

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM072820\
 Data File : BM026895.D
 Acq On : 28 Jul 2020 17:44
 Operator : JU/CG
 Sample : L3388-06
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AD2

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-BM072720MA.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	2.772	4	6	11	rVB2	12630	15249	0.67%	0.083%
2	2.907	24	29	37	rBV	312988	462019	20.40%	2.520%
3	2.978	37	41	56	rVB	1954402	2264778	100.00%	12.354%
4	3.172	72	74	79	rBV5	2177	3975	0.18%	0.022%
5	3.219	79	82	87	rVB	23488	29842	1.32%	0.163%
6	3.860	185	191	195	rBV2	15574	24092	1.06%	0.131%
7	3.943	202	205	208	rBV3	5309	6274	0.28%	0.034%
8	4.160	236	242	245	rBV4	4270	7478	0.33%	0.041%
9	4.290	261	264	266	rBV4	2410	3225	0.14%	0.018%
10	4.390	279	281	286	rVB5	1952	2678	0.12%	0.015%
11	4.554	306	309	312	rBV3	2847	3283	0.14%	0.018%
12	4.825	350	355	362	rBV	15873	23100	1.02%	0.126%
13	5.648	492	495	497	rVB3	1884	2333	0.10%	0.013%
14	6.842	688	698	707	rBV	216417	339428	14.99%	1.852%
15	7.007	720	726	736	rVB	153857	240974	10.64%	1.315%
16	7.207	754	760	766	rBV	203965	313131	13.83%	1.708%
17	7.331	776	781	785	rVB4	2662	4482	0.20%	0.024%
18	7.672	833	839	846	rBV	305693	464481	20.51%	2.534%
19	7.754	848	853	855	rVB6	1506	2360	0.10%	0.013%
20	8.372	947	958	970	rBV	262574	410048	18.11%	2.237%
21	8.484	973	977	984	rVB2	10234	16501	0.73%	0.090%
22	8.813	1025	1033	1042	rVV	199400	319365	14.10%	1.742%
23	8.995	1061	1064	1067	rBV3	1844	2678	0.12%	0.015%
24	9.531	1149	1155	1166	rBV	131300	218088	9.63%	1.190%
25	9.883	1211	1215	1221	rBV5	2525	3888	0.17%	0.021%
26	10.066	1236	1246	1255	rBV	249404	410112	18.11%	2.237%
27	10.448	1304	1311	1318	rVB	365673	605982	26.76%	3.306%
28	10.578	1326	1333	1341	rBV	239602	408739	18.05%	2.230%
29	12.936	1729	1734	1737	rBV3	1757	3067	0.14%	0.017%
30	13.219	1775	1782	1787	rBV6	2032	4551	0.20%	0.025%
31	13.601	1839	1847	1850	rBV5	3182	5209	0.23%	0.028%
32	13.648	1850	1855	1860	rVV5	9349	15449	0.68%	0.084%
33	13.719	1861	1867	1872	rVV	458752	604283	26.68%	3.296%
34	13.766	1872	1875	1881	rVB	62436	78550	3.47%	0.428%

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35	13.995	1908	1914	1922	rBV	500562	710964	31.39%	3.878%
36	14.177	1941	1945	1948	rVB2	1538	2282	0.10%	0.012%
37	14.307	1960	1967	1976	rBV	558179	785408	34.68%	4.284%
38	14.389	1978	1981	1985	rBV3	2430	2930	0.13%	0.016%
39	14.489	1988	1998	2008	rBV	189611	273396	12.07%	1.491%
40	14.583	2013	2014	2018	rVB3	3097	2312	0.10%	0.013%
41	14.742	2037	2041	2047	rBV4	6790	9676	0.43%	0.053%
42	14.871	2059	2063	2065	rBV2	2067	2652	0.12%	0.014%
43	15.301	2130	2136	2143	rVB	614245	845263	37.32%	4.611%
44	15.407	2149	2154	2164	rBV2	114781	165762	7.32%	0.904%
45	15.542	2171	2177	2182	rBV5	6150	9156	0.40%	0.050%
46	16.324	2307	2310	2313	rVB3	2614	2413	0.11%	0.013%
47	16.748	2378	2382	2384	rVB3	1831	2274	0.10%	0.012%
48	17.048	2427	2433	2443	rVB	661050	889802	39.29%	4.854%
49	17.148	2443	2450	2458	rBV	645041	893131	39.44%	4.872%
50	17.459	2498	2503	2509	rBV6	2583	6039	0.27%	0.033%
51	17.636	2529	2533	2537	rVB5	1892	2619	0.12%	0.014%
52	17.754	2550	2553	2556	rBV2	2838	2969	0.13%	0.016%
53	17.854	2566	2570	2574	rVB4	1789	2301	0.10%	0.013%
54	17.971	2586	2590	2595	rBV2	35667	46605	2.06%	0.254%
55	18.283	2639	2643	2648	rVB6	4091	5564	0.25%	0.030%
56	18.936	2747	2754	2758	rBV8	7006	11604	0.51%	0.063%
57	19.012	2762	2767	2769	rBV5	2886	4310	0.19%	0.024%
58	19.106	2780	2783	2787	rBV3	4933	5722	0.25%	0.031%
59	19.230	2801	2804	2805	rVB3	3383	2946	0.13%	0.016%
60	19.259	2805	2809	2817	rBV5	22323	35347	1.56%	0.193%
61	19.330	2818	2821	2829	rVB5	5621	12522	0.55%	0.068%
62	19.442	2834	2840	2849	rVV	766292	1023303	45.18%	5.582%
63	19.530	2853	2855	2859	rVB5	3584	4752	0.21%	0.026%
64	19.577	2859	2863	2864	rBV3	2436	3290	0.15%	0.018%
65	19.730	2886	2889	2891	rVB4	3261	2776	0.12%	0.015%
66	19.948	2923	2926	2928	rBV4	2919	4340	0.19%	0.024%
67	20.000	2933	2935	2938	rVB4	5642	4440	0.20%	0.024%
68	20.036	2938	2941	2943	rBV3	4360	5716	0.25%	0.031%
69	20.177	2962	2965	2969	rBV3	14164	15712	0.69%	0.086%
70	20.347	2989	2994	2998	rBV	123125	161007	7.11%	0.878%
71	20.389	2998	3001	3010	rVB8	36160	73715	3.25%	0.402%

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72	20.847	3076	3079	3082	rBV5	7183	9166	0.40%	0.050%
73	20.989	3100	3103	3110	rVV4	46610	62350	2.75%	0.340%
74	21.065	3112	3116	3122	rVB8	24958	38664	1.71%	0.211%
75	21.242	3141	3146	3151	rBV	817820	997721	44.05%	5.443%
76	21.794	3238	3240	3244	rVB5	20256	21747	0.96%	0.119%
77	21.877	3251	3254	3260	rBV4	38185	49957	2.21%	0.273%
78	22.847	3414	3419	3429	rBV	144718	261609	11.55%	1.427%
79	23.318	3493	3499	3506	rVB2	534833	958962	42.34%	5.231%
80	23.459	3512	3523	3534	rVB	542632	1065627	47.05%	5.813%
81	24.065	3620	3626	3633	rBV	209392	391443	17.28%	2.135%
82	24.135	3634	3638	3647	rVB9	40946	75299	3.32%	0.411%
83	25.688	3896	3902	3914	rVB	83975	218459	9.65%	1.192%
84	26.277	3997	4002	4012	rVB3	56897	141757	6.26%	0.773%
85	26.529	4038	4045	4059	rVB3	184742	533812	23.57%	2.912%
86	28.082	4304	4309	4321	rVB3	45447	150704	6.65%	0.822%

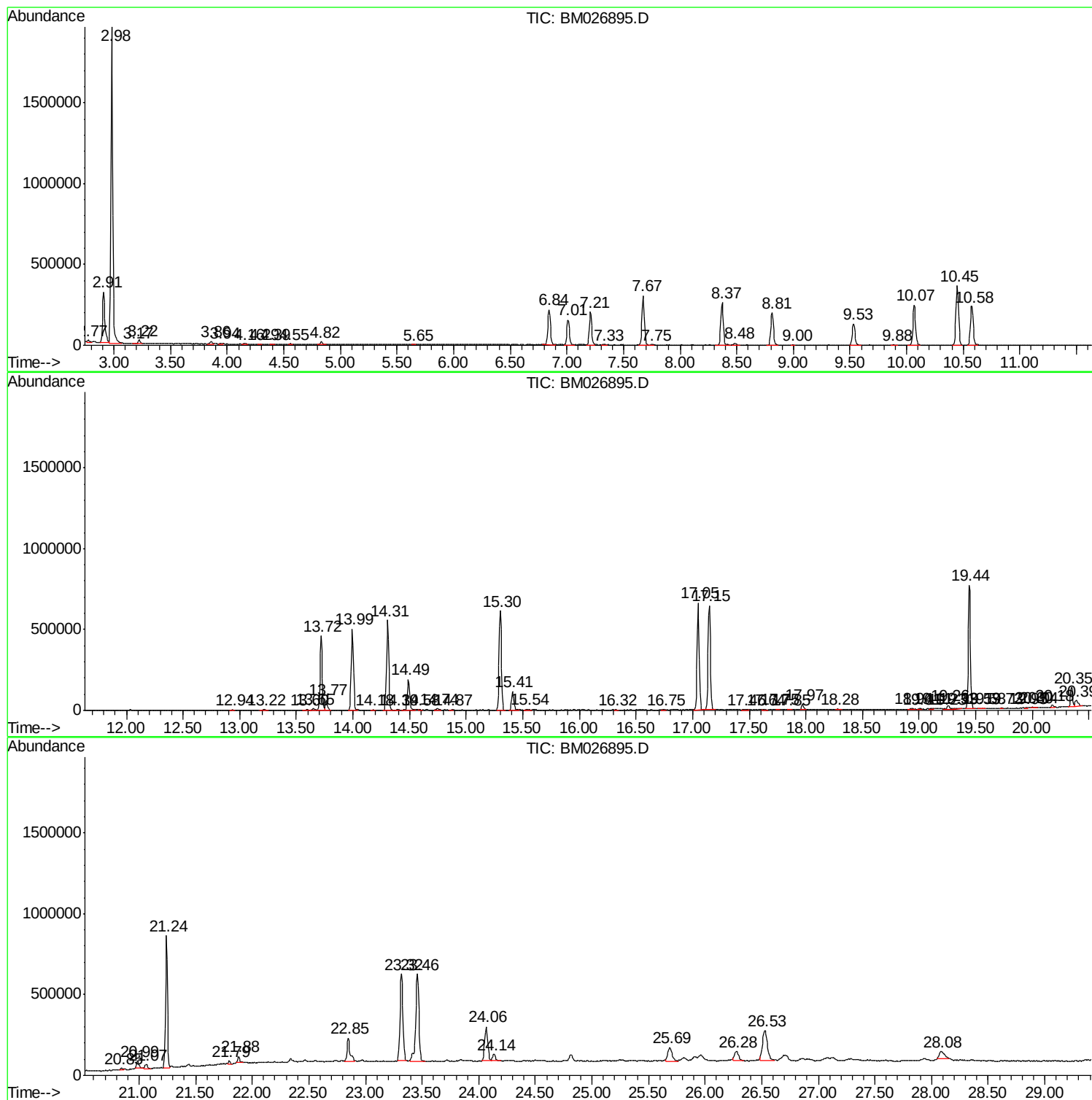
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM072720MA.M
Quant Title : SVOA CALIBRATION

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



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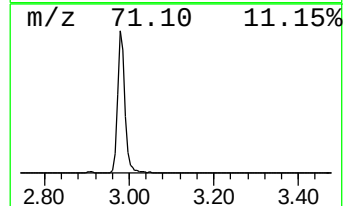
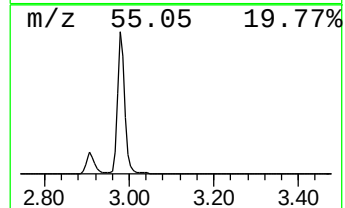
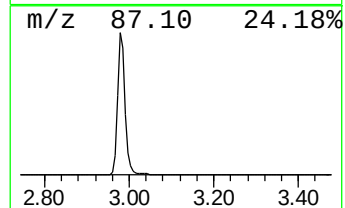
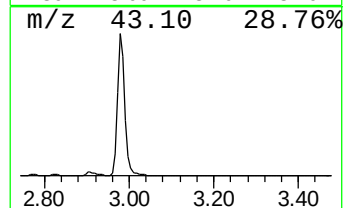
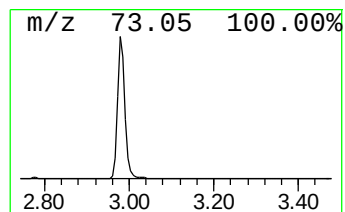
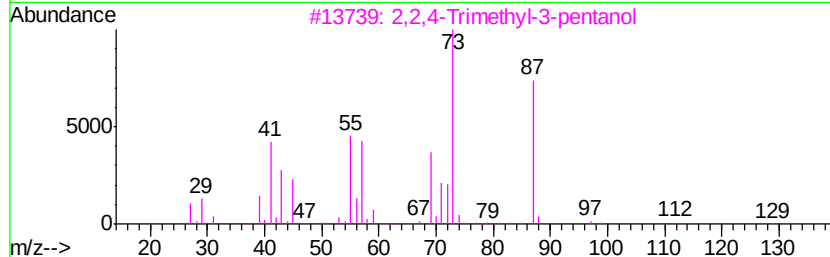
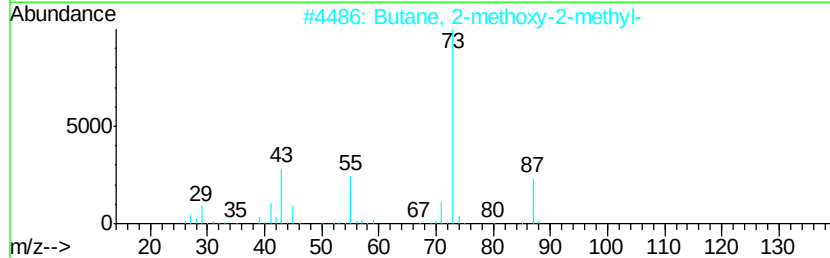
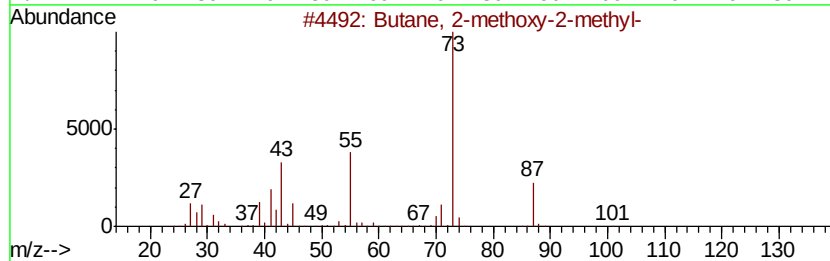
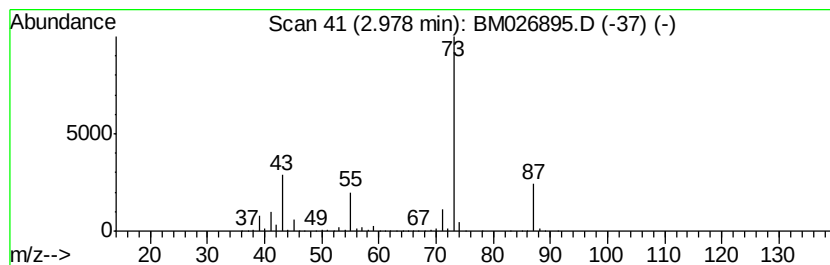
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM072720MA.M
Quant Title : SVOA CALIBRATION

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.
2.98	97.52 ng/ul	2264780	1,4-Dichlorobenzene-d4	7.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	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	39
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
5		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9



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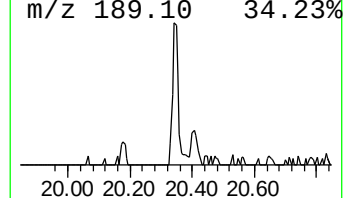
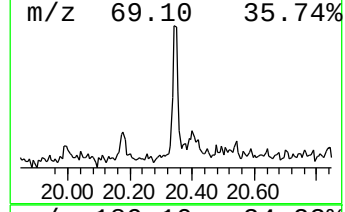
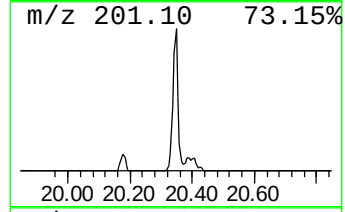
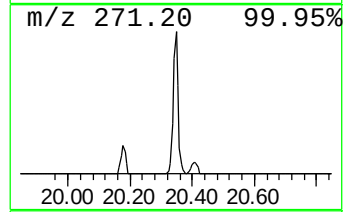
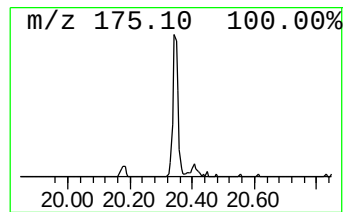
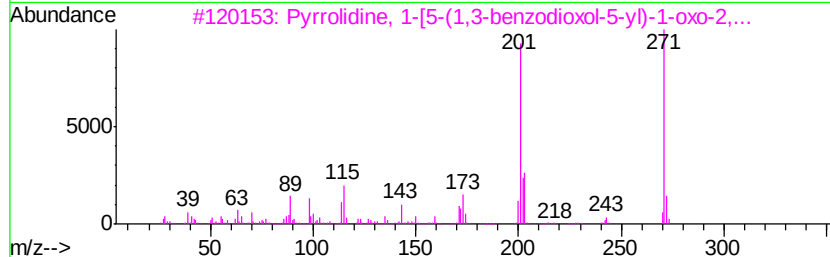
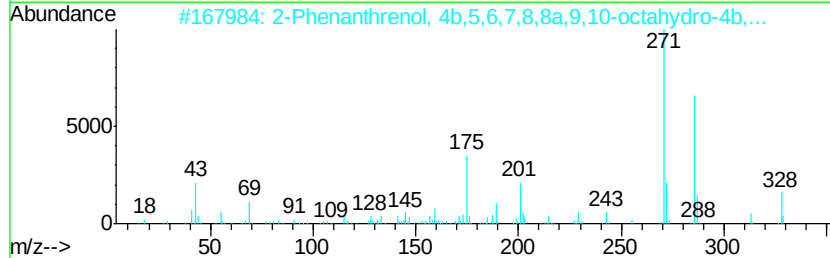
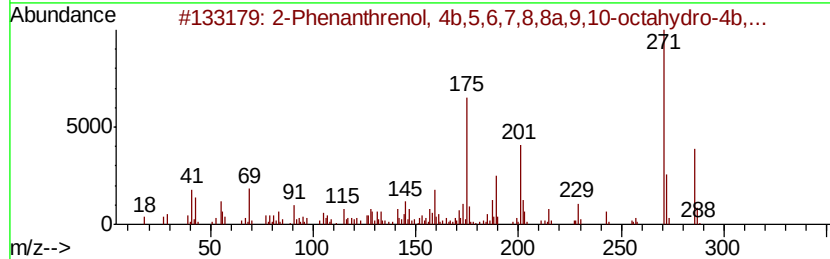
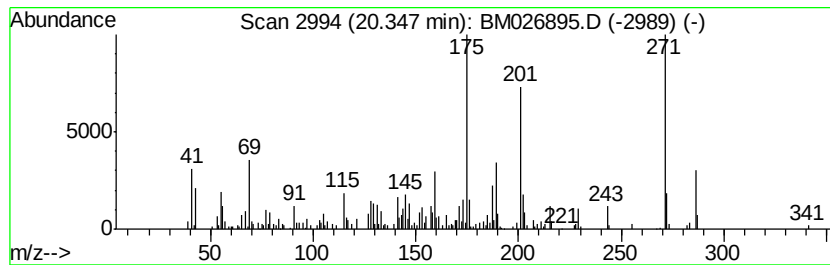
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 2-Phenanthrenol, 4b,5,6,7,8... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.35	3.23 ng/ul	161007	Chrysene-d12	21.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Phenanthrenol, 4b,5,6,7,8,8a,9...	286	C20H30O	000511-15-9	91
2		2-Phenanthrenol, 4b,5,6,7,8,8a,9...	328	C22H32O2	015340-82-6	87
3		Pyrrolidine, 1-[5-(1,3-benzodiox...	271	C16H17NO3	025924-78-1	38
4		Dihydronormorphinone	271	C16H17NO3	014696-23-2	15
5		4-Methyl-2,6-dihydroxyquinoline	175	C10H9NO2	034982-01-9	15



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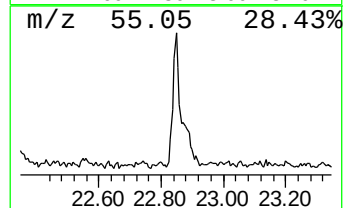
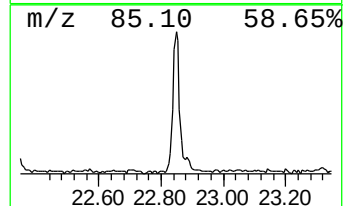
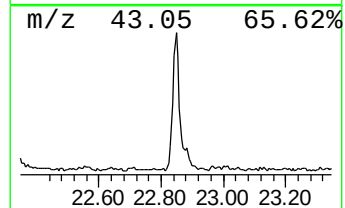
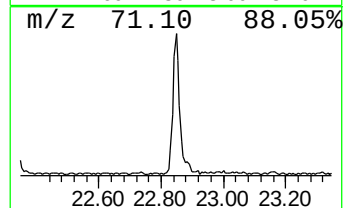
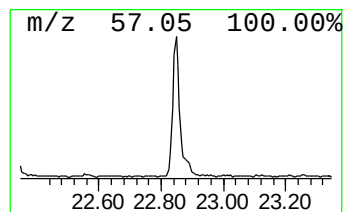
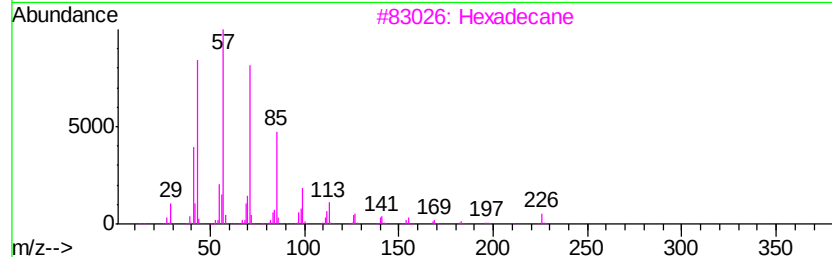
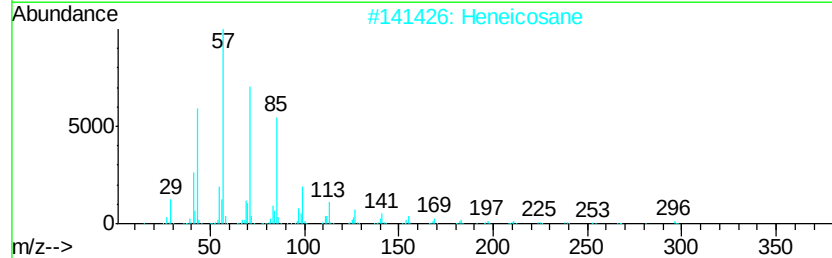
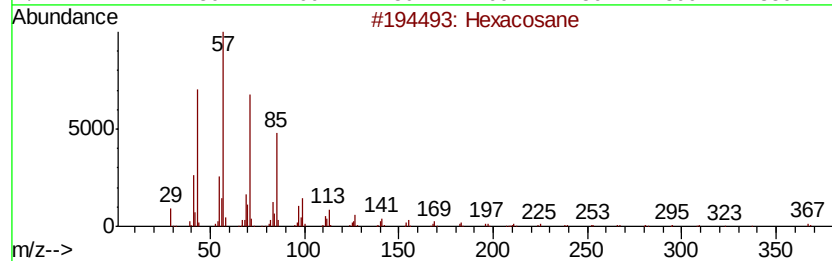
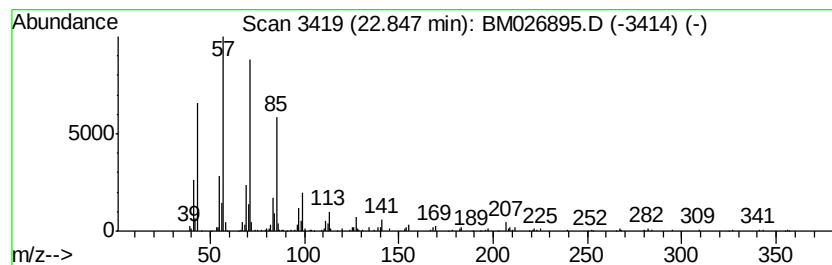
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.85	4.91 ng/ul	261609	Perylene-d12	23.46

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	91
2		Heneicosane	296	C21H44	000629-94-7	91
3		Hexadecane	226	C16H34	000544-76-3	91
4		Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91
5		Heneicosane	296	C21H44	000629-94-7	91



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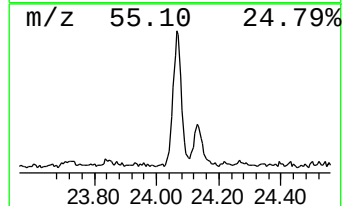
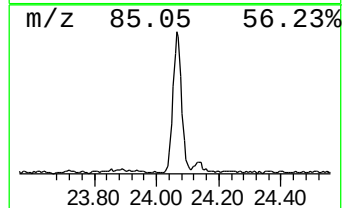
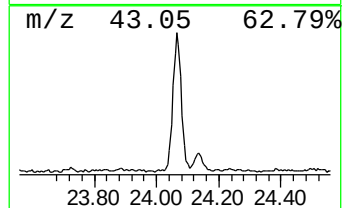
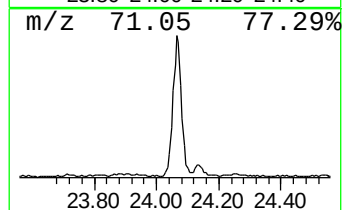
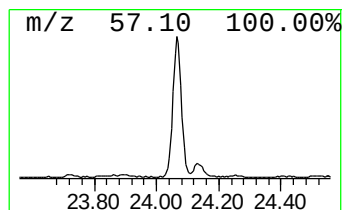
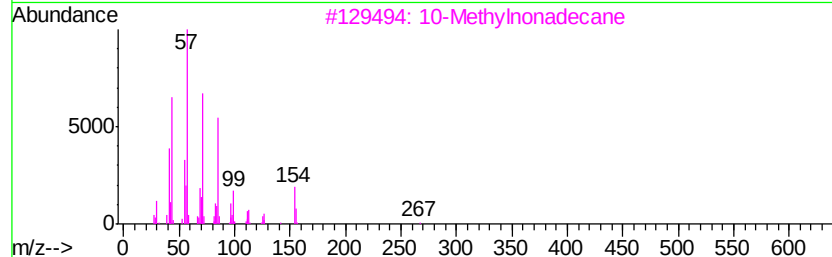
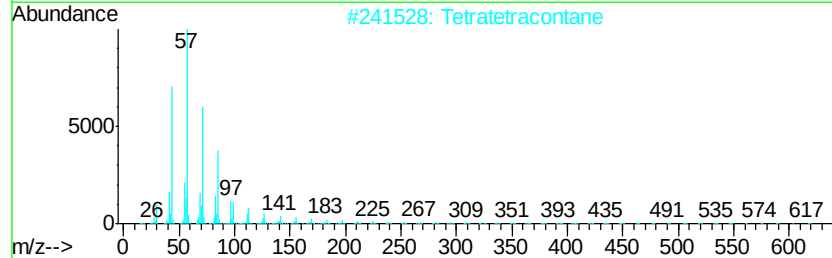
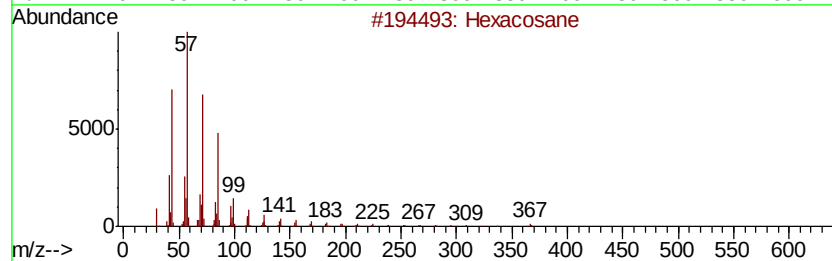
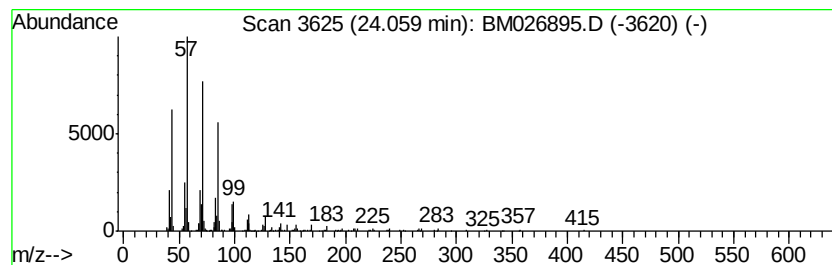
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.06	7.35 ng/ul	391443	Perylene-d12	23.46

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	87
2		Tetratetracontane	619	C44H90	007098-22-8	87
3		10-Methylnonadecane	282	C20H42	056862-62-5	86
4		Heneicosane	296	C21H44	000629-94-7	83
5		Docosane	310	C22H46	000629-97-0	81



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM072820\
Data File : BM026895.D
Acq On : 28 Jul 2020 17:44
Operator : JU/CG
Sample : L3388-06
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0AD2

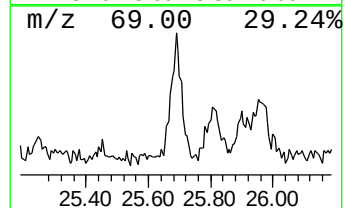
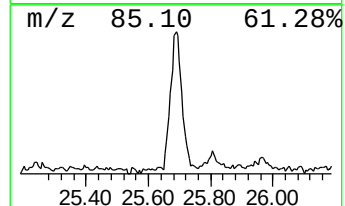
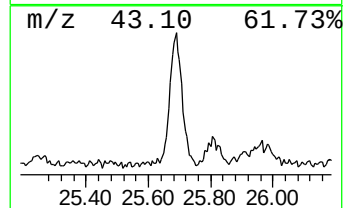
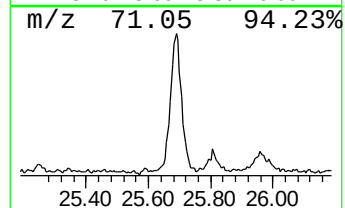
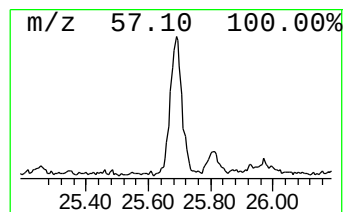
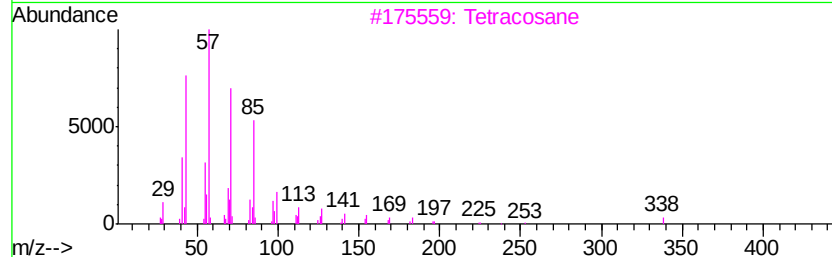
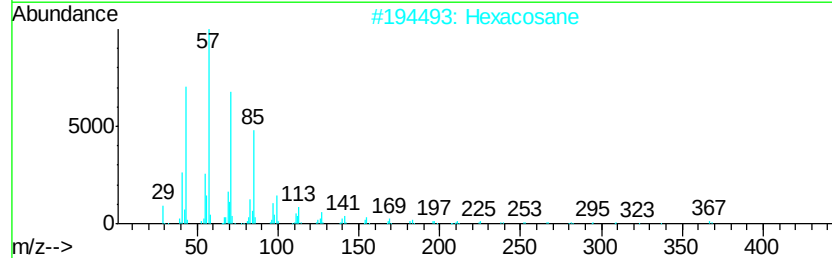
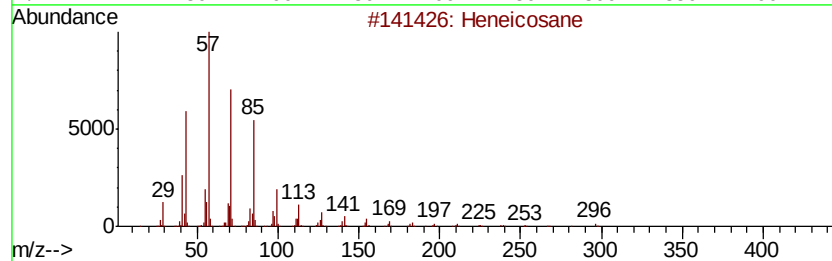
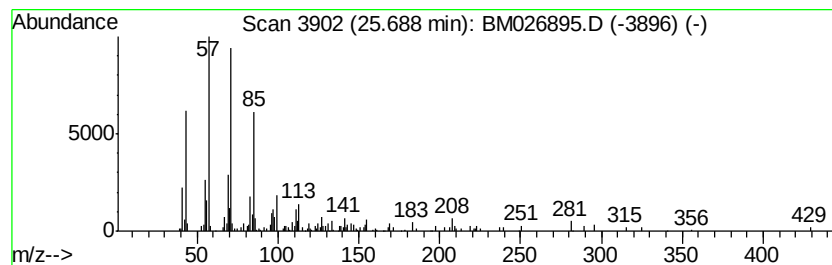
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.69	4.10 ng/ul	218459	Perylene-d12	23.46

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	87
2		Hexacosane	366	C26H54	000630-01-3	87
3		Tetracosane	338	C24H50	000646-31-1	87
4		triacontane	422	C30H62	000638-68-6	83
5		Eicosane	282	C20H42	000112-95-8	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM072820\
Data File : BM026895.D
Acq On : 28 Jul 2020 17:44
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Sample : L3388-06
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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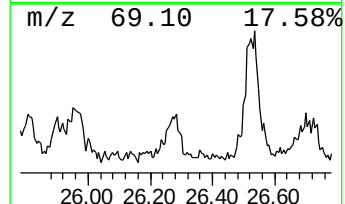
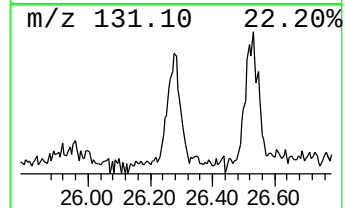
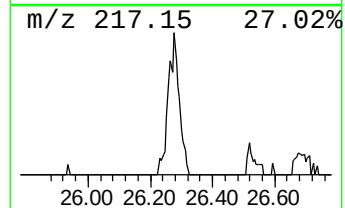
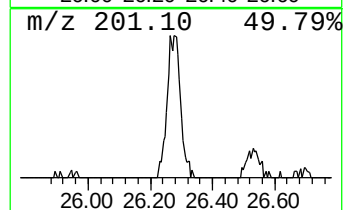
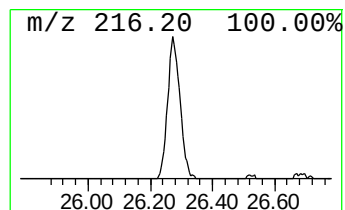
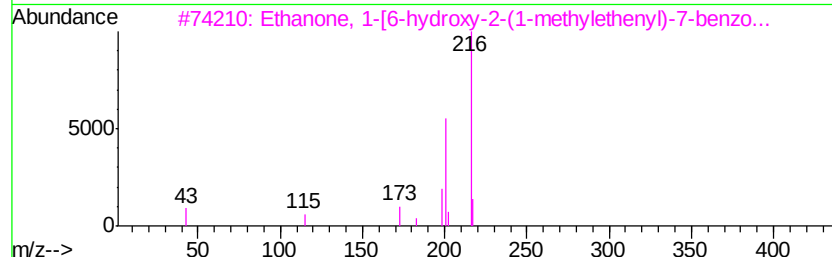
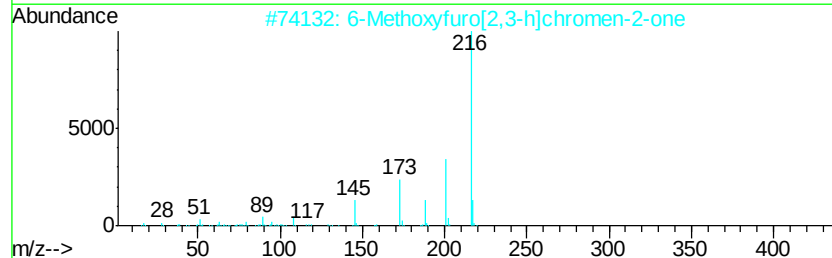
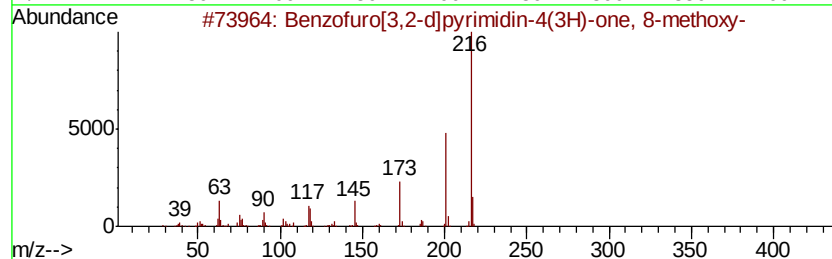
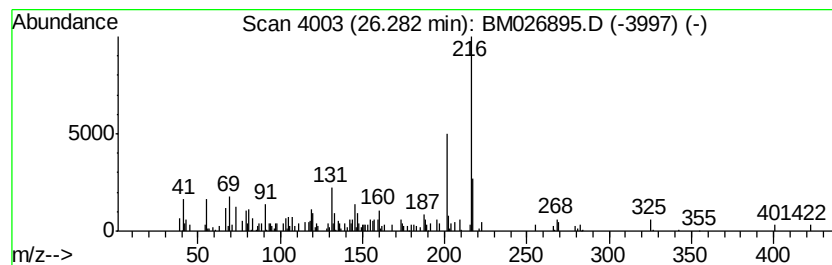
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzo-furo[3,2-d]pyrimidin-4... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.28	2.66 ng/ul	141757	Perylene-d12	23.46

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzo-furo[3,2-d]pyrimidin-4(3H)-...	216	C11H8N2O3	329210-45-9	64
2		6-Methoxyfuro[2,3-h]chromen-2-one	216	C12H8O4	000483-66-9	59
3		Ethanone, 1-[6-hydroxy-2-(1-meth...	216	C13H12O3	055682-75-2	53
4		Cyclobutane, tetrakis(1-methylet...	216	C16H24	088919-66-8	53
5		Methoxsalen	216	C12H8O4	000298-81-7	50



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM072820\
Data File : BM026895.D
Acq On : 28 Jul 2020 17:44
Operator : JU/CG
Sample : L3388-06
Misc :
ALS Vial : 14 Sample Multiplier: 1

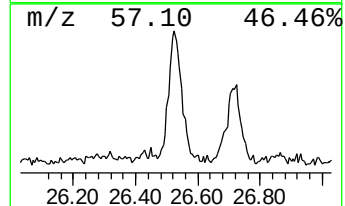
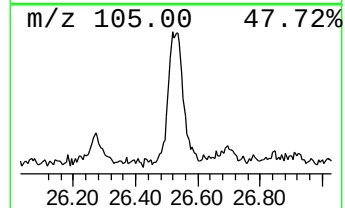
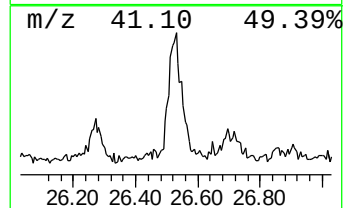
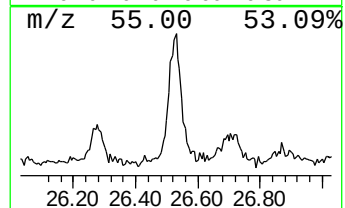
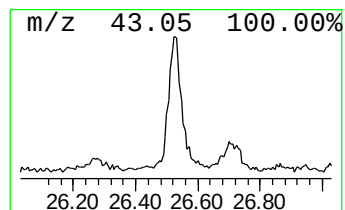
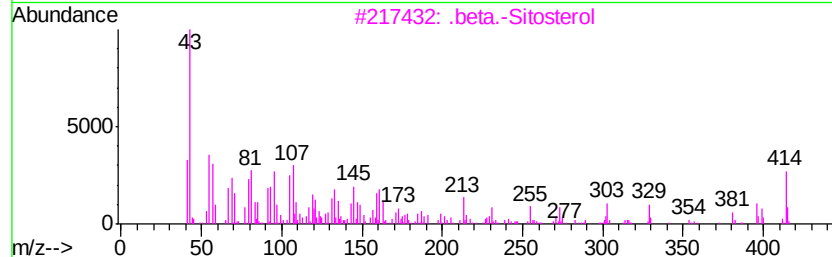
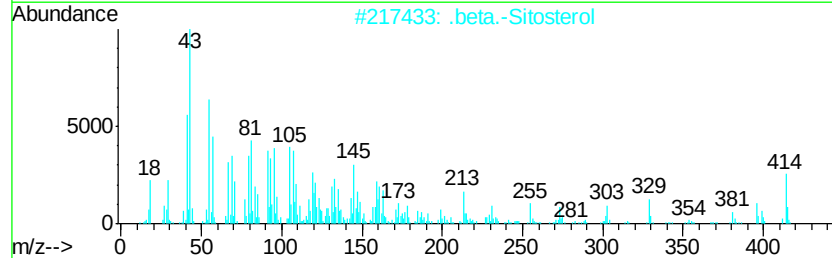
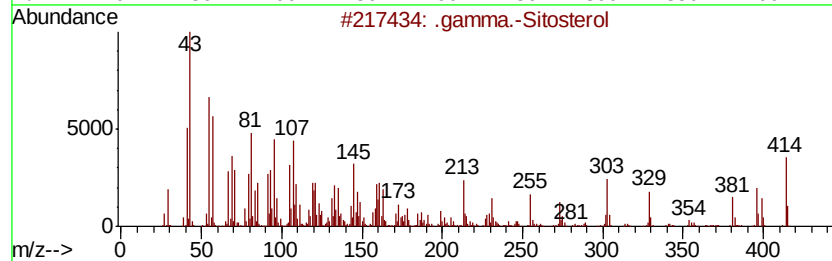
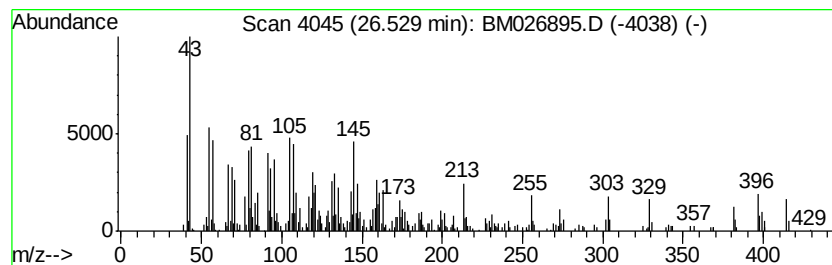
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 .gamma.-Sitosterol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.53	10.02 ng/ul	533812	Perylene-d12	23.46
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		.gamma.-Sitosterol	414 C29H50O	000083-47-6 99
2		.beta.-Sitosterol	414 C29H50O	000083-46-5 98
3		.beta.-Sitosterol	414 C29H50O	000083-46-5 53
4		Stigmastan-3,5-diene	396 C29H48	1000214-16-4 45
5		Campesterol	400 C28H48O	000474-62-4 43



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM072820\
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Acq On : 28 Jul 2020 17:44
Operator : JU/CG
Sample : L3388-06
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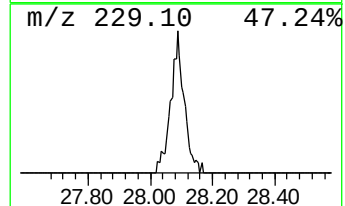
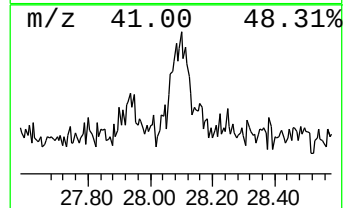
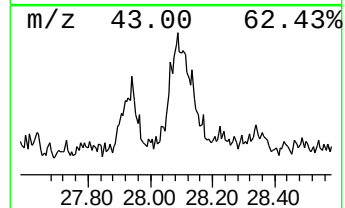
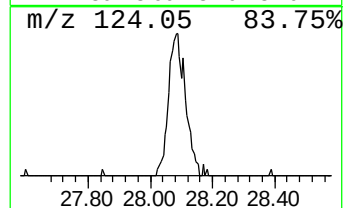
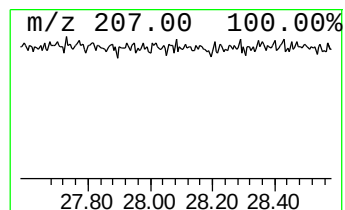
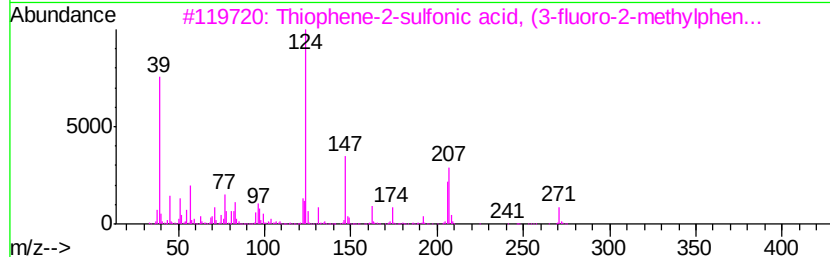
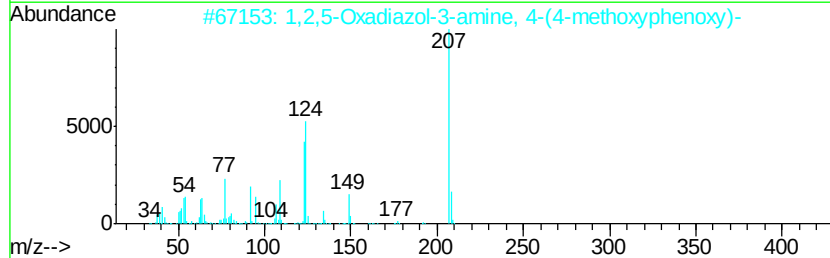
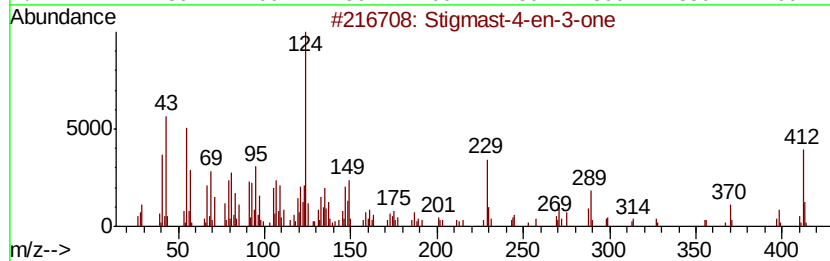
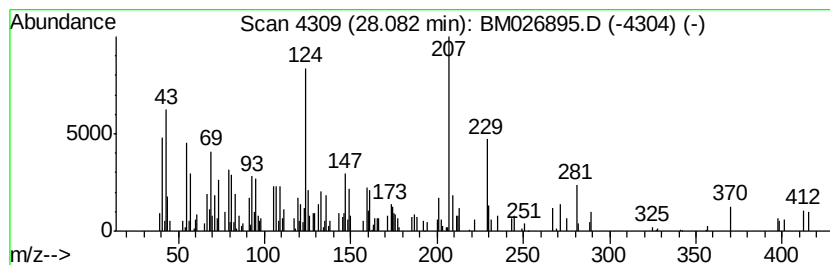
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 unknown-01 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
28.08	2.83 ng/ul	150704	Perylene-d12	23.46
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Stigmast-4-en-3-one	412 C29H48O	001058-61-3 46
2		1,2,5-Oxadiazol-3-amine, 4-(4-me...	207 C9H9N3O3	1000351-72-7 27
3		Thiophene-2-sulfonic acid, (3-fl...	271 C11H10FN02S2	1000311-21-1 25
4		Benz[bl]-1,4-oxazepine-4(5H)-thio...	207 C11H13NOS	1000258-63-4 18
5		3-(3-Fluoroanilino)-1-(3-nitroph...	288 C15H13FN2O3	300726-64-1 18



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM072820\
Data File : BM026895.D
Acq On : 28 Jul 2020 17:44
Operator : JU/CG
Sample : L3388-06
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM072720MA.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...	2.98	97.5	ng/ul	2264780	1	7.67	464481	20.0
2-Phenanthrenol, ...	20.35	3.2	ng/ul	161007	5	21.24	997721	20.0
(DEL) Alkane: Str...	22.85	4.9	ng/ul	261609	6	23.46	1065630	20.0
(DEL) Alkane: Str...	24.06	7.3	ng/ul	391443	6	23.46	1065630	20.0
(DEL) Alkane: Str...	25.69	4.1	ng/ul	218459	6	23.46	1065630	20.0
Benzofuro[3,2-d]p...	26.28	2.7	ng/ul	141757	6	23.46	1065630	20.0
.gamma.-Sitosterol	26.53	10.0	ng/ul	533812	6	23.46	1065630	20.0
unknown-01	28.08	2.8	ng/ul	150704	6	23.46	1065630	20.0