

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
 Data File : BM036570.D
 Acq On : 16 Sep 2022 19:34
 Operator : CG/JU
 Sample : N4601-19
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A5760

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: BM036570.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	32	rVB	1333847	2074783	38.83%	3.740%
2	4.163	179	183	189	rVB	21563	30786	0.58%	0.055%
3	6.728	615	619	622	rBV5	3604	6032	0.11%	0.011%
4	7.169	687	694	707	rVB	343404	604363	11.31%	1.089%
5	7.351	718	725	733	rBV	575023	879393	16.46%	1.585%
6	7.551	752	759	769	rBV	336952	536671	10.05%	0.967%
7	8.028	833	840	847	rBV	1124442	1794164	33.58%	3.234%
8	8.087	847	850	855	rVB4	10437	15730	0.29%	0.028%
9	8.722	951	958	971	rBV	436800	698470	13.07%	1.259%
10	9.187	1030	1037	1047	rBV	590650	972288	18.20%	1.753%
11	9.275	1050	1052	1060	rVV8	6544	12899	0.24%	0.023%
12	9.357	1060	1066	1071	rVB9	5378	11788	0.22%	0.021%
13	9.587	1102	1105	1106	rVV3	7087	8983	0.17%	0.016%
14	9.628	1108	1112	1119	rVB	23431	40357	0.76%	0.073%
15	9.916	1152	1161	1168	rBV	134052	232788	4.36%	0.420%
16	10.451	1244	1252	1261	rBV	264314	443373	8.30%	0.799%
17	10.610	1275	1279	1284	rBV3	17534	29313	0.55%	0.053%
18	10.839	1310	1318	1328	rVB	1658864	2691315	50.37%	4.851%
19	10.969	1335	1340	1346	rVB2	18633	34326	0.64%	0.062%
20	12.122	1529	1536	1542	rVB4	10058	25021	0.47%	0.045%
21	12.257	1556	1559	1563	rVB5	9380	14763	0.28%	0.027%
22	12.416	1581	1586	1592	rVB	16929	29428	0.55%	0.053%
23	13.210	1715	1721	1724	rVB8	5400	6838	0.13%	0.012%
24	13.392	1746	1752	1756	rVB9	3902	7243	0.14%	0.013%
25	13.486	1763	1768	1773	rBV	30905	42428	0.79%	0.076%
26	13.551	1776	1779	1784	rVV6	5853	10334	0.19%	0.019%
27	13.616	1788	1790	1797	rVB7	5278	10733	0.20%	0.019%
28	13.886	1831	1836	1841	rBV9	8542	13316	0.25%	0.024%
29	14.057	1859	1865	1877	rBV	1078904	1456129	27.25%	2.625%
30	14.139	1877	1879	1883	rVB5	4477	5488	0.10%	0.010%
31	14.298	1901	1906	1907	rBV4	5260	7458	0.14%	0.013%
32	14.345	1907	1914	1927	rVB	1693794	2431894	45.52%	4.383%
33	14.551	1945	1949	1953	rVB3	9943	11653	0.22%	0.021%
34	14.651	1959	1966	1975	rBV	2557837	3718954	69.61%	6.703%
35	14.816	1986	1994	2000	rVB8	19119	35599	0.67%	0.064%
36	14.880	2000	2005	2006	rBV5	5131	6543	0.12%	0.012%

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Integrator: RTE

Smoothing : OFF

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Min Area: 0.1 % of largest Peak

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Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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

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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
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37	15.351	2083	2085	2088	rBV4	4494	5983	0.11%	0.011%
38	15.451	2098	2102	2104	rBV5	7995	12022	0.23%	0.022%
39	15.498	2106	2110	2114	rVB3	10960	16477	0.31%	0.030%
40	15.633	2127	2133	2145	rBV	2323381	3360991	62.91%	6.058%

41	15.739	2145	2151	2159	rVB5	19135	30188	0.57%	0.054%
42	15.874	2169	2174	2178	rBV4	21644	32372	0.61%	0.058%
43	16.063	2200	2206	2209	rBV6	6517	12048	0.23%	0.022%
44	16.180	2221	2226	2229	rBV7	4603	6857	0.13%	0.012%
45	16.363	2253	2257	2259	rBV5	7756	9077	0.17%	0.016%

46	16.416	2263	2266	2270	rVB5	11562	13255	0.25%	0.024%
47	16.627	2298	2302	2307	rBV5	4888	9840	0.18%	0.018%
48	16.704	2310	2315	2319	rBV7	10679	17884	0.33%	0.032%
49	16.892	2340	2347	2350	rBV4	17621	30981	0.58%	0.056%
50	17.151	2387	2391	2393	rBV3	14859	19118	0.36%	0.034%

51	17.310	2412	2418	2420	rBV6	7380	9619	0.18%	0.017%
52	17.386	2423	2431	2441	rBV	3239564	4517165	84.55%	8.142%
53	17.480	2441	2447	2462	rVV	2450654	3607681	67.53%	6.503%
54	17.892	2514	2517	2519	rBV4	6490	6422	0.12%	0.012%
55	17.962	2526	2529	2534	rVB	17125	19261	0.36%	0.035%

56	18.062	2542	2546	2550	rVB	61352	70099	1.31%	0.126%
57	18.280	2579	2583	2586	rBV6	13761	20176	0.38%	0.036%
58	18.357	2593	2596	2600	rVB2	10769	12975	0.24%	0.023%
59	18.580	2631	2634	2639	rBV7	11593	15506	0.29%	0.028%
60	18.904	2685	2689	2694	rVB3	18285	25834	0.48%	0.047%

61	19.209	2738	2741	2747	rVB2	21163	26509	0.50%	0.048%
62	19.333	2758	2762	2767	rVB2	29568	37421	0.70%	0.067%
63	19.398	2770	2773	2777	rVV	24735	33399	0.63%	0.060%
64	19.704	2821	2825	2828	rBV3	16065	21110	0.40%	0.038%
65	19.762	2829	2835	2844	rBV	3160599	4222076	79.03%	7.610%

66	20.280	2920	2923	2926	rBV4	25726	36331	0.68%	0.065%
67	20.315	2926	2929	2933	rVV	84948	98514	1.84%	0.178%
68	20.768	3002	3006	3010	rBV	174966	216699	4.06%	0.391%
69	20.809	3010	3013	3017	rVB2	66000	68358	1.28%	0.123%
70	21.262	3086	3090	3099	rBV	637454	869100	16.27%	1.567%

71	21.550	3133	3139	3144	rBV	4107998	5181544	96.98%	9.340%
72	21.721	3164	3168	3176	rBV2	131064	176762	3.31%	0.319%
73	21.909	3197	3200	3204	rVB2	94591	102540	1.92%	0.185%
74	22.203	3244	3250	3260	rBV	740793	1191651	22.30%	2.148%
75	22.956	3375	3378	3386	rVB5	126591	178215	3.34%	0.321%

76	23.274	3428	3432	3435	rBV	182735	279909	5.24%	0.505%
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Integration Parameters: LSCINT.P
Integrator: RTE
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Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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Title : SVOA CALIBRATION

77	23.315	3436	3439	3450	rVB	275576	465455	8.71%	0.839%
78	23.833	3520	3527	3537	rVB2	2022181	3896764	72.94%	7.024%
79	23.992	3547	3554	3568	rVB2	2674186	5342659	100.00%	9.630%
80	24.674	3665	3670	3677	rVB2	185188	368643	6.90%	0.664%
81	26.850	4032	4040	4057	rVB3	588316	1832144	34.29%	3.302%

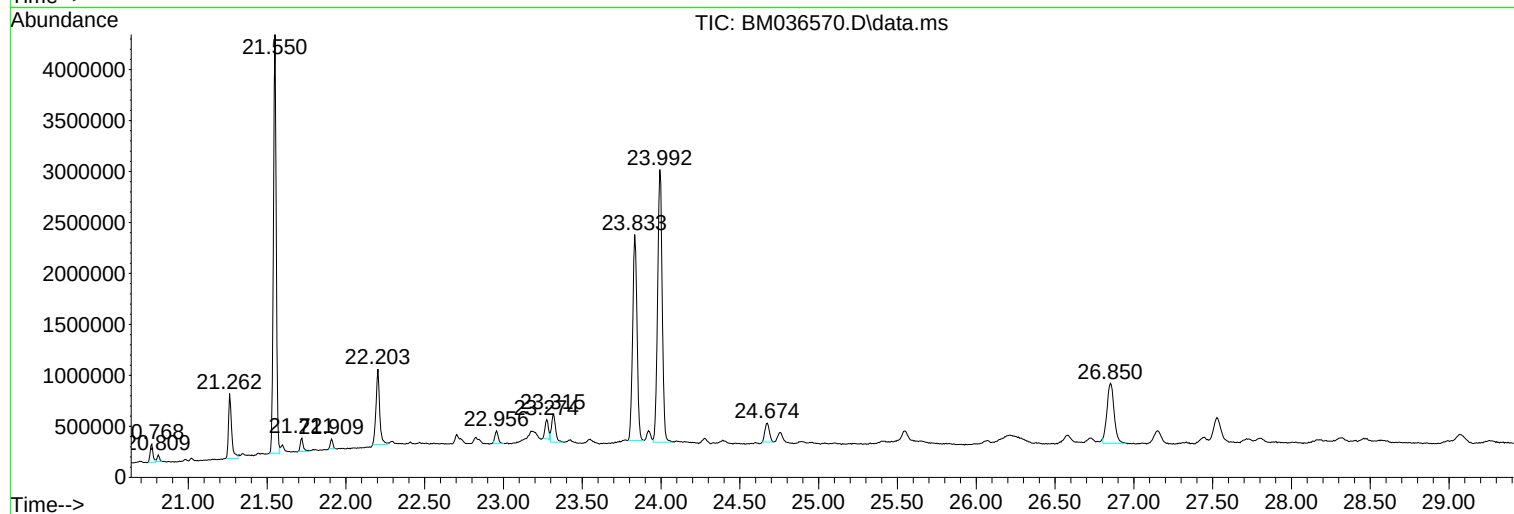
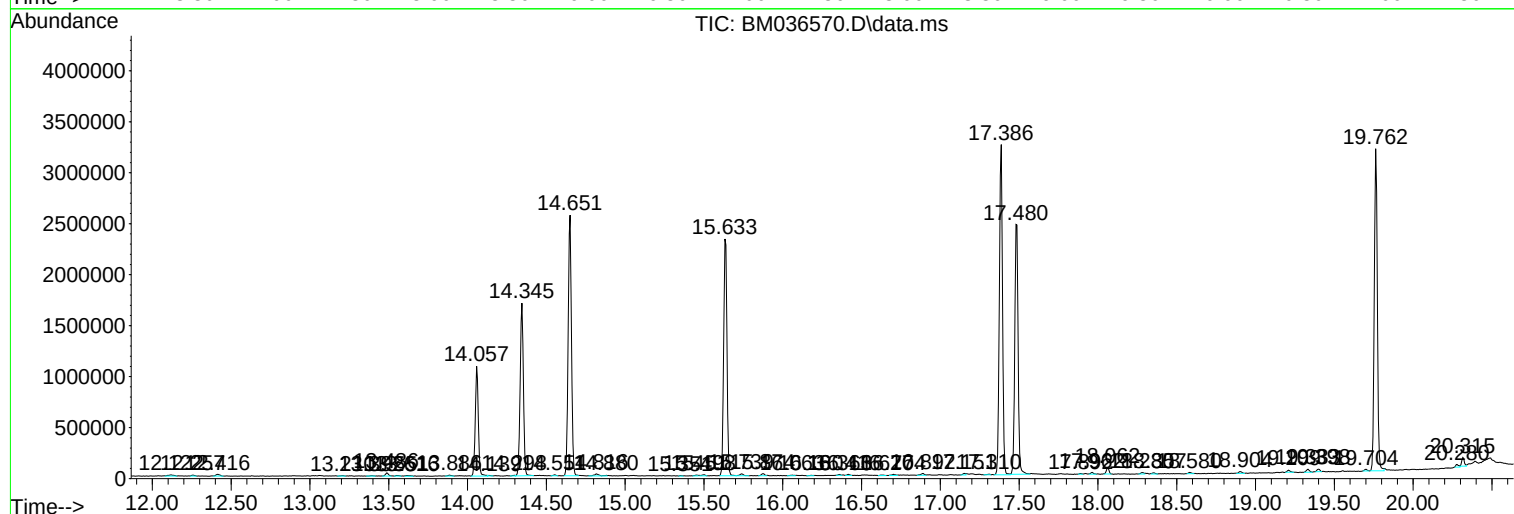
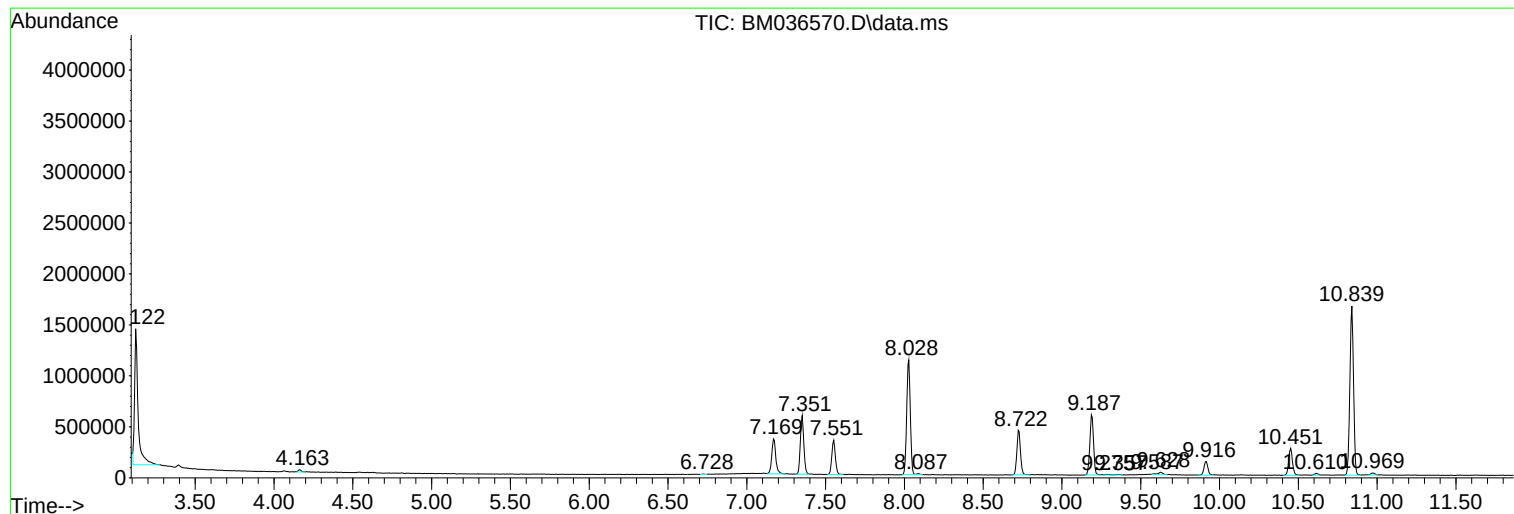
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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
Data File : BM036570.D
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Operator : CG/JU
Sample : N4601-19
Misc :
ALS Vial : 16 Sample Multiplier: 1

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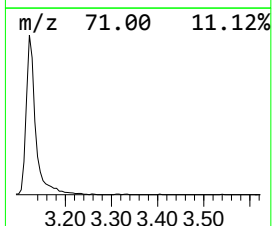
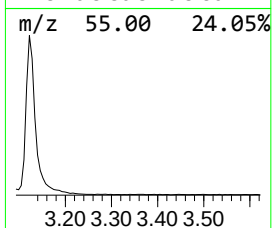
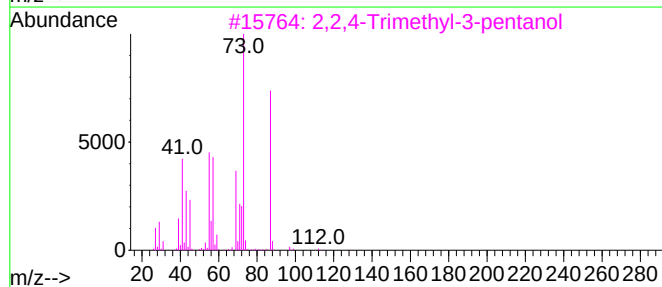
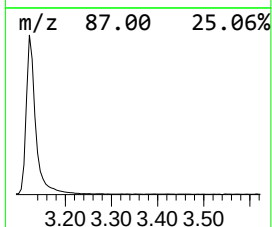
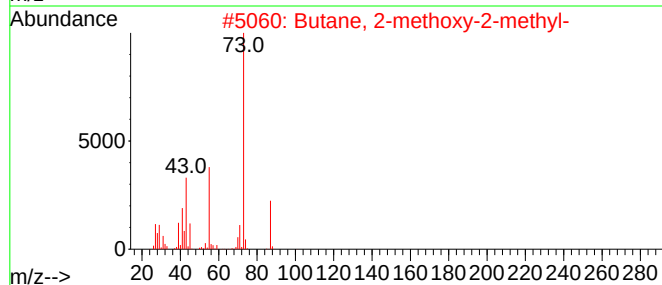
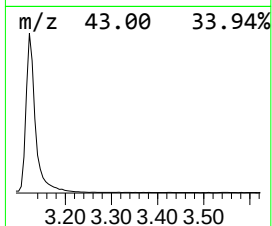
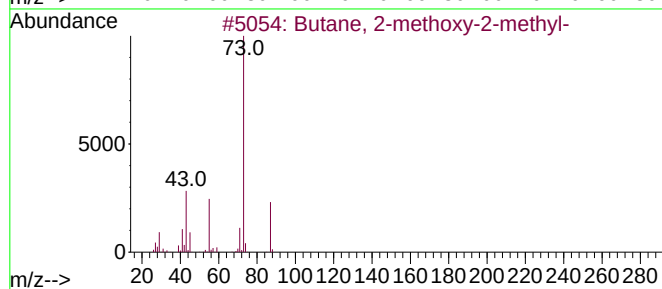
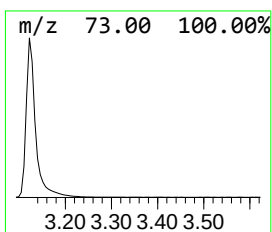
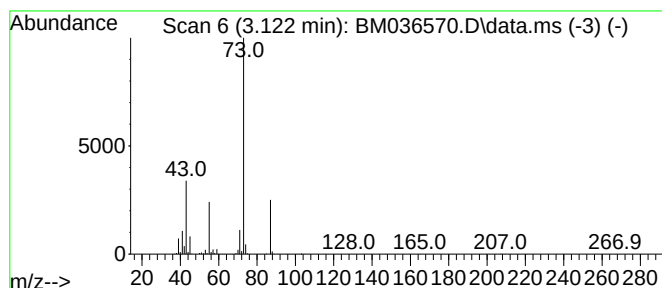
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.122	23.13 ng/ul	2074780	1,4-Dichlorobenzene-d4	8.028

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	74
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	23



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ALS Vial : 16 Sample Multiplier: 1

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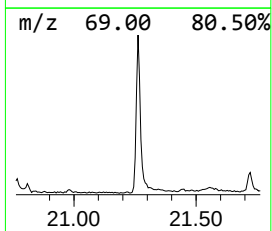
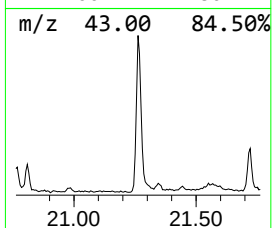
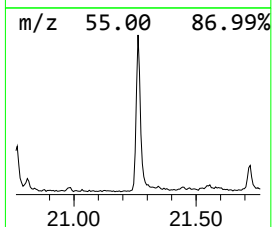
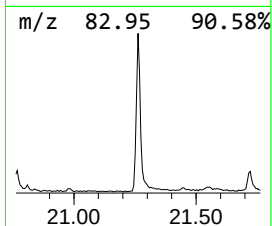
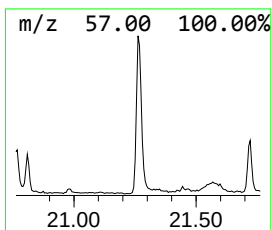
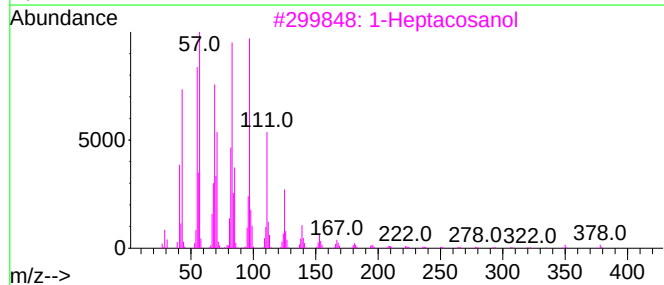
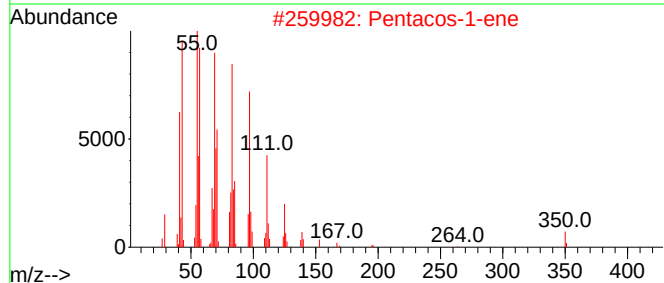
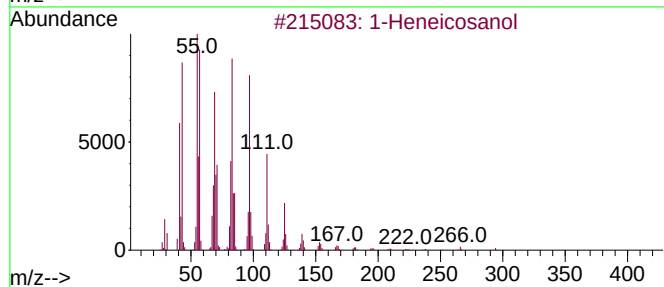
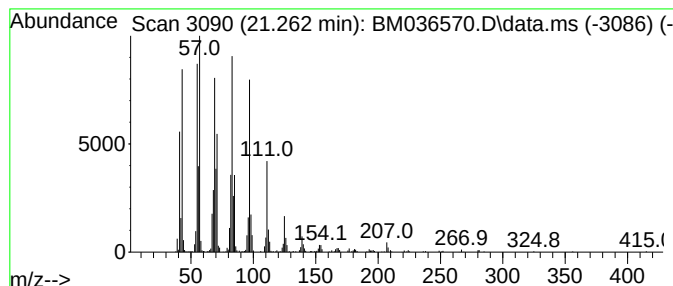
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 1-Heneicosanol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.262	3.35 ng/ul	869100	Chrysene-d12	21.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heneicosanol	312	C21H44O	015594-90-8	95
2			Pentacos-1-ene	350	C25H50	016980-85-1	94
3			1-Heptacosanol	396	C27H56O	002004-39-9	94
4			n-Tetracosanol-1	354	C24H50O	000506-51-4	94
5			Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	94



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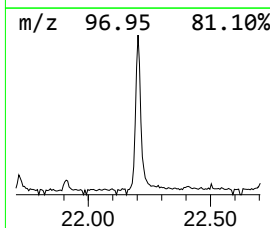
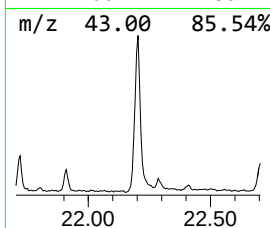
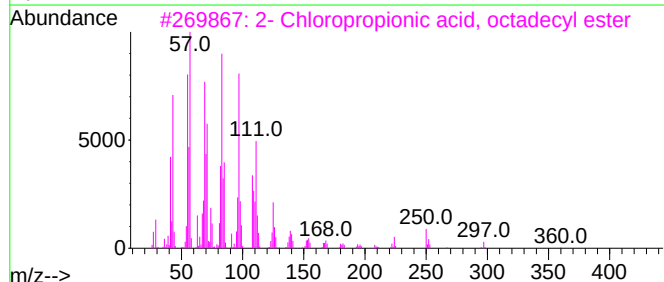
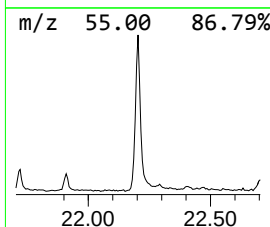
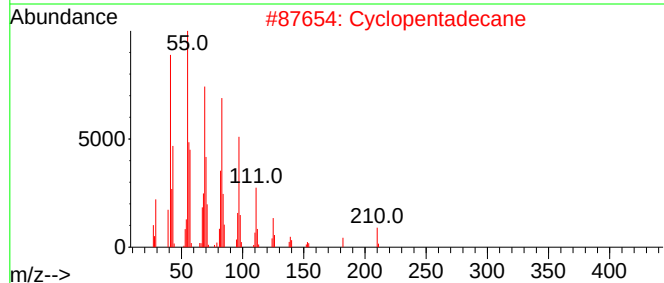
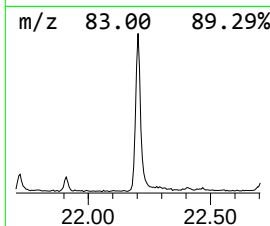
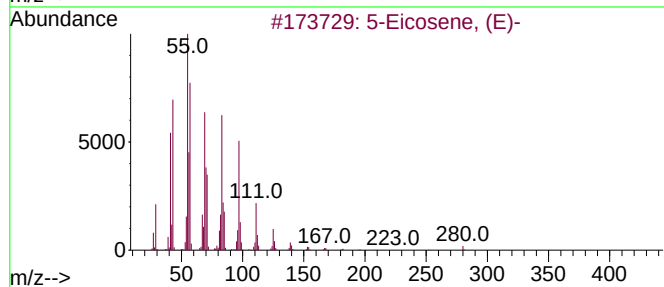
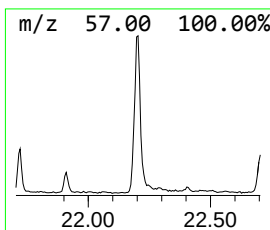
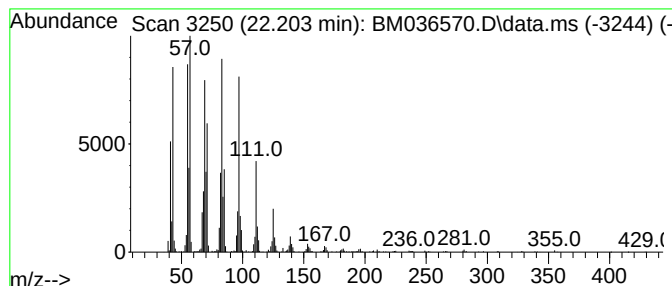
Instrument :
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.203	4.60 ng/ul	1191650	Chrysene-d12	21.551	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Eicosene, (E)-	280	C20H40	074685-30-6	96
2	Cyclopentadecane	210	C15H30	000295-48-7	94
3	2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	94
4	1-Heneicosyl formate	340	C22H44O2	077899-03-7	94
5	1-Heptacosanol	396	C27H56O	002004-39-9	94



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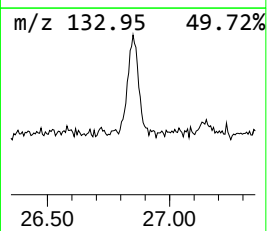
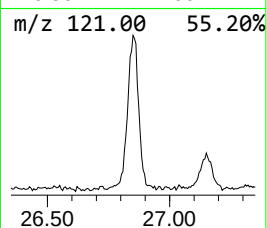
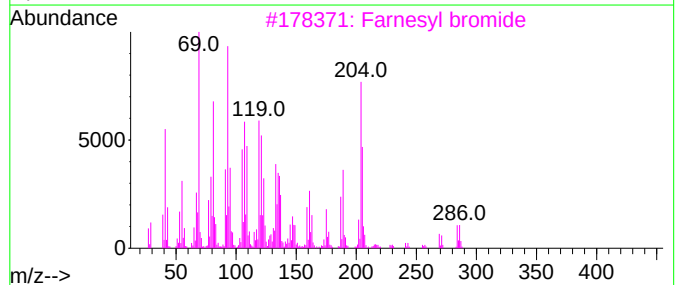
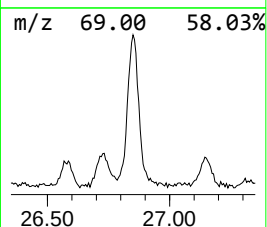
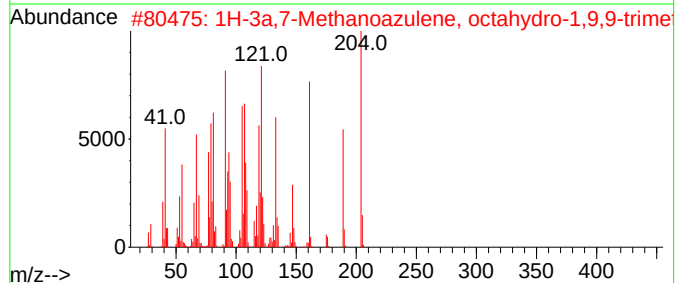
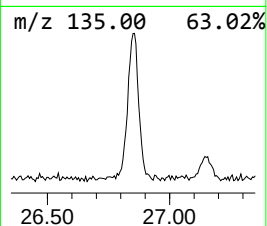
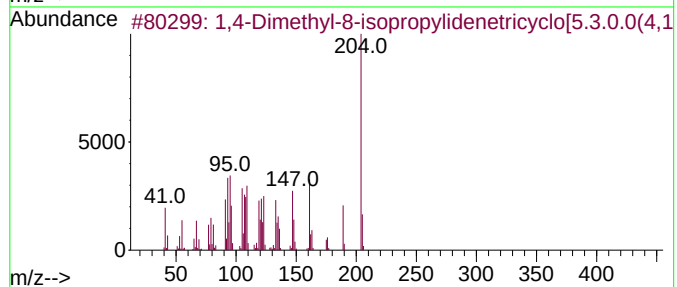
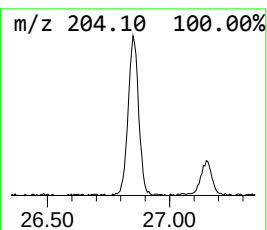
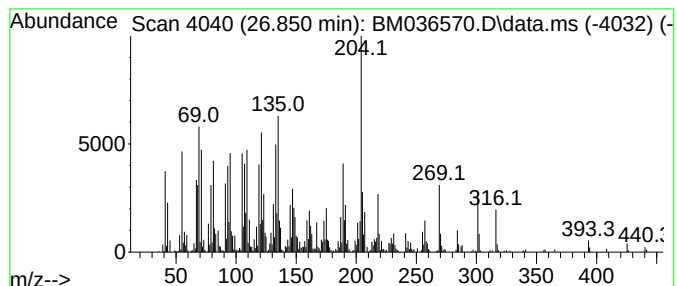
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 1,4-Dimethyl-8-isopropylide... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.850	6.86 ng/ul	1832140	Perylene-d12	23.991

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,4-Dimethyl-8-isopropylidenetri...	204	C15H24	1010140-07-7	59
2			1H-3a,7-Methanoazulene, octahydr...	204	C15H24	000508-55-4	55
3			Farnesyl bromide	284	C15H25Br	006874-67-5	49
4			1H-3a,7-Methanoazulene, 2,3,6,7,...	204	C15H24	000560-32-7	49
5			Guaia-3,9-diene	204	C15H24	000489-83-8	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
Data File : BM036570.D
Acq On : 16 Sep 2022 19:34
Operator : CG/JU
Sample : N4601-19
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
A5760

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	3.122	23.1	ng/ul	2074780	1	8.028	1794160	20.0
1-Heneicosanol	21.262	3.4	ng/ul	869100	5	21.551	5181540	20.0
(DEL) Alkane: S...	22.203	4.6	ng/ul	1191650	5	21.551	5181540	20.0
1,4-Dimethyl-8-...	26.850	6.9	ng/ul	1832140	6	23.991	5342660	20.0