

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM041421\
 Data File : BM029464.D
 Acq On : 14 Apr 2021 12:34
 Operator : CG/JU
 Sample : M1906-08
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 X3313

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 0.1 % of largest Peak
 Max Peaks: 100
 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-BM041221.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.922	3	6	20	rVB	539619	727120	11.49%	0.899%
2	3.187	48	51	59	rVB	20352	30697	0.49%	0.038%
3	3.599	118	121	146	rBV	179695	370416	5.85%	0.458%
4	3.911	170	174	178	rVB	9733	14090	0.22%	0.017%
5	4.128	207	211	215	rVB3	12126	17095	0.27%	0.021%
6	4.228	224	228	245	rVB	221150	328899	5.20%	0.407%
7	4.534	276	280	283	rBV2	9588	11506	0.18%	0.014%
8	4.822	325	329	338	rBV	151305	221646	3.50%	0.274%
9	6.828	663	670	687	rBV2	1108994	2349301	37.13%	2.905%
10	7.016	696	702	716	rBV	260105	437119	6.91%	0.541%
11	7.210	729	735	737	rBV	405566	611907	9.67%	0.757%
12	7.246	737	741	754	rVB	887094	1420519	22.45%	1.757%
13	7.675	807	814	824	rBV	424620	658017	10.40%	0.814%
14	8.381	924	934	939	rBV	470064	722608	11.42%	0.894%
15	8.440	939	944	949	rVV	476596	734202	11.60%	0.908%
16	8.499	949	954	975	rVB	982584	1567555	24.77%	1.939%
17	8.752	989	997	1004	rBV	629405	1047964	16.56%	1.296%
18	8.828	1004	1010	1021	rVV	354475	589047	9.31%	0.728%
19	9.263	1080	1084	1089	rBV2	7244	8877	0.14%	0.011%
20	9.546	1126	1132	1143	rBV	303715	499346	7.89%	0.618%
21	10.104	1216	1227	1239	rBV2	1209212	2683718	42.41%	3.319%
22	10.457	1280	1287	1291	rBV	572981	961233	15.19%	1.189%
23	10.510	1291	1296	1303	rVB	1474005	2348616	37.12%	2.905%
24	10.598	1303	1311	1322	rBV	361222	662578	10.47%	0.819%
25	11.416	1442	1450	1462	rBV5	18984	66614	1.05%	0.082%
26	11.998	1543	1549	1554	rBV2	21021	36069	0.57%	0.045%
27	12.051	1554	1558	1563	rVB3	6830	11379	0.18%	0.014%
28	12.128	1563	1571	1581	rBV	1794642	2759384	43.61%	3.413%
29	12.345	1604	1608	1617	rVB	19519	30825	0.49%	0.038%
30	12.798	1679	1685	1690	rBV2	25013	40809	0.64%	0.050%
31	13.187	1735	1751	1766	rBV	1918998	2881196	45.54%	3.563%
32	13.728	1837	1843	1848	rBV	879293	1196919	18.92%	1.480%
33	13.775	1848	1851	1859	rVB	63062	90958	1.44%	0.112%
34	14.010	1879	1891	1903	rBV	892029	1333446	21.07%	1.649%

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Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM041221.M
 Title : SVOA CALIBRATION

35	14.316	1937	1943	1948	rVV	828996	1215504	19.21%	1.503%
36	14.381	1948	1954	1961	rVB	1318201	1865472	29.48%	2.307%
37	14.516	1970	1977	1994	rBV	430426	654621	10.35%	0.810%
38	14.739	2011	2015	2020	rVB6	4996	8753	0.14%	0.011%
39	15.128	2074	2081	2086	rBV3	11544	25958	0.41%	0.032%
40	15.181	2086	2090	2096	rVB2	16726	22540	0.36%	0.028%
41	15.310	2105	2112	2117	rBV	1208132	1672232	26.43%	2.068%
42	15.369	2117	2122	2127	rVV	1532506	2113987	33.41%	2.614%
43	15.445	2127	2135	2146	rVV2	1700095	2706660	42.78%	3.347%
44	15.551	2149	2153	2161	rVV4	14145	30673	0.48%	0.038%
45	15.680	2169	2175	2181	rVB6	7730	12919	0.20%	0.016%
46	15.880	2203	2209	2215	rBV9	5811	11797	0.19%	0.015%
47	15.998	2224	2229	2232	rBV5	7120	11362	0.18%	0.014%
48	16.045	2232	2237	2243	rVV6	6318	15091	0.24%	0.019%
49	16.098	2243	2246	2252	rVB5	8189	12640	0.20%	0.016%
50	16.216	2259	2266	2275	rBV6	10248	20269	0.32%	0.025%
51	16.304	2277	2281	2286	rVB	16000	20932	0.33%	0.026%
52	16.475	2302	2310	2314	rBV5	17023	34038	0.54%	0.042%
53	16.533	2314	2320	2325	rBV	1969975	2598712	41.07%	3.214%
54	16.710	2347	2350	2356	rVB3	8149	12124	0.19%	0.015%
55	16.804	2361	2366	2368	rVV6	6508	10233	0.16%	0.013%
56	16.839	2368	2372	2375	rVV4	11024	18887	0.30%	0.023%
57	16.880	2375	2379	2386	rVB	36709	53672	0.85%	0.066%
58	16.992	2394	2398	2403	rVB2	16732	20695	0.33%	0.026%
59	17.057	2403	2409	2413	rBV	912719	1410025	22.28%	1.744%
60	17.104	2413	2417	2422	rVV	2258944	3206501	50.68%	3.965%
61	17.157	2422	2426	2432	rVB	1374140	1850294	29.24%	2.288%
62	17.351	2455	2459	2464	rBV3	12969	21731	0.34%	0.027%
63	17.392	2464	2466	2470	rVB5	7656	10521	0.17%	0.013%
64	17.463	2470	2478	2490	rBV	3027485	4231151	66.87%	5.233%
65	17.710	2512	2520	2522	rBV7	14960	25897	0.41%	0.032%
66	17.857	2538	2545	2550	rBV3	11209	24474	0.39%	0.030%
67	17.963	2558	2563	2568	rBV	114804	161660	2.55%	0.200%
68	18.033	2572	2575	2578	rVB	31710	38346	0.61%	0.047%
69	18.127	2588	2591	2599	rVB7	6848	11742	0.19%	0.015%
70	18.463	2643	2648	2654	rBV	37401	61204	0.97%	0.076%
71	18.727	2686	2693	2698	rBV	271570	336911	5.32%	0.417%

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72	18.816	2704	2708	2716	rBV	67589	98384	1.55%	0.122%
73	18.886	2716	2720	2727	rVB3	28481	48291	0.76%	0.060%
74	19.086	2748	2754	2759	rBV	20580	32659	0.52%	0.040%
75	19.133	2760	2762	2769	rVB4	12712	23277	0.37%	0.029%
76	19.245	2777	2781	2789	rBV3	40436	66470	1.05%	0.082%
77	19.333	2791	2796	2797	rVV5	19446	27755	0.44%	0.034%
78	19.363	2797	2801	2805	rVV	82193	118603	1.87%	0.147%
79	19.451	2809	2816	2818	rVV	1454371	2018079	31.89%	2.496%
80	19.480	2818	2821	2827	rVV	3133340	4034850	63.77%	4.990%
81	19.527	2827	2829	2832	rVB2	14647	14020	0.22%	0.017%
82	19.904	2889	2893	2897	rVB	33120	40944	0.65%	0.051%
83	20.115	2926	2929	2935	rVB4	16633	23832	0.38%	0.029%
84	20.386	2970	2975	2980	rBV	3015785	3450050	54.53%	4.267%
85	20.586	3005	3009	3014	rBV	360994	423944	6.70%	0.524%
86	20.633	3014	3017	3021	rVB	68516	84894	1.34%	0.105%
87	20.698	3026	3028	3032	rVB2	28086	27546	0.44%	0.034%
88	20.968	3069	3074	3087	rBV2	423758	743207	11.75%	0.919%
89	21.074	3090	3092	3096	rVB	29425	32485	0.51%	0.040%
90	21.168	3103	3108	3113	rBV	3430963	3656917	57.79%	4.523%
91	21.239	3115	3120	3126	rVB	997455	1197228	18.92%	1.481%
92	21.833	3218	3221	3230	rBV4	48592	99672	1.58%	0.123%
93	22.292	3295	3299	3305	rVB2	87526	104228	1.65%	0.129%
94	22.792	3380	3384	3389	rVB	135168	200582	3.17%	0.248%
95	23.309	3465	3472	3476	rBV	904882	1672530	26.43%	2.068%
96	23.451	3489	3496	3506	rVB2	596533	1146908	18.13%	1.418%
97	23.986	3582	3587	3595	rVB	189825	338274	5.35%	0.418%
98	24.721	3708	3712	3719	rVB3	147646	301756	4.77%	0.373%
99	25.680	3864	3875	3897	rVB2	2176932	6327416	100.00%	7.825%
100	26.350	3979	3989	4004	rVB	815221	2473467	39.09%	3.059%

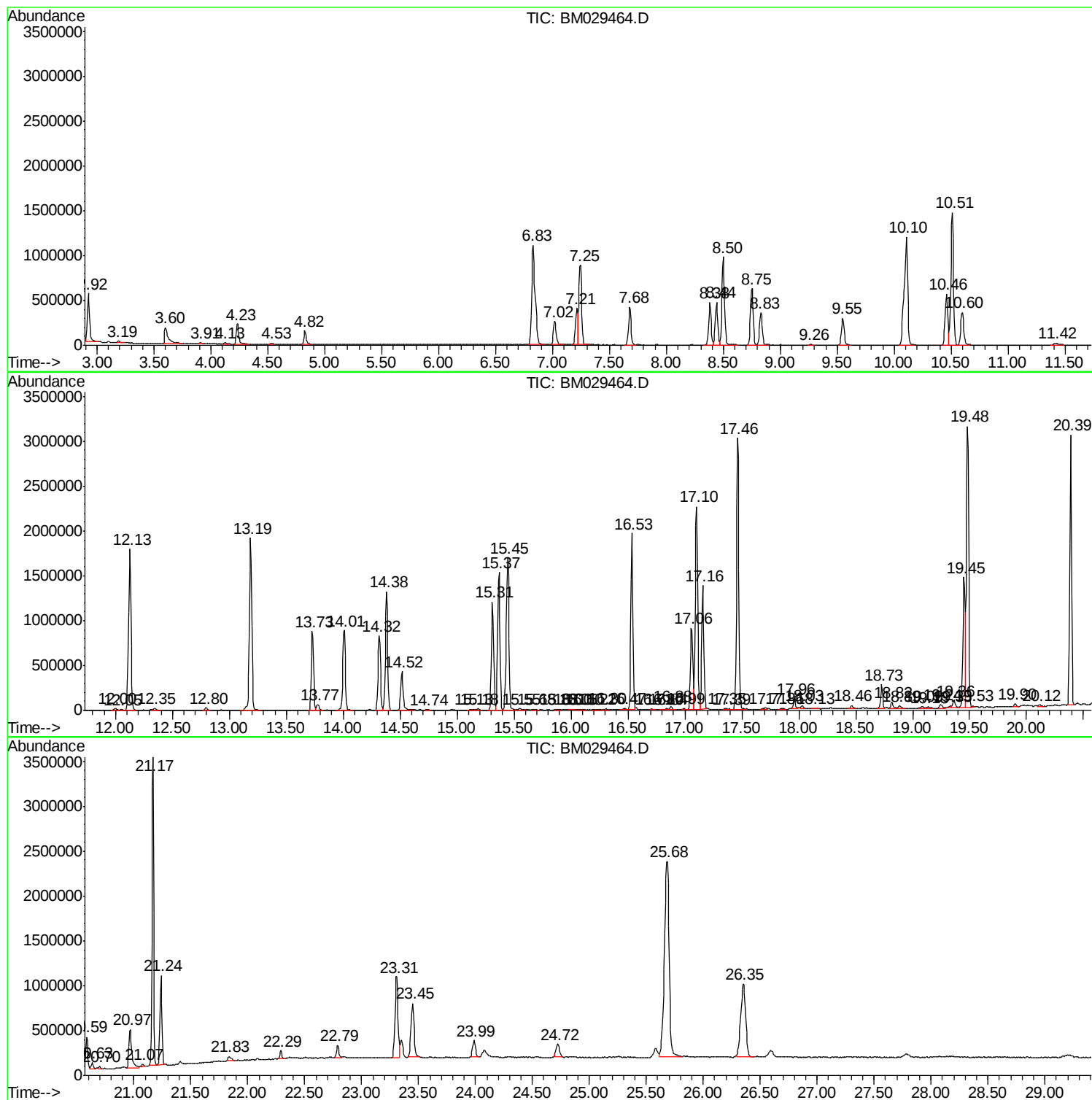
Sum of corrected areas: 80860171

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM041221.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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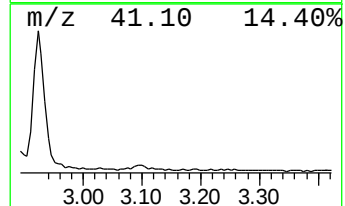
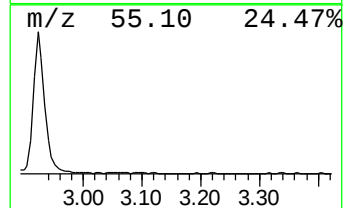
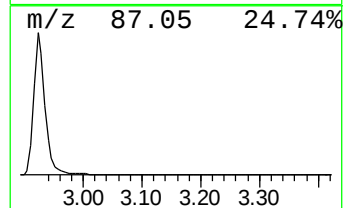
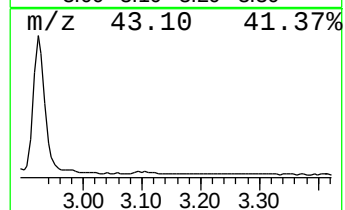
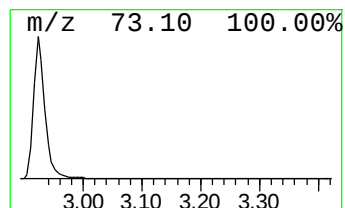
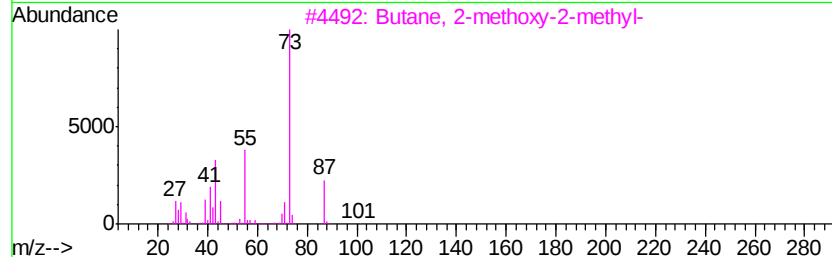
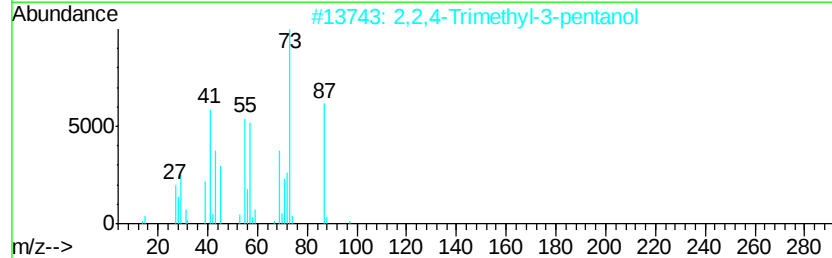
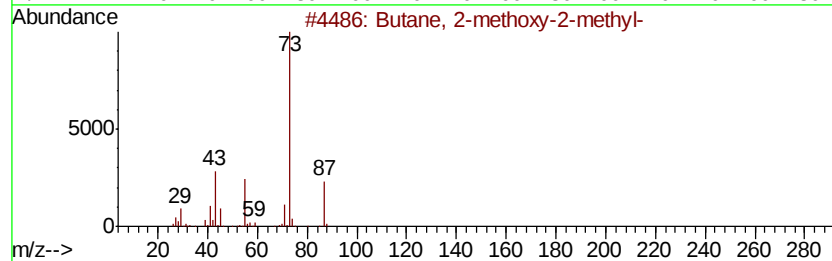
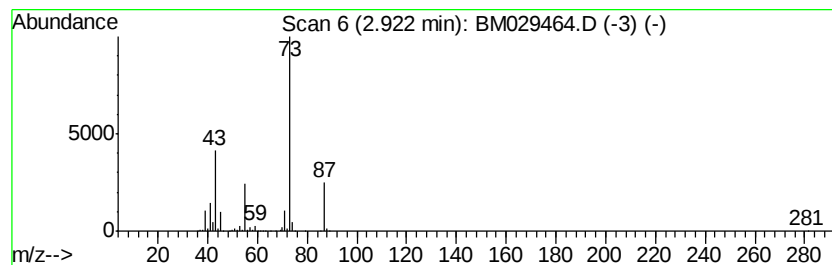
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM041221.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.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.
2.92	22.10 ng/ul	727120	1,4-Dichlorobenzene-d4	7.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	64
3		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
4		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
5		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	23



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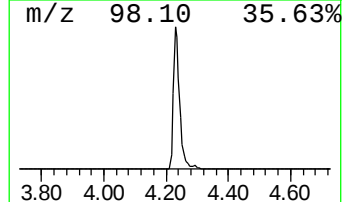
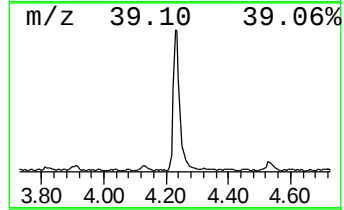
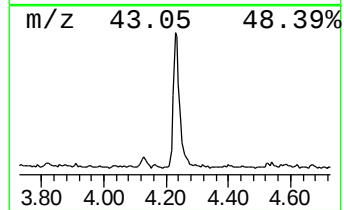
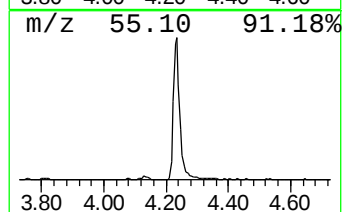
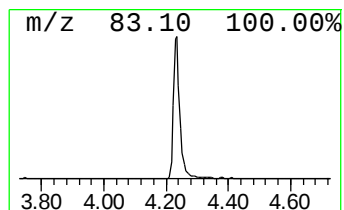
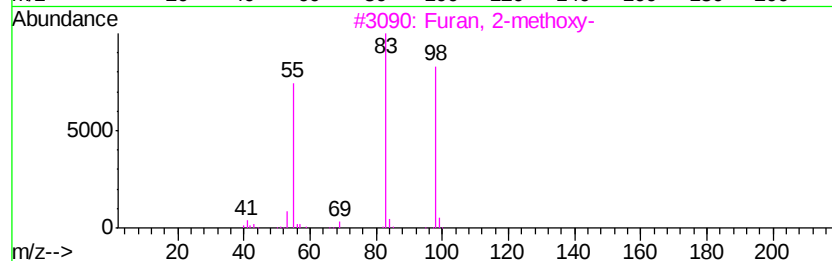
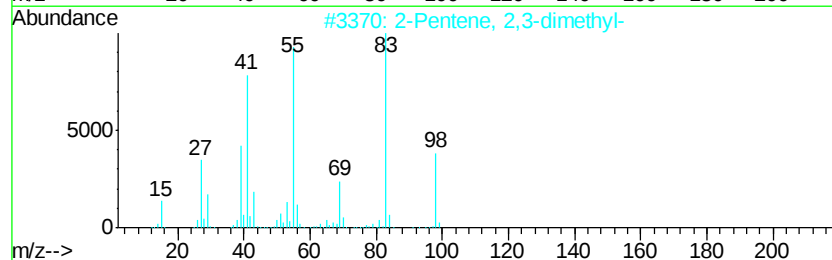
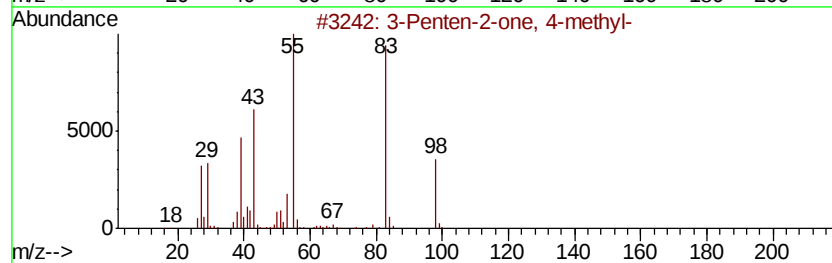
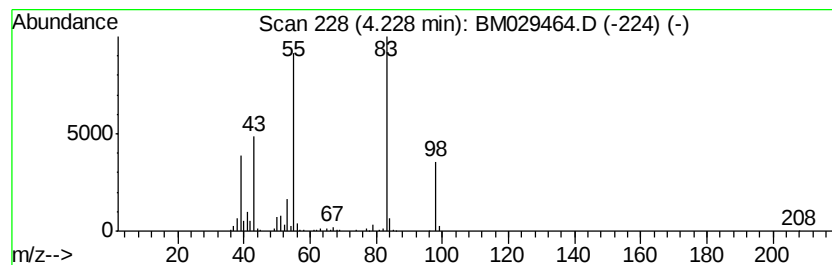
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM041221.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 3-Penten-2-one, 4-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.23	10.00 ng/ul	328899	1,4-Dichlorobenzene-d4	7.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	94
2		2-Pentene, 2,3-dimethyl-	98	C7H14	010574-37-5	90
3		Furan, 2-methoxy-	98	C5H6O2	025414-22-6	83
4		3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	76
5		3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	76



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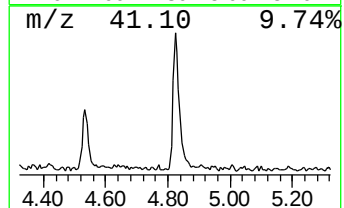
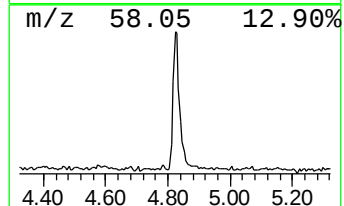
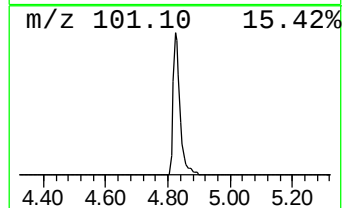
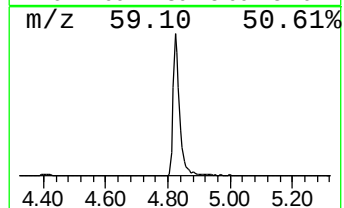
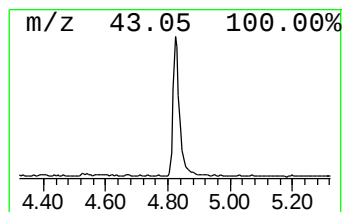
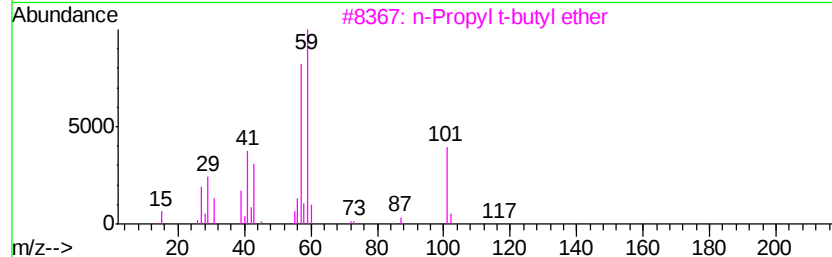
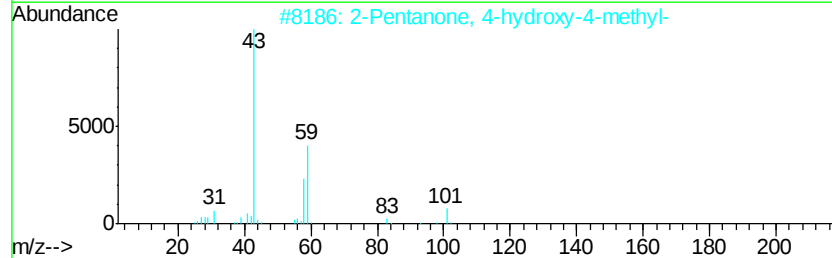
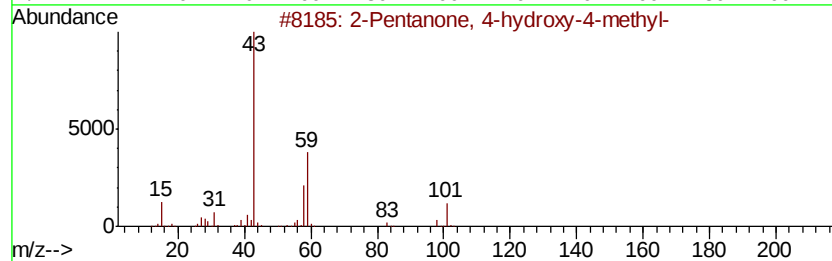
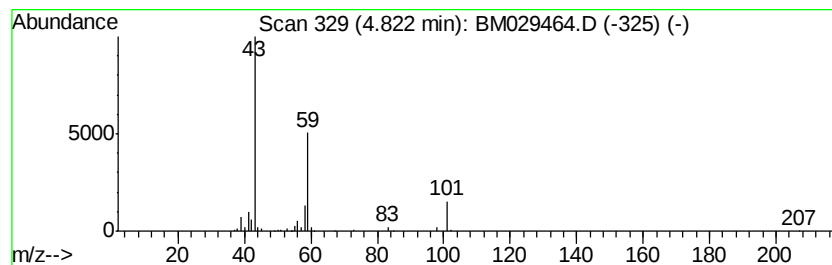
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.82	6.74 ng/ul	221646	1,4-Dichlorobenzene-d4	7.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
3		n-Propyl t-butyl ether	116	C7H16O	1000334-81-9	33
4		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	28
5		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM041421\
Data File : BM029464.D
Acq On : 14 Apr 2021 12:34
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Sample : M1906-08
Misc :
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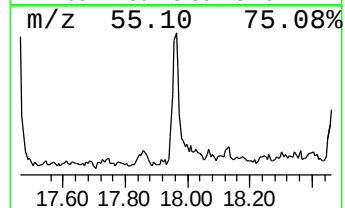
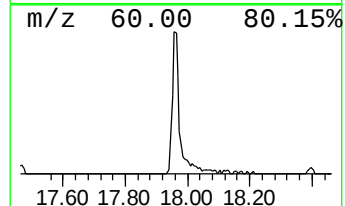
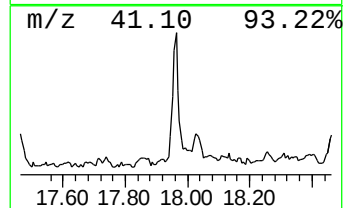
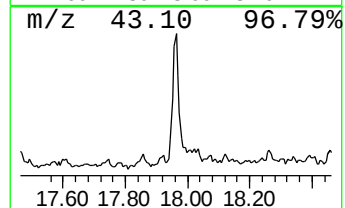
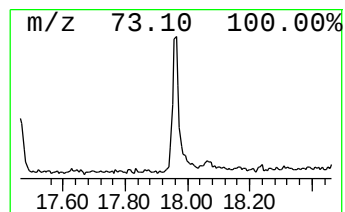
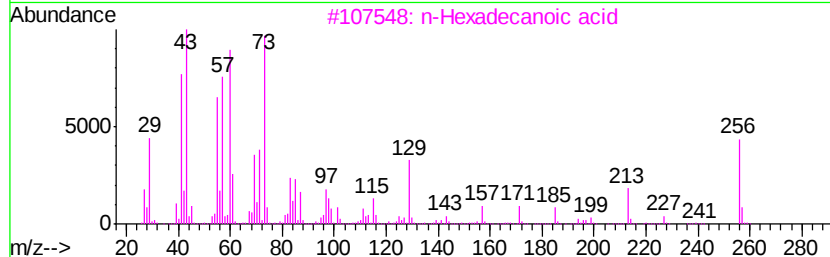
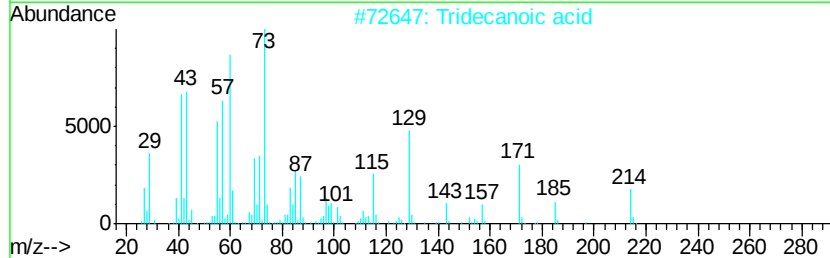
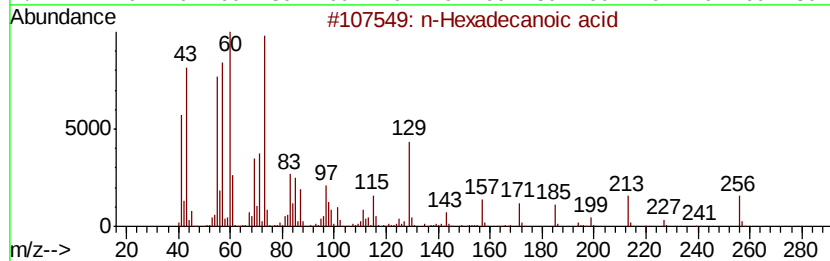
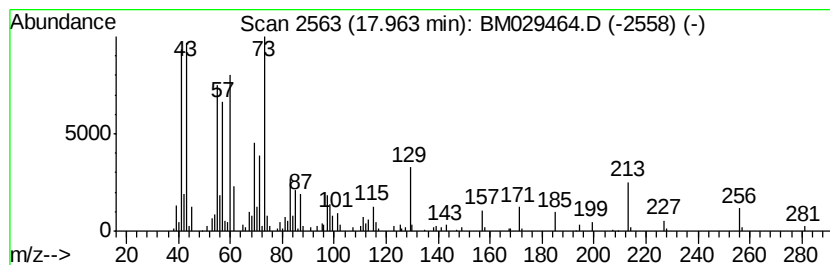
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 n-Hexadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.96	2.29 ng/ul	161660	Phenanthrene-d10	17.06

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tridecanoic acid	214	C13H26O2	000638-53-9	90
3		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	76
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	76
5		Tetradecanoic acid	228	C14H28O2	000544-63-8	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM041421\
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ClientSampleId :
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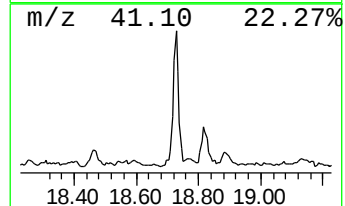
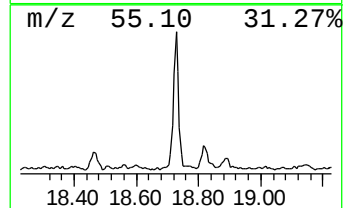
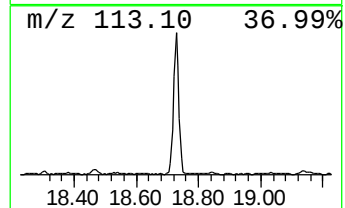
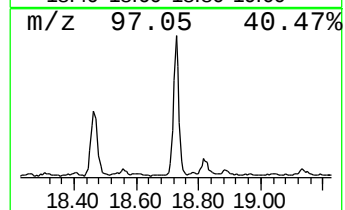
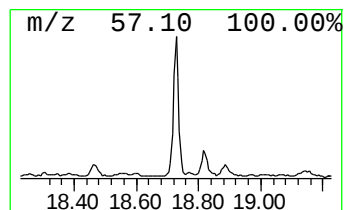
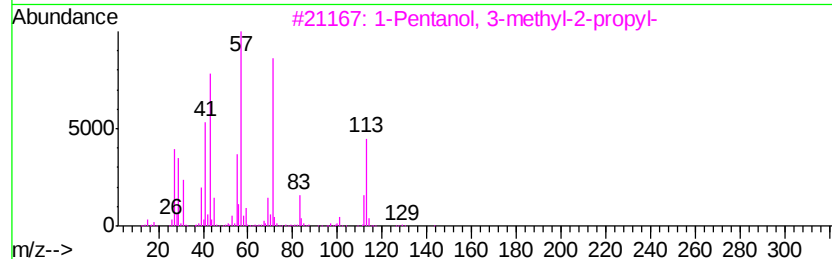
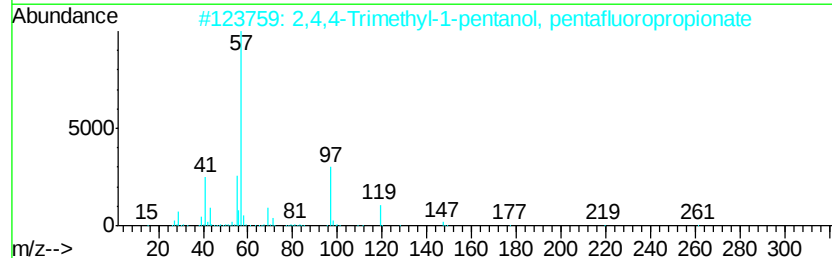
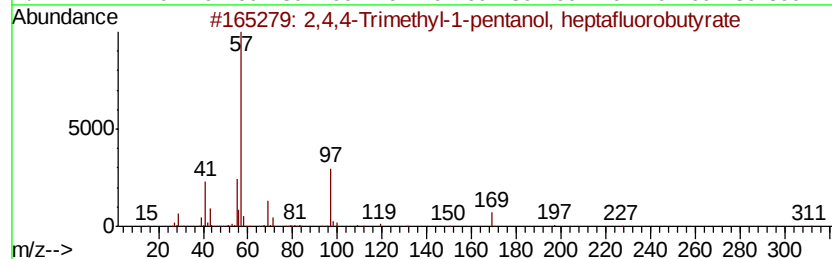
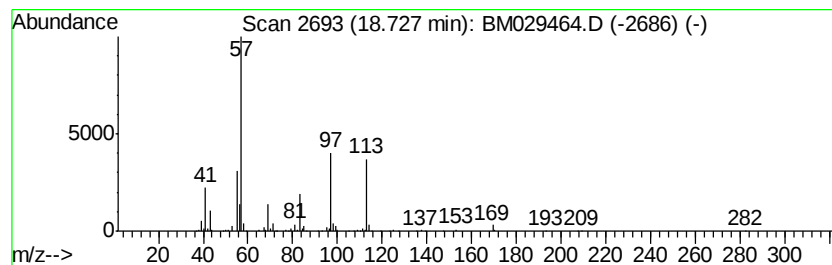
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 2,4,4-Trimethyl-1-pentanol,... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.73	4.78 ng/ul	336911	Phenanthrene-d10	17.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,4,4-Trimethyl-1-pentanol, hept...	326	C12H17F7O2	1000365-01-4	50
2		2,4,4-Trimethyl-1-pentanol, pent...	276	C11H17F5O2	1000365-51-0	47
3		1-Pentanol, 3-methyl-2-propyl-	144	C9H20O	054004-40-9	47
4		2,4,4,6,6,8,8-Heptamethyl-1-nonene	224	C16H32	015796-04-0	45
5		2,4,4-Trimethyl-1-pentanol, trif...	226	C10H17F3O2	1000365-19-5	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM041421\
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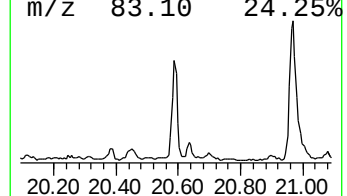
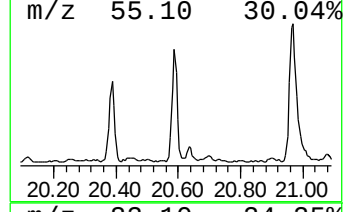
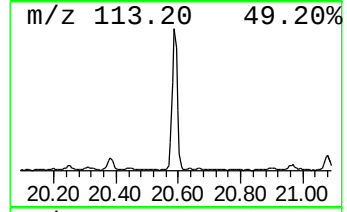
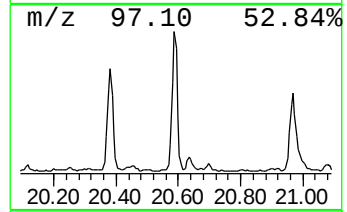
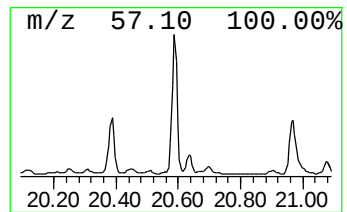
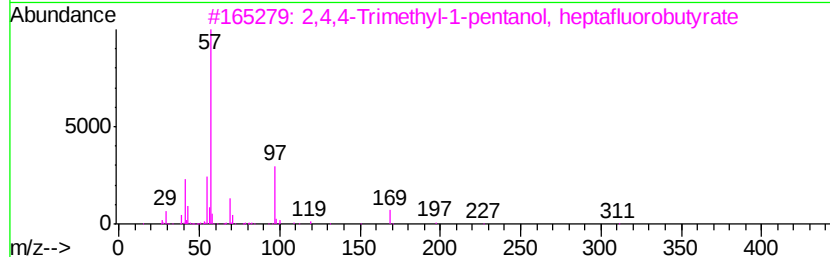
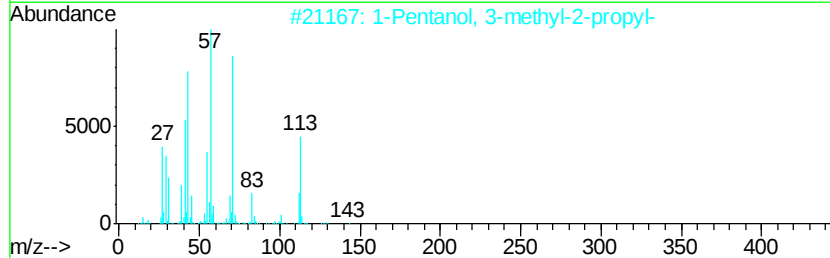
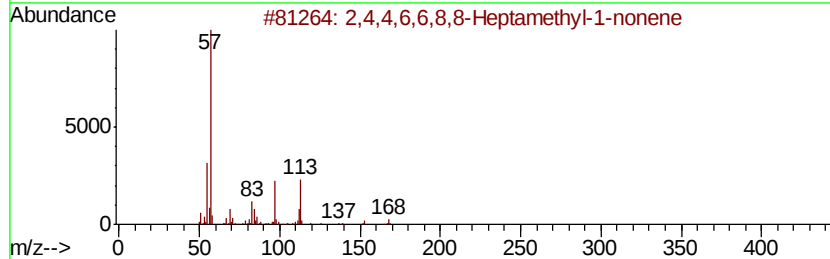
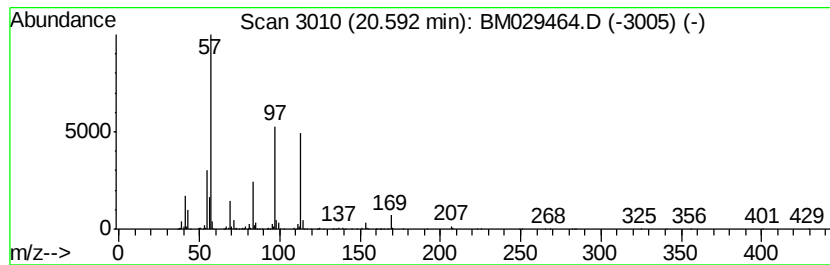
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.59	7.08 ng/ul	423944	Chrysene-d12	21.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,4,4,6,6,8,8-Heptamethyl-1-nonene	224	C16H32	015796-04-0	59
2		1-Pentanol, 3-methyl-2-propyl-	144	C9H20O	054004-40-9	40
3		2,4,4-Trimethyl-1-pentanol, hept...	326	C12H17F7O2	1000365-01-4	32
4		Eicosyl heptafluorobutylrate	494	C24H41F7O2	1000351-83-5	27
5		2,4,4-Trimethyl-1-pentanol, pent...	276	C11H17F5O2	1000365-51-0	27



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM041421\
Data File : BM029464.D
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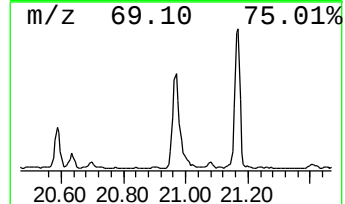
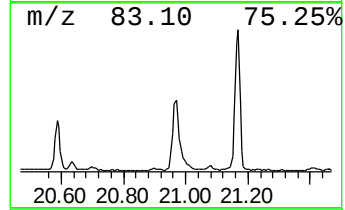
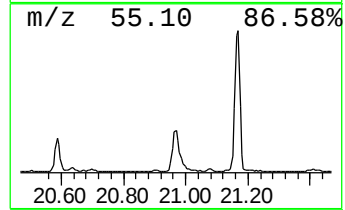
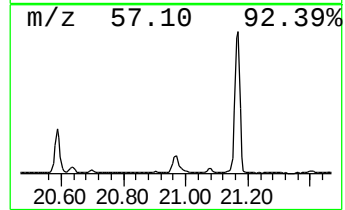
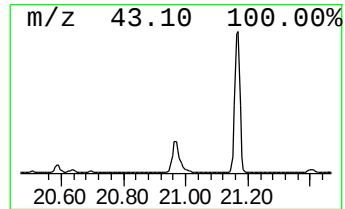
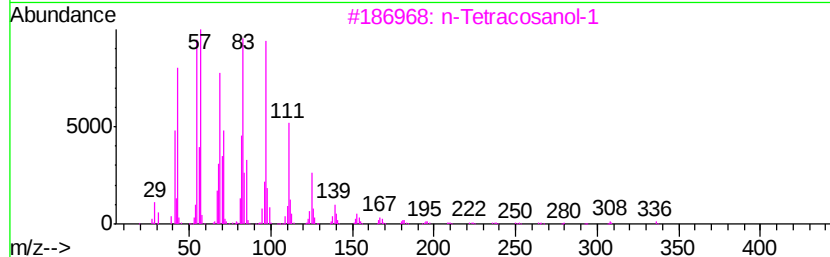
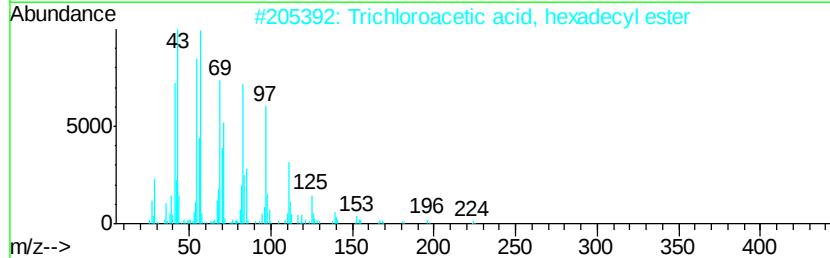
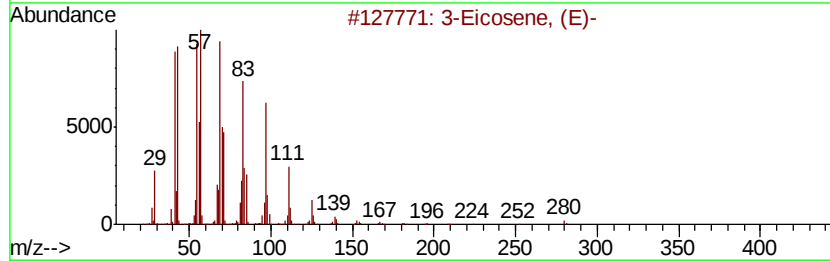
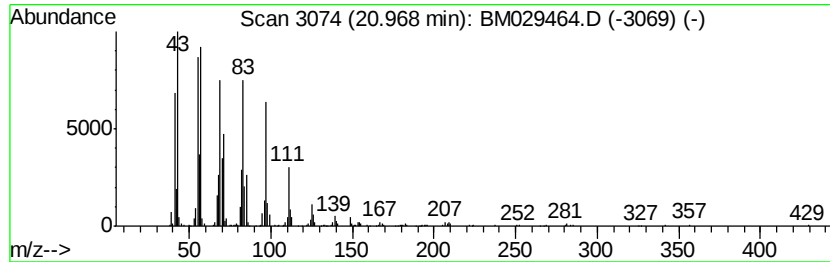
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.97	12.42 ng/ul	743207	Chrysene-d12	21.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Eicosene, (E)-	280	C20H40	074685-33-9	94
2		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
3		n-Tetracosanol-1	354	C24H50O	000506-51-4	94
4		1-Docosene	308	C22H44	001599-67-3	93
5		1-Nonadecene	266	C19H38	018435-45-5	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM041421\
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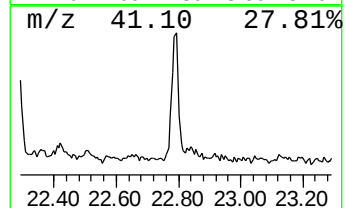
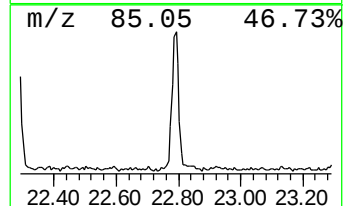
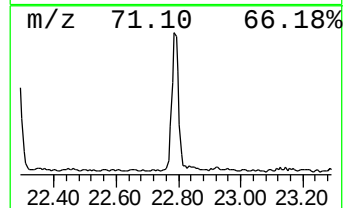
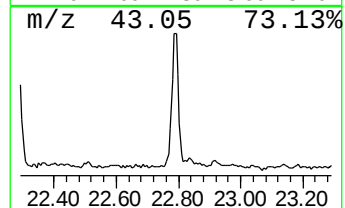
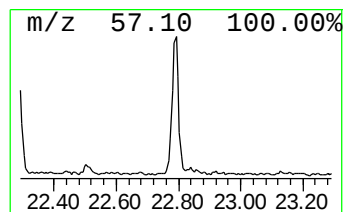
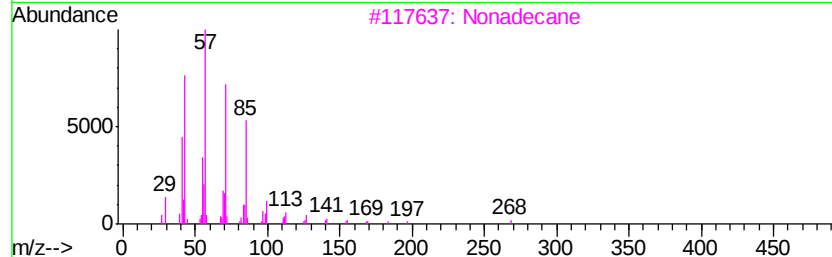
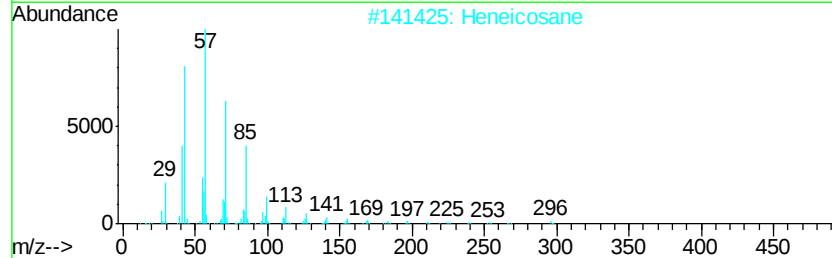
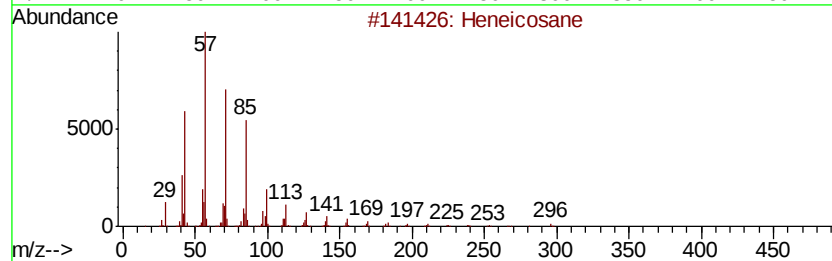
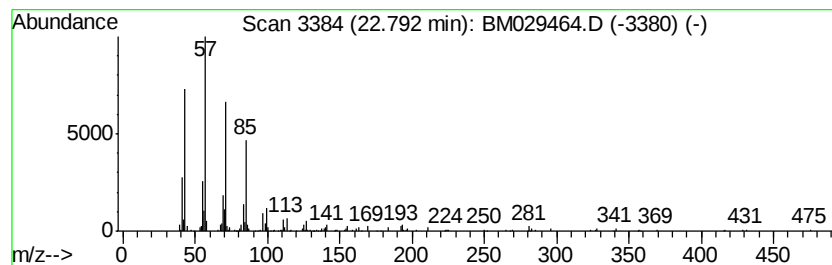
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TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.79	3.50 ng/ul	200582	Perylene-d12	23.45

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	95
2		Heneicosane	296	C21H44	000629-94-7	91
3		Nonadecane	268	C19H40	000629-92-5	90
4		Heptacosane	380	C27H56	000593-49-7	87
5		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM041421\
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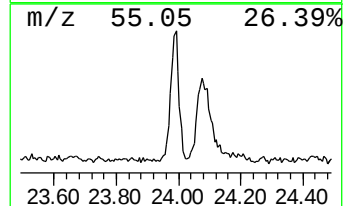
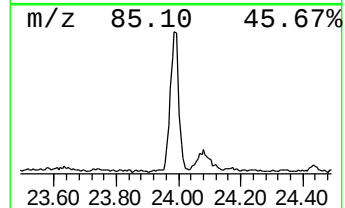
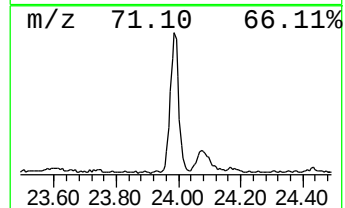
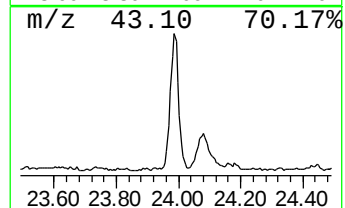
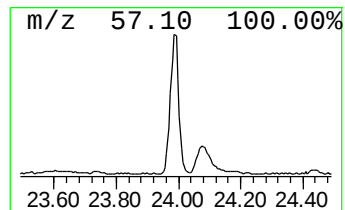
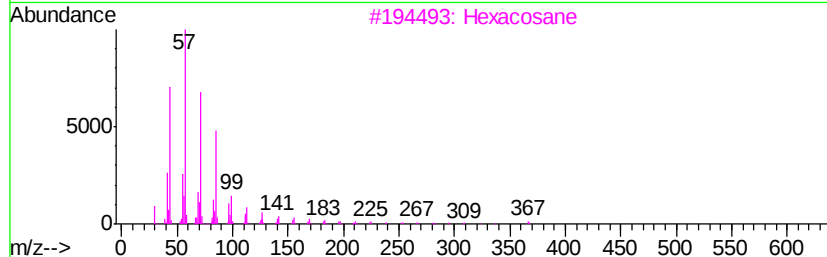
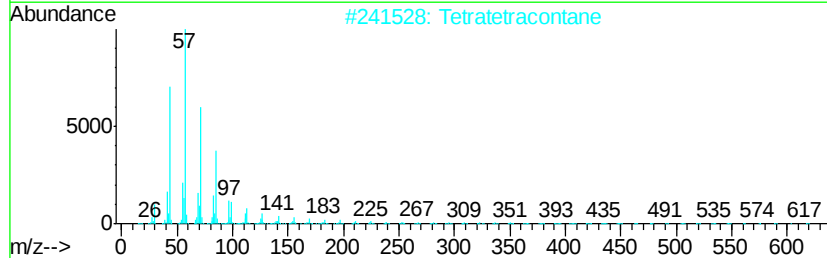
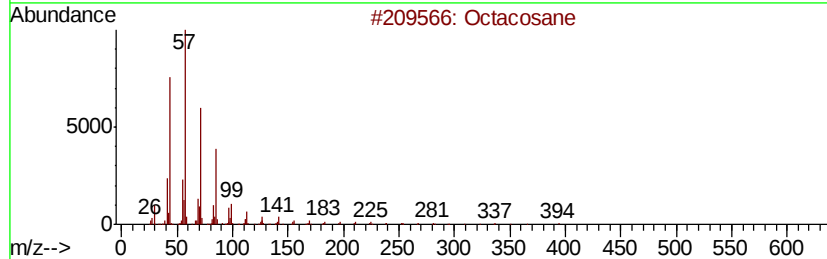
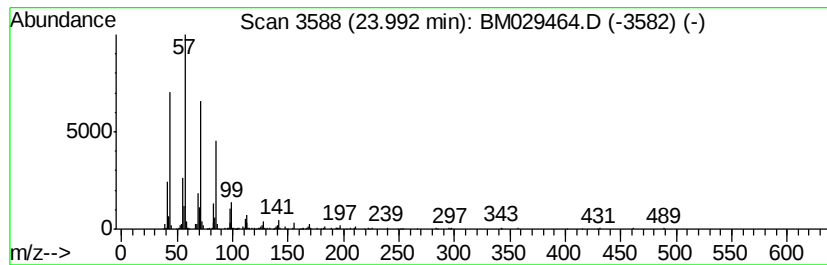
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.99	5.90 ng/ul	338274	Perylene-d12	23.45

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	90
2		Tetratetracontane	619	C44H90	007098-22-8	90
3		Hexacosane	366	C26H54	000630-01-3	90
4		Heptacosane	380	C27H56	000593-49-7	87
5		Eicosane	282	C20H42	000112-95-8	87



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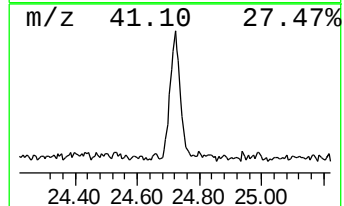
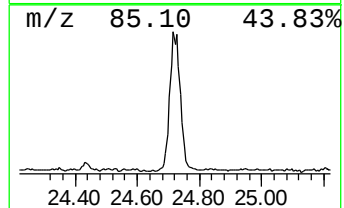
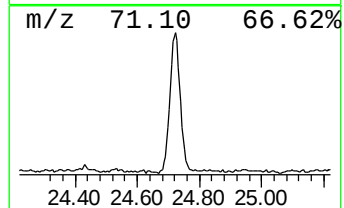
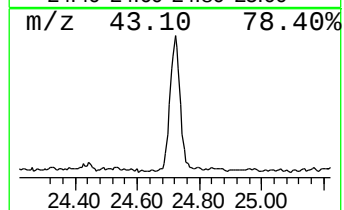
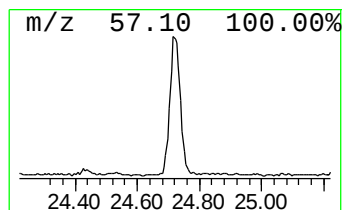
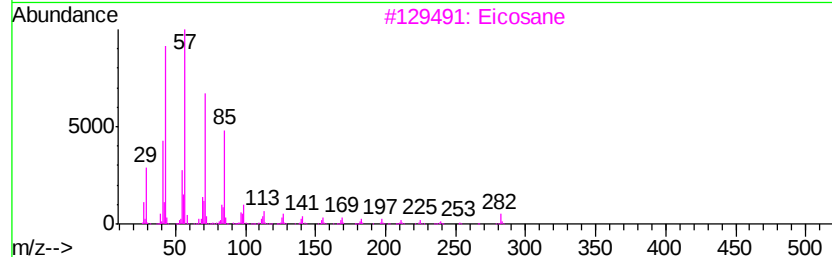
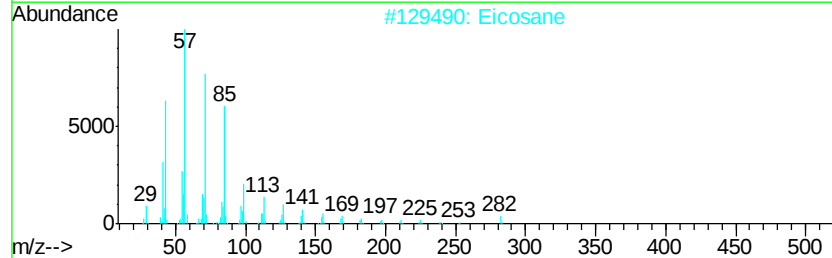
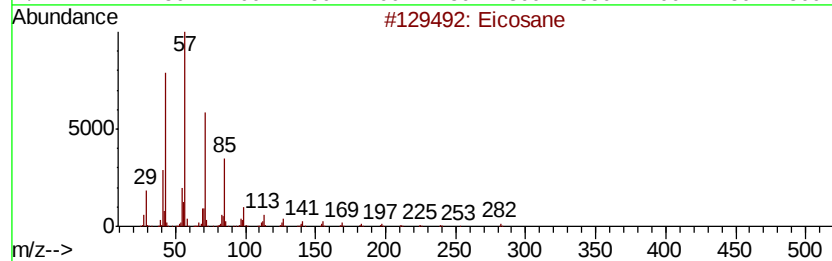
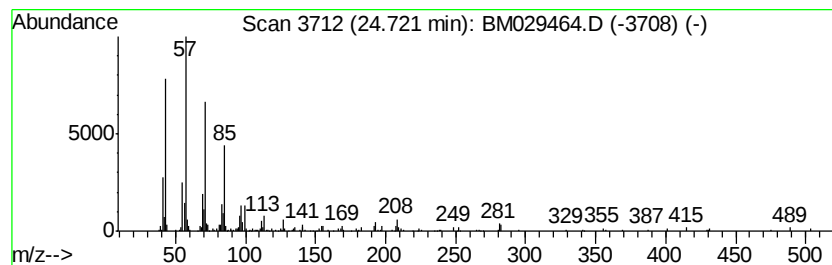
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM041221.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.72	5.26 ng/ul	301756	Perylene-d12	23.45

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	94
2		Eicosane	282	C20H42	000112-95-8	90
3		Eicosane	282	C20H42	000112-95-8	90
4		Hexacosane	366	C26H54	000630-01-3	76
5		Tetratetracontane	619	C44H90	007098-22-8	76



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM041421\
 Data File : BM029464.D
 Acq On : 14 Apr 2021 12:34
 Operator : CG/JU
 Sample : M1906-08
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 X3313

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM041221.M
 Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	2.92	22.1	ng/ul	727120	1	7.68	658017	20.0
3-Penten-2-one, 4...	4.23	10.0	ng/ul	328899	1	7.68	658017	20.0
2-Pentanone, 4-hy...	4.82	6.7	ng/ul	221646	1	7.68	658017	20.0
n-Hexadecanoic acid	17.96	2.3	ng/ul	161660	4	17.06	1410030	20.0
2,4,4-Trimethyl-1...	18.73	4.8	ng/ul	336911	4	17.06	1410030	20.0
(DEL) Alkane: Str...	20.59	7.1	ng/ul	423944	5	21.24	1197230	20.0
(DEL) Alkane: Str...	20.97	12.4	ng/ul	743207	5	21.24	1197230	20.0
(DEL) Alkane: Str...	22.79	3.5	ng/ul	200582	6	23.45	1146910	20.0
(DEL) Alkane: Str...	23.99	5.9	ng/ul	338274	6	23.45	1146910	20.0
(DEL) Alkane: Str...	24.72	5.3	ng/ul	301756	6	23.45	1146910	20.0