

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092322\
 Data File : BM036678.D
 Acq On : 23 Sep 2022 15:42
 Operator : CG/JU
 Sample : N4706-05
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 CB8K8

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-BM091522.M
 Title : SVOA CALIBRATION

Signal : TIC: BM036678.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.122	3	6	15	rVB	6781678	8321379	92.65%	6.852%
2	3.193	15	18	30	rVB	236095	364368	4.06%	0.300%
3	3.393	48	52	55	rBV	88032	120405	1.34%	0.099%
4	3.428	55	58	65	rVB	85145	109186	1.22%	0.090%
5	3.675	97	100	106	rBV	48590	65532	0.73%	0.054%
6	3.822	122	125	139	rBV	360653	541448	6.03%	0.446%
7	4.057	160	165	171	rVB	44060	70204	0.78%	0.058%
8	4.157	177	182	186	rVB	31662	43618	0.49%	0.036%
9	4.604	253	258	268	rVB	166955	257104	2.86%	0.212%
10	5.699	438	444	450	rBV4	12907	29378	0.33%	0.024%
11	7.175	688	695	708	rVB	574273	967721	10.77%	0.797%
12	7.345	717	724	732	rBV	1476403	2309552	25.71%	1.902%
13	7.545	748	758	766	rBV	1571158	2455832	27.34%	2.022%
14	7.616	766	770	776	rVB7	16716	34020	0.38%	0.028%
15	7.887	810	816	821	rVB8	8729	17372	0.19%	0.014%
16	8.016	830	838	845	rBV	1712635	2655007	29.56%	2.186%
17	8.081	845	849	853	rBV5	11250	17852	0.20%	0.015%
18	8.392	895	902	911	rVV2	50629	95052	1.06%	0.078%
19	8.728	952	959	970	rVV	1481329	2439774	27.16%	2.009%
20	8.810	970	973	978	rVV6	11267	22569	0.25%	0.019%
21	8.922	987	992	995	rVV6	11265	18999	0.21%	0.016%
22	9.034	1004	1011	1016	rVB10	12781	27868	0.31%	0.023%
23	9.128	1022	1027	1029	rBV3	19960	25686	0.29%	0.021%
24	9.181	1029	1036	1046	rVB	1780551	2813832	31.33%	2.317%
25	9.292	1051	1055	1060	rVB6	22423	39466	0.44%	0.032%
26	9.351	1060	1065	1068	rBV6	13112	23095	0.26%	0.019%
27	9.404	1070	1074	1080	rVB9	15529	33689	0.38%	0.028%
28	9.475	1080	1086	1090	rBV3	14298	28262	0.31%	0.023%
29	9.545	1091	1098	1104	rBV2	122713	211919	2.36%	0.174%
30	9.622	1104	1111	1114	rBV2	35551	81118	0.90%	0.067%
31	9.769	1132	1136	1141	rVV8	22875	36380	0.41%	0.030%
32	9.822	1141	1145	1149	rVB6	11107	19666	0.22%	0.016%
33	9.904	1152	1159	1171	rBV	1302792	2137690	23.80%	1.760%
34	10.134	1191	1198	1203	rVB9	19008	45521	0.51%	0.037%
35	10.228	1203	1214	1220	rBV6	33235	80700	0.90%	0.066%
36	10.357	1232	1236	1240	rVB7	12350	17864	0.20%	0.015%

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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
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37	10.445	1244	1251	1260	rBV	2181394	3573474	39.79%	2.942%
38	10.604	1273	1278	1282	rBV	47076	89947	1.00%	0.074%
39	10.828	1309	1316	1323	rBV	2456611	4095837	45.60%	3.372%
40	10.963	1331	1339	1350	rBV	897525	1580740	17.60%	1.302%
41	11.186	1373	1377	1381	rBV5	19120	25834	0.29%	0.021%
42	11.369	1404	1408	1413	rBV3	18105	32318	0.36%	0.027%
43	11.422	1413	1417	1424	rVB10	13111	28736	0.32%	0.024%
44	11.533	1431	1436	1442	rVB	51511	86681	0.97%	0.071%
45	11.622	1445	1451	1455	rBV7	27325	58828	0.65%	0.048%
46	11.669	1455	1459	1464	rVV6	22039	47936	0.53%	0.039%
47	11.728	1466	1469	1476	rVB9	19823	38626	0.43%	0.032%
48	11.804	1476	1482	1486	rBV8	17966	35874	0.40%	0.030%
49	12.022	1514	1519	1525	rBV8	23321	52521	0.58%	0.043%
50	12.086	1525	1530	1535	rVV6	33488	60518	0.67%	0.050%
51	12.163	1535	1543	1548	rVV	183421	317344	3.53%	0.261%
52	12.204	1548	1550	1555	rVB4	25137	35205	0.39%	0.029%
53	12.322	1566	1570	1574	rBV3	15508	23435	0.26%	0.019%
54	12.404	1579	1584	1598	rVB4	39765	163716	1.82%	0.135%
55	12.580	1611	1614	1619	rBV5	26724	37419	0.42%	0.031%
56	12.692	1631	1633	1639	rVB6	18021	22147	0.25%	0.018%
57	12.751	1639	1643	1648	rVB7	19220	26161	0.29%	0.022%
58	13.016	1683	1688	1695	rBV3	43716	72899	0.81%	0.060%
59	13.133	1706	1708	1718	rVB3	13725	24411	0.27%	0.020%
60	13.263	1726	1730	1734	rVB	69251	104568	1.16%	0.086%
61	13.374	1736	1749	1753	rBV	28017	94447	1.05%	0.078%
62	13.551	1775	1779	1785	rBV3	34555	54653	0.61%	0.045%
63	13.604	1785	1788	1792	rVV4	20336	29176	0.32%	0.024%
64	13.739	1808	1811	1816	rBV4	15302	27008	0.30%	0.022%
65	13.822	1822	1825	1828	rBV5	26319	39472	0.44%	0.032%
66	13.863	1828	1832	1835	rVV4	50613	80283	0.89%	0.066%
67	13.898	1835	1838	1844	rVB	54483	84786	0.94%	0.070%
68	14.057	1859	1865	1871	rVB	4096814	5597428	62.32%	4.609%
69	14.257	1895	1899	1902	rBV5	36764	53153	0.59%	0.044%
70	14.286	1902	1904	1906	rVV2	26885	30492	0.34%	0.025%
71	14.339	1906	1913	1924	rVB	4539006	6698824	74.59%	5.516%
72	14.445	1926	1931	1943	rVB	180619	309465	3.45%	0.255%
73	14.569	1948	1952	1957	rVB5	66720	94813	1.06%	0.078%
74	14.639	1957	1964	1972	rBV	3700158	5476819	60.98%	4.509%
75	14.833	1992	1997	2005	rBV	517795	755325	8.41%	0.622%
76	15.057	2032	2035	2038	rVB4	38177	47766	0.53%	0.039%

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77	15.386	2087	2091	2095	rBV2	59474	93218	1.04%	0.077%
78	15.445	2097	2101	2106	rVB3	50308	98882	1.10%	0.081%
79	15.627	2126	2132	2144	rVB	5947548	8421030	93.76%	6.934%
80	15.739	2145	2151	2159	rBV	1556873	2144632	23.88%	1.766%
81	15.857	2167	2171	2183	rVB	273124	504208	5.61%	0.415%
82	16.145	2214	2220	2230	rBV	4334812	5896105	65.65%	4.855%
83	16.468	2271	2275	2280	rBV	227805	281983	3.14%	0.232%
84	16.651	2300	2306	2311	rBV	2589532	3409585	37.96%	2.807%
85	17.139	2386	2389	2391	rBV3	40826	48693	0.54%	0.040%
86	17.380	2424	2430	2437	rBV	4145545	5982937	66.61%	4.926%
87	17.474	2440	2446	2455	rVB	5972735	8488723	94.51%	6.989%
88	18.051	2540	2544	2549	rVB	219231	280521	3.12%	0.231%
89	18.274	2577	2582	2590	rBV	1164066	1698855	18.92%	1.399%
90	19.021	2706	2709	2712	rBV	108339	123330	1.37%	0.102%
91	19.398	2770	2773	2776	rBV	105537	135441	1.51%	0.112%
92	19.480	2784	2787	2793	rVB	407261	450376	5.01%	0.371%
93	19.545	2794	2798	2805	rBV	566104	791886	8.82%	0.652%
94	19.756	2829	2834	2839	rBV	6406841	8981416	100.00%	7.395%
95	19.803	2839	2842	2850	rVB	334851	439641	4.90%	0.362%
96	20.756	3001	3004	3006	rBV	313767	317010	3.53%	0.261%
97	21.250	3085	3088	3096	rBV	1173485	1329068	14.80%	1.094%
98	21.545	3133	3138	3149	rBV	3846902	4890771	54.45%	4.027%
99	23.827	3520	3526	3533	rBV2	3106872	5928883	66.01%	4.882%
100	23.986	3547	3553	3563	rVB2	2041945	4027431	44.84%	3.316%

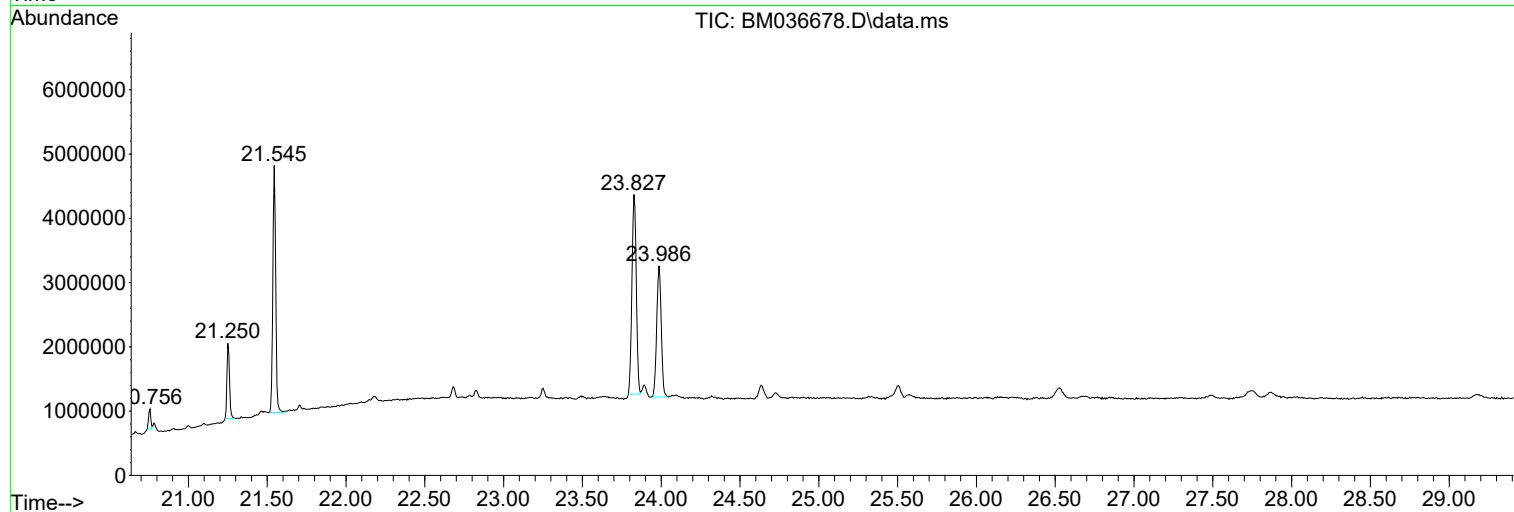
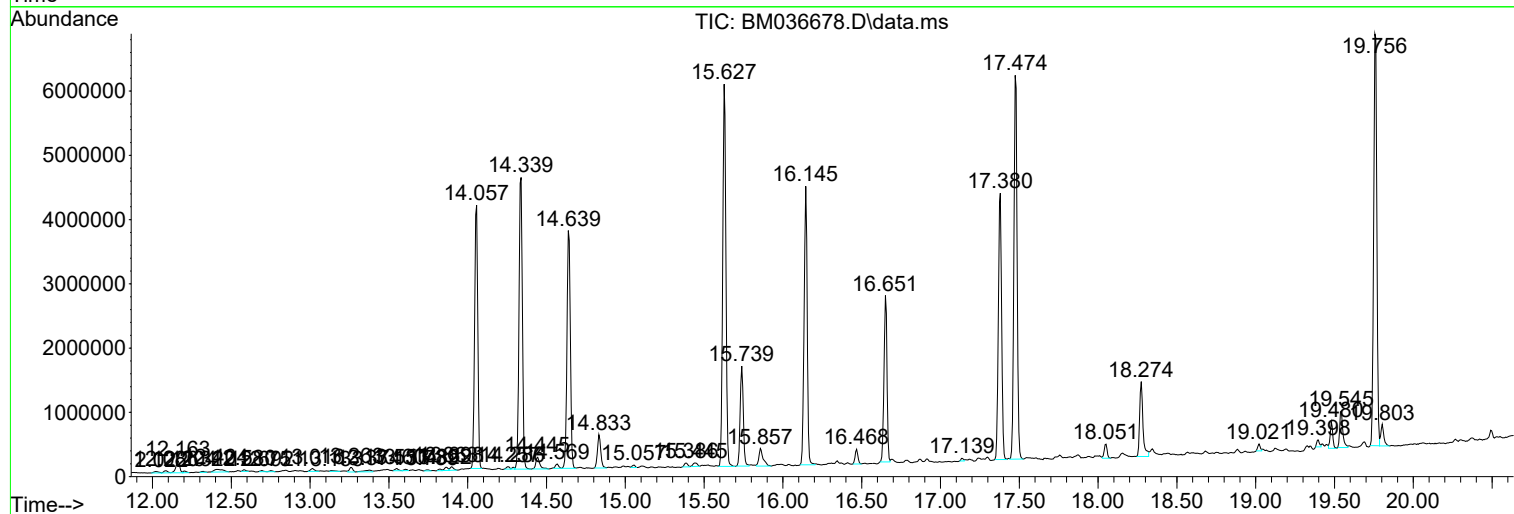
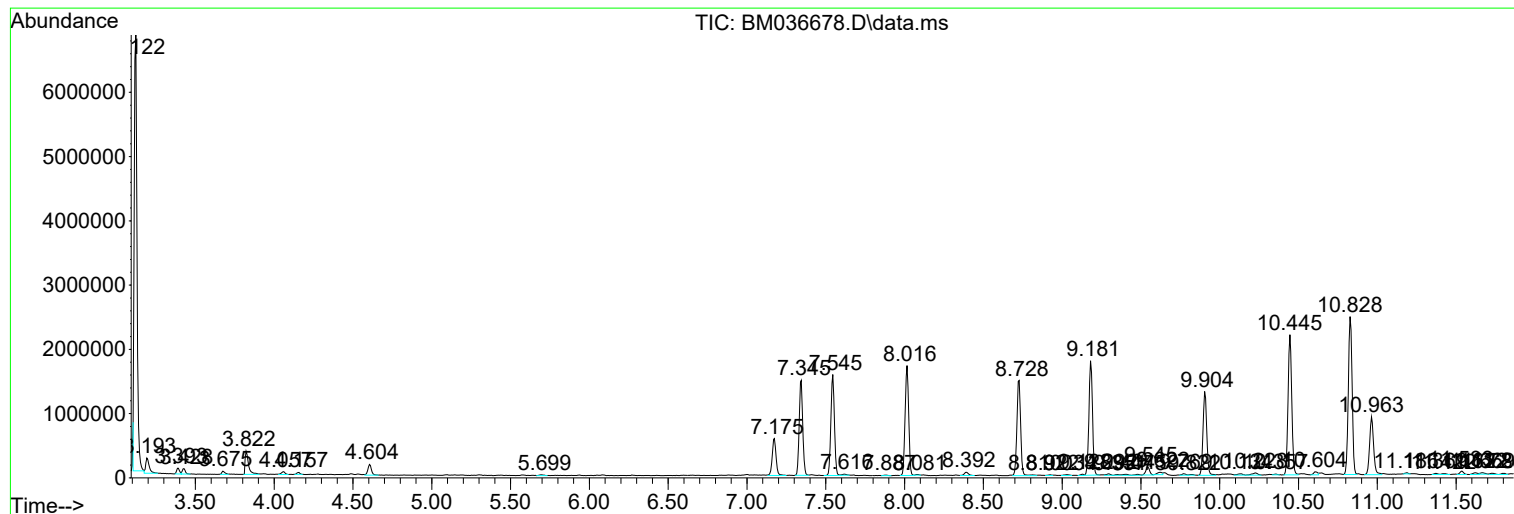
Sum of corrected areas: 121452838

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

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



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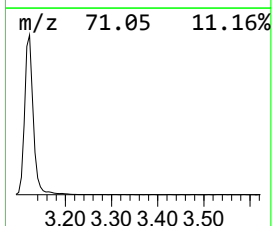
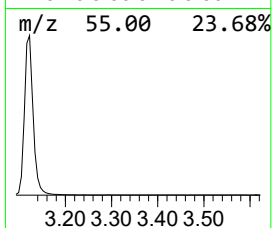
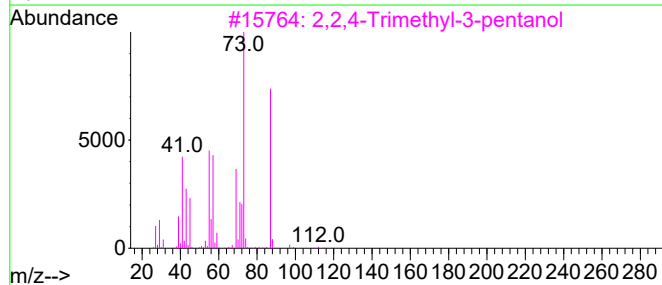
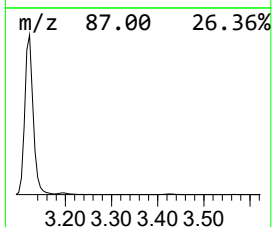
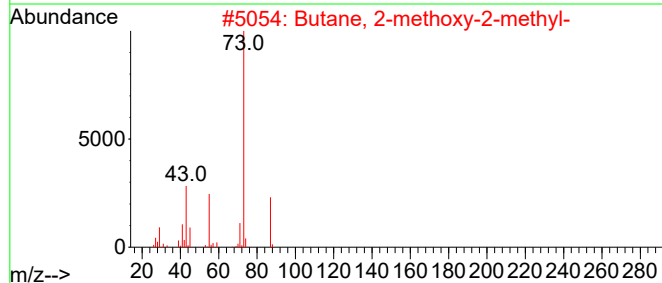
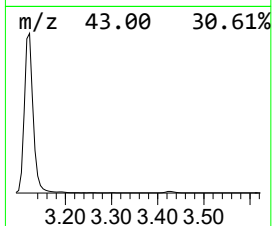
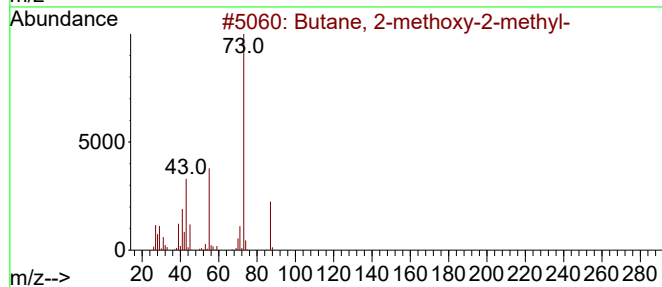
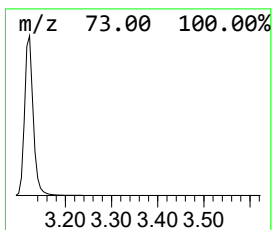
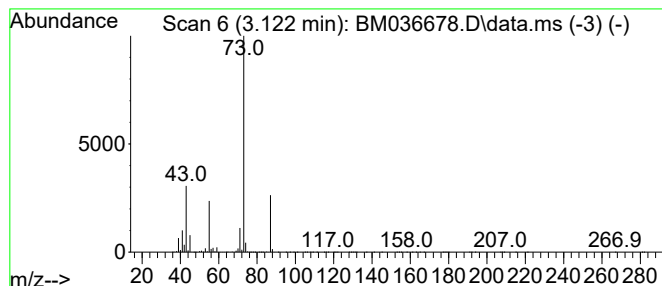
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.122	62.68 ng/ul	8321380	1,4-Dichlorobenzene-d4	8.016

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	78
3			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
4			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
5			Borane, dimethoxy-	74	C2H7BO2	004542-61-4	9



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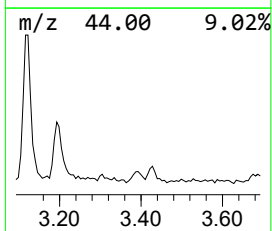
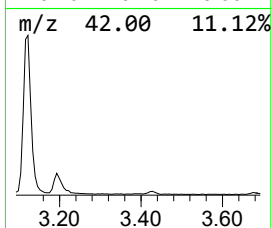
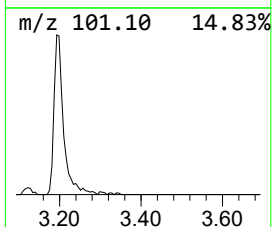
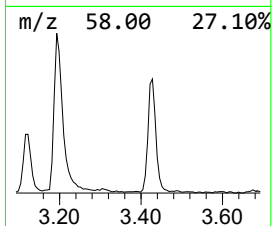
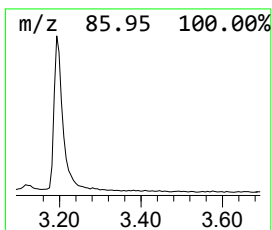
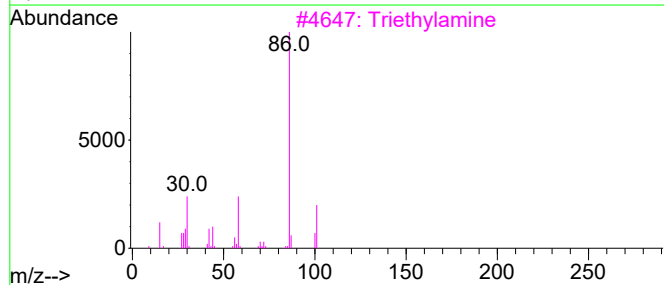
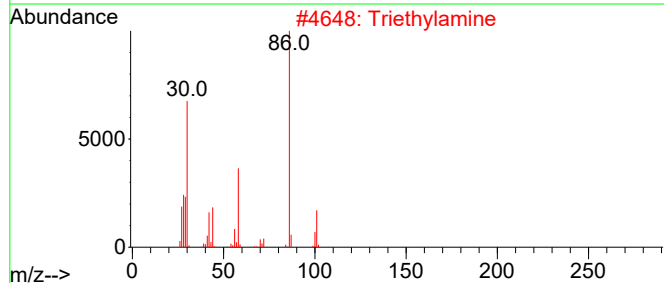
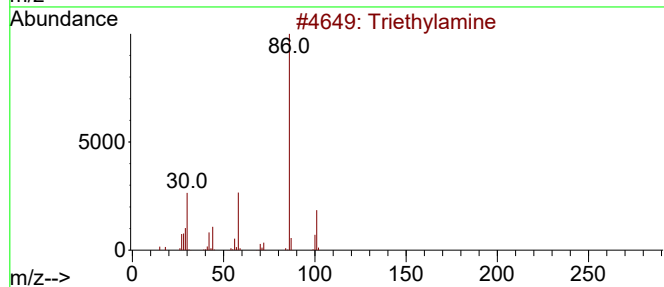
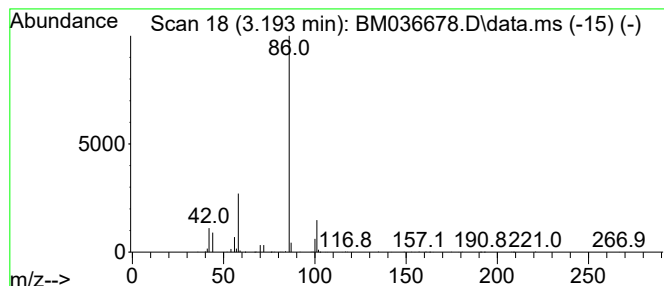
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Triethylamine Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.193	2.74 ng/ul	364368	1,4-Dichlorobenzene-d4	8.016

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triethylamine	101	C6H15N	000121-44-8	91
2			Triethylamine	101	C6H15N	000121-44-8	90
3			Triethylamine	101	C6H15N	000121-44-8	87
4			Tetraethylammonium bromide	209	C8H20BrN	000071-91-0	83
5			Tetraethylammonium chloride	165	C8H20ClN	000056-34-8	83



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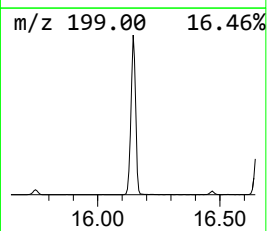
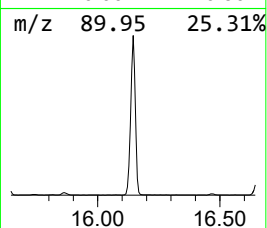
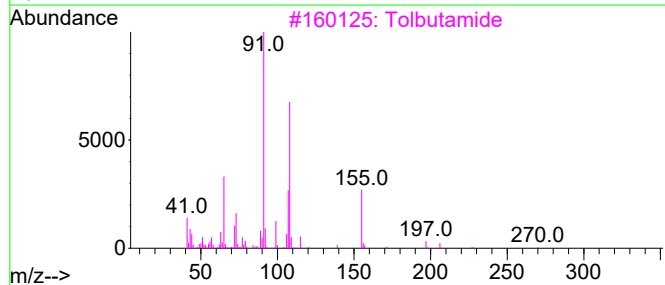
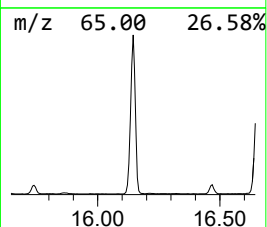
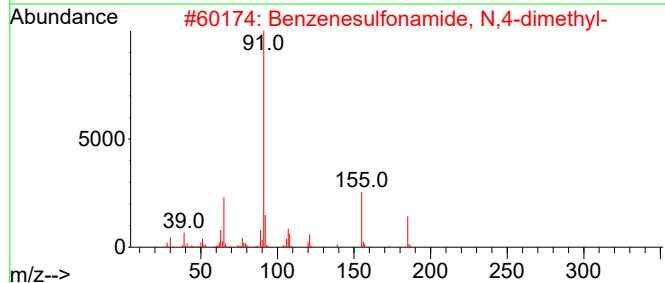
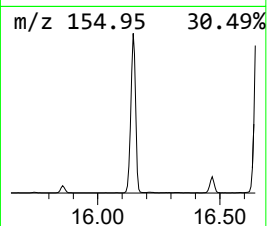
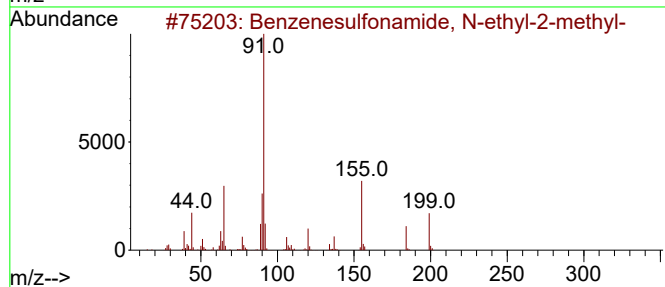
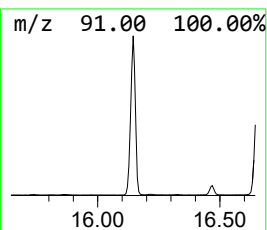
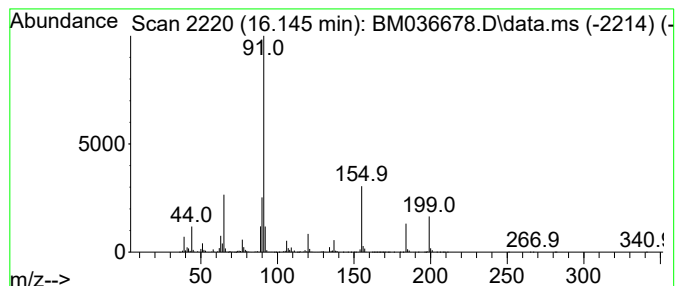
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzenesulfonamide, N-ethyl... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.145	19.71 ng/ul	5896110	Phenanthrene-d10	17.380

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	98
2			Benzenesulfonamide, N,4-dimethyl-	185	C8H11NO2S	000640-61-9	50
3			Tolbutamide	270	C12H18N2O3S	000064-77-7	50
4			Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	49
5			Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092322\
Data File : BM036678.D
Acq On : 23 Sep 2022 15:42
Operator : CG/JU
Sample : N4706-05
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB8K8

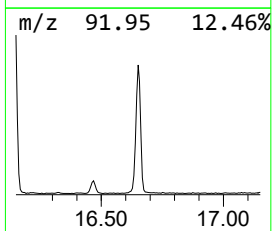
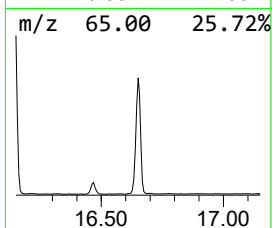
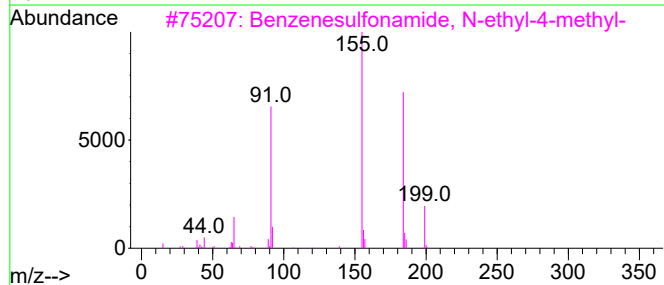
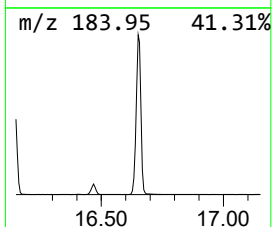
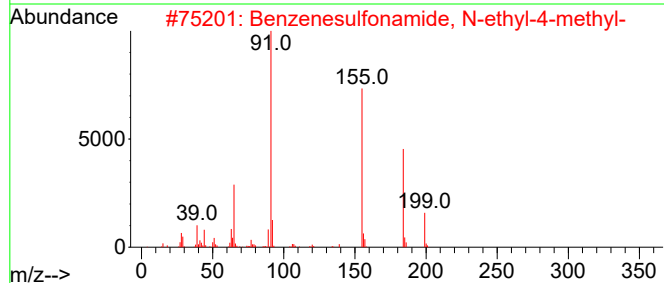
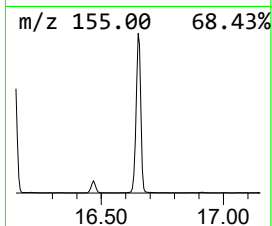
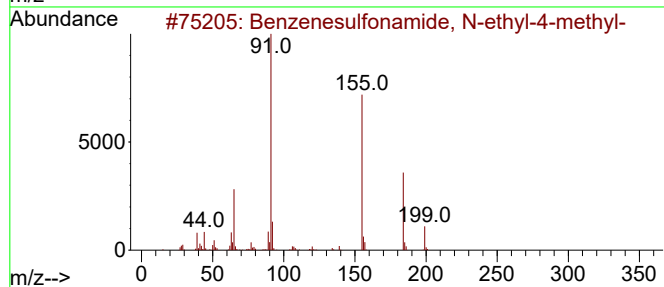
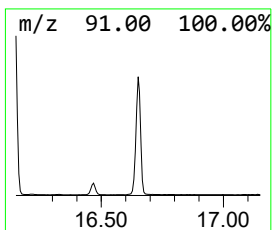
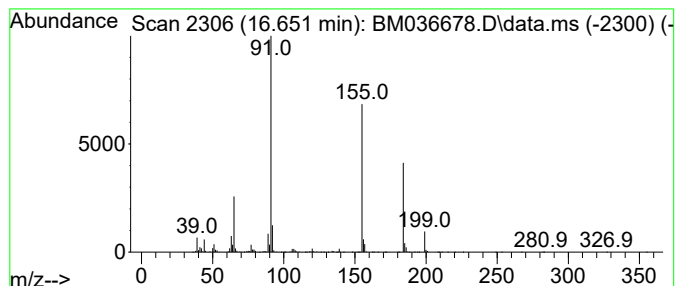
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Benzenesulfonamide, N-ethyl... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.651	11.40 ng/ul	3409590	Phenanthrene-d10	17.380

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	98
2			Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	96
3			Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	91
4			4-Methyl-N-(2-oxobutyl)benzenesu...	241	C11H15NO3S	1000192-14-5	86
5			N-isobutyl-4-methylbenzenesulfon...	227	C11H17NO2S	023705-38-6	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092322\
Data File : BM036678.D
Acq On : 23 Sep 2022 15:42
Operator : CG/JU
Sample : N4706-05
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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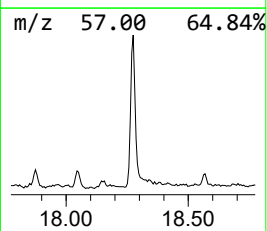
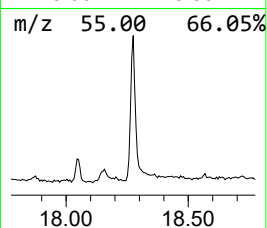
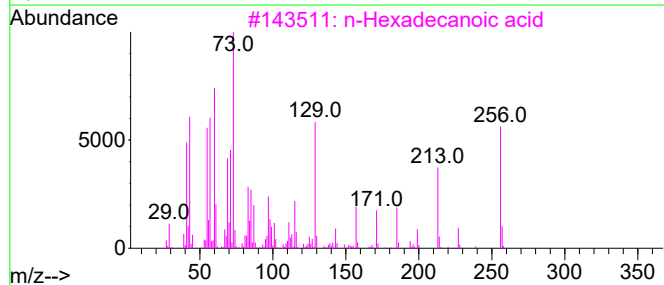
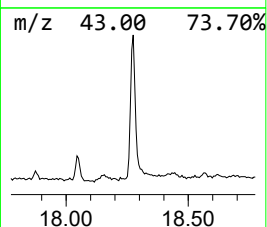
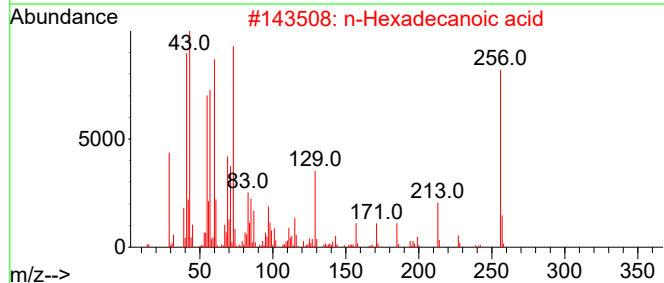
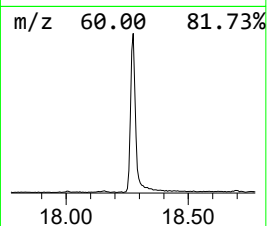
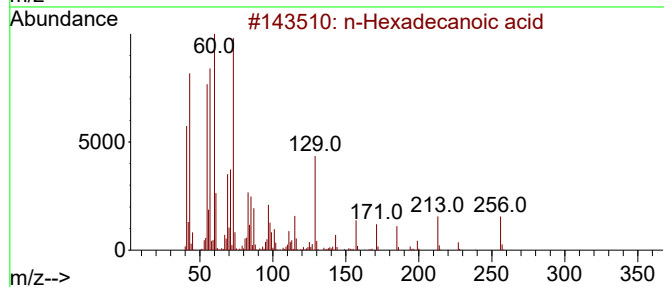
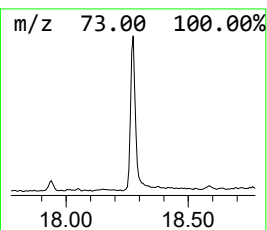
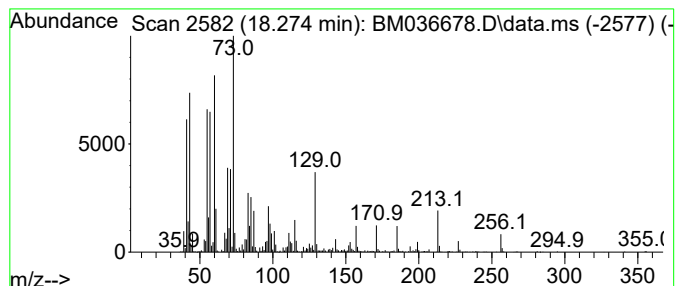
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.274	5.68 ng/ul	1698860	Phenanthrene-d10	17.380

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	99
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5			Tridecanoic acid	214	C13H26O2	000638-53-9	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092322\
Data File : BM036678.D
Acq On : 23 Sep 2022 15:42
Operator : CG/JU
Sample : N4706-05
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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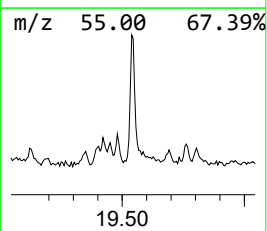
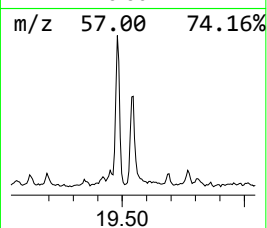
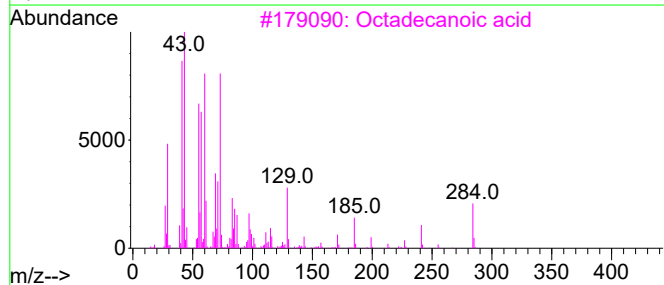
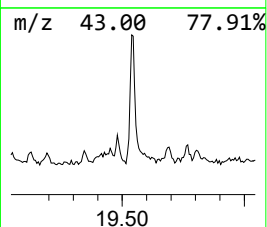
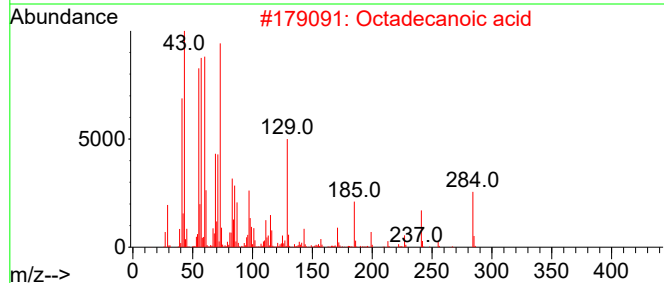
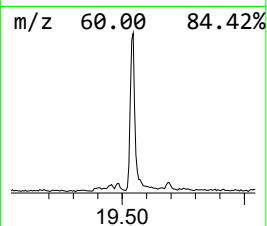
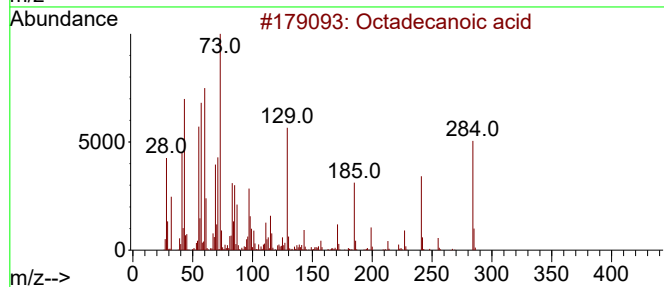
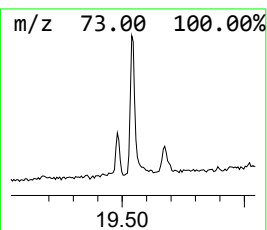
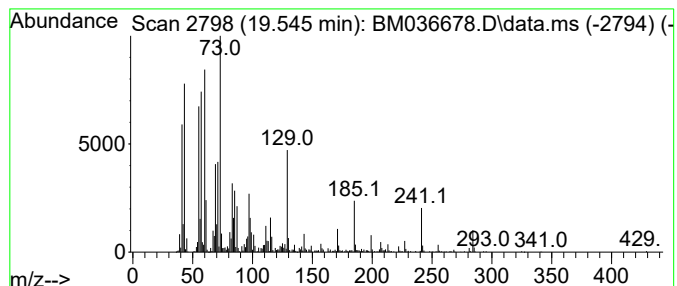
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Octadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.545	3.24 ng/ul	791886	Chrysene-d12	21.544

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	95
5	Octadecanoic acid			284	C18H36O2	000057-11-4	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092322\
Data File : BM036678.D
Acq On : 23 Sep 2022 15:42
Operator : CG/JU
Sample : N4706-05
Misc :
ALS Vial : 10 Sample Multiplier: 1

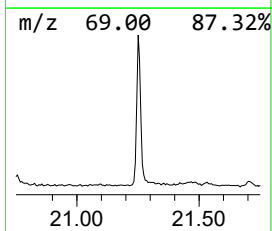
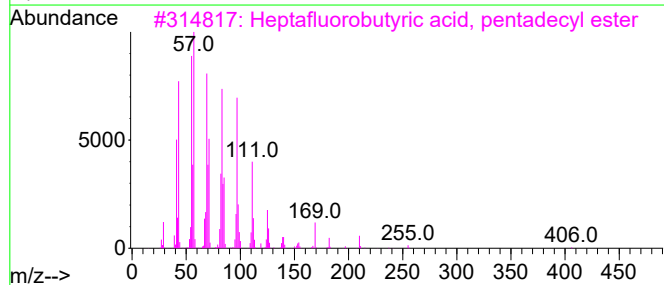
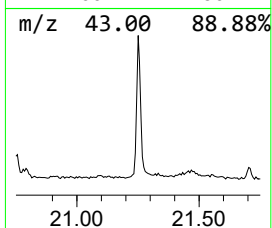
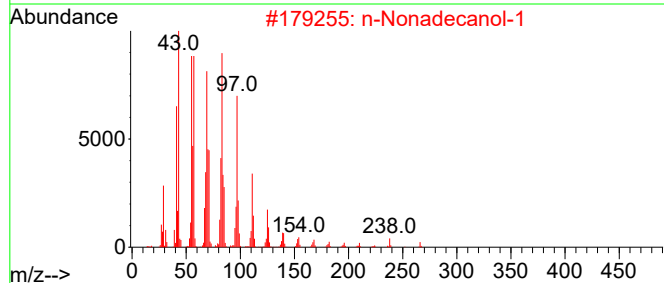
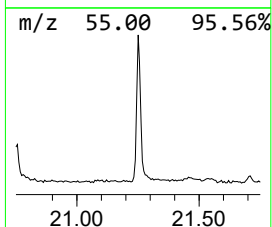
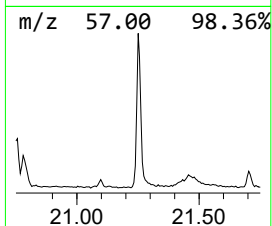
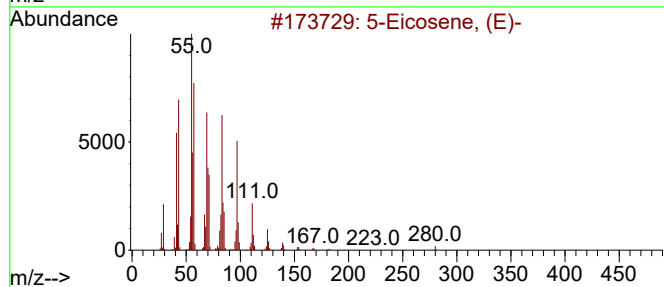
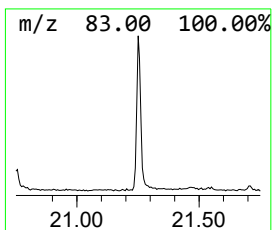
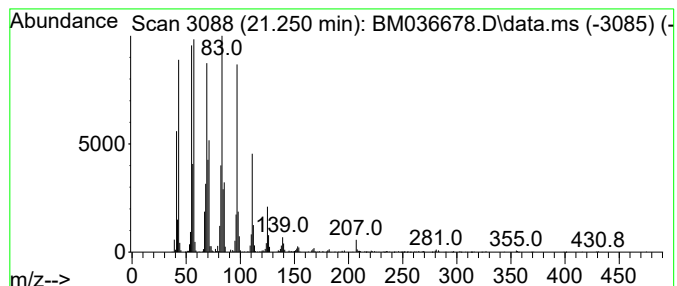
Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.250	5.43 ng/ul	1329070	Chrysene-d12	21.544	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Eicosene, (E)-	280	C20H40	074685-30-6	99
2	n-Nonadecanol-1	284	C19H40O	001454-84-8	95
3	Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	94
4	1-Heneicosanol	312	C21H44O	015594-90-8	94
5	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092322\
Data File : BM036678.D
Acq On : 23 Sep 2022 15:42
Operator : CG/JU
Sample : N4706-05
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.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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Triethylamine	3.193	2.7	ng/ul	364368	1	8.016	2655010	20.0
Benzenesulfonam...	16.145	19.7	ng/ul	5896110	4	17.380	5982940	20.0
Benzenesulfonam...	16.651	11.4	ng/ul	3409590	4	17.380	5982940	20.0
n-Hexadecanoic ...	18.274	5.7	ng/ul	1698860	4	17.380	5982940	20.0
Octadecanoic acid	19.545	3.2	ng/ul	791886	5	21.544	4890770	20.0
(DEL) Alkane: S...	21.250	5.4	ng/ul	1329070	5	21.544	4890770	20.0