

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM111920\
 Data File : BM027844.D
 Acq On : 19 Nov 2020 14:59
 Operator : JU/CG
 Sample : L4710-16
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0B45

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\SOM-EPA-BM111220MA.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	3.104	14	20	25	rBV	78947	145605	18.30%	1.168%
2	3.193	31	35	49	rVB	539977	795839	100.00%	6.382%
3	3.493	83	86	94	rVV	5027	7486	0.94%	0.060%
4	3.969	163	167	171	rVB3	2462	3795	0.48%	0.030%
5	4.081	180	186	192	rBV	11271	16826	2.11%	0.135%
6	4.163	195	200	203	rVV4	2329	3622	0.46%	0.029%
7	4.222	206	210	219	rVB2	4749	8481	1.07%	0.068%
8	4.322	224	227	232	rVB2	2741	4166	0.52%	0.033%
9	4.563	264	268	274	rVB5	3113	5192	0.65%	0.042%
10	5.322	391	397	401	rBV2	3089	4907	0.62%	0.039%
11	5.416	410	413	418	rBV2	1935	3031	0.38%	0.024%
12	7.457	751	760	770	rVV	126774	210187	26.41%	1.686%
13	7.639	785	791	798	rVB	92297	143762	18.06%	1.153%
14	7.851	818	827	835	rBV	129252	203301	25.55%	1.630%
15	7.916	835	838	840	rBV2	2775	3340	0.42%	0.027%
16	8.334	902	909	915	rBV	202033	319858	40.19%	2.565%
17	9.034	1021	1028	1035	rVB	138058	220317	27.68%	1.767%
18	9.510	1102	1109	1115	rBV	123717	210155	26.41%	1.685%
19	9.910	1170	1177	1183	rBV	4980	7616	0.96%	0.061%
20	10.239	1225	1233	1239	rVB	103493	172476	21.67%	1.383%
21	10.775	1317	1324	1331	rVV	157901	263393	33.10%	2.112%
22	11.169	1384	1391	1398	rBV	273702	455561	57.24%	3.653%
23	11.292	1404	1412	1423	rBV	100309	166107	20.87%	1.332%
24	13.321	1752	1757	1761	rVV2	2742	4009	0.50%	0.032%
25	13.563	1793	1798	1803	rVV3	3331	5142	0.65%	0.041%
26	13.827	1840	1843	1848	rVV3	5089	6172	0.78%	0.049%
27	14.339	1924	1930	1935	rBV	270323	378045	47.50%	3.032%
28	14.386	1935	1938	1944	rVB	132750	170133	21.38%	1.364%
29	14.568	1965	1969	1975	rBV4	4629	7370	0.93%	0.059%
30	14.645	1976	1982	1992	rVB	302178	444721	55.88%	3.566%
31	14.821	2009	2012	2018	rVB4	2316	3515	0.44%	0.028%
32	14.951	2027	2034	2040	rBV	418084	592903	74.50%	4.755%
33	15.104	2054	2060	2072	rBV	120422	165561	20.80%	1.328%
34	15.257	2083	2086	2091	rBV2	3041	3428	0.43%	0.027%

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Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111220MA.M
 Title : SVOA CALIBRATION

35	15.333	2097	2099	2103	rVV3	3159	4085	0.51%	0.033%
36	15.715	2159	2164	2166	rBV	3840	6337	0.80%	0.051%
37	15.757	2168	2171	2176	rVB4	2962	4367	0.55%	0.035%
38	15.927	2194	2200	2206	rVV	369185	521931	65.58%	4.186%
39	16.027	2212	2217	2225	rVB	107318	143178	17.99%	1.148%
40	16.162	2236	2240	2244	rVB3	3727	4832	0.61%	0.039%
41	16.598	2308	2314	2321	rVV6	5676	10517	1.32%	0.084%
42	16.792	2340	2347	2350	rVV6	2408	4351	0.55%	0.035%
43	16.892	2359	2364	2368	rVV	6873	9541	1.20%	0.077%
44	17.115	2396	2402	2406	rBV3	1891	3151	0.40%	0.025%
45	17.192	2410	2415	2417	rVB5	2262	3016	0.38%	0.024%
46	17.392	2445	2449	2452	rBV4	4994	6449	0.81%	0.052%
47	17.662	2489	2495	2500	rVV	471172	679303	85.36%	5.448%
48	17.704	2500	2502	2507	rVV	83746	107581	13.52%	0.863%
49	17.762	2507	2512	2523	rVV	400380	577940	72.62%	4.635%
50	17.915	2535	2538	2543	rVB2	4684	6239	0.78%	0.050%
51	18.051	2557	2561	2565	rBV2	8059	10316	1.30%	0.083%
52	18.256	2592	2596	2599	rBV4	2844	4061	0.51%	0.033%
53	18.292	2599	2602	2608	rVB5	3820	5733	0.72%	0.046%
54	18.403	2618	2621	2624	rBV4	1917	3390	0.43%	0.027%
55	18.509	2634	2639	2645	rBV2	79779	110999	13.95%	0.890%
56	18.574	2646	2650	2658	rVB2	16830	28709	3.61%	0.230%
57	18.721	2668	2675	2679	rBV2	17869	34027	4.28%	0.273%
58	18.751	2679	2680	2686	rVB3	6659	6092	0.77%	0.049%
59	18.803	2686	2689	2694	rVB4	4238	5709	0.72%	0.046%
60	18.856	2694	2698	2703	rBV6	3599	6556	0.82%	0.053%
61	19.015	2720	2725	2728	rBV2	16877	23931	3.01%	0.192%
62	19.050	2728	2731	2735	rVB2	17761	22405	2.82%	0.180%
63	19.280	2762	2770	2778	rBV	56714	88090	11.07%	0.706%
64	19.362	2780	2784	2788	rVB	13563	18920	2.38%	0.152%
65	19.421	2790	2794	2801	rVV7	13785	25460	3.20%	0.204%
66	19.509	2807	2809	2812	rVV3	8366	8230	1.03%	0.066%
67	19.562	2814	2818	2824	rVB3	9342	15233	1.91%	0.122%
68	19.680	2833	2838	2844	rVB	325727	440855	55.39%	3.535%
69	19.756	2846	2851	2858	rBV4	35544	51645	6.49%	0.414%
70	19.815	2859	2861	2865	rBV5	3567	4158	0.52%	0.033%
71	19.898	2872	2875	2883	rVB5	17372	32896	4.13%	0.264%

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72	20.009	2888	2894	2897	rVV	486056	718427	90.27%	5.762%
73	20.039	2897	2899	2905	rVV	255551	312742	39.30%	2.508%
74	20.103	2906	2910	2915	rVV2	11956	16313	2.05%	0.131%
75	20.209	2925	2928	2933	rVB	17187	21802	2.74%	0.175%
76	20.339	2946	2950	2953	rBV2	60681	81111	10.19%	0.650%
77	20.480	2970	2974	2976	rBV4	17572	25794	3.24%	0.207%
78	20.515	2976	2980	2985	rVB	42570	70046	8.80%	0.562%
79	20.615	2994	2997	3001	rBV	33064	38774	4.87%	0.311%
80	20.680	3003	3008	3011	rVV2	16617	25660	3.22%	0.206%
81	20.815	3029	3031	3034	rBV2	10485	9350	1.17%	0.075%
82	21.250	3102	3105	3110	rBV2	24546	35291	4.43%	0.283%
83	21.415	3131	3133	3134	rBV	19056	15843	1.99%	0.127%
84	21.444	3134	3138	3143	rVB2	195000	234458	29.46%	1.880%
85	21.491	3144	3146	3150	rVV2	20221	22624	2.84%	0.181%
86	21.533	3151	3153	3160	rVB2	24794	34335	4.31%	0.275%
87	21.650	3169	3173	3177	rBV	81210	97114	12.20%	0.779%
88	21.768	3186	3193	3197	rBV2	446995	777529	97.70%	6.235%
89	21.803	3197	3199	3209	rVB	146009	184187	23.14%	1.477%
90	21.886	3211	3213	3217	rVB3	30924	34101	4.28%	0.273%
91	21.933	3218	3221	3224	rBV2	25494	32078	4.03%	0.257%
92	22.280	3277	3280	3285	rBV5	39882	51401	6.46%	0.412%
93	22.350	3289	3292	3299	rVV4	49260	87917	11.05%	0.705%
94	22.850	3373	3377	3380	rBV	45210	59666	7.50%	0.478%
95	23.403	3467	3471	3474	rBV	51072	64852	8.15%	0.520%
96	23.450	3474	3479	3485	rBV2	150101	297708	37.41%	2.388%
97	24.033	3573	3578	3584	rBV2	182838	335768	42.19%	2.693%
98	24.191	3599	3605	3611	rBV2	187845	358856	45.09%	2.878%
99	24.438	3644	3647	3658	rVB2	41481	83167	10.45%	0.667%
100	24.738	3694	3698	3706	rVB2	71786	146272	18.38%	1.173%

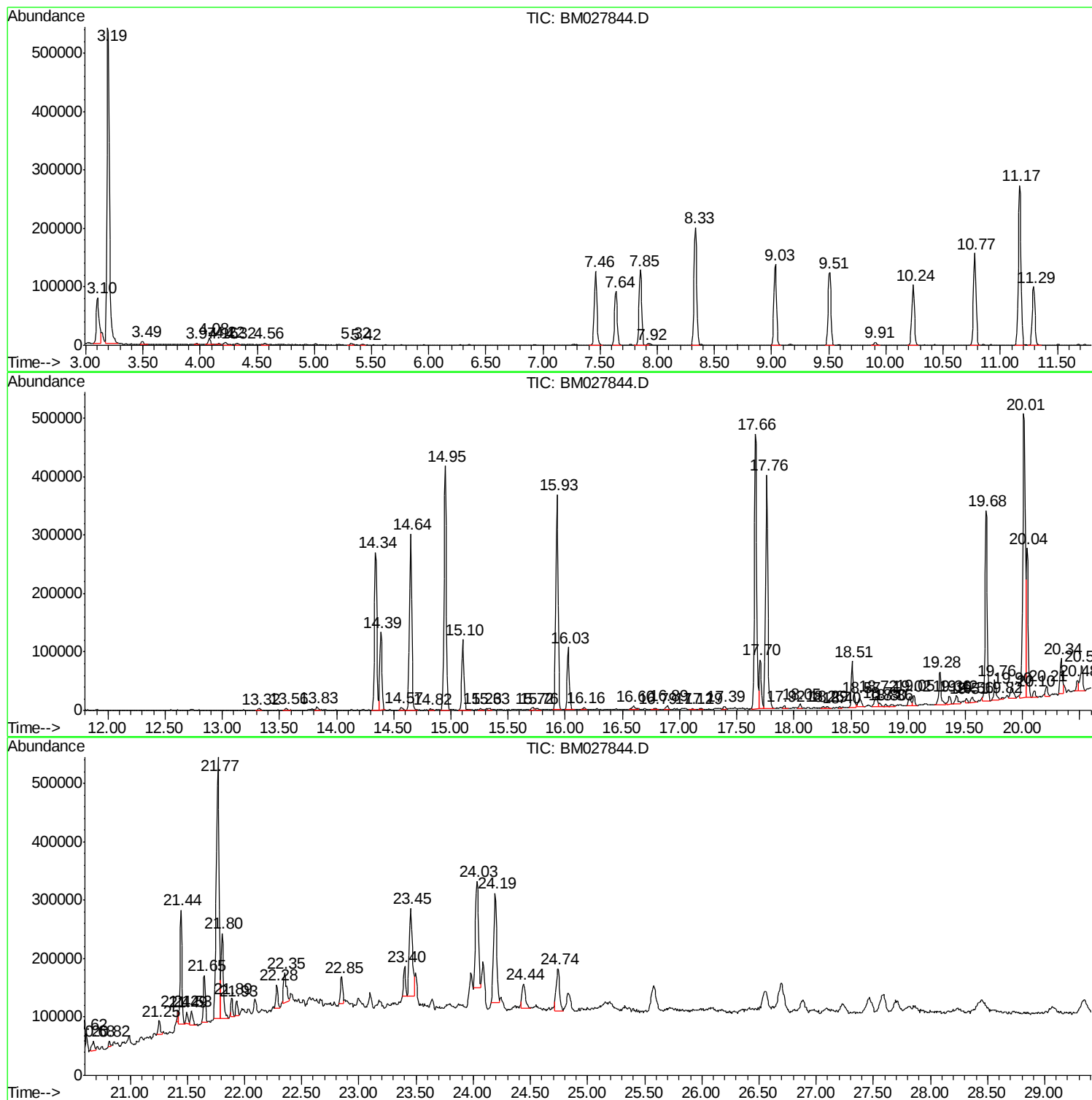
Sum of corrected areas: 12469441

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111220MA.M
Quant Title : SVOA CALIBRATION

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



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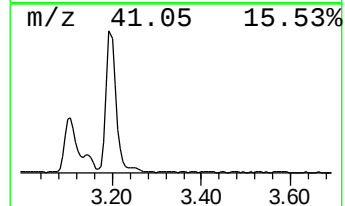
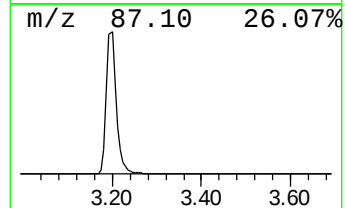
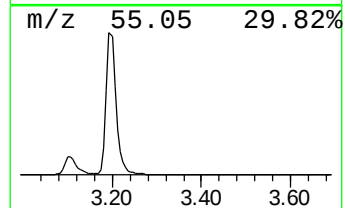
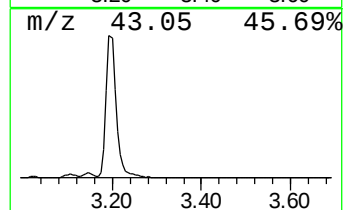
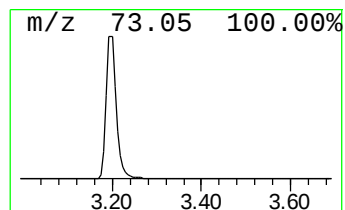
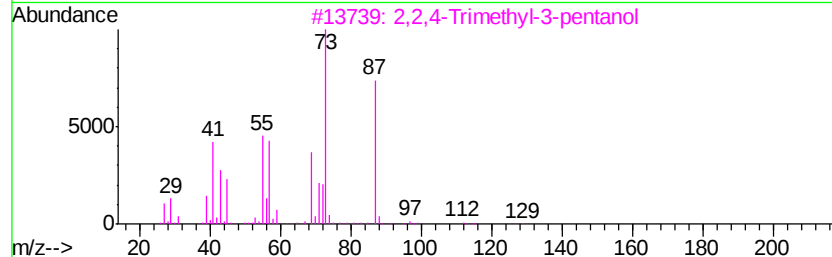
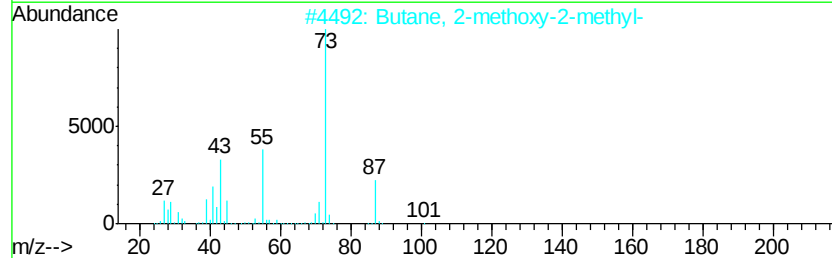
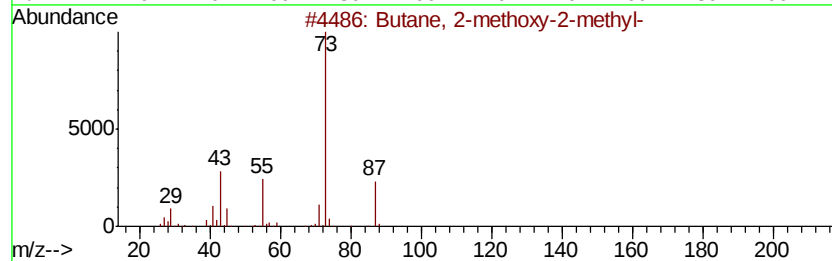
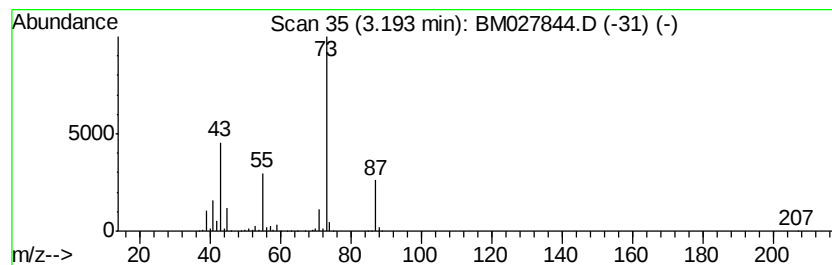
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111220MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.19	49.76 ng/ul	795839	1,4-Dichlorobenzene-d4	8.33

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	64	
3	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40	
4	Silane, tetramethyl-	88	C4H12Si	000075-76-3	17	
5	N-Ethylformamide	73	C3H7NO	000627-45-2	9	



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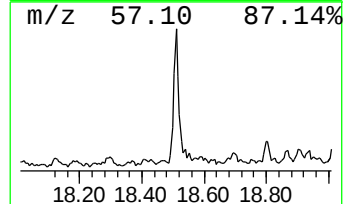
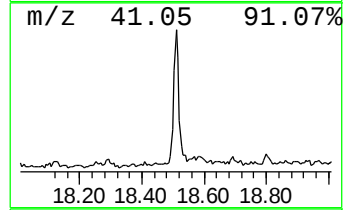
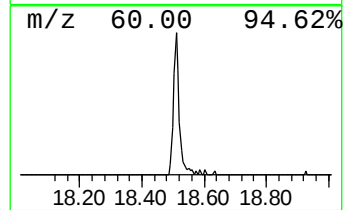
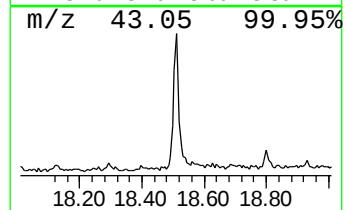
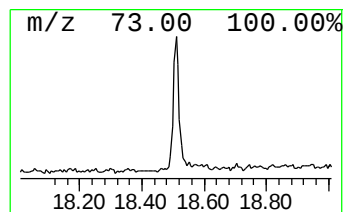
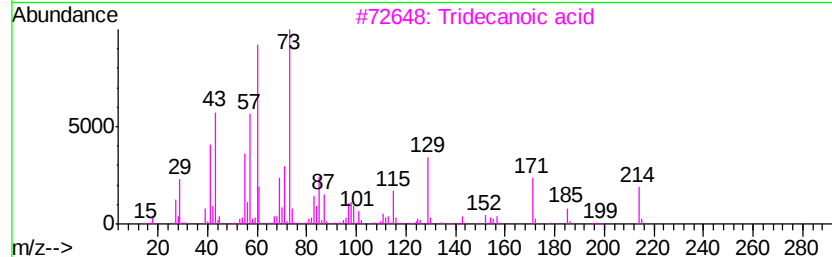
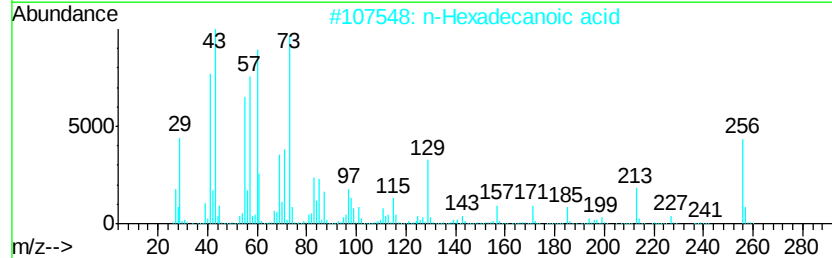
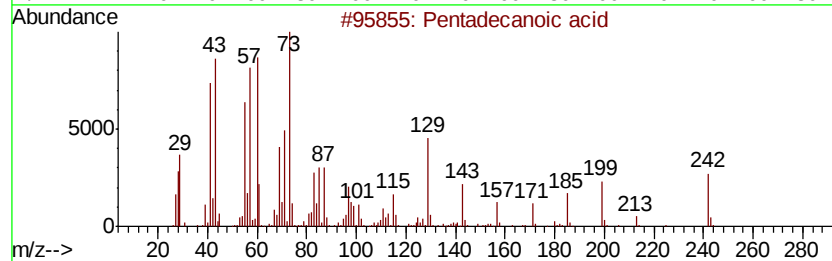
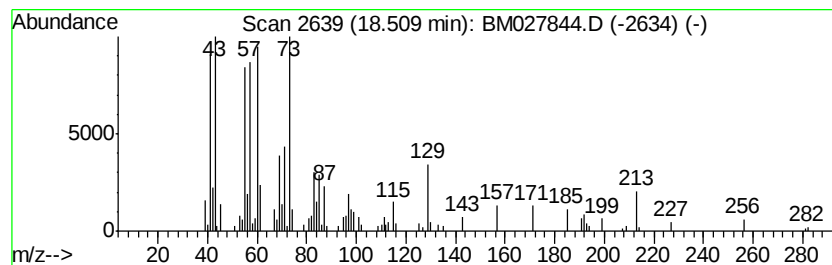
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111220MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Pentadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.51	3.27 ng/ul	110999	Phenanthrene-d10	17.66

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecanoic acid	242	C15H30O2	001002-84-2	87	
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	87	
3	Tridecanoic acid	214	C13H26O2	000638-53-9	76	
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	74	
5	Pentadecanoic acid	242	C15H30O2	001002-84-2	74	



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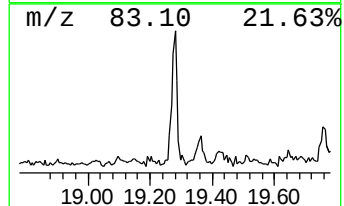
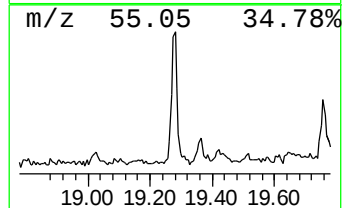
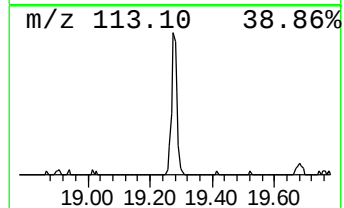
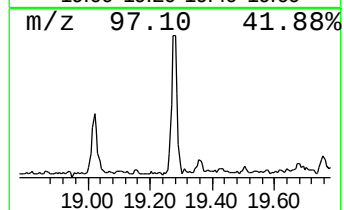
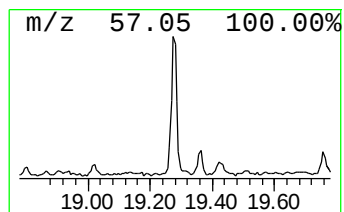
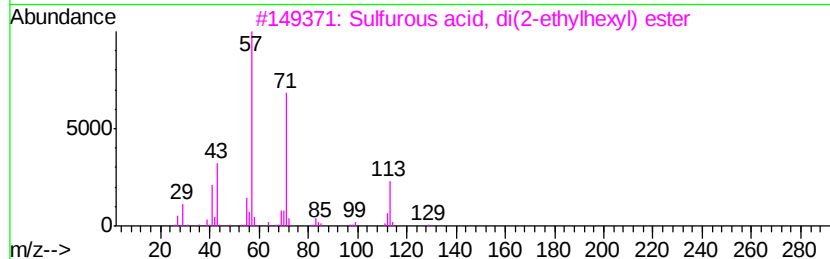
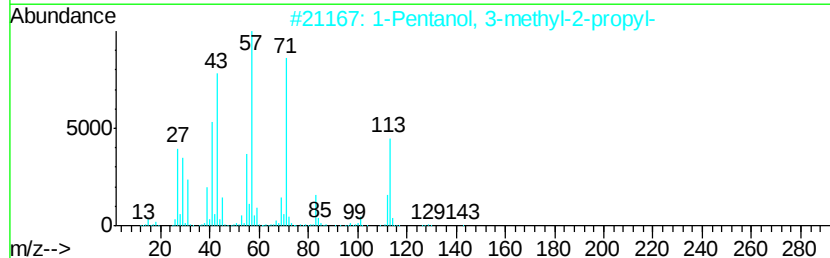
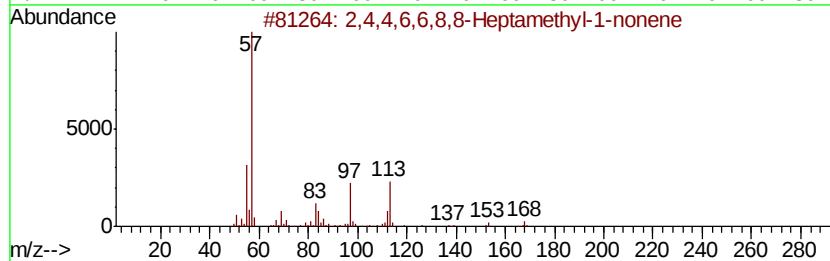
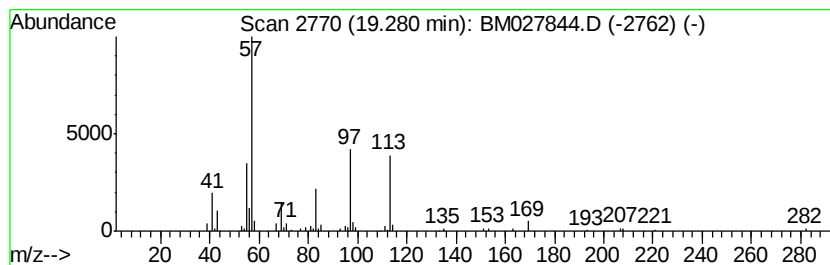
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Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.28	2.59 ng/ul	88090	Phenanthrene-d10	17.66

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	64
2		1-Pentanol, 3-methyl-2-propyl-	144	C9H20O	054004-40-9	43
3		Sulfurous acid, di(2-ethylhexyl)...	306	C16H34O3S	1000309-19-1	35
4		2,4,4-Trimethyl-1-pentanol, hept...	326	C12H17F7O2	1000365-01-4	35
5		4-Methylnonanoic acid	172	C10H20O2	045019-28-1	28



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM111920\
Data File : BM027844.D
Acq On : 19 Nov 2020 14:59
Operator : JU/CG
Sample : L4710-16
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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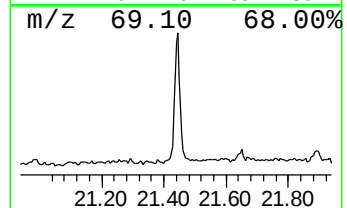
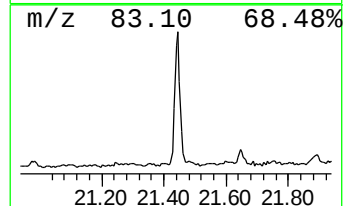
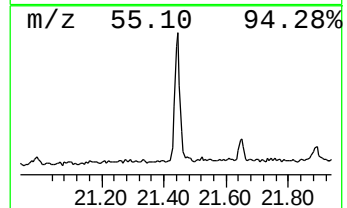
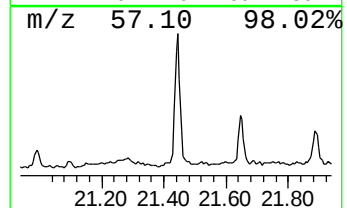
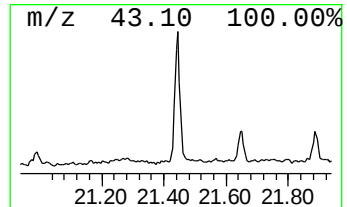
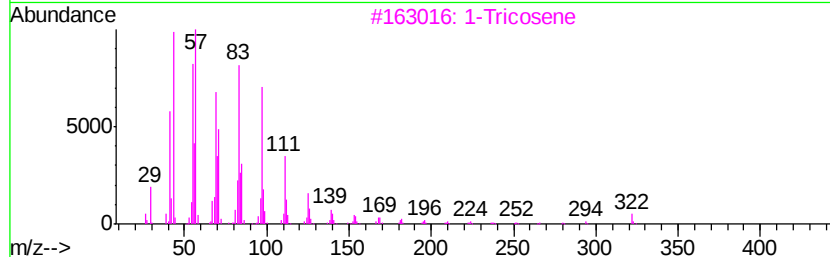
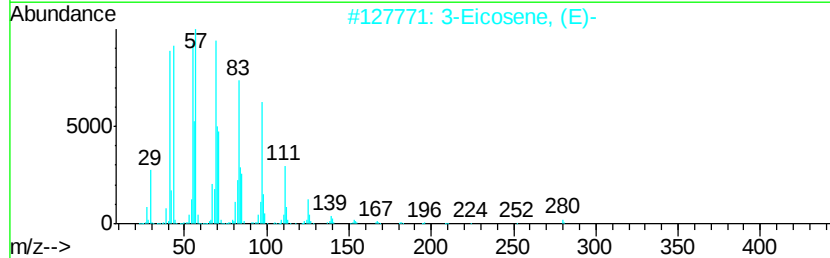
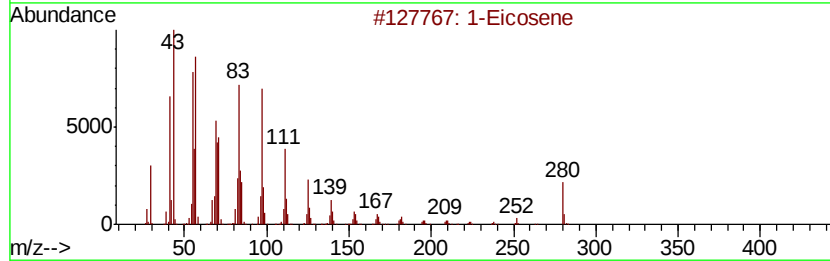
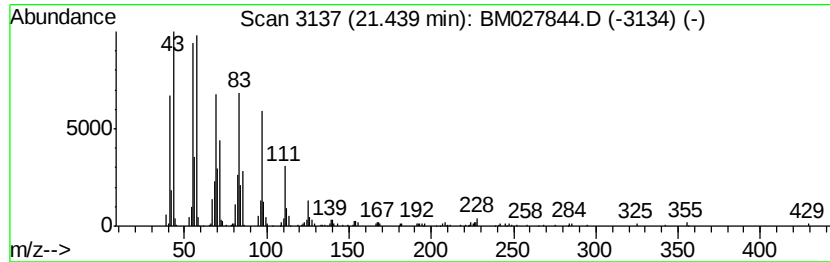
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.44	6.03 ng/ul	234458	Chrysene-d12	21.77

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Eicosene	280	C20H40	003452-07-1	94
2			3-Eicosene, (E)-	280	C20H40	074685-33-9	91
3			1-Tricosene	322	C23H46	018835-32-0	90
4			Z-8-Hexadecene	224	C16H32	1000130-87-5	90
5			Nonadecyl trifluoroacetate	380	C21H39F3O2	1000351-76-3	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM111920\
Data File : BM027844.D
Acq On : 19 Nov 2020 14:59
Operator : JU/CG
Sample : L4710-16
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0B45

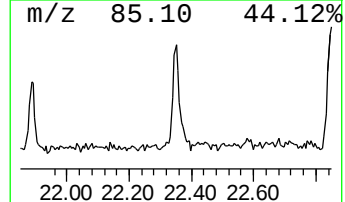
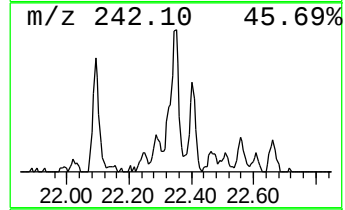
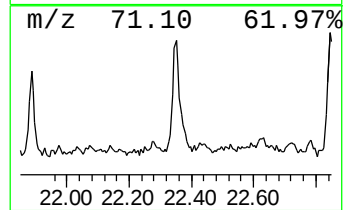
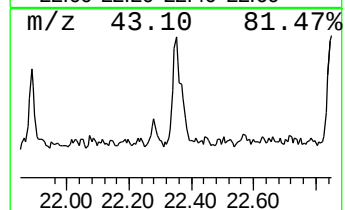
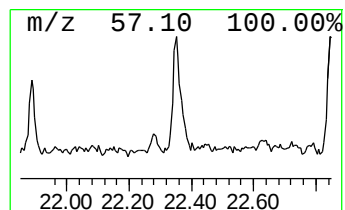
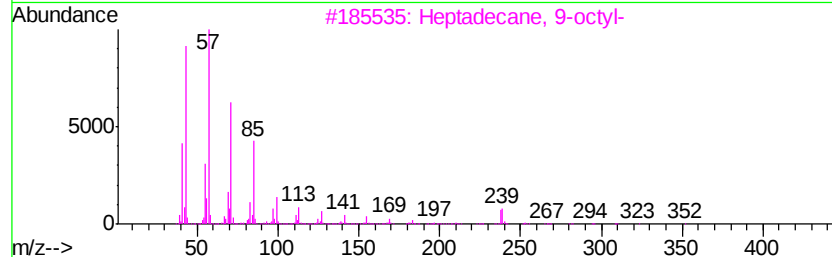
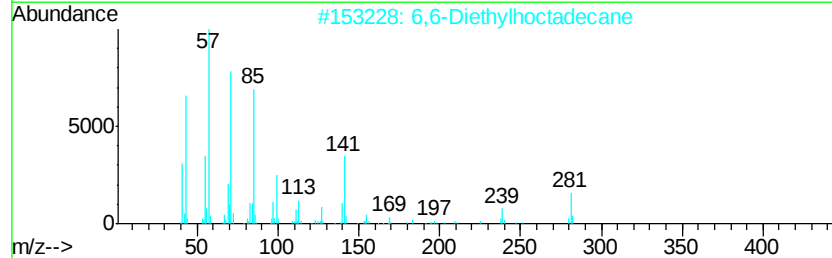
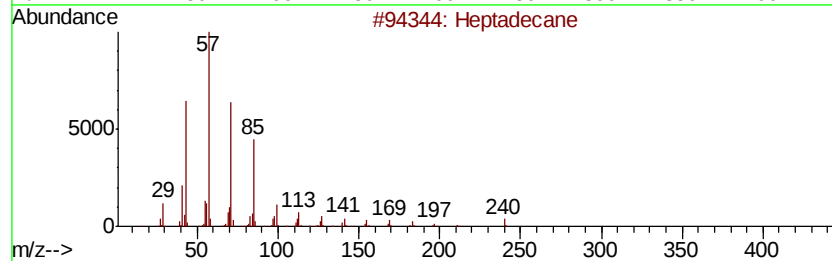
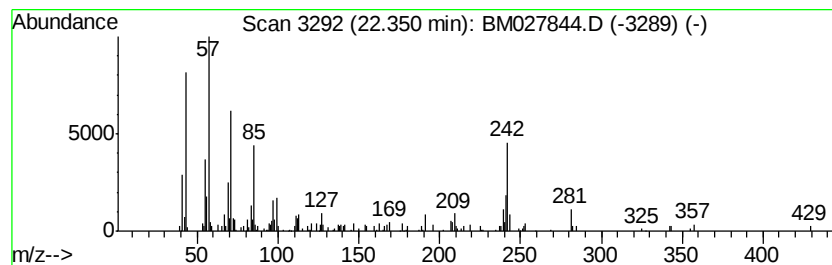
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.35	2.26 ng/ul	87917	Chrysene-d12	21.77

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane	240	C17H36	000629-78-7	50
2		6,6-Diethyloctadecane	310	C22H46	1000360-41-8	46
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	43
4		Tridecane, 6-propyl-	226	C16H34	055045-10-8	41
5		Hexadecane	226	C16H34	000544-76-3	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM111920\
Data File : BM027844.D
Acq On : 19 Nov 2020 14:59
Operator : JU/CG
Sample : L4710-16
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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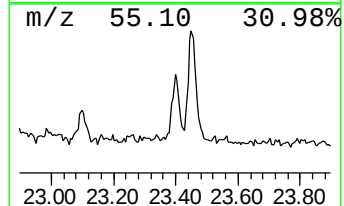
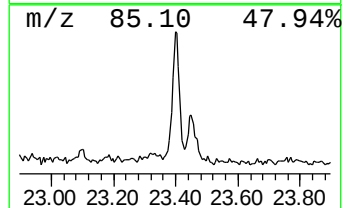
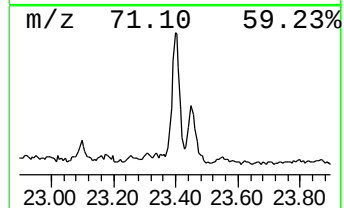
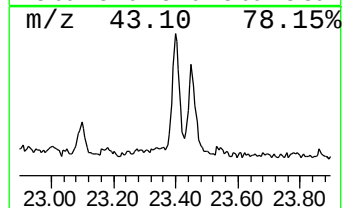
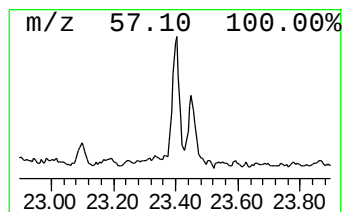
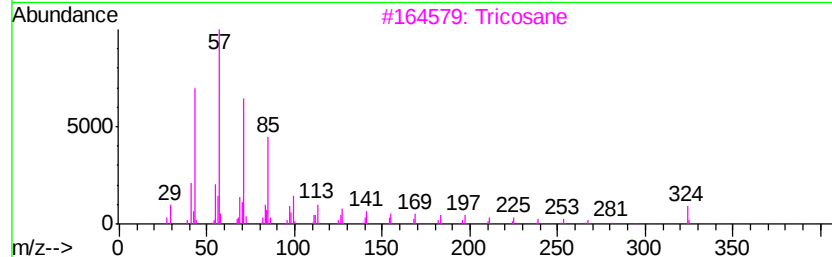
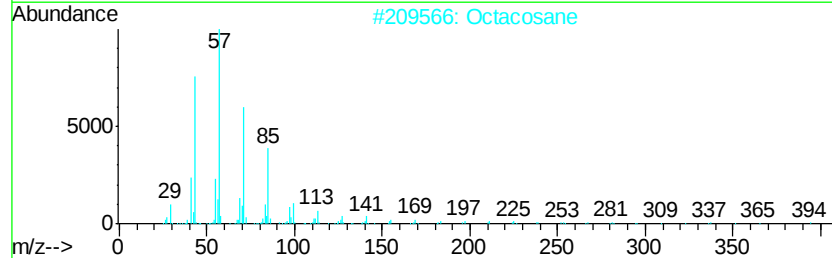
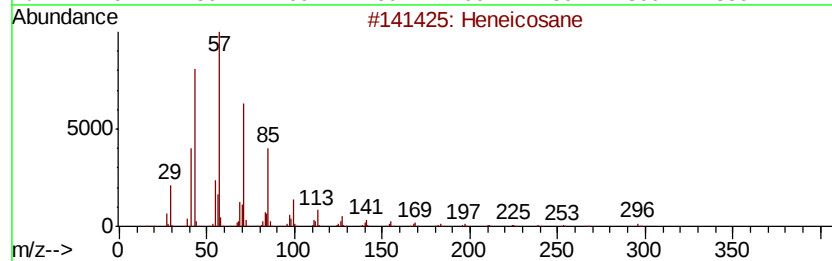
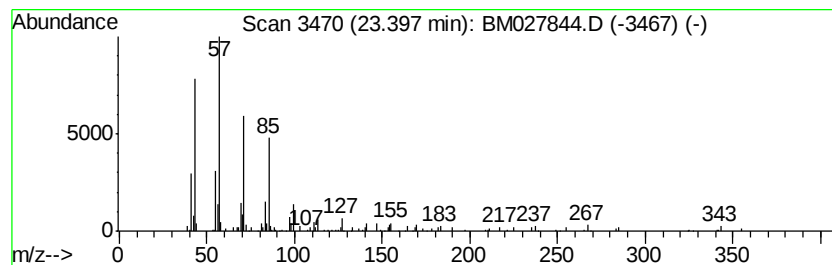
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.40	3.61 ng/ul	64852	Perylene-d12	24.19

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	96
2			Octacosane	394	C28H58	000630-02-4	90
3			Tricosane	324	C23H48	000638-67-5	90
4			Tricosane	324	C23H48	000638-67-5	90
5			Hexacosane	366	C26H54	000630-01-3	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM111920\
Data File : BM027844.D
Acq On : 19 Nov 2020 14:59
Operator : JU/CG
Sample : L4710-16
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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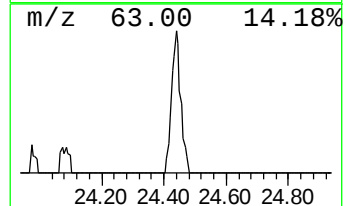
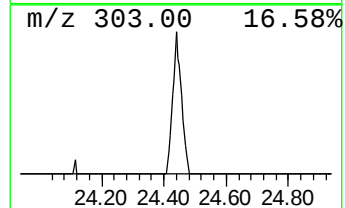
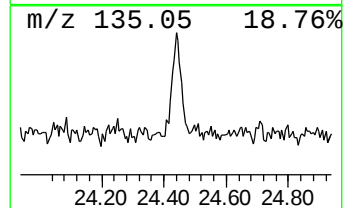
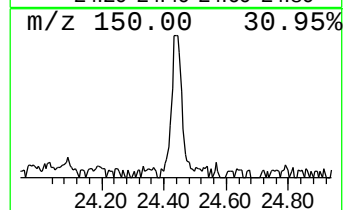
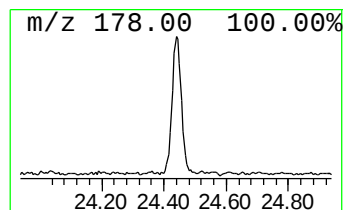
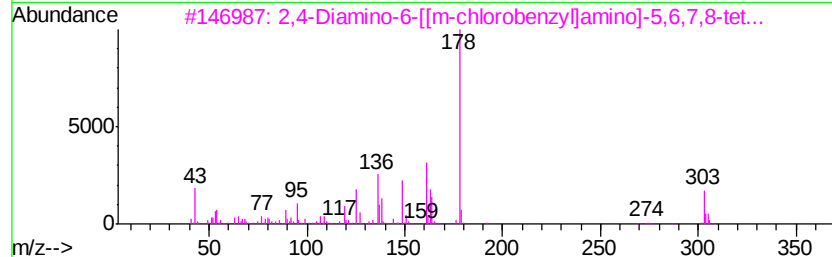
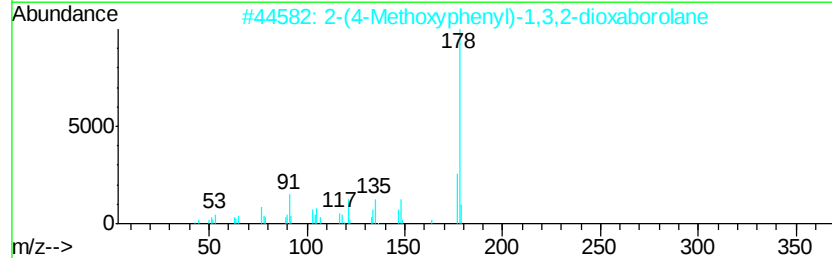
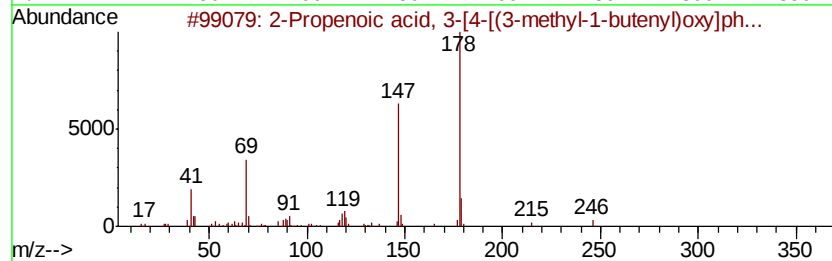
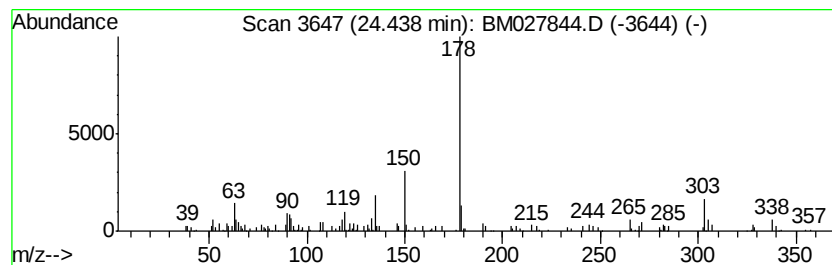
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 2-Propenoic acid, 3-[4-[(3-... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.44	4.64 ng/ul	83167	Perylene-d12	24.19

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Propenoic acid, 3-[4-[(3-methy...	246	C15H18O3	055256-10-5	53
2		2-(4-Methoxyphenyl)-1,3,2-dioxab...	178	C9H11BO3	069519-11-5	52
3		2,4-Diamino-6-[m-chlorobenzyl]a...	303	C15H18ClN5	1000253-34-5	50
4		Phenol, 2-ethoxy-4-(2-propenyl)-	178	C11H14O2	001755-54-0	49
5		2,4(1H,3H)-Pteridinedione, 7-met...	178	C7H6N4O2	013401-38-2	47



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM111920\
Data File : BM027844.D
Acq On : 19 Nov 2020 14:59
Operator : JU/CG
Sample : L4710-16
Misc :
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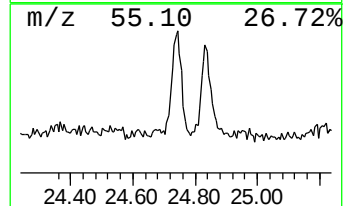
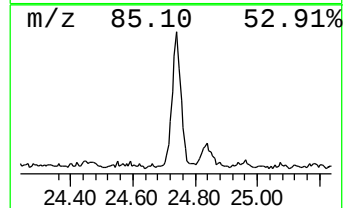
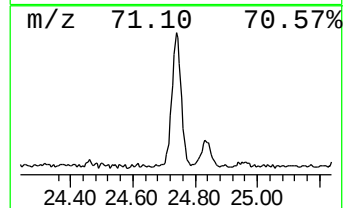
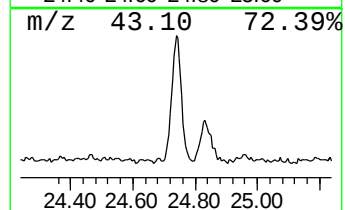
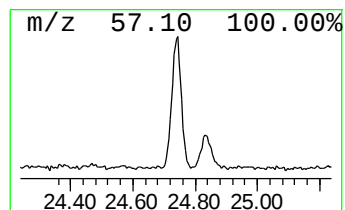
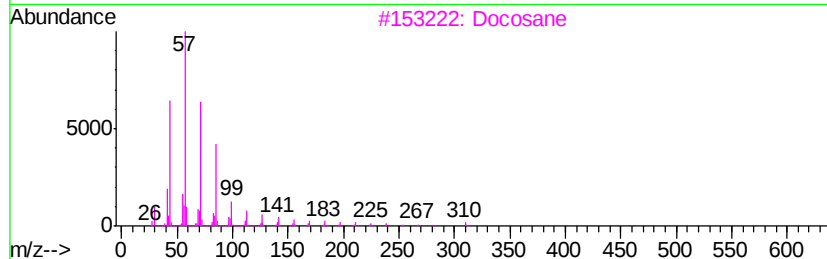
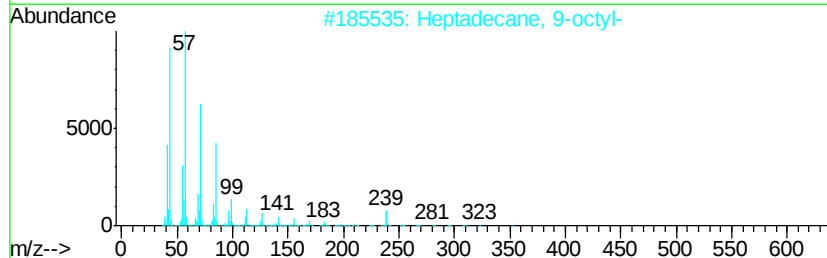
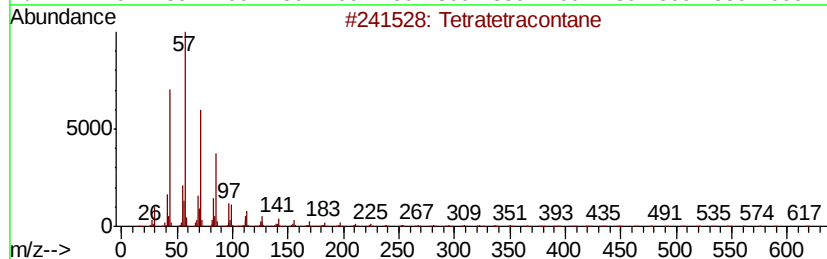
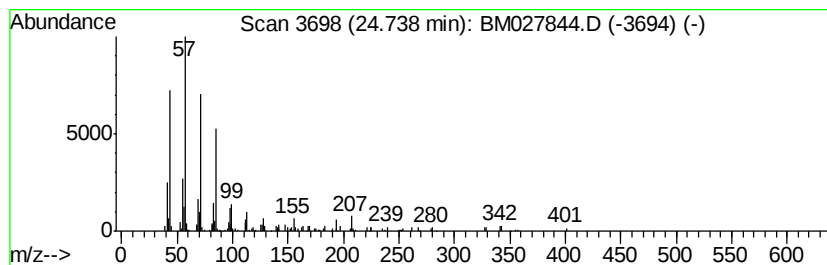
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.74	8.15 ng/ul	146272	Perylene-d12	24.19

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetratetracontane	619	C44H90	007098-22-8	90
2			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87
3			Docosane	310	C22H46	000629-97-0	87
4			Triacontane	422	C30H62	000638-68-6	87
5			Tridecane, 7-hexyl-	268	C19H40	007225-66-3	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM111920\
Data File : BM027844.D
Acq On : 19 Nov 2020 14:59
Operator : JU/CG
Sample : L4710-16
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0B45

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111220MA.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...	3.19	49.8	ng/ul	795839	1	8.33	319858	20.0
Pentadecanoic acid	18.51	3.3	ng/ul	110999	4	17.66	679303	20.0
(DEL) Alkane: Str...	19.28	2.6	ng/ul	88090	4	17.66	679303	20.0
(DEL) Alkane: Str...	21.44	6.0	ng/ul	234458	5	21.77	777529	20.0
(DEL) Alkane: Str...	22.35	2.3	ng/ul	87917	5	21.77	777529	20.0
(DEL) Alkane: Str...	23.40	3.6	ng/ul	64852	6	24.19	358856	20.0
2-Propenoic acid,...	24.44	4.6	ng/ul	83167	6	24.19	358856	20.0
(DEL) Alkane: Str...	24.74	8.2	ng/ul	146272	6	24.19	358856	20.0