

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121719\
 Data File : BM024171.D
 Acq On : 19 Dec 2019 02:22
 Operator : JU
 Sample : K6171-14
 Misc :
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0BX1

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-BM121719MA.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	3.010	19	21	30	rVB	37364	59488	0.92%	0.066%
2	3.299	66	70	77	rVB	36377	61196	0.94%	0.067%
3	3.634	121	127	134	rBV2	60107	141749	2.19%	0.156%
4	3.704	134	139	148	rVB	312546	487722	7.52%	0.538%
5	4.410	251	259	267	rBV	185712	365433	5.63%	0.403%
6	4.551	275	283	289	rBV	160235	286332	4.41%	0.316%
7	4.610	289	293	302	rVB	64975	107280	1.65%	0.118%
8	4.998	352	359	368	rBV	60283	126206	1.95%	0.139%
9	5.898	508	512	521	rVB2	25516	44423	0.68%	0.049%
10	6.028	528	534	543	rBV	115960	203861	3.14%	0.225%
11	6.522	612	618	626	rBV4	23797	50541	0.78%	0.056%
12	6.645	632	639	652	rBV	763313	1429052	22.03%	1.576%
13	6.798	659	665	677	rBV	596830	989817	15.26%	1.092%
14	6.987	690	697	709	rBV	838073	1386502	21.38%	1.529%
15	7.445	768	775	785	rBV	1619544	2576399	39.72%	2.841%
16	7.539	786	791	794	rVV5	15415	34314	0.53%	0.038%
17	7.663	806	812	817	rBV	59483	93612	1.44%	0.103%
18	7.745	823	826	832	rVB7	15358	24145	0.37%	0.027%
19	8.075	878	882	888	rBV4	15525	34497	0.53%	0.038%
20	8.169	892	898	903	rBV	1043511	1651253	25.46%	1.821%
21	8.228	903	908	914	rVV	1736875	2828812	43.62%	3.120%
22	8.275	914	916	924	rVB	86754	141460	2.18%	0.156%
23	8.545	956	962	965	rBV6	24149	41416	0.64%	0.046%
24	8.598	965	971	981	rVV	767492	1309688	20.19%	1.444%
25	8.686	981	986	995	rVV7	21598	62006	0.96%	0.068%
26	9.034	1036	1045	1052	rBV	44012	80159	1.24%	0.088%
27	9.316	1086	1093	1110	rBV	626706	1084972	16.73%	1.197%
28	9.610	1138	1143	1145	rVV6	11121	22125	0.34%	0.024%
29	9.633	1145	1147	1152	rVV4	17318	27377	0.42%	0.030%
30	9.857	1176	1185	1198	rBV	1028183	1887185	29.10%	2.081%
31	10.216	1238	1246	1263	rBV	2477281	4163603	64.20%	4.592%
32	10.369	1263	1272	1290	rBV	630719	1332148	20.54%	1.469%
33	10.810	1340	1347	1356	rBV2	44177	120513	1.86%	0.133%
34	11.657	1486	1491	1497	rVB9	20039	40516	0.62%	0.045%

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Integration Parameters: LSCINT.P

Integrator: RTE
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Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM121719MA.M
 Title : SVOA CALIBRATION

35	11.739	1497	1505	1514	rBV	176533	330481	5.10%	0.364%
36	11.816	1514	1518	1527	rVB3	21195	44550	0.69%	0.049%
37	12.080	1557	1563	1576	rBV2	59020	150964	2.33%	0.166%
38	12.651	1655	1660	1664	rBV6	10649	23181	0.36%	0.026%
39	12.957	1706	1712	1722	rBV	212575	364903	5.63%	0.402%
40	13.427	1786	1792	1799	rVB8	17946	34760	0.54%	0.038%
41	13.533	1804	1810	1814	rVV	2093719	2862992	44.14%	3.158%
42	13.580	1814	1818	1828	rVV	454284	665678	10.26%	0.734%
43	13.798	1848	1855	1867	rBV	2350273	3562262	54.92%	3.929%
44	13.945	1875	1880	1891	rBV	202492	390805	6.03%	0.431%
45	14.033	1893	1895	1899	rVV4	20632	21796	0.34%	0.024%
46	14.110	1899	1908	1916	rVV	3854536	5760782	88.82%	6.353%
47	14.174	1916	1919	1924	rVB6	20685	29961	0.46%	0.033%
48	14.363	1945	1951	1972	rBV	354886	951688	14.67%	1.050%
49	14.515	1972	1977	1981	rVV5	41360	96376	1.49%	0.106%
50	14.563	1981	1985	1995	rVV5	99208	197843	3.05%	0.218%
51	14.874	2032	2038	2043	rBV3	26984	53569	0.83%	0.059%
52	14.933	2045	2048	2051	rVV	24753	34114	0.53%	0.038%
53	15.110	2072	2078	2085	rBV	3293787	4548009	70.12%	5.016%
54	15.263	2099	2104	2113	rBV	360889	559294	8.62%	0.617%
55	15.351	2113	2119	2125	rVV3	41405	100687	1.55%	0.111%
56	15.510	2142	2146	2150	rVB3	28958	37321	0.58%	0.041%
57	15.557	2150	2154	2159	rBV	52708	66405	1.02%	0.073%
58	15.857	2202	2205	2209	rVV4	22967	35210	0.54%	0.039%
59	16.033	2228	2235	2241	rBV8	21634	49459	0.76%	0.055%
60	16.098	2242	2246	2249	rBV6	13515	22930	0.35%	0.025%
61	16.215	2263	2266	2273	rVB5	31457	49838	0.77%	0.055%
62	16.304	2273	2281	2288	rBV2	73407	131076	2.02%	0.145%
63	16.527	2316	2319	2326	rVB7	30814	56466	0.87%	0.062%
64	16.615	2326	2334	2337	rBV9	12915	33690	0.52%	0.037%
65	16.651	2337	2340	2344	rVV2	38874	53309	0.82%	0.059%
66	16.680	2344	2345	2352	rVB5	15372	21686	0.33%	0.024%
67	16.792	2360	2364	2370	rBV7	34446	54925	0.85%	0.061%
68	16.862	2370	2376	2381	rBV	4619618	6485744	100.00%	7.153%
69	16.904	2381	2383	2387	rVV	245470	309445	4.77%	0.341%
70	16.962	2387	2393	2402	rVV2	3366252	4719022	72.76%	5.204%
71	17.068	2407	2411	2417	rVB9	27204	50121	0.77%	0.055%

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72	17.180	2426	2430	2438	rVB3	49754	81784	1.26%	0.090%
73	17.280	2440	2447	2457	rVV2	144337	290841	4.48%	0.321%
74	17.392	2463	2466	2470	rBV4	19559	21840	0.34%	0.024%
75	17.521	2486	2488	2491	rBV3	17637	21754	0.34%	0.024%
76	17.668	2509	2513	2520	rBV9	50423	84088	1.30%	0.093%
77	17.739	2521	2525	2529	rBV4	53583	81118	1.25%	0.089%
78	17.792	2529	2534	2553	rBV	842795	1414541	21.81%	1.560%
79	17.933	2555	2558	2563	rVB2	50223	85977	1.33%	0.095%
80	18.104	2582	2587	2591	rBV8	23452	48346	0.75%	0.053%
81	18.233	2604	2609	2610	rBV4	29718	53343	0.82%	0.059%
82	18.298	2617	2620	2626	rVB5	38203	54120	0.83%	0.060%
83	18.668	2680	2683	2686	rBV3	30005	42676	0.66%	0.047%
84	18.709	2686	2690	2698	rVV3	207956	309335	4.77%	0.341%
85	18.821	2705	2709	2717	rBV9	44216	104069	1.60%	0.115%
86	18.933	2724	2728	2735	rBV	676678	883120	13.62%	0.974%
87	19.086	2750	2754	2760	rBV2	188332	292729	4.51%	0.323%
88	19.268	2780	2785	2797	rBV2	3806037	5978048	92.17%	6.593%
89	19.856	2882	2885	2889	rVB	109760	120051	1.85%	0.132%
90	20.309	2958	2962	2966	rBV	873502	945302	14.58%	1.043%
91	20.815	3043	3048	3054	rBV	1417788	1927457	29.72%	2.126%
92	21.074	3087	3092	3097	rBV	4533359	6020895	92.83%	6.640%
93	21.674	3191	3194	3206	rVB2	692716	1347964	20.78%	1.487%
94	22.115	3266	3269	3274	rVB	480615	575000	8.87%	0.634%
95	22.591	3346	3350	3354	rVB2	1050696	1408957	21.72%	1.554%
96	23.074	3426	3432	3436	rBV	1898173	3261566	50.29%	3.597%
97	23.121	3437	3440	3447	rVB2	683594	1057882	16.31%	1.167%
98	23.203	3449	3454	3462	rVB	2596365	4381102	67.55%	4.832%
99	23.727	3539	3543	3549	rVB	709330	1210509	18.66%	1.335%
100	24.409	3654	3659	3671	rVB4	854001	2310660	35.63%	2.548%

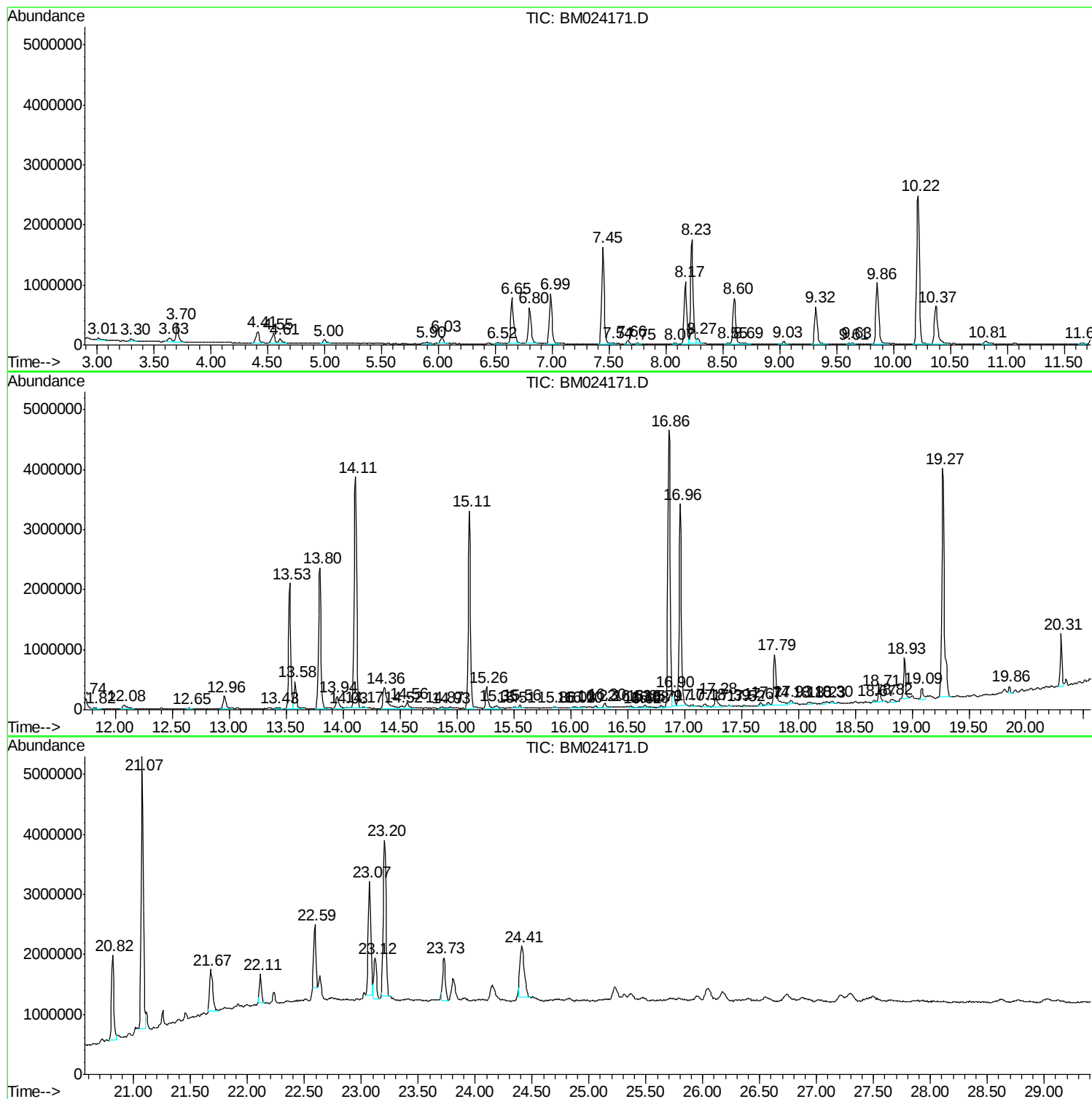
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ALS Vial : 45 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampled :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM121719MA.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
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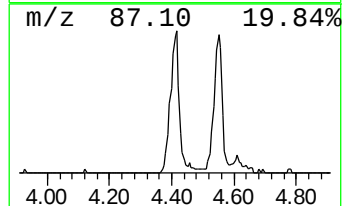
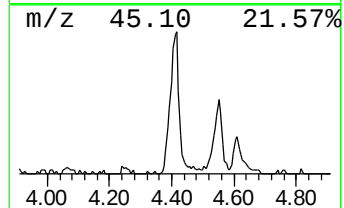
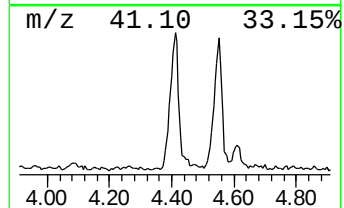
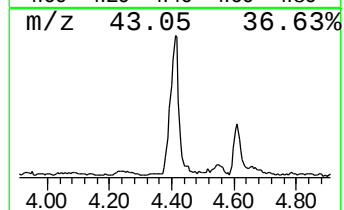
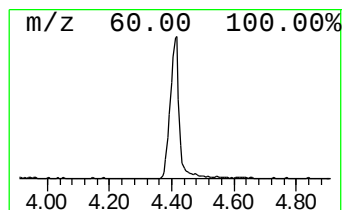
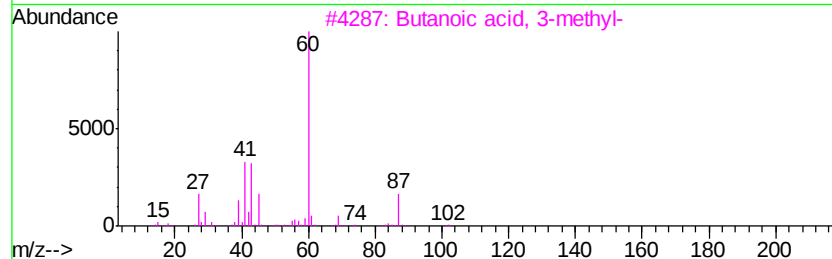
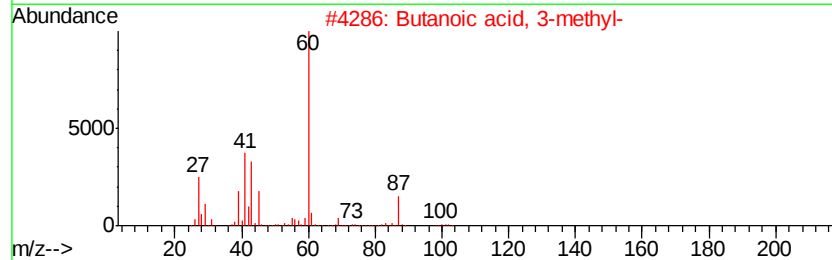
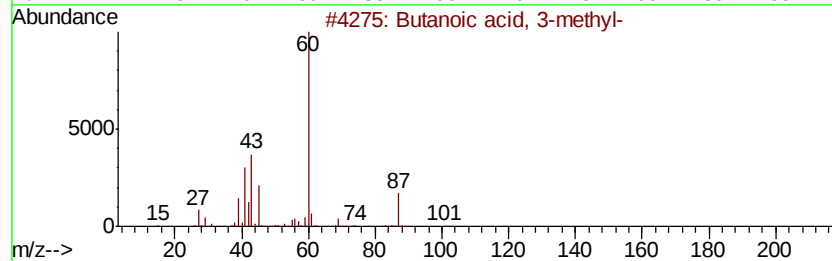
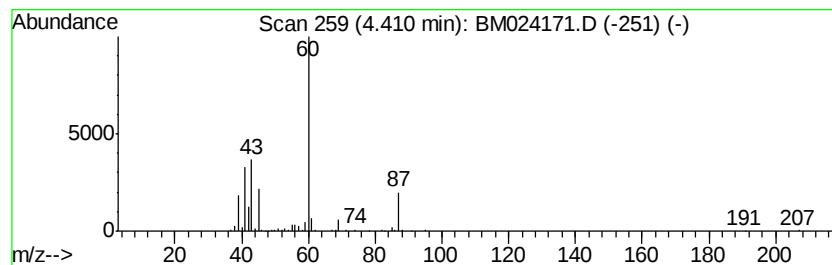
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM121719MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Butanoic acid, 3-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.41	2.84 ng/ul	365433	1,4-Dichlorobenzene-d4	7.45

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butanoic acid, 3-methyl-	102	C5H10O2	000503-74-2	90
2		Butanoic acid, 3-methyl-	102	C5H10O2	000503-74-2	78
3		Butanoic acid, 3-methyl-	102	C5H10O2	000503-74-2	78
4		Pentanoic acid, 3-methyl-	116	C6H12O2	000105-43-1	56
5		1,2-Hydrazinedicarboxaldehyde	88	C2H4N2O2	000628-36-4	38



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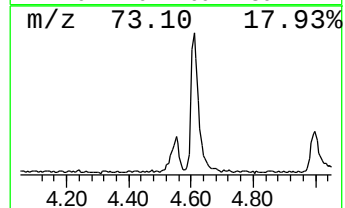
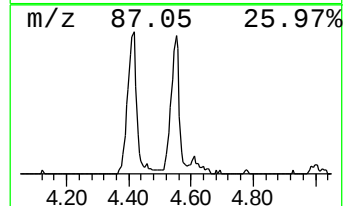
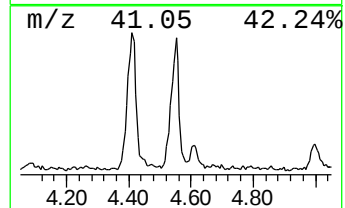
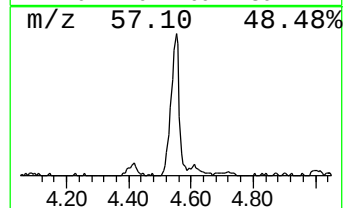
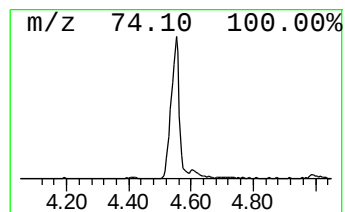
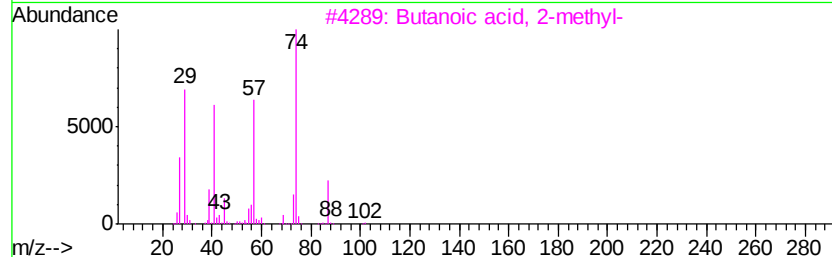
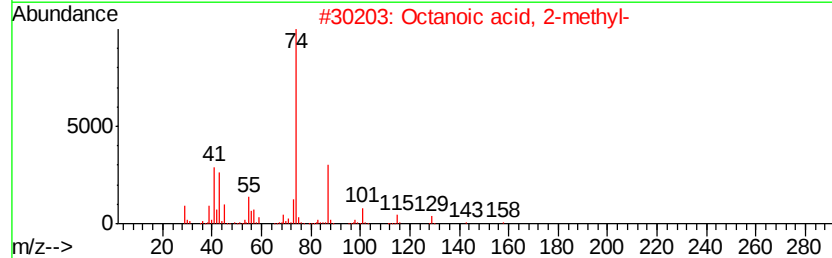
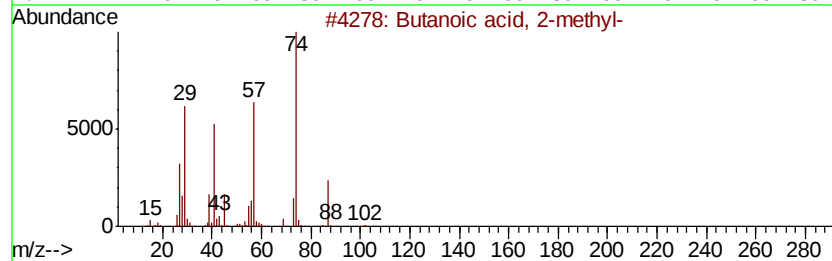
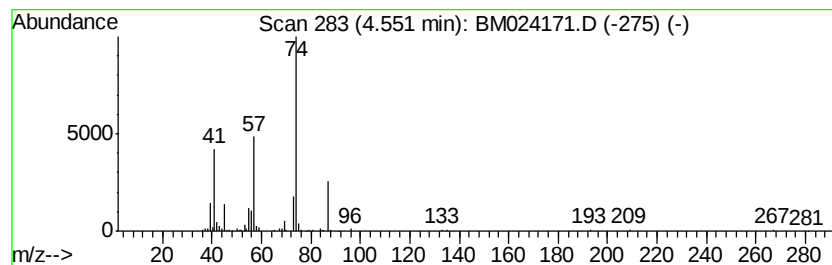
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Butanoic acid, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.		
4.55	2.22 ng/ul	286332	1,4-Dichlorobenzene-d4		7.45		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butanoic acid, 2-methyl-	102	C5H10O2	000116-53-0	83
2			Octanoic acid, 2-methyl-	158	C9H18O2	003004-93-1	78
3			Butanoic acid, 2-methyl-	102	C5H10O2	000116-53-0	78
4			Pentanoic acid, 2-methyl-	116	C6H12O2	000097-61-0	64
5			Hexanoic acid, 2-methyl-	130	C7H14O2	004536-23-6	64



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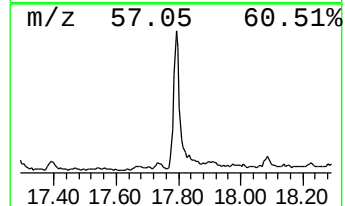
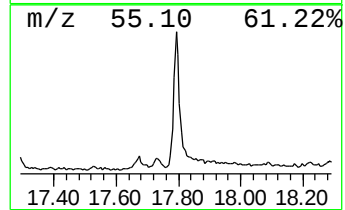
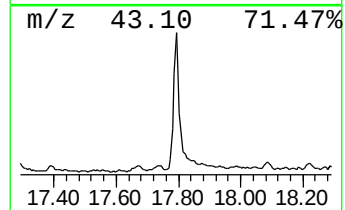
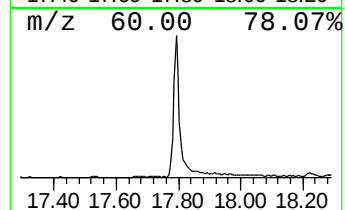
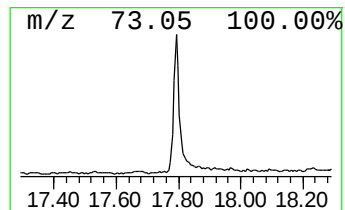
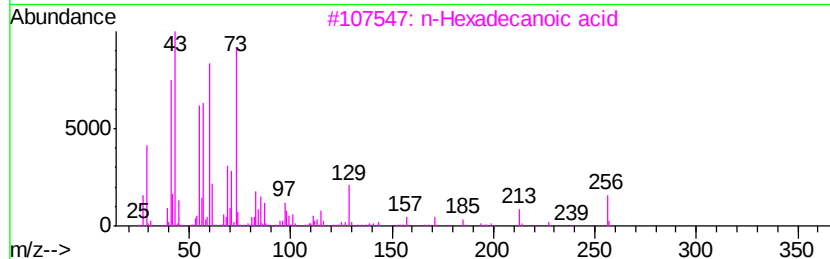
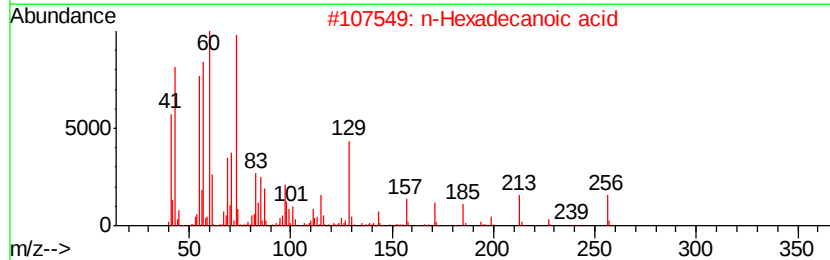
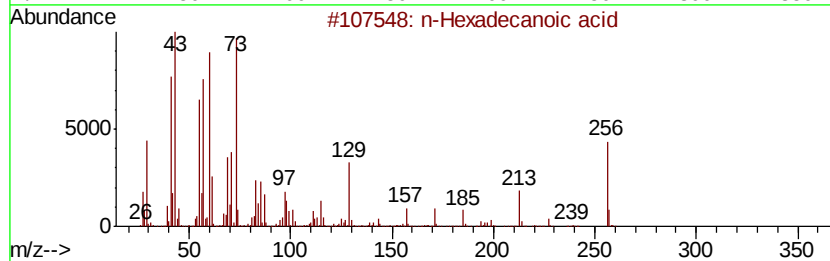
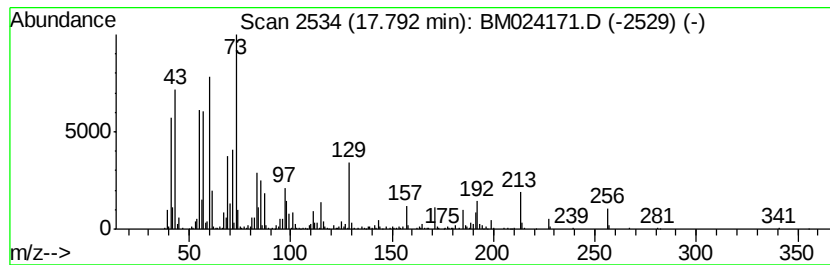
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.79	4.36 ng/ul	1414540	Phenanthrene-d10	16.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
4		Tridecanoic acid	214	C13H26O2	000638-53-9	91
5		Octadecanoic acid	284	C18H36O2	000057-11-4	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121719\
Data File : BM024171.D
Acq On : 19 Dec 2019 02:22
Operator : JU
Sample : K6171-14
Misc :
ALS Vial : 45 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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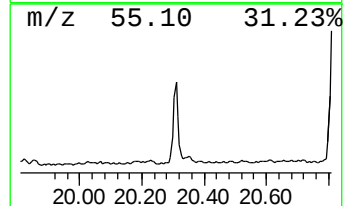
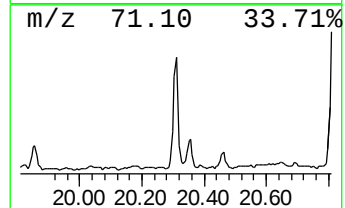
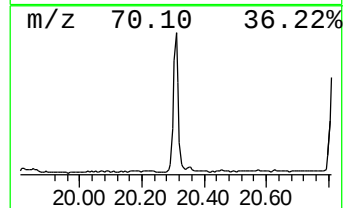
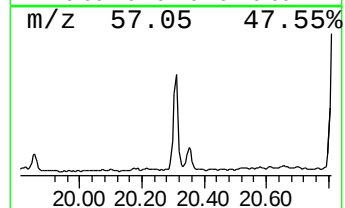
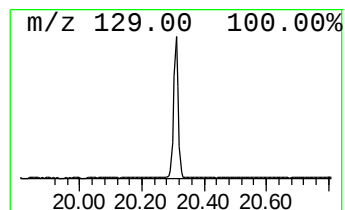
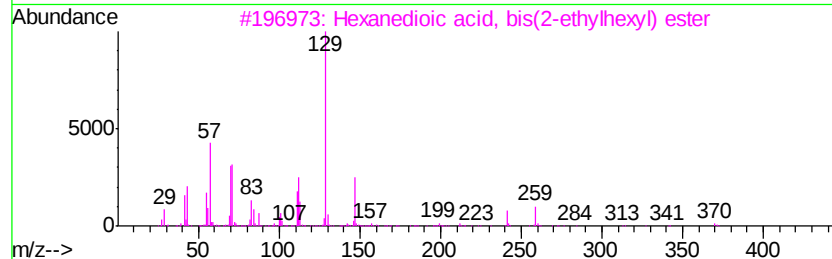
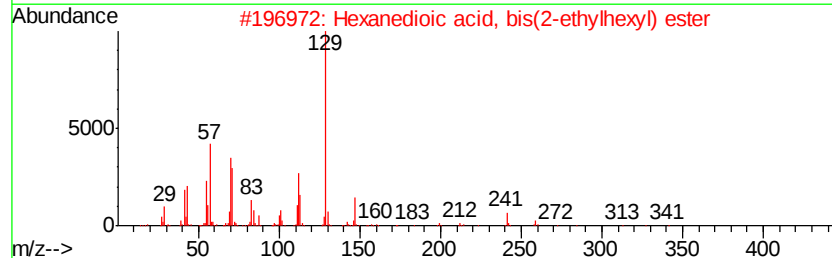
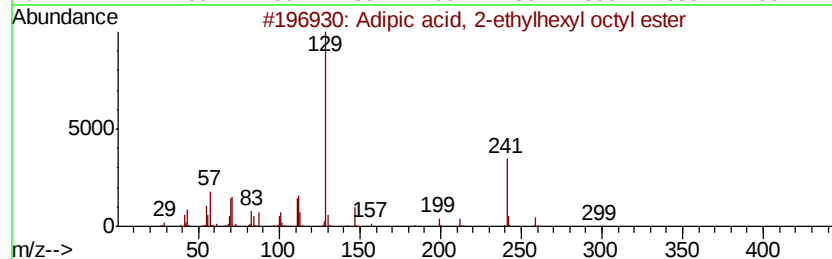
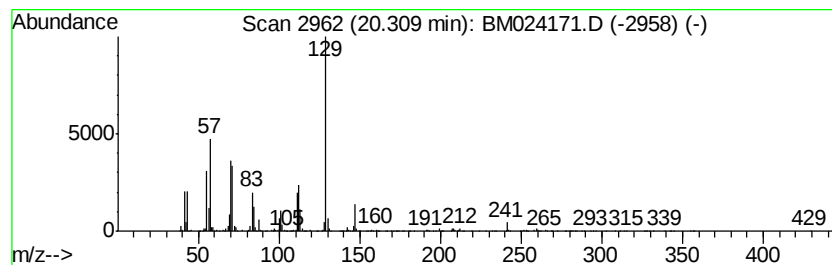
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Adipic acid, 2-ethylhexyl o... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.31	3.14 ng/ul	945302	Chrysene-d12	21.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Adipic acid, 2-ethylhexyl octyl ...	370	C22H42O4	1000324-48-6	91
2		Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	87
3		Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	86
4		Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	83
5		Diisooctyl adipate	370	C22H42O4	001330-86-5	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121719\
Data File : BM024171.D
Acq On : 19 Dec 2019 02:22
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ClientSampleId :
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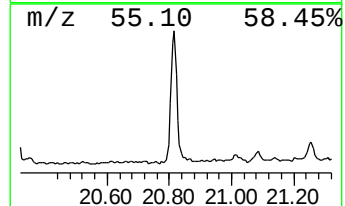
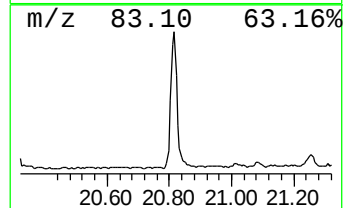
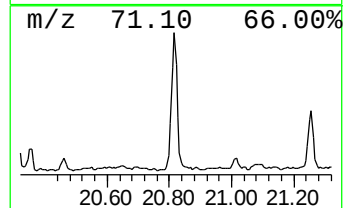
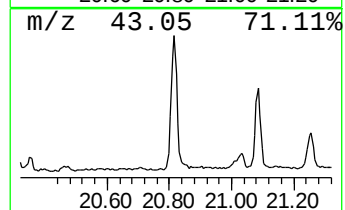
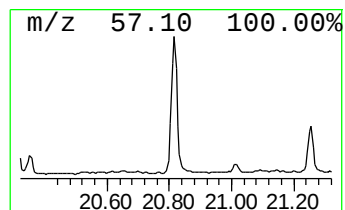
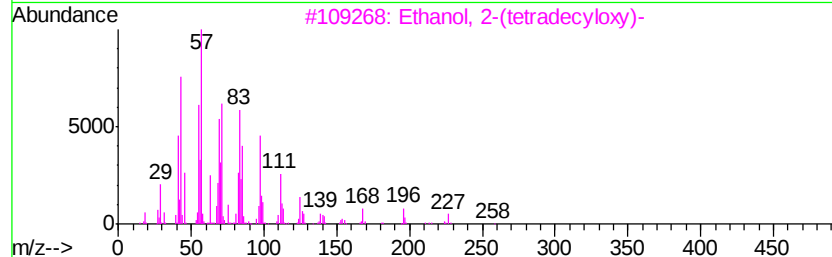
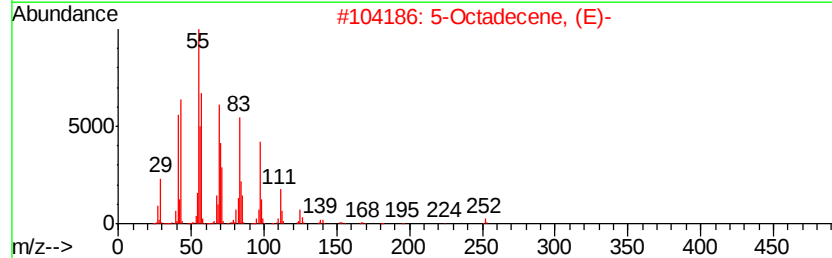
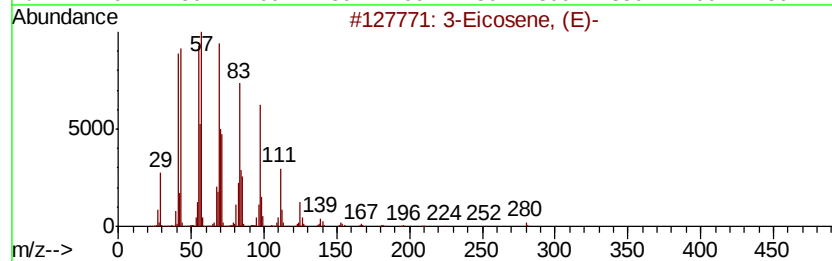
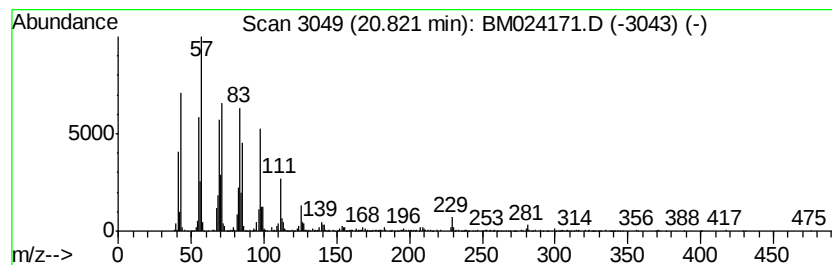
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Cyclic20.82 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.82	6.40 ng/ul	1927460	Chrysene-d12	21.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Eicosene, (E)-	280	C20H40	074685-33-9	97
2		5-Octadecene, (E)-	252	C18H36	007206-21-5	95
3		Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	91
4		Oxalic acid, isobutyl hexadecyl ...	370	C22H42O4	1000309-38-1	91
5		Oxalic acid, isobutyl heptadecyl...	384	C23H44O4	1000309-38-2	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121719\
Data File : BM024171.D
Acq On : 19 Dec 2019 02:22
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Sample : K6171-14
Misc :
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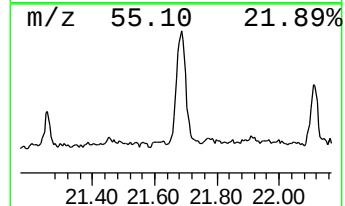
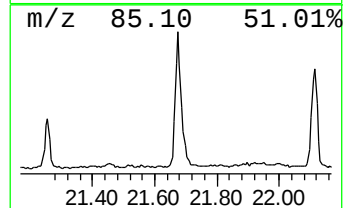
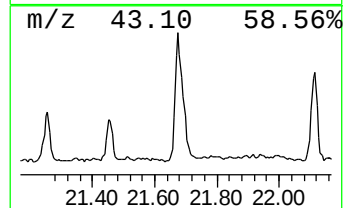
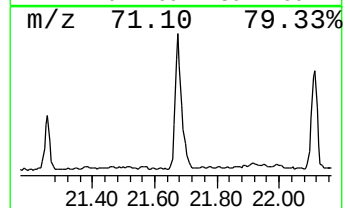
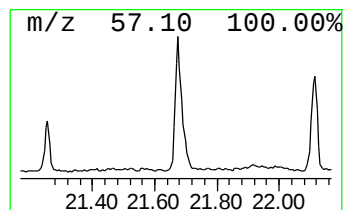
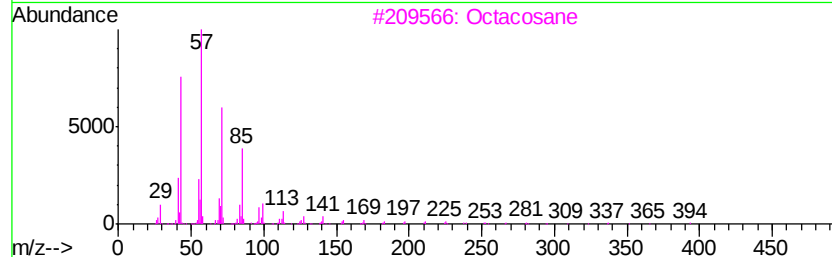
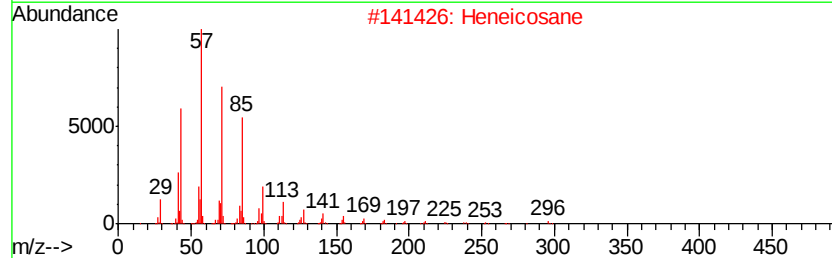
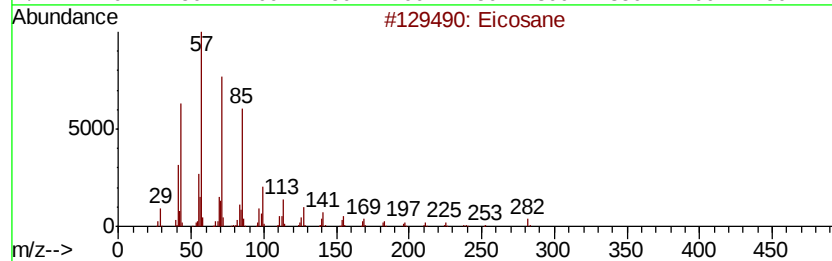
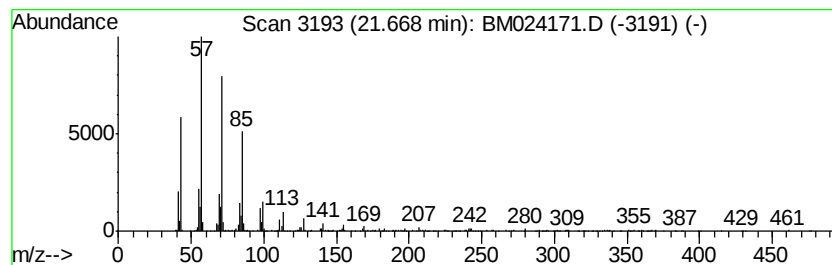
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.67	4.48 ng/ul	1347960	Chrysene-d12	21.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	95
2		Heneicosane	296	C21H44	000629-94-7	91
3		Octacosane	394	C28H58	000630-02-4	91
4		Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91
5		Hexacosane	366	C26H54	000630-01-3	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121719\
Data File : BM024171.D
Acq On : 19 Dec 2019 02:22
Operator : JU
Sample : K6171-14
Misc :
ALS Vial : 45 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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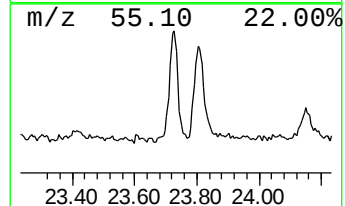
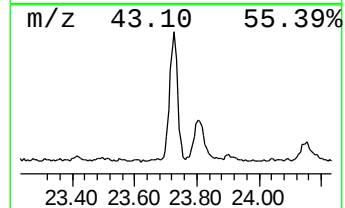
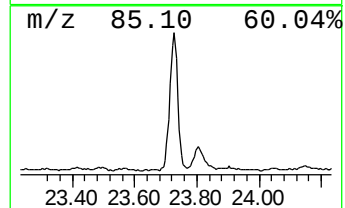
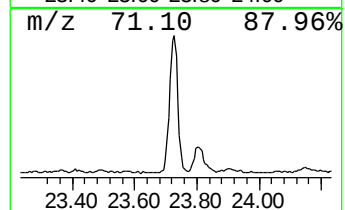
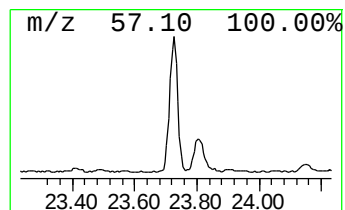
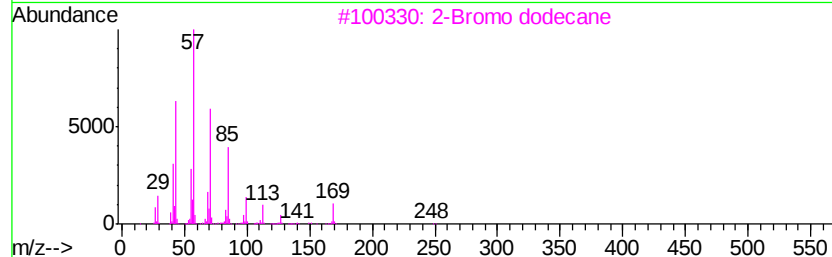
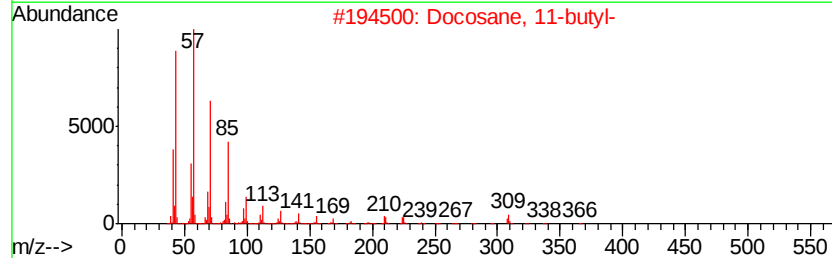
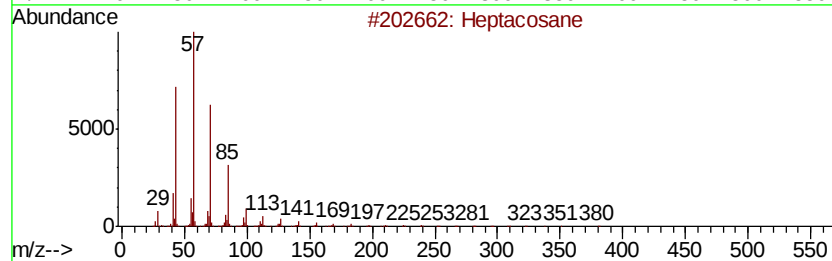
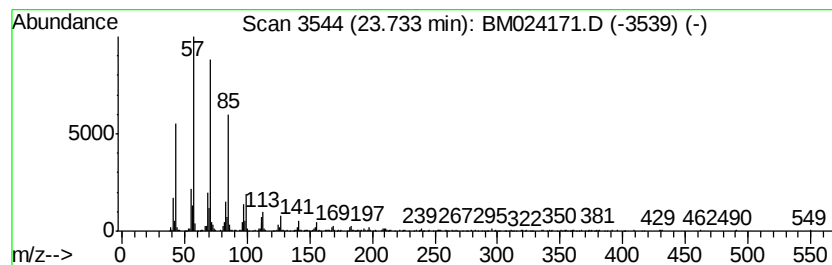
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.73	5.53 ng/ul	1210510	Perylene-d12	23.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptacosane	380	C27H56	000593-49-7	94
2		Docosane, 11-butyl-	366	C26H54	013475-76-8	93
3		2-Bromo dodecane	248	C12H25Br	013187-99-0	93
4		Tetracosane	338	C24H50	000646-31-1	91
5		2-methyloctacosane	408	C29H60	1000376-72-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121719\
Data File : BM024171.D
Acq On : 19 Dec 2019 02:22
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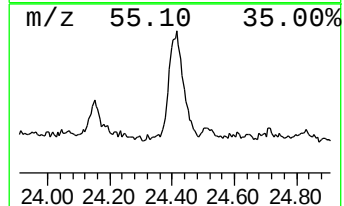
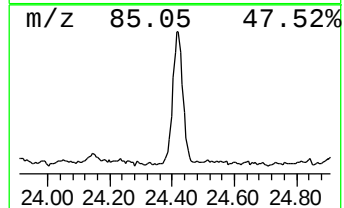
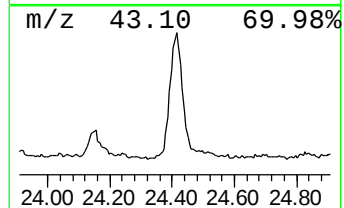
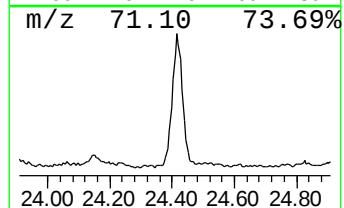
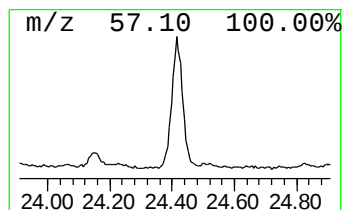
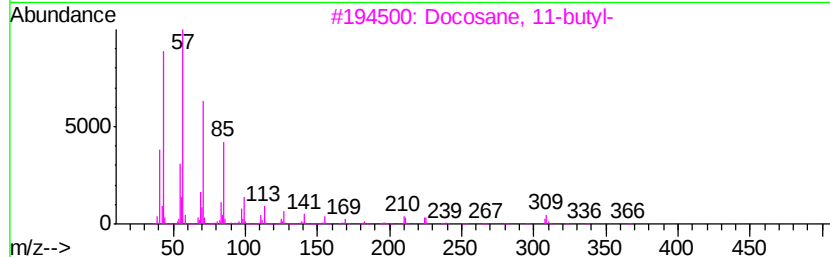
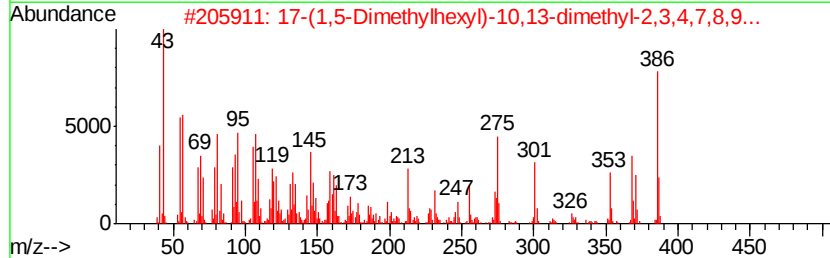
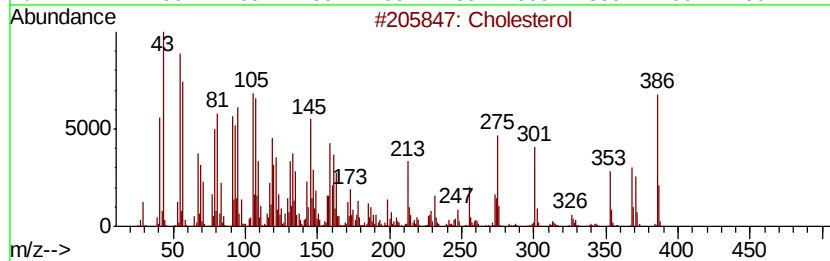
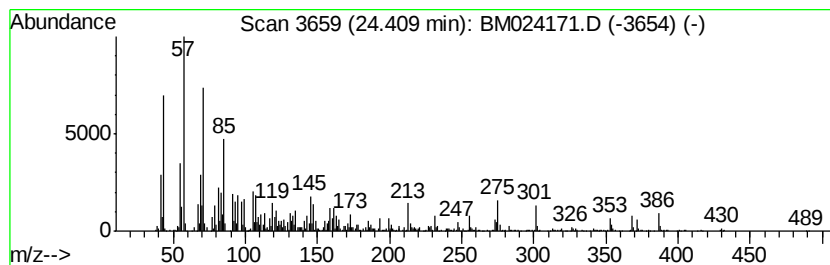
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Cholesterol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.41	10.55 ng/ul	2310660	Perylene-d12	23.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cholesterol	386	C27H46O	000057-88-5	91
2		17-(1,5-Dimethylhexyl)-10,13-dim...	386	C27H46O	1000210-38-4	91
3		Docosane, 11-butyl-	366	C26H54	013475-76-8	86
4		Sulfurous acid, dodecyl 2-propyl...	292	C15H32O3S	1000309-12-3	70
5		Dodecane	170	C12H26	000112-40-3	70



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM121719\
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Sample : K6171-14
Misc :
ALS Vial : 45 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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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
Butanoic acid, 3-...	4.41	2.8	ng/ul	365433	1	7.45	2576400	20.0
Butanoic acid, 2-...	4.55	2.2	ng/ul	286332	1	7.45	2576400	20.0
n-Hexadecanoic acid	17.79	4.4	ng/ul	1414540	4	16.86	6485740	20.0
Adipic acid, 2-et...	20.31	3.1	ng/ul	945302	5	21.07	6020900	20.0
(DEL) Alkane: Cyc...	20.82	6.4	ng/ul	1927460	5	21.07	6020900	20.0
(DEL) Alkane: Str...	21.67	4.5	ng/ul	1347960	5	21.07	6020900	20.0
(DEL) Alkane: Str...	23.73	5.5	ng/ul	1210510	6	23.20	4381100	20.0
Cholesterol	24.41	10.6	ng/ul	2310660	6	23.20	4381100	20.0