

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042223\
 Data File : BM039582.D
 Acq On : 22 Apr 2023 18:41
 Operator : CG/JU
 Sample : 02440-12
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 COMQ7

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_M\Methods\SFAM-EPA-BM041823.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BM039582.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.213	26	30	41	rVB2	56755	93274	1.88%	0.174%
2	3.649	102	104	114	rBV	56194	91507	1.84%	0.170%
3	3.866	137	141	145	rVB	16540	24434	0.49%	0.046%
4	3.960	154	157	165	rVB2	21972	34390	0.69%	0.064%
5	4.707	280	284	290	rBV2	21256	45120	0.91%	0.084%
6	4.843	302	307	313	rBV4	15469	32664	0.66%	0.061%
7	5.372	395	397	402	rVB3	7140	9899	0.20%	0.018%
8	5.501	415	419	425	rVB	13071	19183	0.39%	0.036%
9	6.684	616	620	625	rVB8	2946	5196	0.10%	0.010%
10	6.884	650	654	659	rBV3	22235	51423	1.03%	0.096%
11	7.007	670	675	691	rVB	163264	376616	7.57%	0.701%
12	7.154	693	700	717	rBV	748807	1259289	25.31%	2.345%
13	7.354	727	734	755	rVB	716121	1242843	24.98%	2.315%
14	7.637	778	782	787	rVB7	4743	7959	0.16%	0.015%
15	7.819	803	813	819	rBV	673927	1124991	22.61%	2.095%
16	7.884	819	824	845	rVV	511309	984812	19.80%	1.834%
17	8.025	845	848	854	rVV4	10495	16298	0.33%	0.030%
18	8.078	854	857	863	rVV5	6420	12186	0.24%	0.023%
19	8.172	872	873	878	rVB5	5143	5940	0.12%	0.011%
20	8.336	896	901	905	rBV5	6764	11941	0.24%	0.022%
21	8.478	918	925	931	rBV3	16532	39090	0.79%	0.073%
22	8.548	931	937	955	rBV	552888	1145992	23.04%	2.134%
23	8.925	994	1001	1006	rBV5	7451	19365	0.39%	0.036%
24	8.989	1006	1012	1031	rVV	857124	1534094	30.84%	2.857%
25	9.148	1034	1039	1054	rVB	124196	265644	5.34%	0.495%
26	9.378	1075	1078	1084	rBV6	10037	19573	0.39%	0.036%
27	9.478	1092	1095	1101	rBV7	6808	13555	0.27%	0.025%
28	9.713	1128	1135	1153	rBV	672967	1239175	24.91%	2.308%
29	9.995	1177	1183	1197	rBV2	75644	249226	5.01%	0.464%
30	10.266	1222	1229	1252	rBV	768903	1677590	33.72%	3.124%
31	10.430	1255	1257	1263	rVV7	10261	19542	0.39%	0.036%
32	10.478	1263	1265	1269	rVB5	7121	11280	0.23%	0.021%
33	10.625	1283	1290	1306	rBV	899969	1542679	31.01%	2.873%
34	10.778	1310	1316	1338	rBV	278059	705182	14.18%	1.313%
35	11.301	1399	1405	1414	rBV2	14212	41073	0.83%	0.076%
36	11.472	1428	1434	1444	rBV4	29615	85774	1.72%	0.160%

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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
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37	11.542	1444	1446	1449	rVV4	5901	6116	0.12%	0.011%
38	11.619	1453	1459	1462	rBV8	5304	9370	0.19%	0.017%
39	11.983	1513	1521	1524	rBV10	3262	8401	0.17%	0.016%
40	12.042	1526	1531	1540	rVB5	11426	19912	0.40%	0.037%
41	12.142	1544	1548	1553	rBV6	6316	11615	0.23%	0.022%
42	12.224	1557	1562	1574	rBV	18371	45510	0.91%	0.085%
43	12.360	1576	1585	1587	rBV4	11373	25003	0.50%	0.047%
44	12.677	1635	1639	1643	rBV6	4347	6084	0.12%	0.011%
45	12.719	1643	1646	1648	rVB3	5177	5379	0.11%	0.010%
46	12.783	1651	1657	1663	rBV	54638	110357	2.22%	0.206%
47	12.854	1666	1669	1674	rVB4	16572	26152	0.53%	0.049%
48	12.995	1690	1693	1697	rVB3	7025	8890	0.18%	0.017%
49	13.071	1697	1706	1712	rBV7	12876	38118	0.77%	0.071%
50	13.289	1737	1743	1746	rBV2	27583	40628	0.82%	0.076%
51	13.330	1747	1750	1752	rVV2	8842	11337	0.23%	0.021%
52	13.371	1753	1757	1761	rVV3	18891	30136	0.61%	0.056%
53	13.448	1768	1770	1780	rVB8	15700	26451	0.53%	0.049%
54	13.648	1800	1804	1805	rBV4	6562	7665	0.15%	0.014%
55	13.713	1805	1815	1816	rBV7	17891	42029	0.84%	0.078%
56	13.889	1837	1845	1858	rBV	1889768	2775846	55.80%	5.170%
57	13.989	1859	1862	1872	rVB6	29020	57647	1.16%	0.107%
58	14.130	1881	1886	1887	rBV	75448	90277	1.81%	0.168%
59	14.171	1887	1893	1903	rVV	2200856	3407030	68.49%	6.345%
60	14.260	1904	1908	1913	rVB	65356	89537	1.80%	0.167%
61	14.366	1923	1926	1929	rVB	14551	15925	0.32%	0.030%
62	14.477	1939	1945	1956	rVB	1274340	1904176	38.28%	3.546%
63	14.707	1980	1984	1988	rBV3	11831	18295	0.37%	0.034%
64	14.771	1989	1995	1998	rBV5	30569	65295	1.31%	0.122%
65	14.871	2008	2012	2017	rBV4	21931	42407	0.85%	0.079%
66	14.995	2028	2033	2038	rVB3	37638	65206	1.31%	0.121%
67	15.054	2040	2043	2046	rVB3	29663	32634	0.66%	0.061%
68	15.236	2070	2074	2078	rBV5	34131	64066	1.29%	0.119%
69	15.401	2096	2102	2107	rVB6	45110	104754	2.11%	0.195%
70	15.471	2107	2114	2124	rBV	2743710	4056574	81.55%	7.555%
71	15.612	2133	2138	2149	rBV	683684	1115996	22.43%	2.078%
72	15.712	2151	2155	2156	rBV3	37659	54917	1.10%	0.102%
73	15.842	2172	2177	2181	rBV	227935	323163	6.50%	0.602%
74	16.071	2210	2216	2221	rBV6	39932	95428	1.92%	0.178%
75	16.124	2221	2225	2231	rVB	115462	177968	3.58%	0.331%
76	16.230	2240	2243	2250	rBV3	35879	65640	1.32%	0.122%

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77	16.354	2261	2264	2269	rBV4	46406	86317	1.74%	0.161%
78	16.530	2290	2294	2296	rBV3	16191	22727	0.46%	0.042%
79	16.642	2311	2313	2317	rVB2	16556	15228	0.31%	0.028%
80	16.754	2327	2332	2339	rBV3	24875	60864	1.22%	0.113%
81	16.977	2367	2370	2373	rBV5	11212	17248	0.35%	0.032%
82	17.230	2407	2413	2423	rBV	1486104	2119538	42.61%	3.947%
83	17.330	2424	2430	2441	rVV	2895847	4233797	85.11%	7.885%
84	17.430	2443	2447	2454	rVV	226515	371787	7.47%	0.692%
85	17.495	2454	2458	2473	rVB	464176	785133	15.78%	1.462%
86	17.895	2523	2526	2535	rVB3	49055	85437	1.72%	0.159%
87	18.130	2561	2566	2575	rBV	318357	545721	10.97%	1.016%
88	18.342	2598	2602	2612	rBV	98249	165047	3.32%	0.307%
89	18.912	2697	2699	2705	rBV6	14436	23889	0.48%	0.044%
90	19.195	2743	2747	2751	rBV3	42749	58838	1.18%	0.110%
91	19.265	2755	2759	2762	rBV2	53692	83792	1.68%	0.156%
92	19.412	2780	2784	2793	rBV3	208138	430456	8.65%	0.802%
93	19.630	2815	2821	2829	rBV	3650852	4974559	100.00%	9.264%
94	20.512	2967	2971	2983	rBV	470934	939251	18.88%	1.749%
95	21.136	3074	3077	3084	rVB4	123277	198569	3.99%	0.370%
96	21.424	3121	3126	3139	rBV	1610169	2232754	44.88%	4.158%
97	23.635	3496	3502	3515	rVB2	2406700	4913409	98.77%	9.150%
98	23.788	3523	3528	3538	rVB2	1077555	2171366	43.65%	4.044%

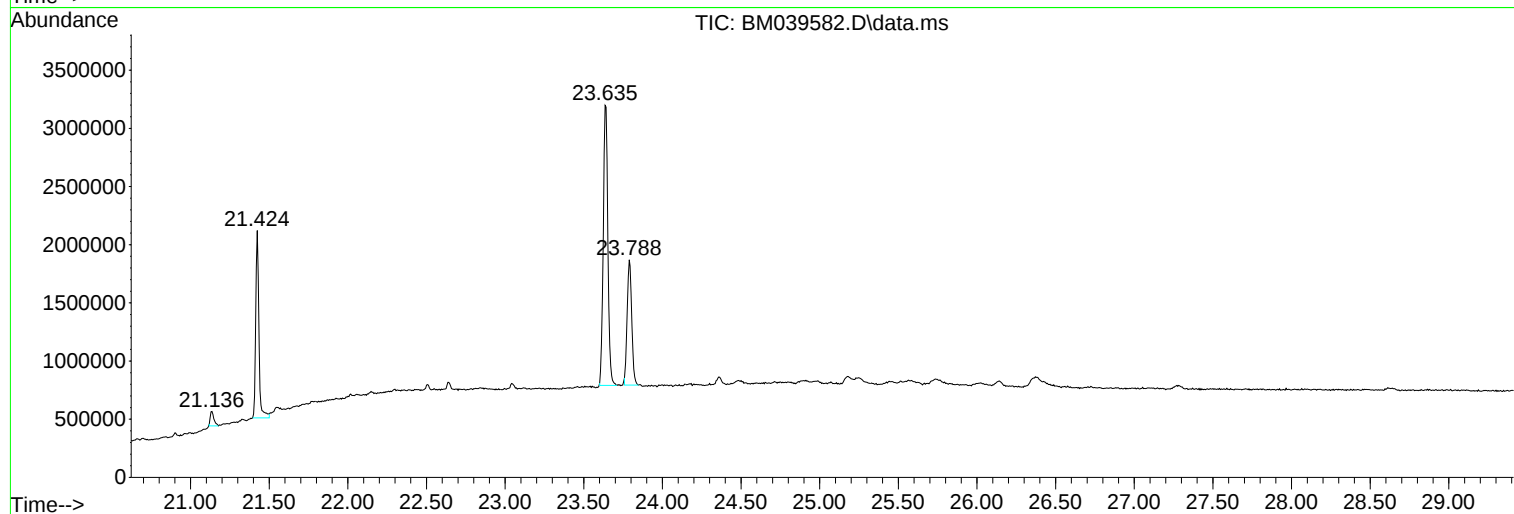
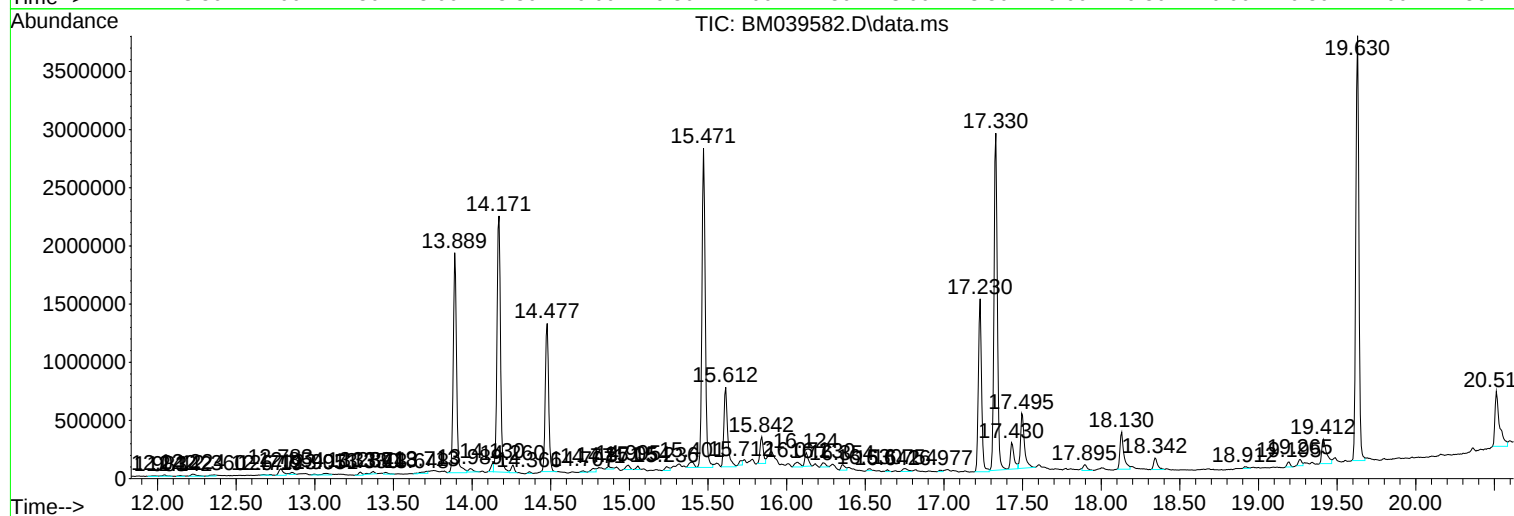
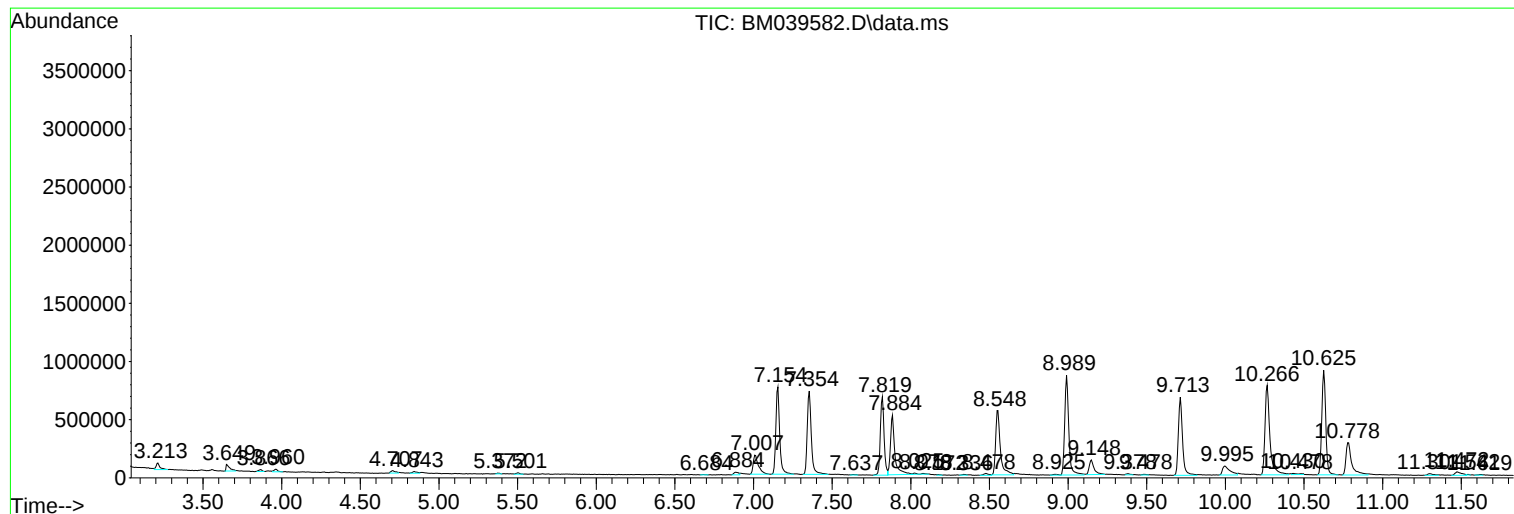
Sum of corrected areas: 53696460

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
Quant Title : SVOA CALIBRATION

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



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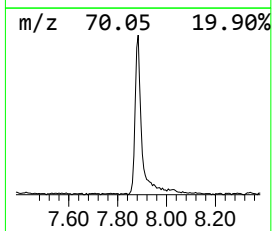
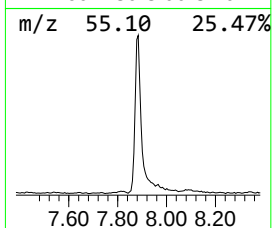
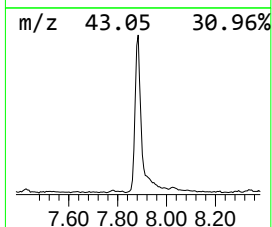
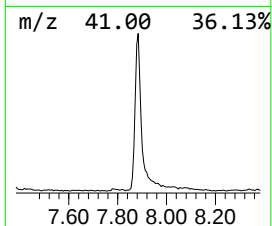
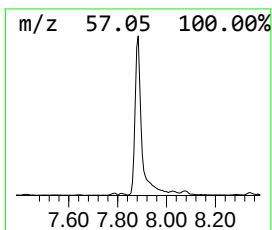
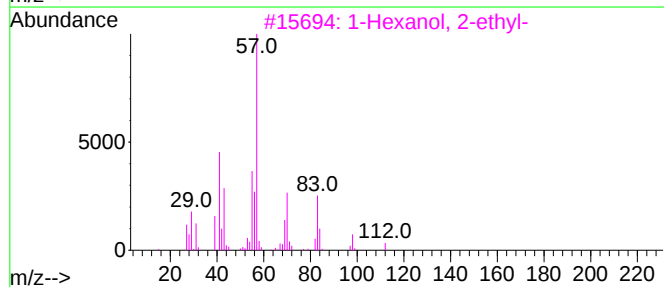
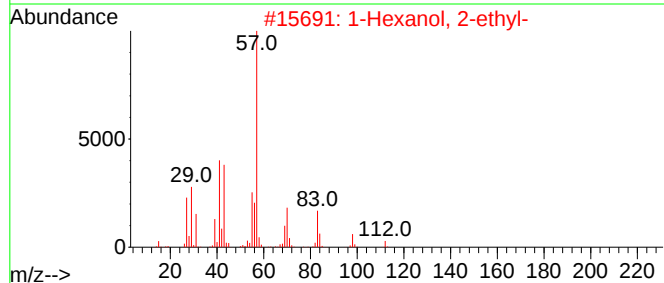
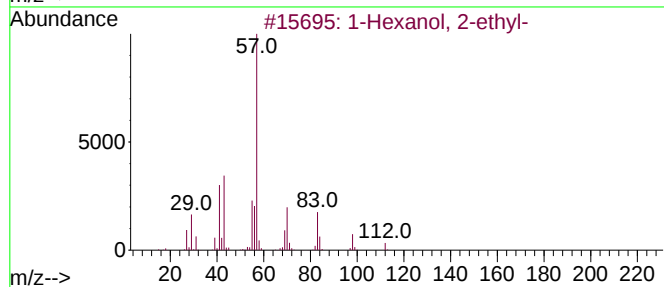
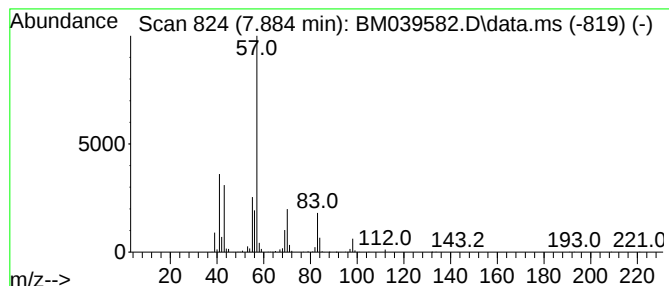
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 1-Hexanol, 2-ethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.884	17.51 ng/ul	984812	1,4-Dichlorobenzene-d4	7.819

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	83
2	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	83
3	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	72
4	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	72
5	1-Isobutoxy-2-ethylhexane			186	C12H26O	1000139-90-3	64



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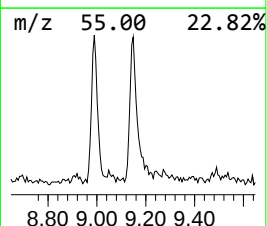
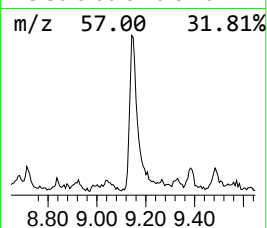
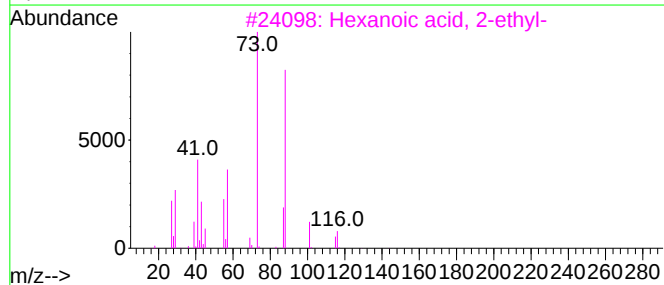
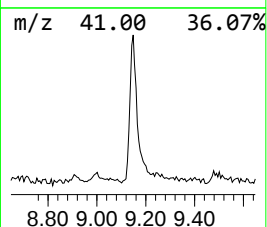
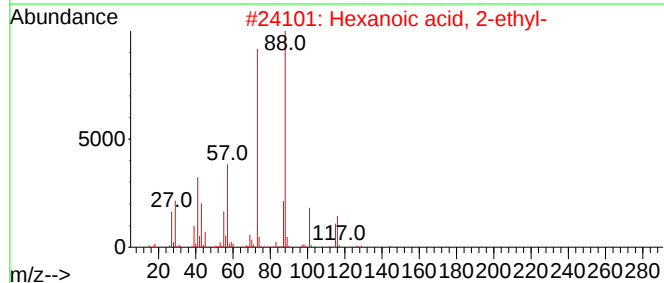
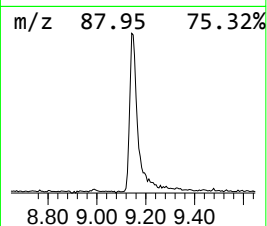
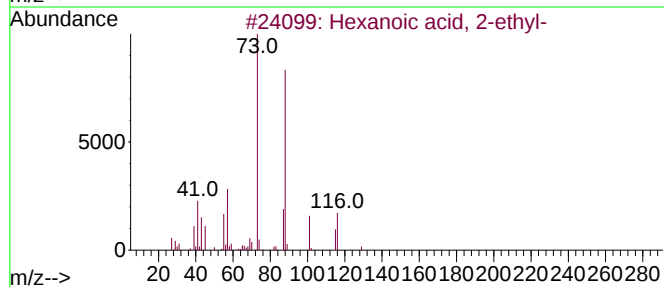
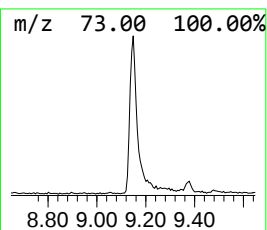
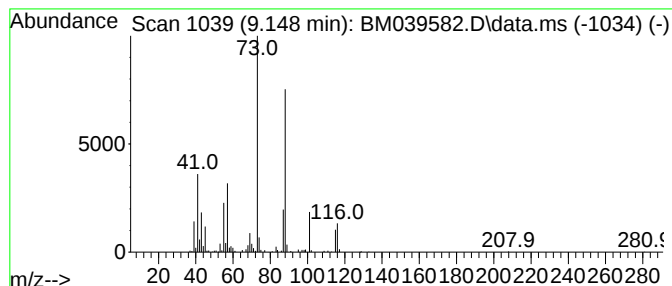
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Hexanoic acid, 2-ethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.148	4.72 ng/ul	265644	1,4-Dichlorobenzene-d4	7.819

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	90
2			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	86
3			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	78
4			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	78
5			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	72



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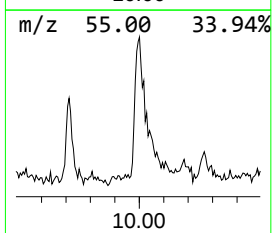
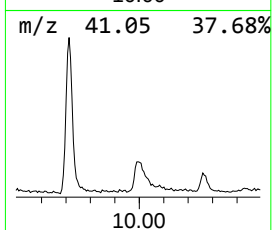
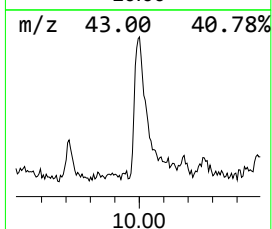
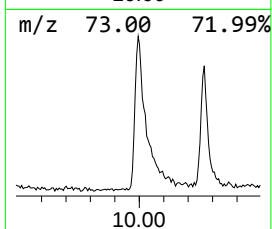
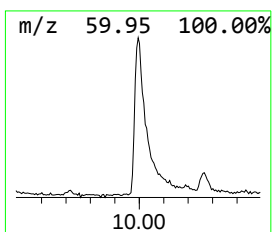
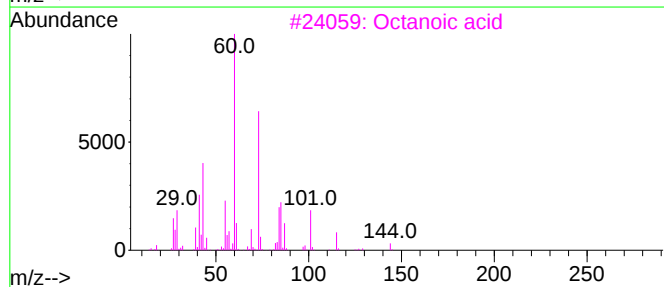
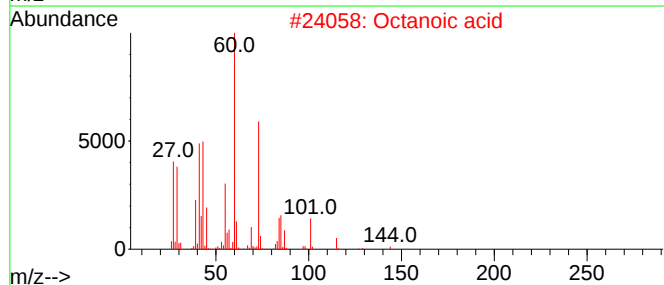
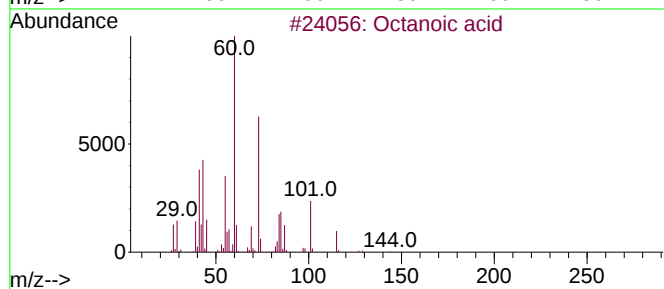
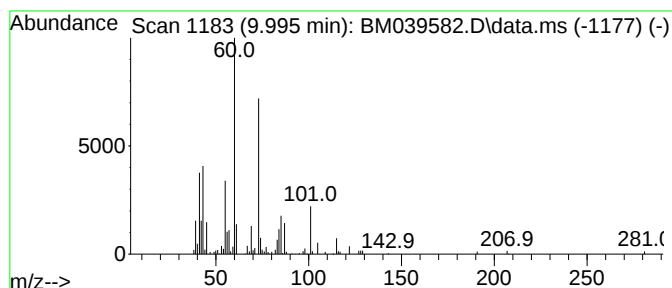
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Octanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.995	3.23 ng/ul	249226	Naphthalene-d8	10.625

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octanoic acid			144	C8H16O2	000124-07-2	86
2	Octanoic acid			144	C8H16O2	000124-07-2	86
3	Octanoic acid			144	C8H16O2	000124-07-2	78
4	Hexanoic acid			116	C6H12O2	000142-62-1	58
5	Hexanoic acid			116	C6H12O2	000142-62-1	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042223\
Data File : BM039582.D
Acq On : 22 Apr 2023 18:41
Operator : CG/JU
Sample : 02440-12
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
COMQ7

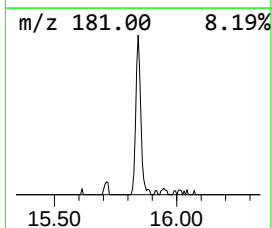
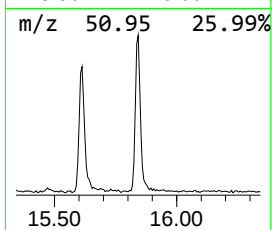
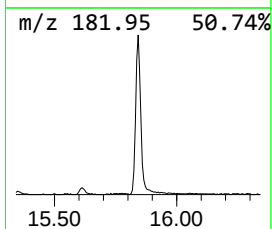
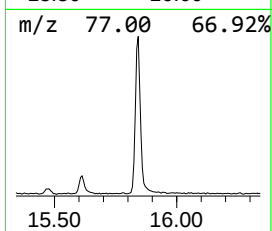
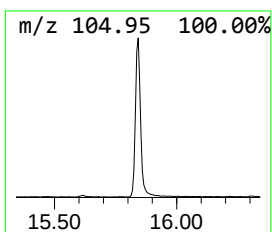
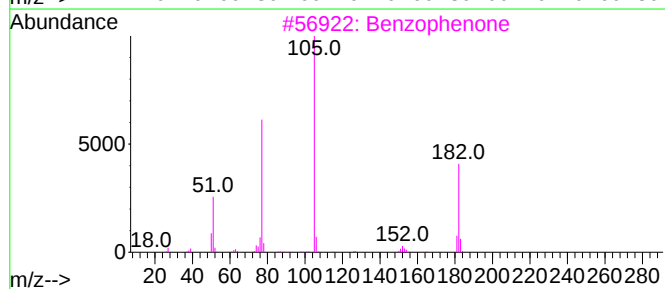
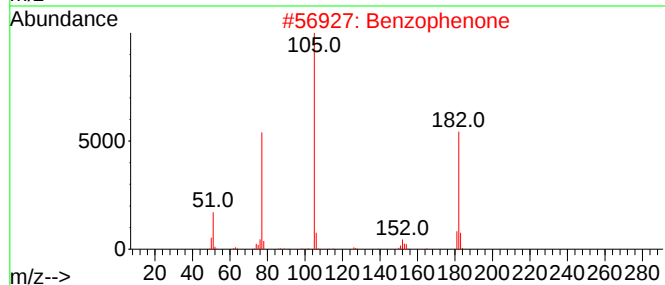
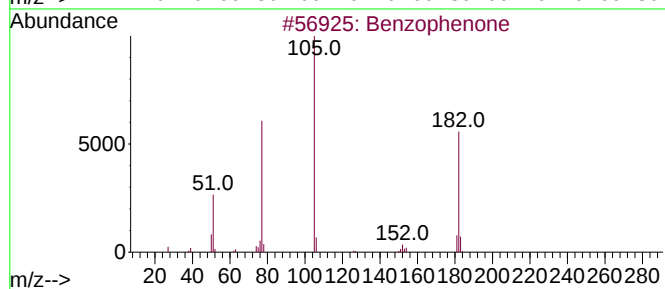
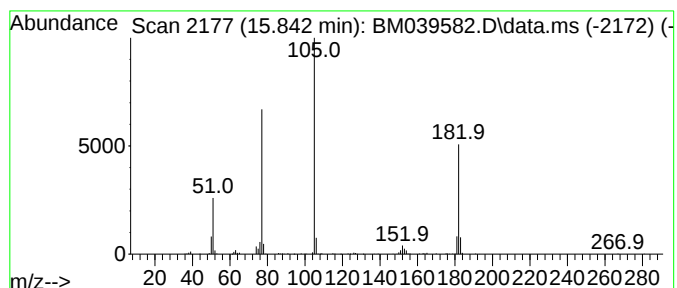
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Benzophenone Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.842	3.39 ng/ul	323163	Acenaphthene-d10	14.477

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzophenone	182	C13H10O	000119-61-9	95
3			Benzophenone	182	C13H10O	000119-61-9	94
4			Benzophenone	182	C13H10O	000119-61-9	93
5			Benzophenone	182	C13H10O	000119-61-9	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042223\
Data File : BM039582.D
Acq On : 22 Apr 2023 18:41
Operator : CG/JU
Sample : 02440-12
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ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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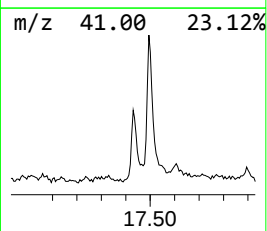
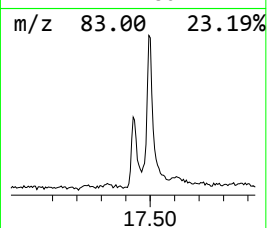
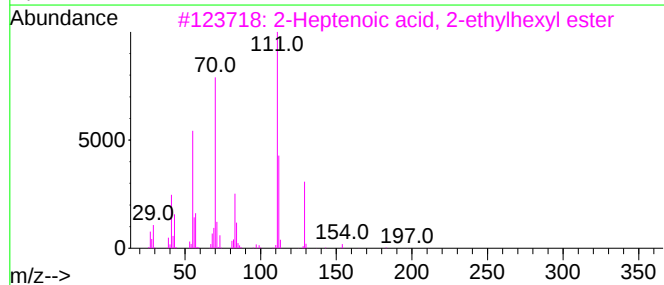
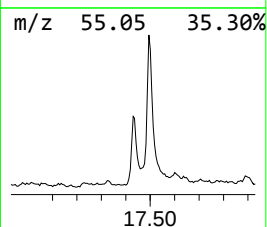
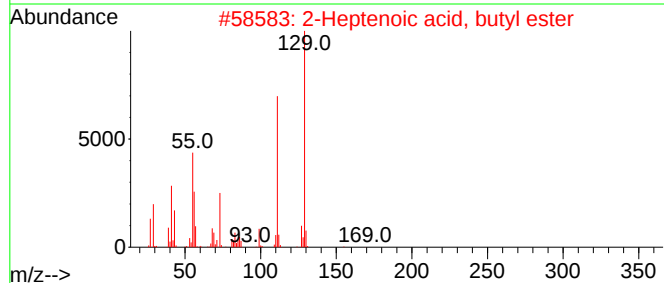
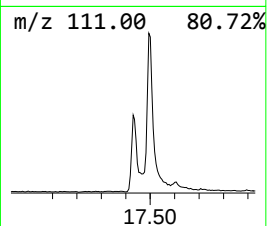
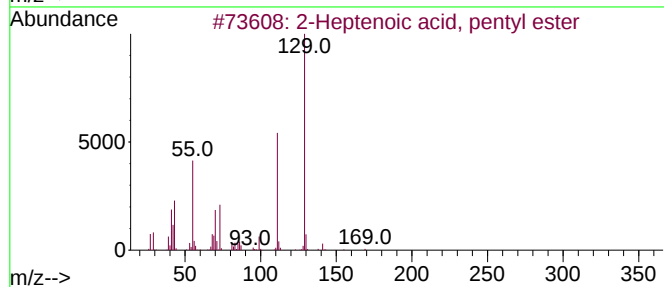
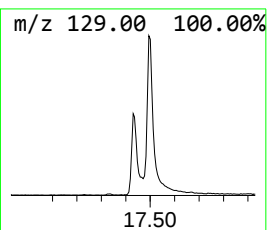
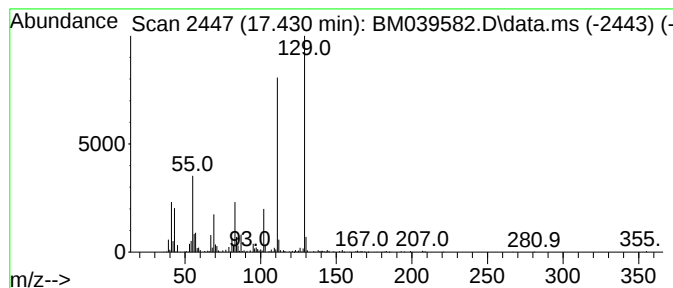
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.430	3.51 ng/ul	371787	Phenanthrene-d10	17.230

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Heptenoic acid, pentyl ester	198	C12H22O2	1000406-09-0	45
2			2-Heptenoic acid, butyl ester	184	C11H20O2	1000406-08-9	45
3			2-Heptenoic acid, 2-ethylhexyl e...	240	C15H28O2	1000406-90-3	43
4			Adipate Dimethylolpropane	232	C11H20O5	1000445-99-1	42
5			2-Heptenoic acid, hexyl ester	212	C13H24O2	1000406-09-2	42



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042223\
Data File : BM039582.D
Acq On : 22 Apr 2023 18:41
Operator : CG/JU
Sample : 02440-12
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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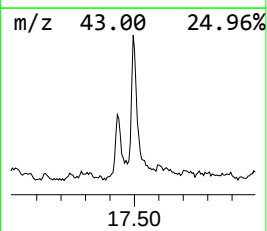
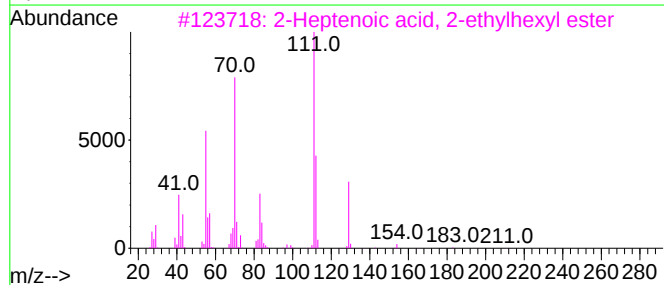
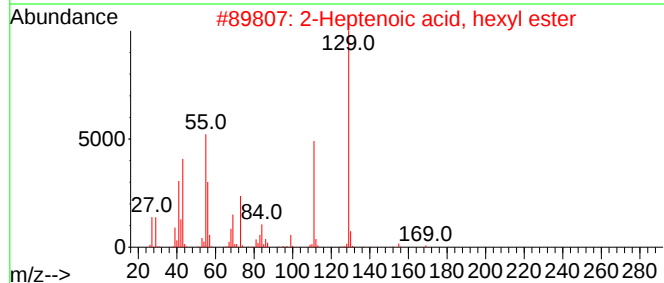
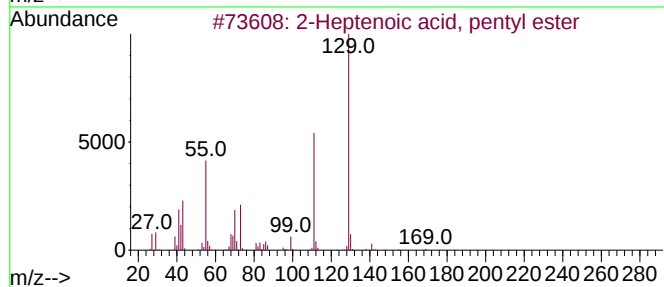
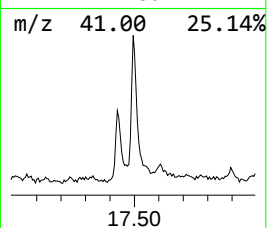
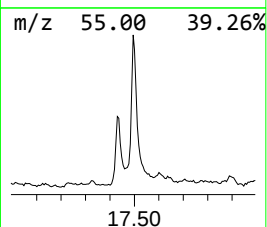
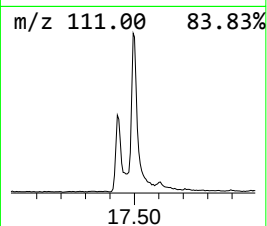
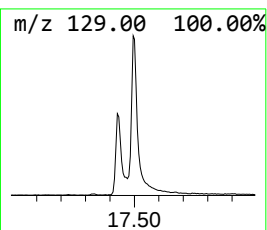
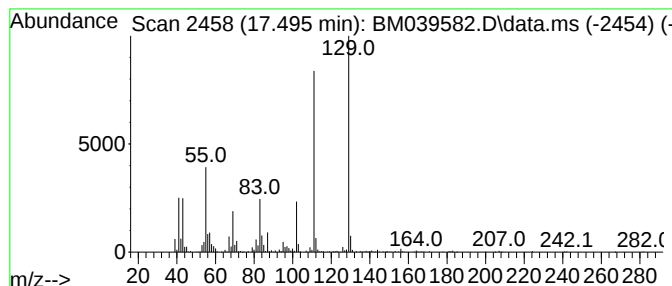
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 2-Heptenoic acid, pentyl ester Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.495	7.41 ng/ul	785133	Phenanthrene-d10	17.230

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Heptenoic acid, pentyl ester	198	C12H22O2	1000406-09-0	59
2			2-Heptenoic acid, hexyl ester	212	C13H24O2	1000406-09-2	53
3			2-Heptenoic acid, 2-ethylhexyl e...	240	C15H28O2	1000406-90-3	43
4			Cyclohexane, 1,1,3-trimethyl-	126	C9H18	003073-66-3	38
5			2-Heptenoic acid, octyl ester	240	C15H28O2	1000406-09-4	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042223\
Data File : BM039582.D
Acq On : 22 Apr 2023 18:41
Operator : CG/JU
Sample : 02440-12
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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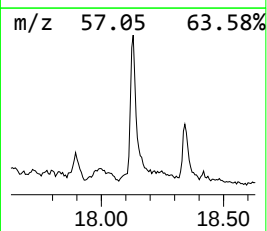
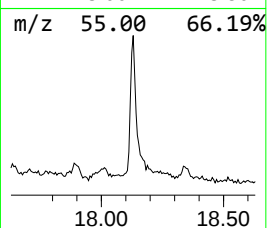
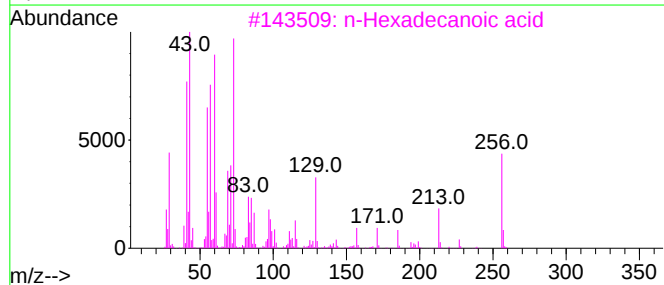
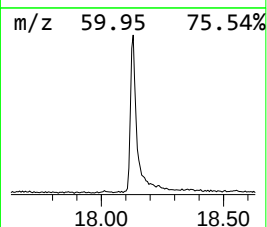
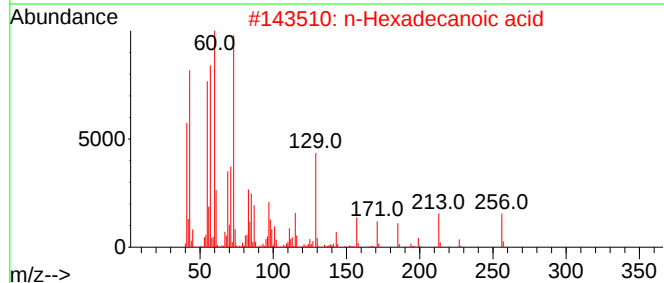
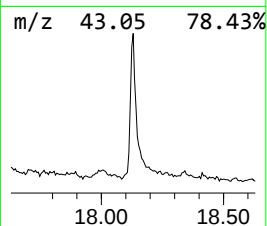
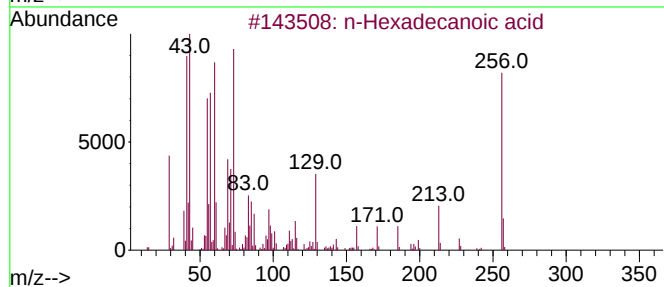
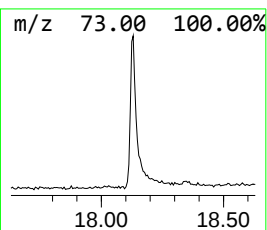
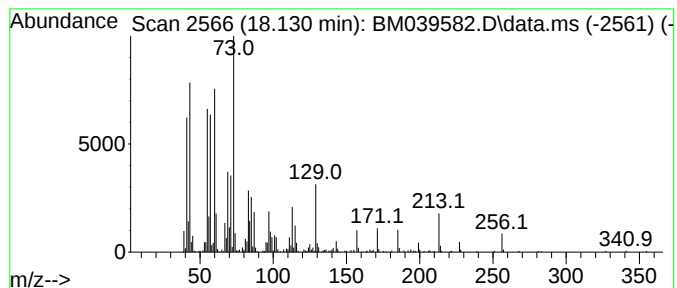
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.130	5.15 ng/ul	545721	Phenanthrene-d10	17.230

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94
4			Pentadecanoic acid	242	C15H30O2	001002-84-2	81
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042223\
Data File : BM039582.D
Acq On : 22 Apr 2023 18:41
Operator : CG/JU
Sample : 02440-12
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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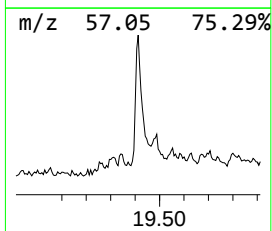
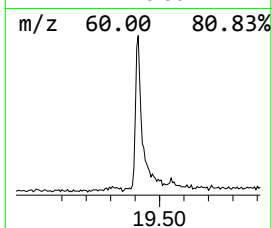
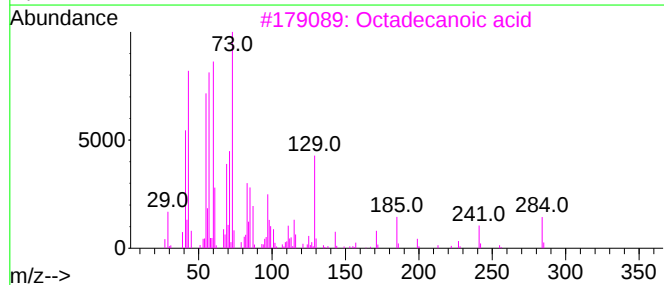
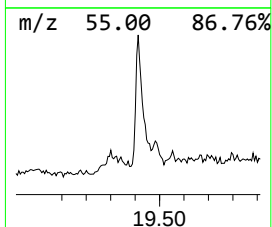
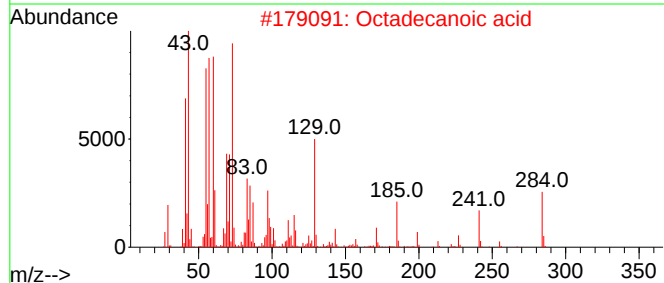
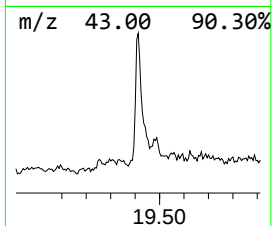
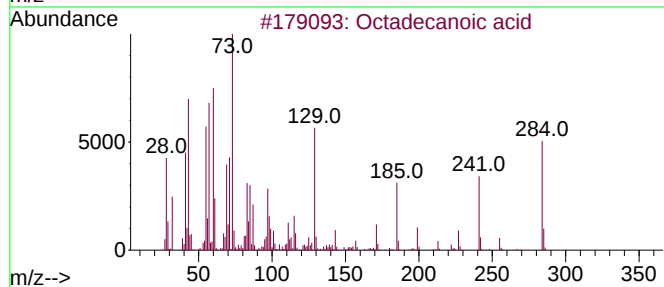
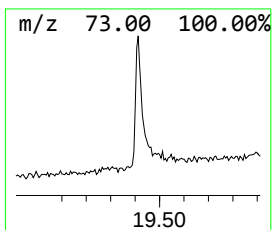
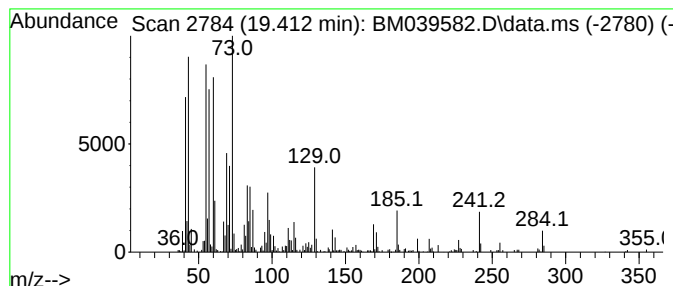
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Octadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.412	3.86 ng/ul	430456	Chrysene-d12	21.424

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Octadecanoic acid	284	C18H36O2	000057-11-4	99
3			Octadecanoic acid	284	C18H36O2	000057-11-4	98
4			Octadecanoic acid	284	C18H36O2	000057-11-4	98
5			Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042223\
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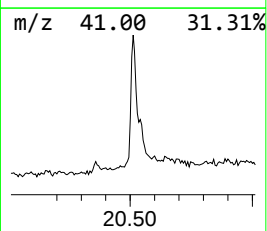
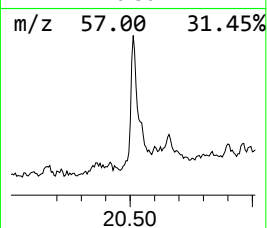
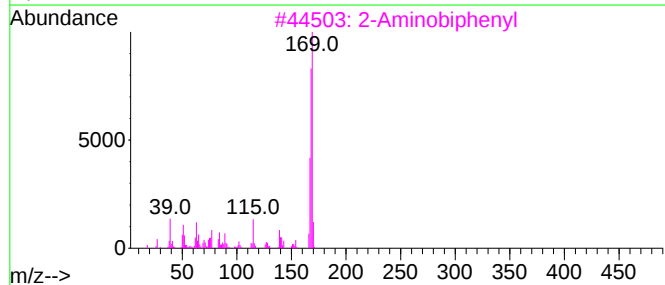
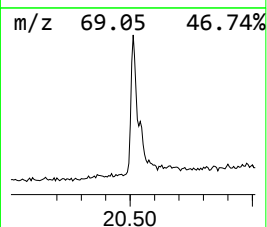
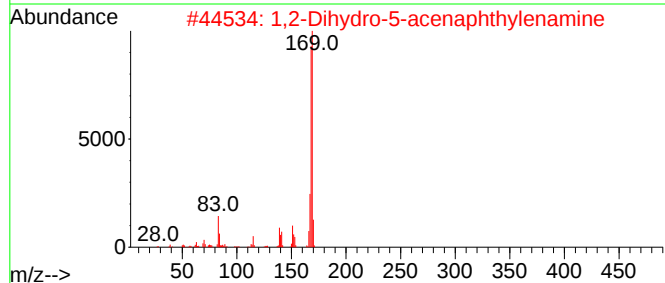
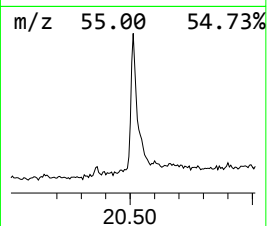
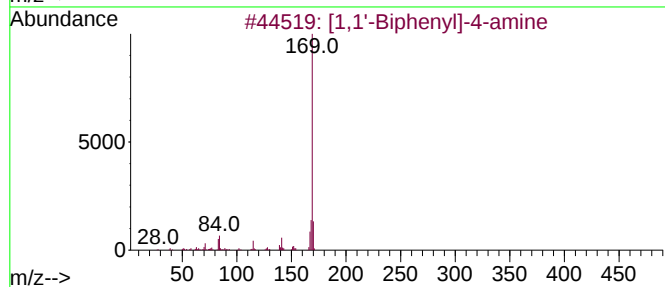
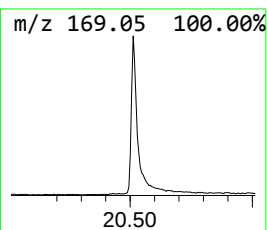
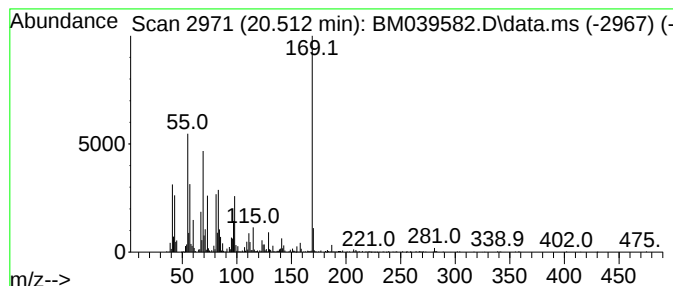
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 unknown-02 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.512	8.41 ng/ul	939251	Chrysene-d12	21.424	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	[1,1'-Biphenyl]-4-amine	169	C12H11N	000092-67-1	46
2	1,2-Dihydro-5-acenaphthylenamine	169	C12H11N	004657-93-6	43
3	2-Aminobiphenyl	169	C12H11N	000090-41-5	43
4	2-p-Tolylpyridine	169	C12H11N	004467-06-5	38
5	[1,1'-Biphenyl]-3-amine	169	C12H11N	002243-47-2	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042223\
Data File : BM039582.D
Acq On : 22 Apr 2023 18:41
Operator : CG/JU
Sample : 02440-12
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
COMQ7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.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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Hexanoic acid, ...	9.148	4.7	ng/ul	265644	1	7.819	1124990	20.0
Octanoic acid	9.995	3.2	ng/ul	249226	2	10.625	1542680	20.0
Benzophenone	15.842	3.4	ng/ul	323163	3	14.477	1904180	20.0
unknown-01	17.430	3.5	ng/ul	371787	4	17.230	2119540	20.0
2-Heptenoic aci...	17.495	7.4	ng/ul	785133	4	17.230	2119540	20.0
n-Hexadecanoic ...	18.130	5.2	ng/ul	545721	4	17.230	2119540	20.0
Octadecanoic acid	19.412	3.9	ng/ul	430456	5	21.424	2232750	20.0
unknown-02	20.512	8.4	ng/ul	939251	5	21.424	2232750	20.0