6-Isopropylquinoline	171	C12H13N	000135-79-5	32
5			Bicyclo[3.2.0]hept-2-ene, 4-isop...	171	C12H13N	1000217-15-6	32



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060722\
Data File : BN020038.D
Acq On : 07 Jun 2022 21:06
Operator : CG/JU
Sample : N3171-08
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBTZ3

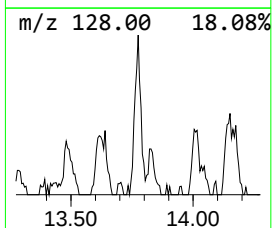
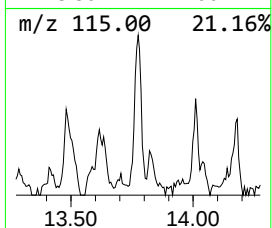
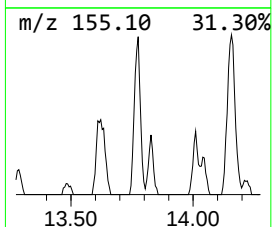
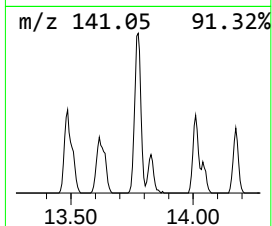
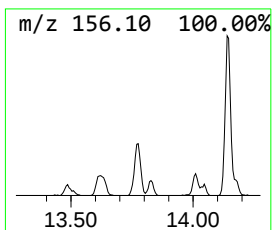
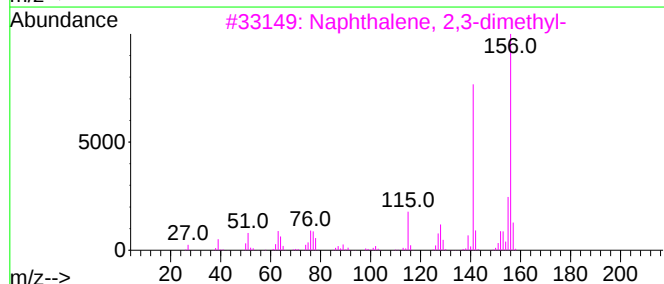
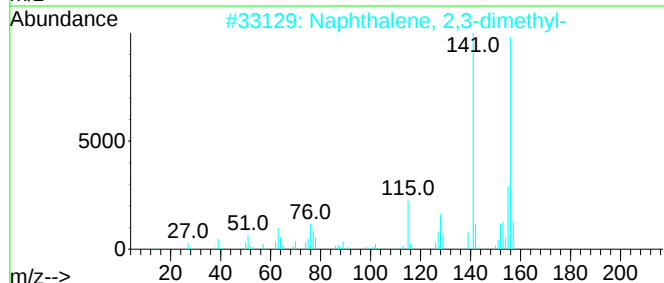
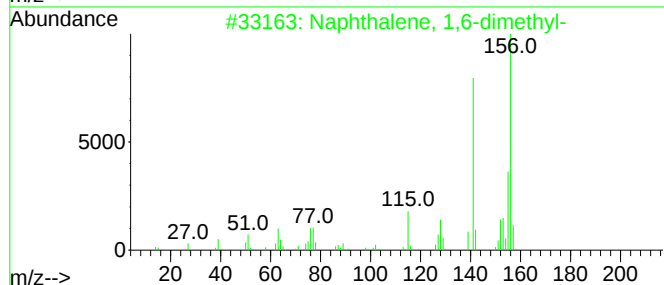
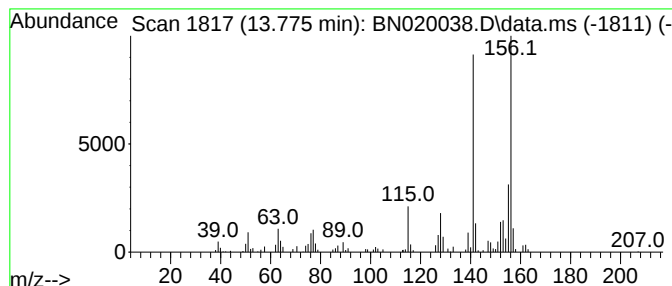
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN060622.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 Naphthalene, 1,6-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.775	2.25 ng/ul	166822	Acenaphthene-d10	14.451

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	98
2			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
3			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
4			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96
5			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060722\
Data File : BN020038.D
Acq On : 07 Jun 2022 21:06
Operator : CG/JU
Sample : N3171-08
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBTZ3

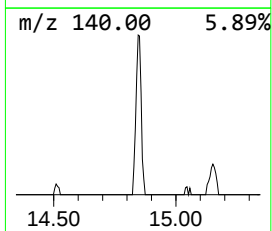
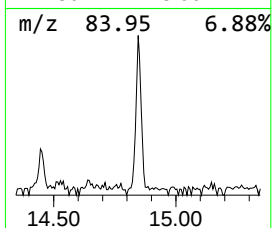
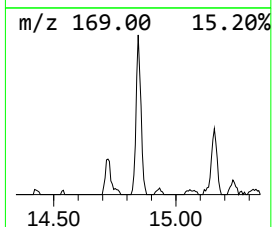
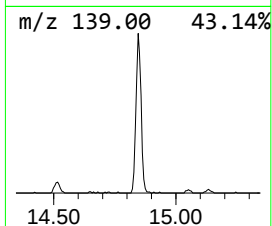
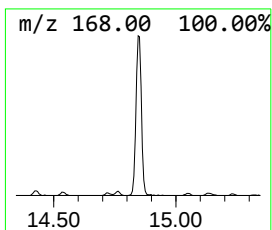
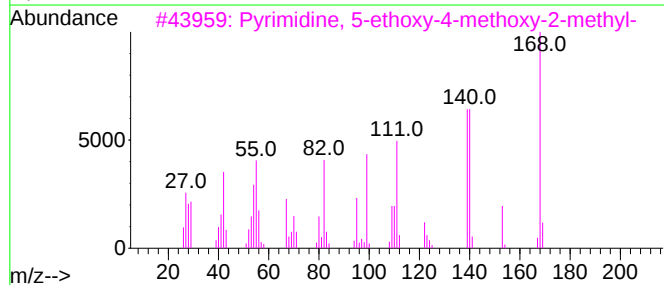
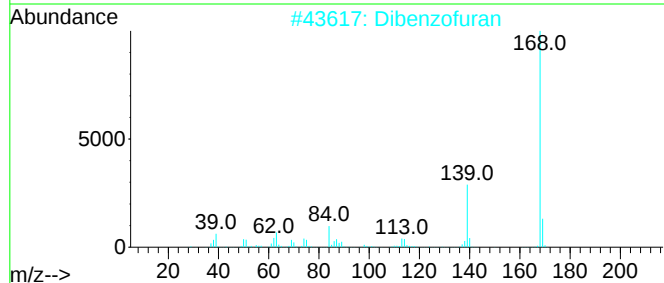
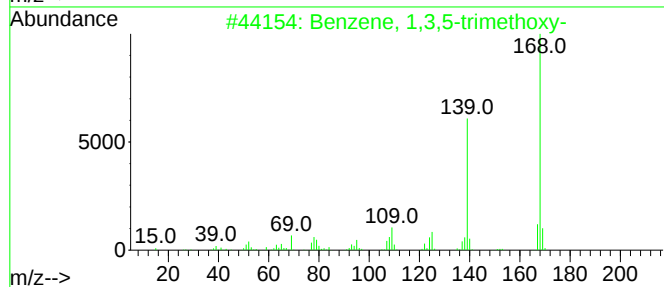
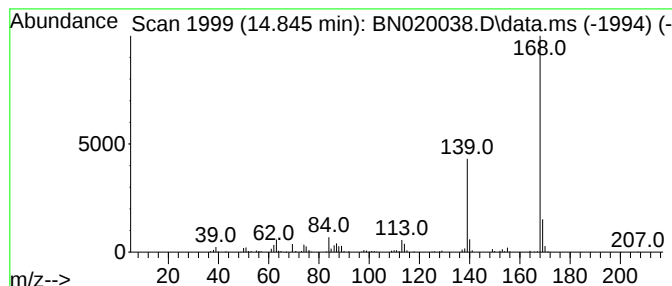
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN060622.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzene, 1,3,5-trimethoxy- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.845	3.79 ng/ul	280255	Acenaphthene-d10	14.451

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3,5-trimethoxy-	168	C9H12O3	000621-23-8	72
2			Dibenzofuran	168	C12H8O	000132-64-9	72
3			Pyrimidine, 5-ethoxy-4-methoxy-2...	168	C8H12N2O2	024614-12-8	64
4			Pyrimidine, 2-(dimethylamino)-5...	168	C6H8N4O2	014233-44-4	56
5			9H-Pyrido[3,4-b]indole	168	C11H8N2	000244-63-3	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060722\
Data File : BN020038.D
Acq On : 07 Jun 2022 21:06
Operator : CG/JU
Sample : N3171-08
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBTZ3

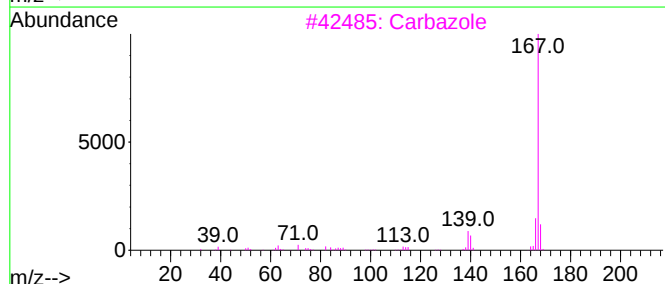
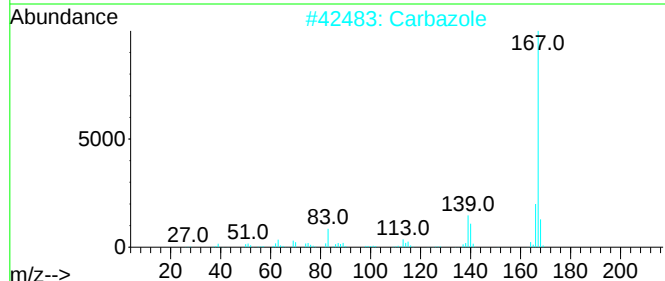
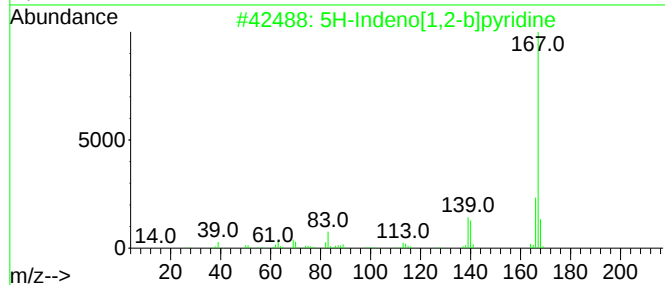
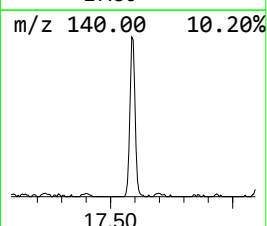
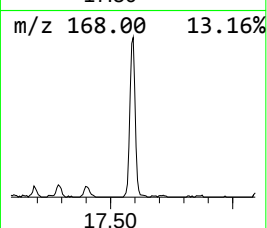
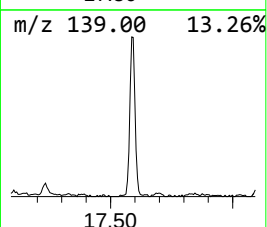
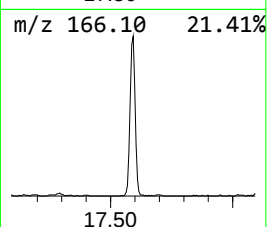
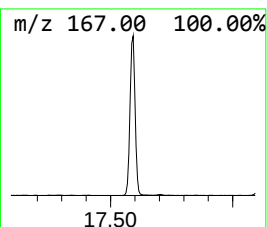
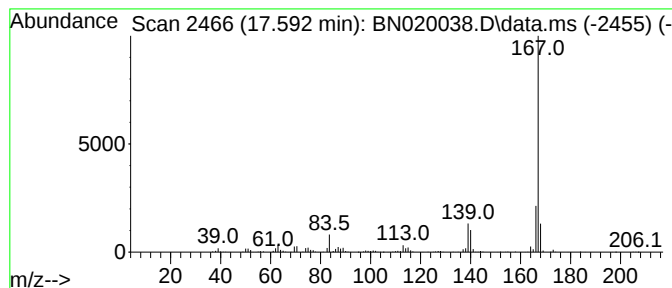
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN060622.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 5H-Indeno[1,2-b]pyridine Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.592	7.03 ng/ul	638761	Phenanthrene-d10	17.192

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5H-Indeno[1,2-b]pyridine	167	C12H9N	000244-99-5	97
2			Carbazole	167	C12H9N	000086-74-8	96
3			Carbazole	167	C12H9N	000086-74-8	91
4			5H-Indeno[1,2-b]pyridine	167	C12H9N	000244-99-5	80
5			5H-Indeno[1,2-b]pyridine	167	C12H9N	000244-99-5	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060722\
Data File : BN020038.D
Acq On : 07 Jun 2022 21:06
Operator : CG/JU
Sample : N3171-08
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBTZ3

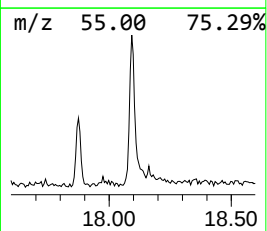
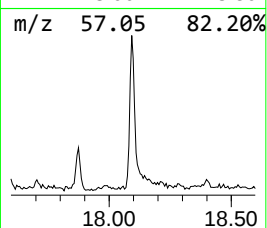
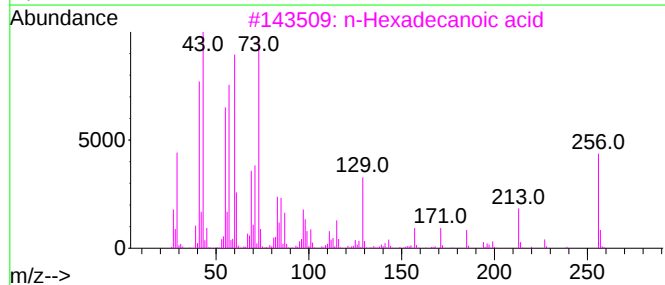
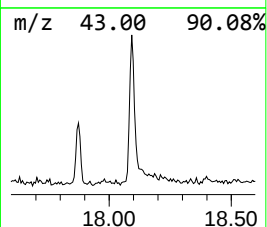
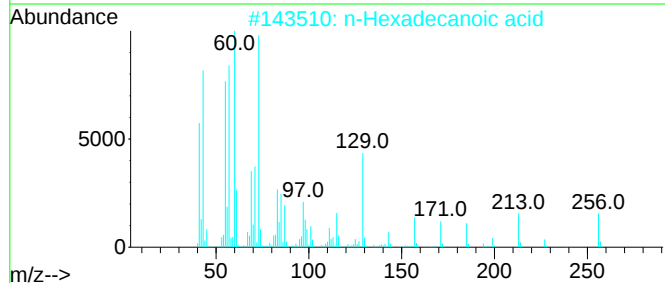
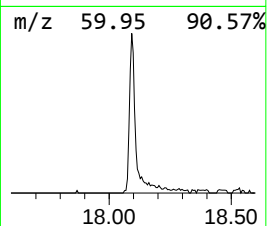
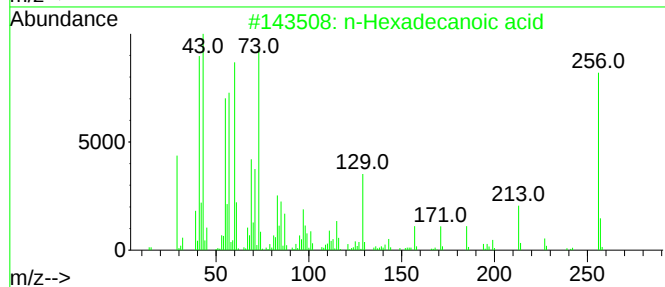
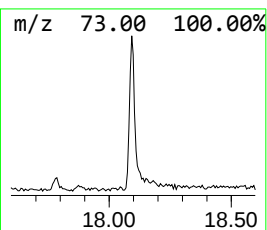
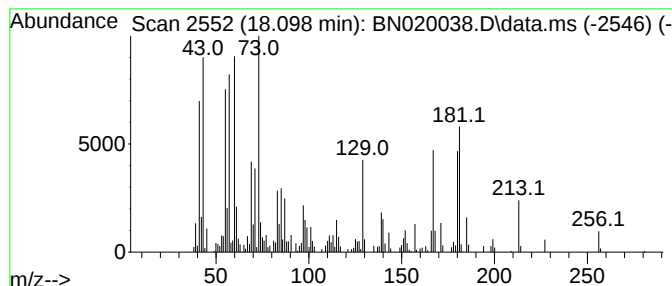
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN060622.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.098	3.56 ng/ul	323708	Phenanthrene-d10	17.192

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
5			Undecanoic acid	186	C11H22O2	000112-37-8	55



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060722\
Data File : BN020038.D
Acq On : 07 Jun 2022 21:06
Operator : CG/JU
Sample : N3171-08
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBTZ3

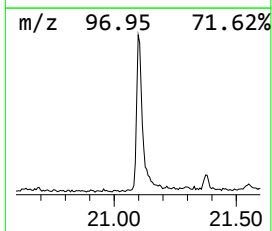
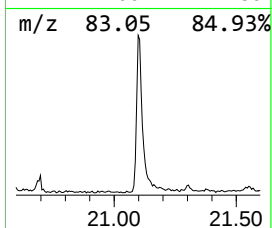
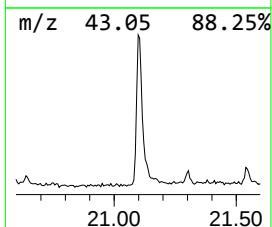
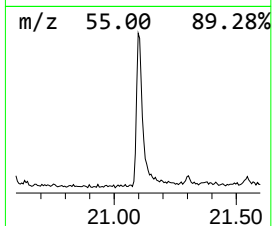
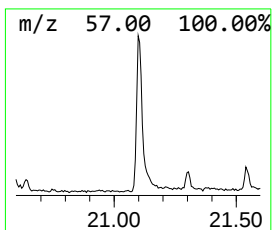
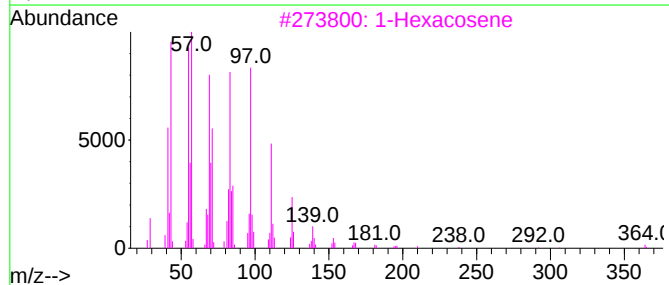
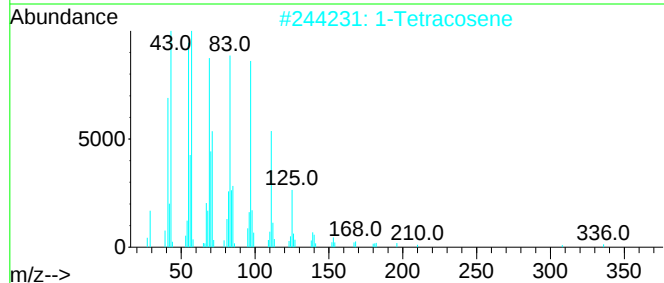
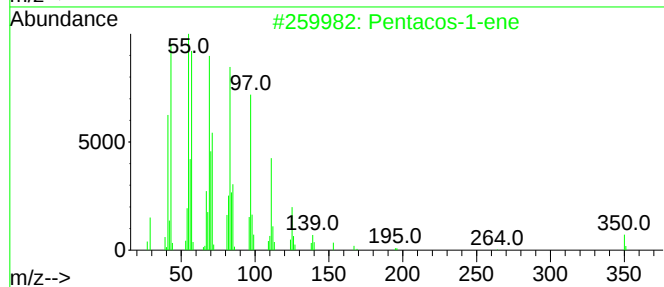
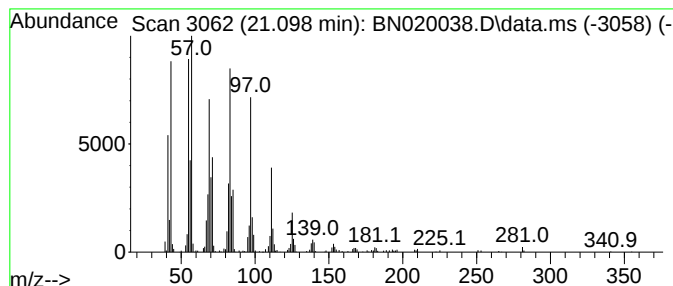
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN060622.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.098	4.59 ng/ul	475187	Chrysene-d12	21.380

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentacos-1-ene	350	C25H50	016980-85-1	94
2			1-Tetracosene	336	C24H48	010192-32-2	94
3			1-Hexacosene	364	C26H52	018835-33-1	94
4			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
5			1-Heneicosanol	312	C21H44O	015594-90-8	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060722\
 Data File : BN020038.D
 Acq On : 07 Jun 2022 21:06
 Operator : CG/JU
 Sample : N3171-08
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBTZ3

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

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Indane	8.199	24.4	ng/ul	821501	1	7.822	674868	20.0
Benzene, 2-ethe...	10.016	2.4	ng/ul	132126	2	10.610	1085620	20.0
unknown-01	12.134	3.9	ng/ul	209730	2	10.610	1085620	20.0
Naphthalene, 1,...	13.775	2.3	ng/ul	166822	3	14.451	1479720	20.0
Benzene, 1,3,5-...	14.845	3.8	ng/ul	280255	3	14.451	1479720	20.0
5H-Indeno[1,2-b...	17.592	7.0	ng/ul	638761	4	17.192	1818120	20.0
n-Hexadecanoic ...	18.098	3.6	ng/ul	323708	4	17.192	1818120	20.0
(DEL) Alkane: S...	21.098	4.6	ng/ul	475187	5	21.380	2068990	20.0