

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041422\
 Data File : BP009833.D
 Acq On : 14 Apr 2022 18:05
 Operator : CG/JU
 Sample : N2293-06
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 ETNN6

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_P\Methods\SFAM-EPA-BP040722.M
 Title : SVOA CALIBRATION

Signal : TIC: BP009833.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.123	3	6	22	rVB	627012	876120	27.93%	2.192%
2	3.381	47	50	60	rVB	12289	19311	0.62%	0.048%
3	3.805	118	122	136	rBV	83171	153260	4.89%	0.383%
4	4.134	174	178	185	rVB3	10087	17351	0.55%	0.043%
5	5.058	330	335	342	rVB3	6335	11389	0.36%	0.028%
6	5.211	357	361	365	rBV2	10279	16326	0.52%	0.041%
7	5.269	365	371	378	rVV	95720	146346	4.67%	0.366%
8	5.340	378	383	391	rVB6	8287	16513	0.53%	0.041%
9	6.640	597	604	613	rVB3	15425	26484	0.84%	0.066%
10	7.111	676	684	696	rBV	256920	434838	13.86%	1.088%
11	7.287	707	714	722	rVV	206447	334958	10.68%	0.838%
12	7.487	740	748	760	rBV	263031	432752	13.80%	1.082%
13	7.599	760	767	773	rVV10	4672	10090	0.32%	0.025%
14	7.963	822	829	837	rBV	438776	723632	23.07%	1.810%
15	8.575	928	933	939	rVB4	22918	35362	1.13%	0.088%
16	8.658	939	947	955	rBV	334889	560330	17.87%	1.402%
17	8.781	961	968	974	rBV	21828	37512	1.20%	0.094%
18	8.846	974	979	986	rVB2	19854	32751	1.04%	0.082%
19	9.116	1018	1025	1037	rVB	283505	467818	14.92%	1.170%
20	9.340	1059	1063	1070	rVB6	7826	12412	0.40%	0.031%
21	9.457	1076	1083	1090	rBV	83277	140881	4.49%	0.352%
22	9.534	1092	1096	1100	rVV2	16341	26643	0.85%	0.067%
23	9.575	1100	1103	1109	rVV3	9665	15942	0.51%	0.040%
24	9.769	1128	1136	1143	rBV	673523	1084435	34.58%	2.713%
25	9.846	1143	1149	1164	rVB	221363	388776	12.40%	0.972%
26	10.199	1202	1209	1216	rBV	5113	11591	0.37%	0.029%
27	10.263	1216	1220	1225	rBV4	7257	11627	0.37%	0.029%
28	10.381	1228	1240	1253	rBV	343303	599730	19.12%	1.500%
29	10.646	1278	1285	1290	rBV4	12599	22724	0.72%	0.057%
30	10.769	1297	1306	1314	rBV	664941	1125308	35.88%	2.815%
31	10.899	1318	1328	1338	rVV	333596	608171	19.39%	1.521%
32	10.993	1338	1344	1358	rVB2	65018	116039	3.70%	0.290%
33	11.116	1359	1365	1372	rVB3	12774	22900	0.73%	0.057%
34	11.316	1388	1399	1408	rBV	391117	669810	21.36%	1.675%
35	11.557	1429	1440	1445	rBV	8378	20829	0.66%	0.052%
36	11.722	1459	1468	1474	rBV6	10864	23936	0.76%	0.060%

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37	11.869	1487	1493	1499	rVB8	5704	8950	0.29%	0.022%
38	12.163	1537	1543	1546	rBV2	25779	41289	1.32%	0.103%
39	12.204	1546	1550	1557	rVB3	44367	73363	2.34%	0.184%
40	12.351	1570	1575	1580	rVB5	9686	15698	0.50%	0.039%
41	12.469	1587	1595	1613	rBV	2125860	3136400	100.00%	7.845%
42	12.604	1613	1618	1622	rVV7	10704	20164	0.64%	0.050%
43	12.828	1647	1656	1660	rBV5	15339	31643	1.01%	0.079%
44	12.875	1660	1664	1671	rVB2	20824	32585	1.04%	0.082%
45	13.322	1734	1740	1748	rVB2	5925	15667	0.50%	0.039%
46	13.434	1748	1759	1764	rBV3	15558	30411	0.97%	0.076%
47	13.522	1767	1774	1780	rVV	1408900	2011063	64.12%	5.031%
48	13.575	1780	1783	1789	rVV4	23947	38088	1.21%	0.095%
49	13.716	1803	1807	1812	rBV6	8323	11867	0.38%	0.030%
50	13.998	1849	1855	1864	rBV	799439	1128853	35.99%	2.824%
51	14.128	1873	1877	1885	rVB3	27913	43469	1.39%	0.109%
52	14.287	1893	1904	1916	rBV	733050	1125907	35.90%	2.816%
53	14.475	1931	1936	1939	rBV5	10874	18405	0.59%	0.046%
54	14.534	1943	1946	1949	rVV5	8365	11926	0.38%	0.030%
55	14.592	1949	1956	1962	rVV	1140972	1632371	52.05%	4.083%
56	14.645	1962	1965	1972	rVB	45195	62465	1.99%	0.156%
57	14.716	1972	1977	1980	rBV6	7063	10403	0.33%	0.026%
58	14.769	1980	1986	1998	rBV	369872	545914	17.41%	1.366%
59	15.004	2022	2026	2030	rVB3	17253	23149	0.74%	0.058%
60	15.240	2061	2066	2074	rVV5	47989	78305	2.50%	0.196%
61	15.404	2088	2094	2099	rBV7	3933	9232	0.29%	0.023%
62	15.451	2099	2102	2109	rVB6	7239	11357	0.36%	0.028%
63	15.587	2117	2125	2133	rBV	1033854	1578053	50.31%	3.947%
64	15.692	2133	2143	2154	rVB	326465	530897	16.93%	1.328%
65	15.828	2161	2166	2169	rBV3	12466	18651	0.59%	0.047%
66	15.875	2169	2174	2180	rVB7	15878	28981	0.92%	0.072%
67	16.316	2245	2249	2251	rBV4	7926	9657	0.31%	0.024%
68	16.416	2262	2266	2271	rVB3	9583	13527	0.43%	0.034%
69	16.769	2322	2326	2334	rVB2	27494	40874	1.30%	0.102%
70	16.957	2353	2358	2363	rVV8	8888	14158	0.45%	0.035%
71	17.016	2363	2368	2374	rVV10	8516	16503	0.53%	0.041%
72	17.110	2380	2384	2391	rBV8	10472	21872	0.70%	0.055%
73	17.163	2391	2393	2398	rVB5	6464	9654	0.31%	0.024%
74	17.339	2417	2423	2434	rVB	1416860	2056190	65.56%	5.143%
75	17.439	2434	2440	2447	rBV2	1313933	1878344	59.89%	4.699%
76	17.528	2452	2455	2462	rVB5	23238	35170	1.12%	0.088%

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77	17.639	2470	2474	2479	rVB8	6463	8920	0.28%	0.022%
78	17.816	2500	2504	2506	rBV2	8241	9493	0.30%	0.024%
79	17.851	2507	2510	2517	rVB8	9135	18946	0.60%	0.047%
80	17.957	2525	2528	2532	rVV6	6229	9545	0.30%	0.024%
81	18.010	2533	2537	2542	rVB8	14921	20241	0.65%	0.051%
82	18.222	2569	2573	2581	rBV3	81037	123309	3.93%	0.308%
83	18.733	2657	2660	2663	rBV5	6854	9286	0.30%	0.023%
84	18.980	2698	2702	2711	rBV2	33820	49338	1.57%	0.123%
85	19.122	2722	2726	2732	rBV2	119085	146114	4.66%	0.365%
86	19.245	2744	2747	2753	rVB6	22567	30451	0.97%	0.076%
87	19.310	2753	2758	2763	rBV3	32357	52432	1.67%	0.131%
88	19.451	2780	2782	2787	rBV6	15678	19371	0.62%	0.048%
89	19.669	2813	2819	2822	rBV	1913412	2580604	82.28%	6.455%
90	19.704	2822	2825	2834	rVB	2712402	3095670	98.70%	7.744%
91	19.957	2864	2868	2871	rBV	91910	108747	3.47%	0.272%
92	19.998	2871	2875	2880	rVB	131230	160338	5.11%	0.401%
93	20.116	2893	2895	2901	rVB5	21890	26045	0.83%	0.065%
94	20.222	2909	2913	2919	rVB	142693	171526	5.47%	0.429%
95	20.869	3019	3023	3027	rBV	136435	165294	5.27%	0.413%
96	21.045	3049	3053	3058	rVB	146227	188967	6.02%	0.473%
97	21.127	3063	3067	3072	rBV	241152	317967	10.14%	0.795%
98	21.422	3112	3117	3123	rBV	1814262	2379341	75.86%	5.952%
99	23.721	3501	3508	3516	rBV	1162595	2354221	75.06%	5.889%
100	23.886	3529	3536	3548	rVB2	1080100	2254846	71.89%	5.640%

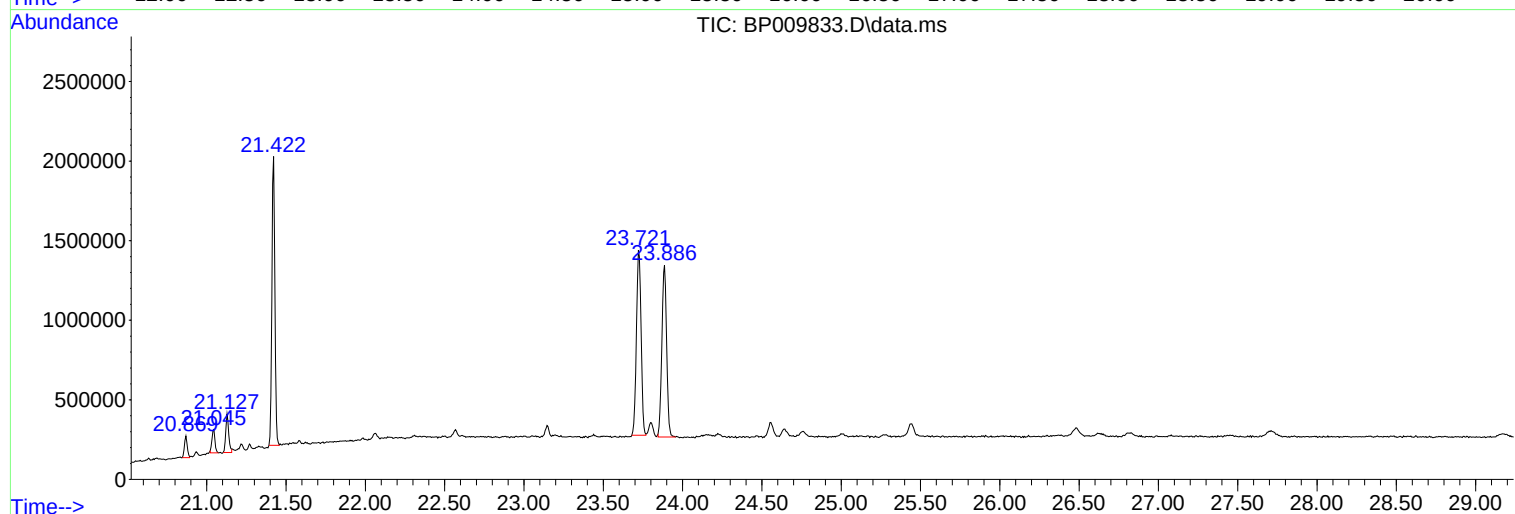
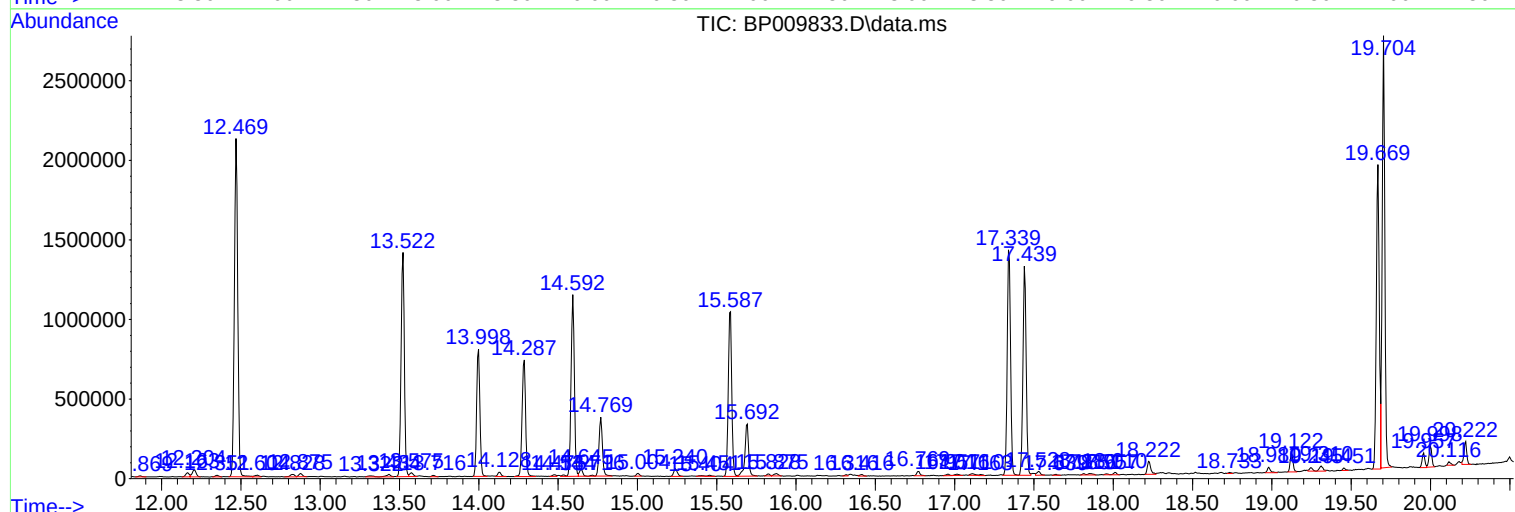
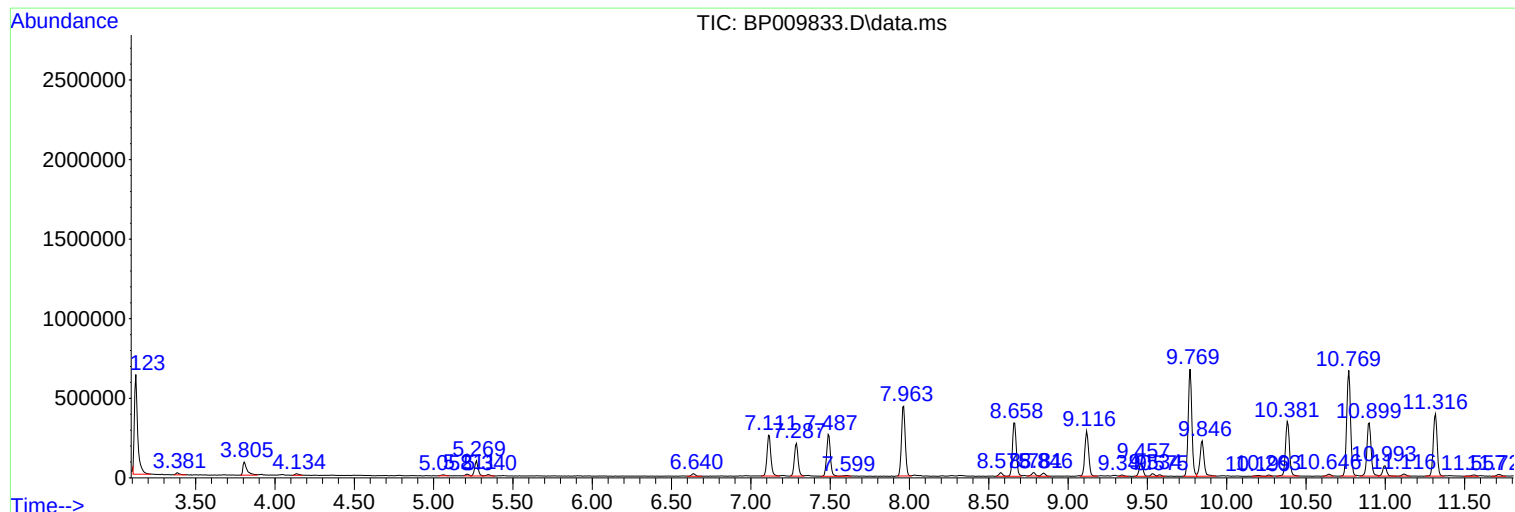
Sum of corrected areas: 39977384

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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



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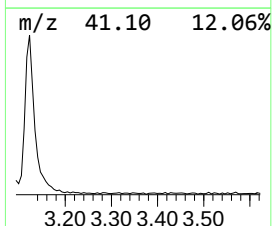
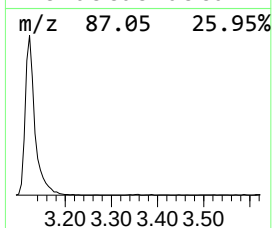
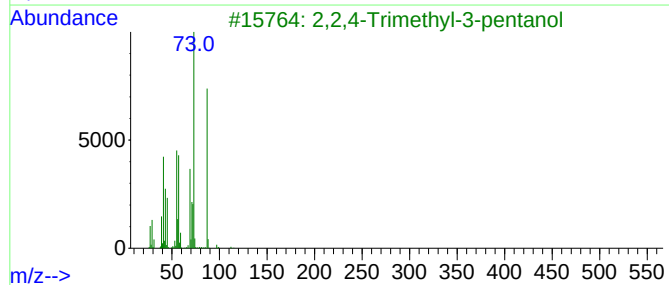
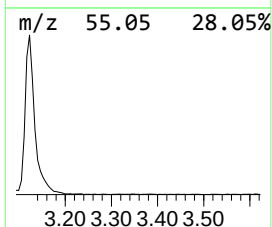
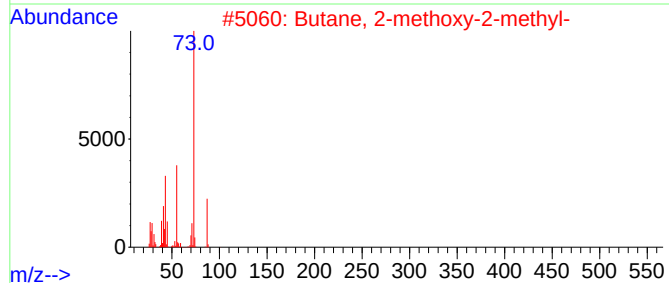
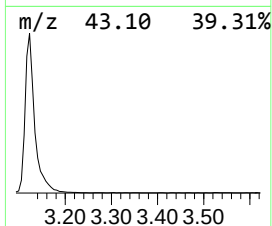
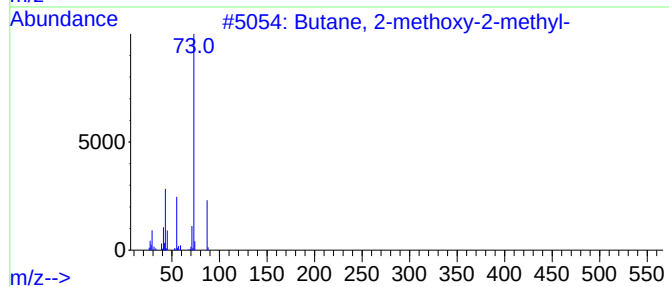
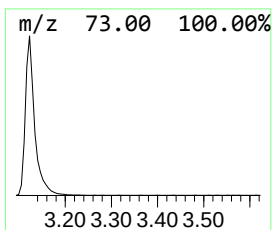
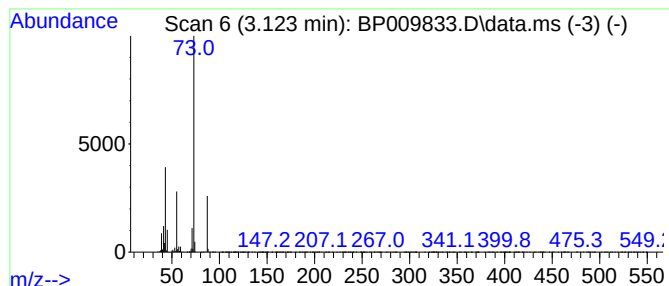
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.123	24.21 ng/ul	876120	1,4-Dichlorobenzene-d4	7.963

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	72
3			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
4			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
5			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12



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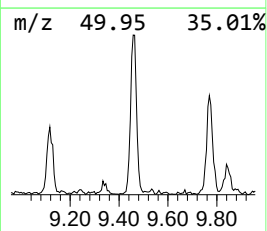
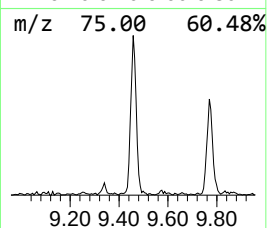
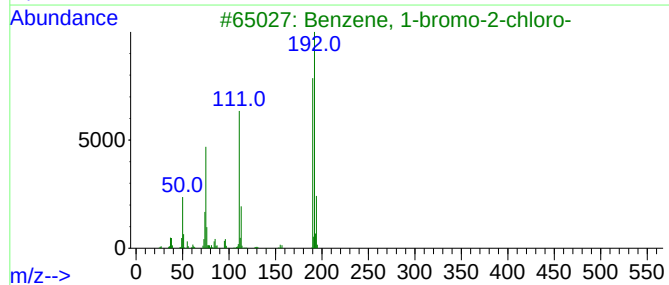
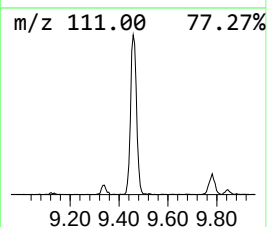
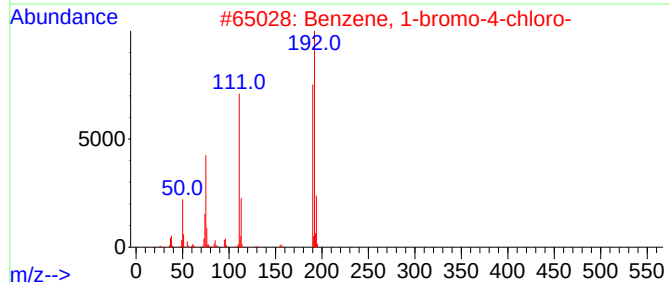
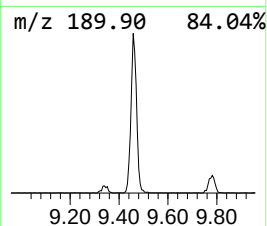
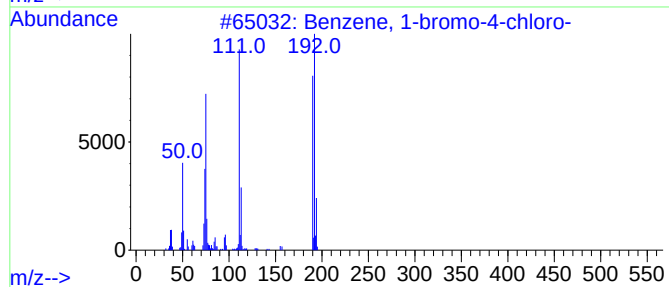
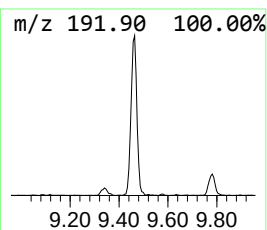
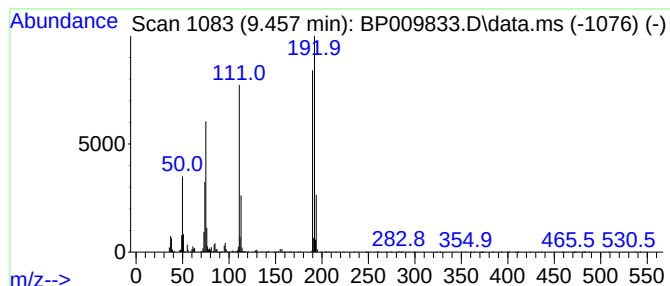
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzene, 1-bromo-4-chloro- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.457	2.50 ng/ul	140881	Naphthalene-d8	10.769

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	99
2			Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	97
3			Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	95
4			4-Chlorophenylselenol	192	C6H5ClSe	016645-10-6	95
5			Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	95



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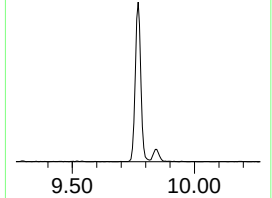
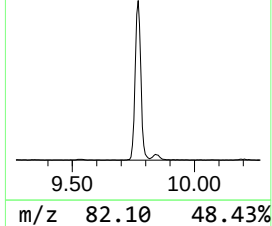
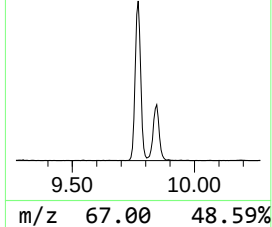
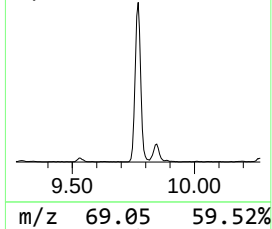
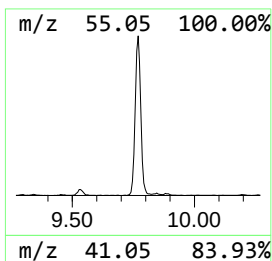
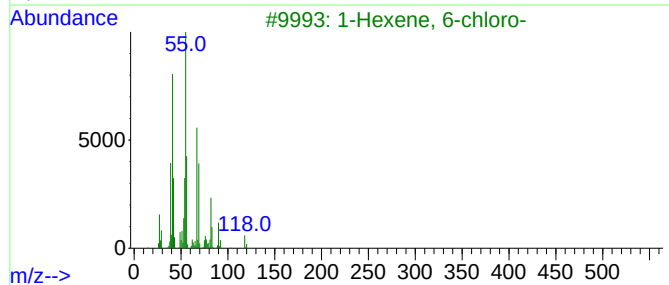
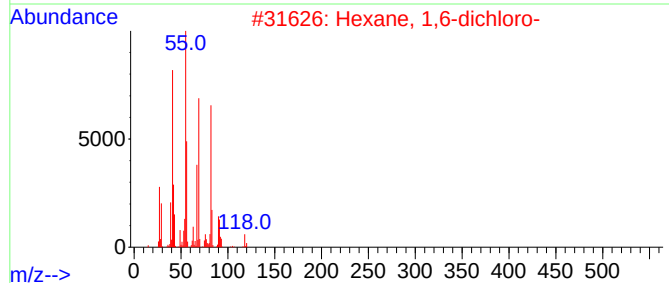
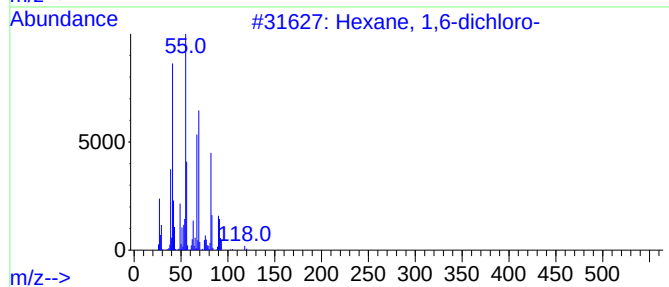
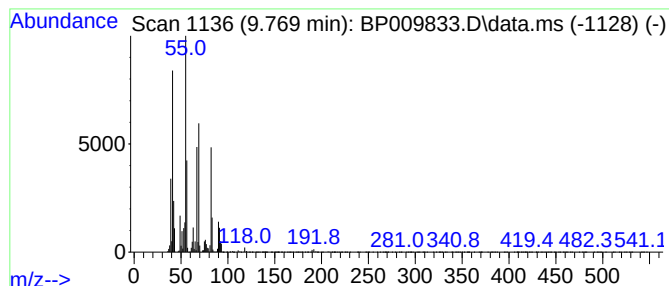
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 Hexane, 1,6-dichloro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.769	19.27 ng/ul	1084440	Naphthalene-d8	10.769

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 1,6-dichloro-	154	C6H12Cl2	002163-00-0	91
2			Hexane, 1,6-dichloro-	154	C6H12Cl2	002163-00-0	80
3			1-Hexene, 6-chloro-	118	C6H11Cl	000928-89-2	50
4			1-Hexanol, 6-chloro-	136	C6H13ClO	002009-83-8	42
5			(E)-6-Methylhept-4-en-1-ol	128	C8H16O	855901-81-4	40



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041422\
Data File : BP009833.D
Acq On : 14 Apr 2022 18:05
Operator : CG/JU
Sample : N2293-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ETNN6

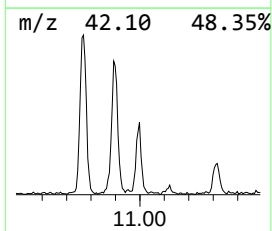
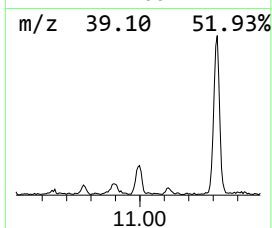
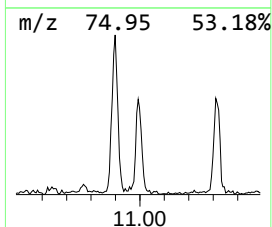
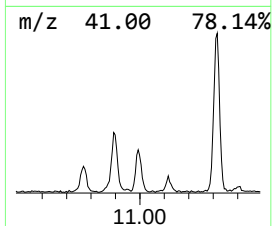
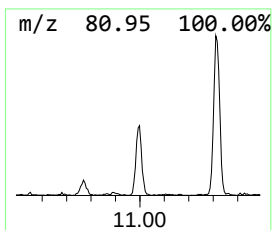
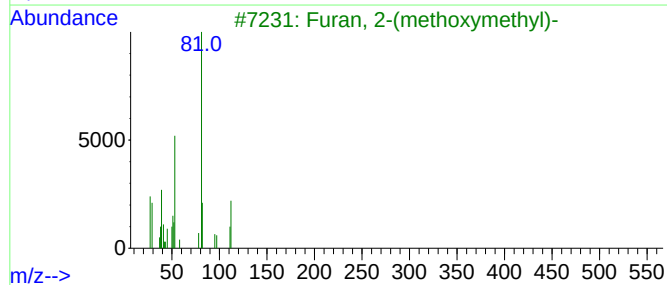
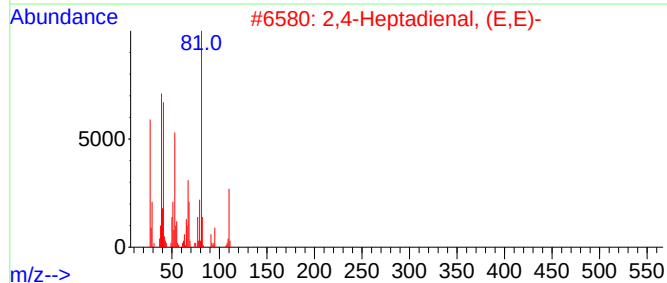
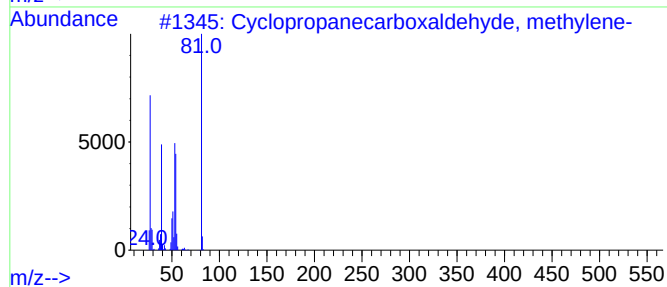
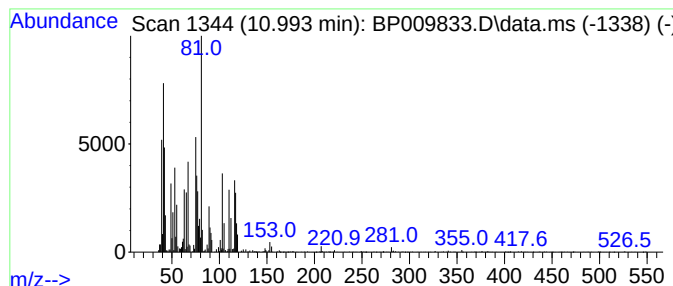
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 unknown-01 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.993	2.06 ng/ul	116039	Naphthalene-d8	10.769

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopropanecarboxaldehyde, meth...	82	C5H6O	142423-24-3	38
2			2,4-Heptadienal, (E,E)-	110	C7H10O	004313-03-5	32
3			Furan, 2-(methoxymethyl)-	112	C6H8O2	013679-46-4	32
4			2-Pentynoic acid ethyl ester	126	C7H10O2	055314-57-3	32
5			Furan, 2-propyl-	110	C7H10O	004229-91-8	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041422\
Data File : BP009833.D
Acq On : 14 Apr 2022 18:05
Operator : CG/JU
Sample : N2293-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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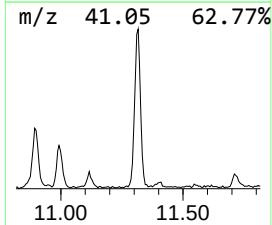
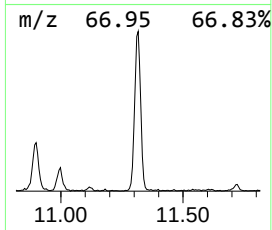
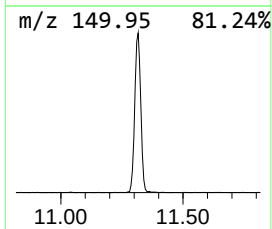
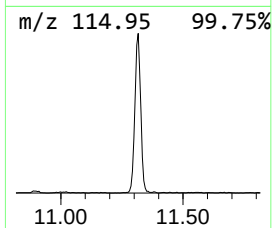
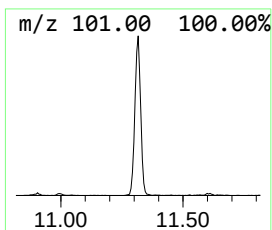
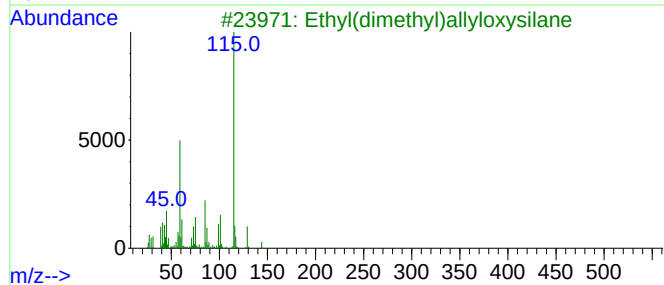
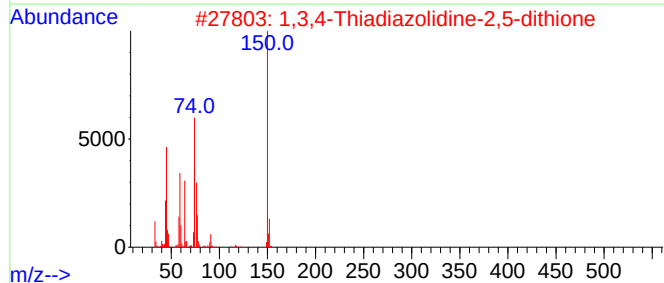
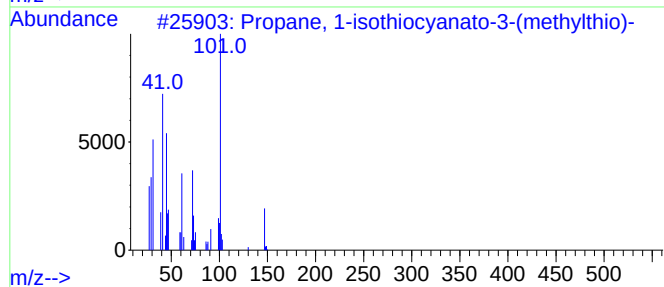
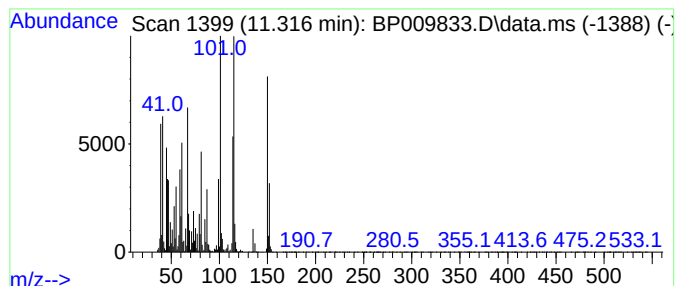
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-02 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.316	11.90 ng/ul	669810	Naphthalene-d8	10.769

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propane, 1-isothiocyanato-3-(met...	147	C5H9NS2	000505-79-3	22
2			1,3,4-Thiadiazolidine-2,5-dithione	150	C2H2N2S3	001072-71-5	14
3			Ethyl(dimethyl)allyloxysilane	144	C7H16OSi	1000278-59-3	14
4			1,2,4-Thiadiazole, 5-amino-	101	C2H3N3S	007552-07-0	14
5			1,3,4-Thiadiazolidine-2,5-dithione	150	C2H2N2S3	001072-71-5	14



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041422\
Data File : BP009833.D
Acq On : 14 Apr 2022 18:05
Operator : CG/JU
Sample : N2293-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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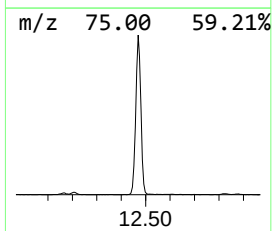
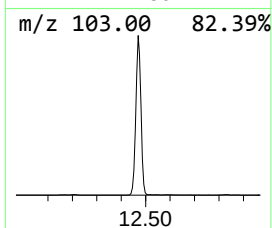
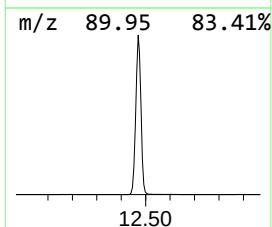
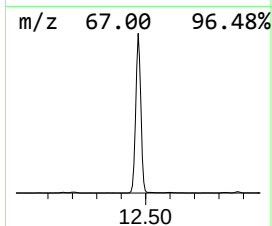
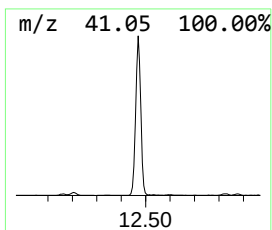
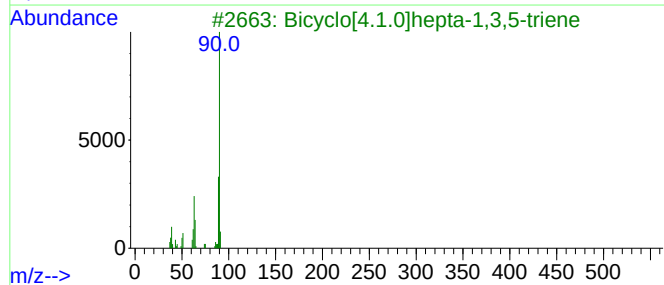
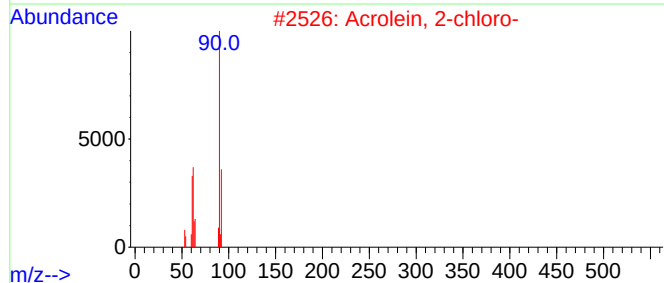
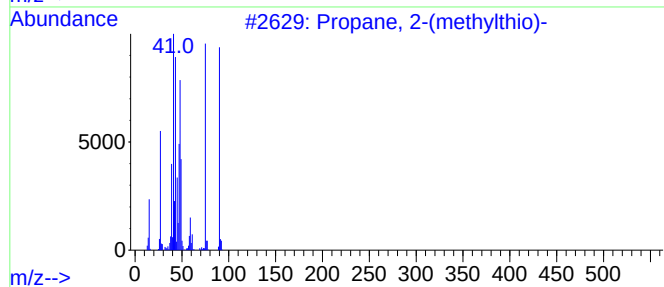
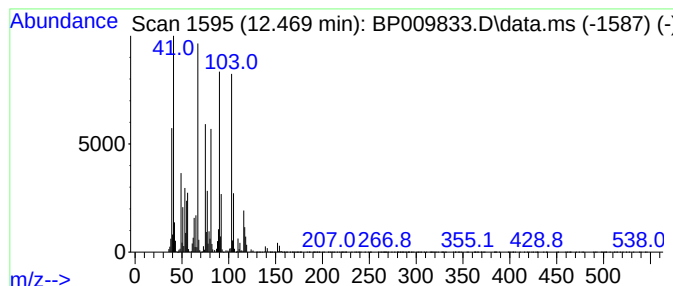
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 Propane, 2-(methylthio)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.469	55.74 ng/ul	3136400	Naphthalene-d8	10.769

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propane, 2-(methylthio)-	90	C4H10S	001551-21-9	50
2			Acrolein, 2-chloro-	90	C3H3ClO	000683-51-2	43
3			Bicyclo[4.1.0]hepta-1,3,5-triene	90	C7H6	004646-69-9	38
4			Propane, 2-(methylthio)-	90	C4H10S	001551-21-9	37
5			Thiourea, methyl-	90	C2H6N2S	000598-52-7	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041422\
Data File : BP009833.D
Acq On : 14 Apr 2022 18:05
Operator : CG/JU
Sample : N2293-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ETNN6

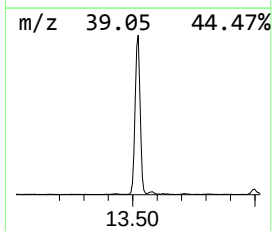
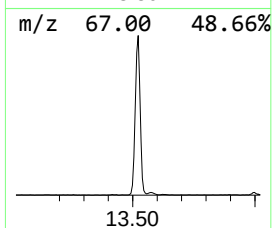
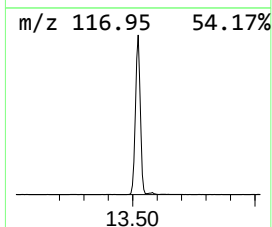
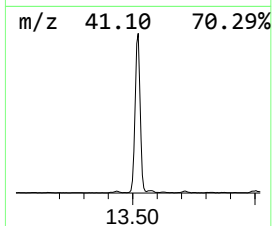
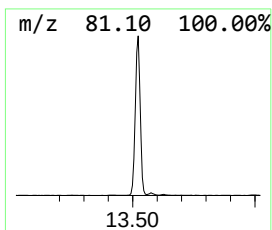
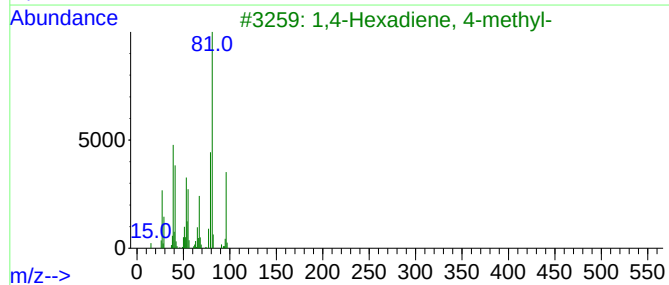
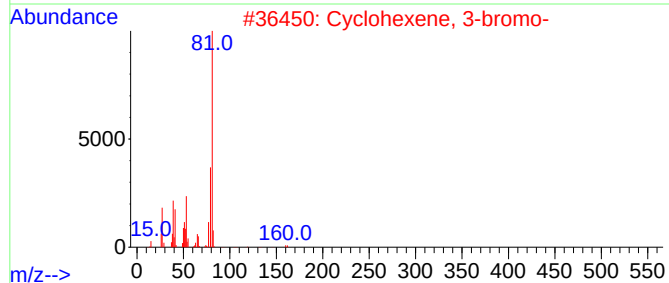
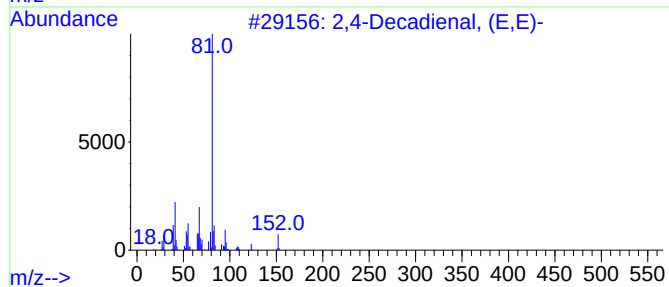
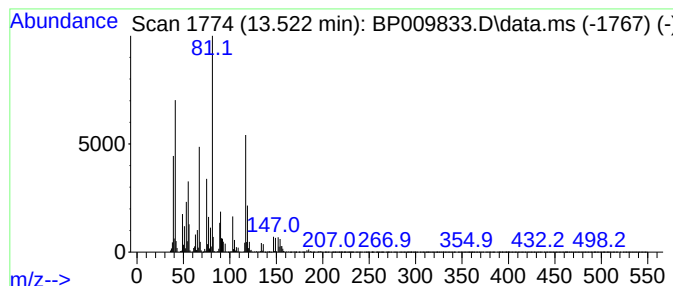
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-03 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.522	24.64 ng/ul	2011060	Acenaphthene-d10	14.592

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,4-Decadienal, (E,E)-	152	C10H16O	025152-84-5	35
2			Cyclohexene, 3-bromo-	160	C6H9Br	001521-51-3	27
3			1,4-Hexadiene, 4-methyl-	96	C7H12	001116-90-1	16
4			2,4-Dimethyl 1,4-pentadiene	96	C7H12	004161-65-3	16
5			Benzonitrile, 4-methyl-	117	C8H7N	000104-85-8	15



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041422\
Data File : BP009833.D
Acq On : 14 Apr 2022 18:05
Operator : CG/JU
Sample : N2293-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

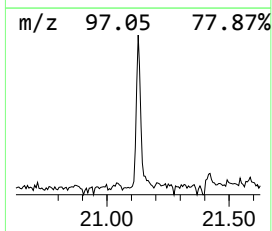
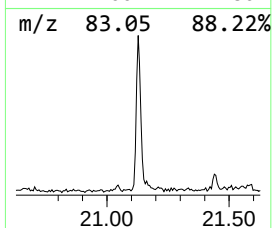
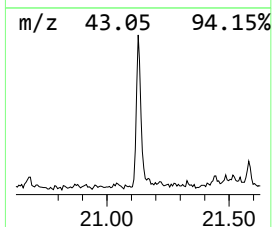
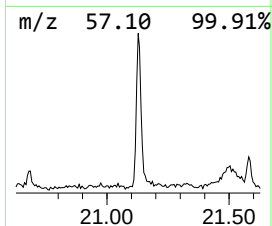
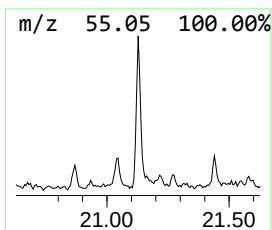
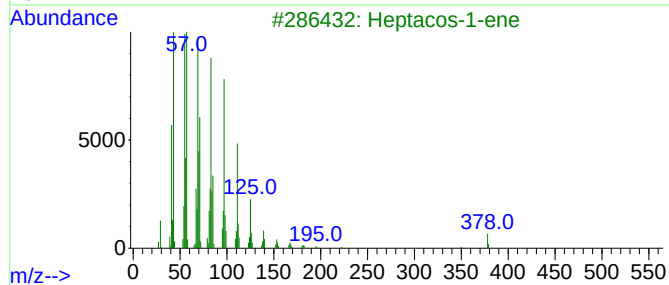
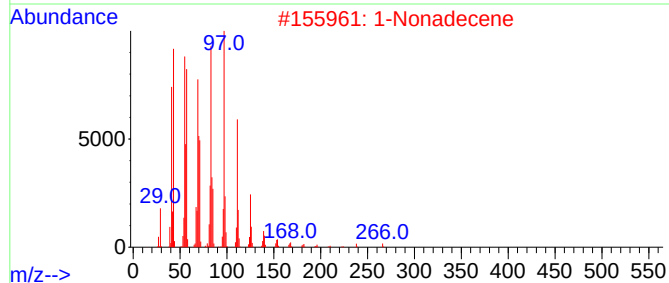
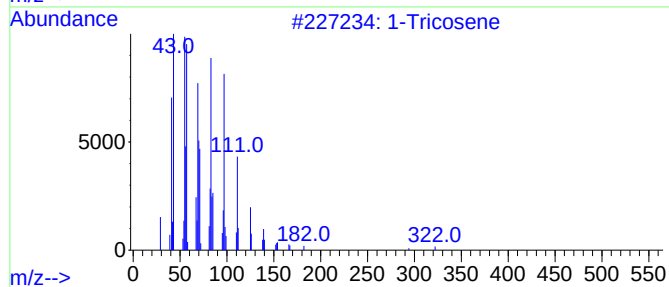
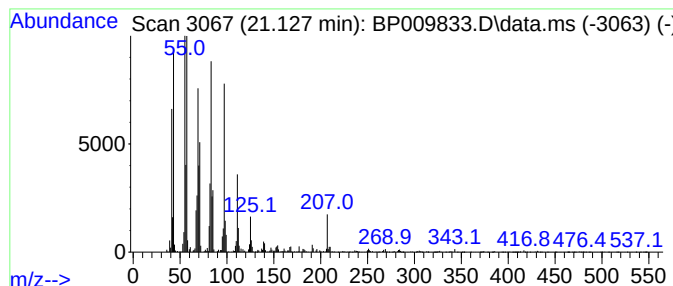
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.127	2.67 ng/ul	317967	Chrysene-d12	21.422	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Tricosene	322	C23H46	018835-32-0	94
2	1-Nonadecene	266	C19H38	018435-45-5	91
3	Heptacos-1-ene	378	C27H54	015306-27-1	91
4	1-Heneicosyl formate	340	C22H44O2	077899-03-7	91
5	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041422\
Data File : BP009833.D
Acq On : 14 Apr 2022 18:05
Operator : CG/JU
Sample : N2293-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ETNN6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.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
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Benzene, 1-brom...	9.457	2.5	ng/ul	140881	2	10.769	1125310	20.0
Hexane, 1,6-dic...	9.769	19.3	ng/ul	1084440	2	10.769	1125310	20.0
unknown-01	10.993	2.1	ng/ul	116039	2	10.769	1125310	20.0
unknown-02	11.316	11.9	ng/ul	669810	2	10.769	1125310	20.0
Propane, 2-(met...	12.469	55.7	ng/ul	3136400	2	10.769	1125310	20.0
unknown-03	13.522	24.6	ng/ul	2011060	3	14.592	1632370	20.0
(DEL) Alkane: S...	21.127	2.7	ng/ul	317967	5	21.422	2379340	20.0