

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082319\
 Data File : BM022291.D
 Acq On : 24 Aug 2019 13:22
 Operator : HP/JU
 Sample : K4242-12
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AJ0

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-BM082219MA.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	3.134	21	25	30	rBV	9678	14526	0.72%	0.064%
2	3.463	77	81	84	rBV3	1566	2475	0.12%	0.011%
3	3.728	120	126	133	rVB2	3640	8703	0.43%	0.038%
4	3.816	139	141	146	rVB3	2215	2691	0.13%	0.012%
5	4.493	251	256	263	rVB3	3270	6088	0.30%	0.027%
6	4.710	290	293	297	rBV2	2487	3675	0.18%	0.016%
7	6.222	545	550	555	rVB	21552	30295	1.51%	0.133%
8	6.751	633	640	655	rBV	236542	434310	21.59%	1.911%
9	6.910	662	667	682	rBV	171241	286959	14.27%	1.263%
10	7.092	692	698	706	rVB	288240	449024	22.32%	1.976%
11	7.557	770	777	787	rBV	191244	306423	15.23%	1.348%
12	7.745	805	809	813	rBV3	5964	5731	0.28%	0.025%
13	8.275	893	899	905	rBV	324925	527986	26.25%	2.323%
14	8.334	905	909	917	rVB	36449	61123	3.04%	0.269%
15	8.728	969	976	985	rBV	256950	434382	21.60%	1.912%
16	9.434	1090	1096	1111	rBV	249788	409244	20.35%	1.801%
17	9.745	1145	1149	1154	rVB2	2212	2915	0.14%	0.013%
18	9.969	1180	1187	1203	rBV	341277	616998	30.68%	2.715%
19	10.328	1241	1248	1256	rBV	292809	467102	23.22%	2.056%
20	10.498	1270	1277	1297	rBV	262112	526757	26.19%	2.318%
21	10.651	1299	1303	1309	rVB3	6183	8634	0.43%	0.038%
22	11.933	1516	1521	1526	rVB2	3829	6224	0.31%	0.027%
23	12.210	1564	1568	1572	rBV3	2492	3233	0.16%	0.014%
24	13.051	1708	1711	1714	rBV2	8902	10955	0.54%	0.048%
25	13.080	1714	1716	1721	rVB	3103	3453	0.17%	0.015%
26	13.498	1784	1787	1789	rVB	3255	3434	0.17%	0.015%
27	13.586	1800	1802	1805	rBV2	4432	4076	0.20%	0.018%
28	13.639	1805	1811	1815	rVV	660965	896995	44.60%	3.947%
29	13.686	1815	1819	1826	rVB	150412	205450	10.21%	0.904%
30	13.904	1848	1856	1867	rBV2	776731	1102470	54.81%	4.851%
31	14.033	1875	1878	1883	rVB2	2134	2466	0.12%	0.011%
32	14.145	1891	1897	1902	rBV	177899	239307	11.90%	1.053%
33	14.210	1902	1908	1915	rVV	455691	675518	33.59%	2.973%
34	14.474	1944	1953	1974	rBV	177050	398711	19.82%	1.755%

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Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM082219MA.M
 Title : SVOA CALIBRATION

35	14.645	1979	1982	1984	rVV3	4572	5831	0.29%	0.026%
36	14.674	1984	1987	1992	rVB4	5483	7112	0.35%	0.031%
37	14.986	2034	2040	2047	rBV	8074	11374	0.57%	0.050%
38	15.215	2073	2079	2091	rBV	983348	1347754	67.01%	5.931%
39	15.380	2102	2107	2116	rBV	222632	311411	15.48%	1.370%
40	15.457	2116	2120	2123	rVB3	7734	10209	0.51%	0.045%
41	15.580	2137	2141	2146	rBV3	15724	19621	0.98%	0.086%
42	15.680	2155	2158	2163	rVB	2482	3689	0.18%	0.016%
43	16.168	2238	2241	2245	rVV	3596	4054	0.20%	0.018%
44	16.380	2271	2277	2282	rBV3	2466	6162	0.31%	0.027%
45	16.598	2312	2314	2318	rBV2	2859	3771	0.19%	0.017%
46	16.745	2336	2339	2345	rVB3	4408	6393	0.32%	0.028%
47	16.874	2358	2361	2365	rVB3	1709	2849	0.14%	0.013%
48	16.974	2372	2378	2384	rBV2	628933	858181	42.67%	3.776%
49	17.074	2389	2395	2402	rBV	1094930	1534243	76.28%	6.752%
50	17.151	2405	2408	2410	rBV4	1743	2100	0.10%	0.009%
51	17.262	2426	2427	2431	rVB2	1911	2138	0.11%	0.009%
52	17.362	2441	2444	2445	rBV2	2025	2056	0.10%	0.009%
53	17.545	2471	2475	2477	rBV4	2254	3035	0.15%	0.013%
54	17.739	2505	2508	2513	rBV4	5649	10119	0.50%	0.045%
55	17.809	2518	2520	2522	rVB2	2332	2191	0.11%	0.010%
56	17.874	2525	2531	2541	rBV	232925	340991	16.95%	1.501%
57	18.151	2575	2578	2582	rBV6	3541	4524	0.22%	0.020%
58	18.227	2589	2591	2593	rVB2	2329	2076	0.10%	0.009%
59	18.292	2599	2602	2607	rBV4	1865	3046	0.15%	0.013%
60	18.421	2620	2624	2626	rBV3	1897	2378	0.12%	0.010%
61	18.668	2664	2666	2670	rBV5	3304	3475	0.17%	0.015%
62	19.015	2722	2725	2728	rBV	25280	36409	1.81%	0.160%
63	19.162	2745	2750	2759	rBV	101442	162461	8.08%	0.715%
64	19.268	2765	2768	2775	rVB2	11010	13976	0.69%	0.062%
65	19.386	2782	2788	2796	rBV	1556787	2011335	100.00%	8.851%
66	19.903	2873	2876	2877	rBV3	9382	10777	0.54%	0.047%
67	20.321	2942	2947	2953	rBV	1183182	1289932	64.13%	5.676%
68	20.745	3014	3019	3024	rBV	415044	476377	23.68%	2.096%
69	20.886	3039	3043	3053	rVB	215484	300737	14.95%	1.323%
70	21.186	3089	3094	3101	rBV	915420	1158575	57.60%	5.098%
71	21.321	3114	3117	3123	rBV	65810	79848	3.97%	0.351%

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Sampling : 1 Min Area: 0.1 % of largest Peak
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Stop Thrs : 0 Peak Location: TOP

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Title : SVOA CALIBRATION

72	21.744	3186	3189	3192	rBV	132694	131469	6.54%	0.579%
73	22.192	3262	3265	3270	rVB	220047	266740	13.26%	1.174%
74	22.680	3343	3348	3355	rVB	527665	727996	36.19%	3.204%
75	23.238	3436	3443	3454	rVB	972628	1977276	98.31%	8.701%
76	23.380	3461	3467	3474	rBV	466609	857308	42.62%	3.773%
77	23.838	3540	3545	3554	rVB	310793	554051	27.55%	2.438%

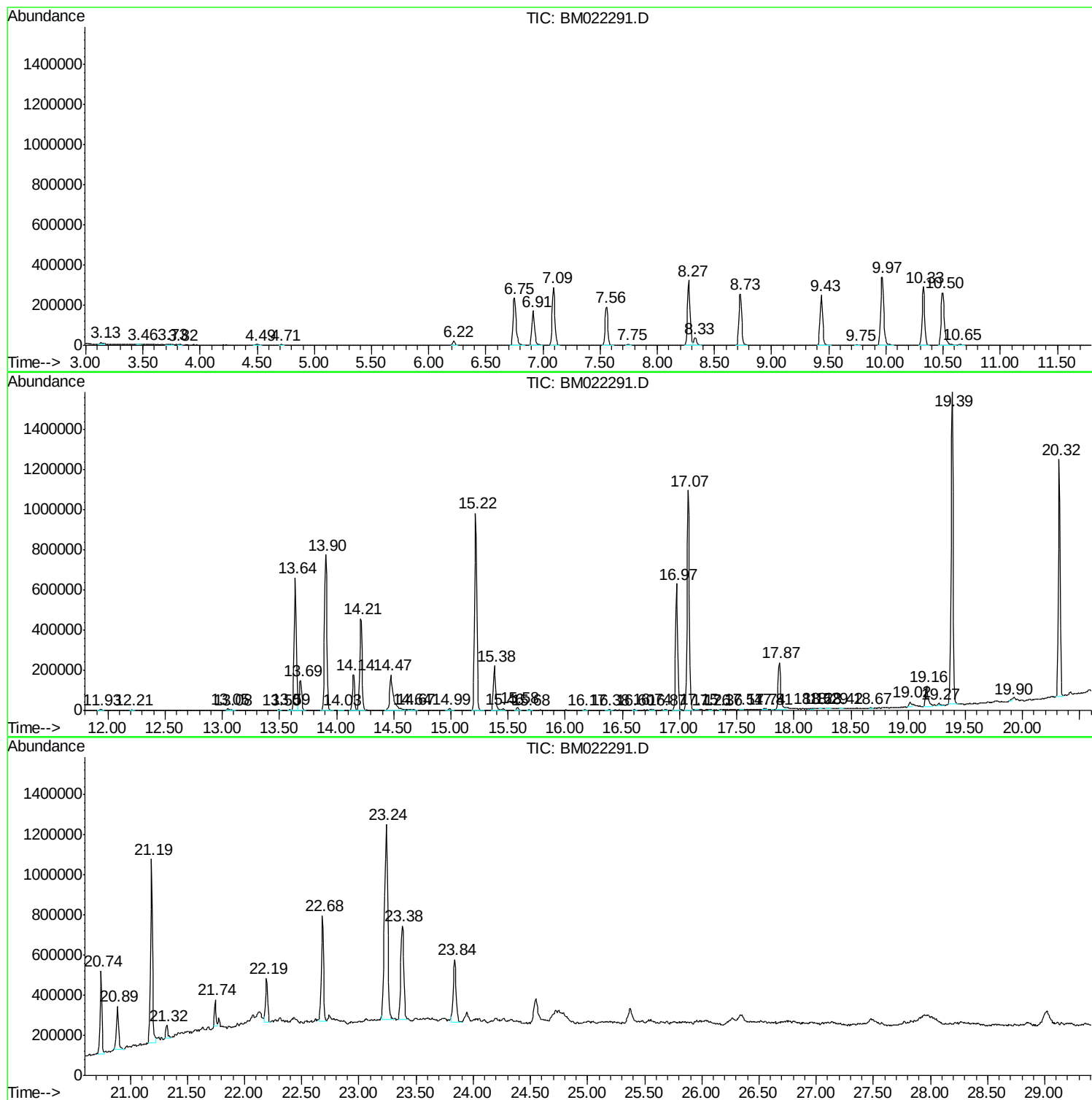
Sum of corrected areas: 22724383

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM082219MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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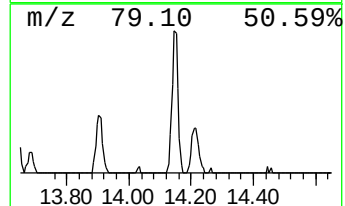
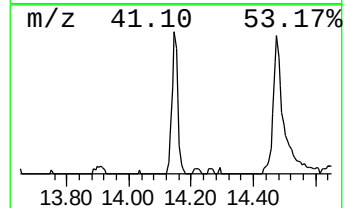
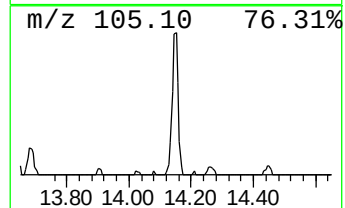
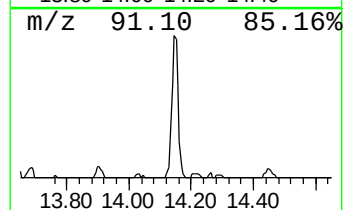
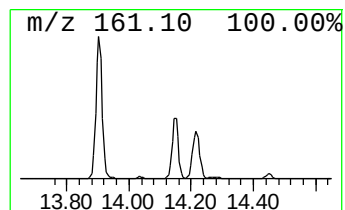
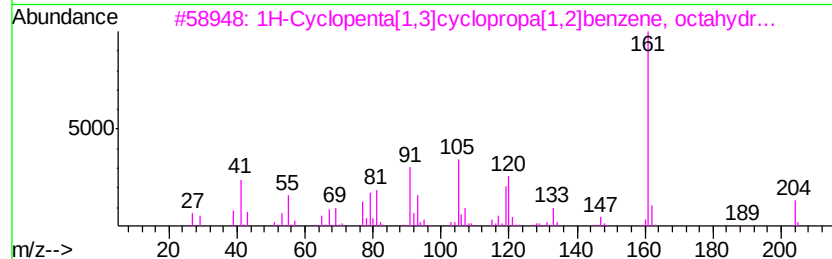
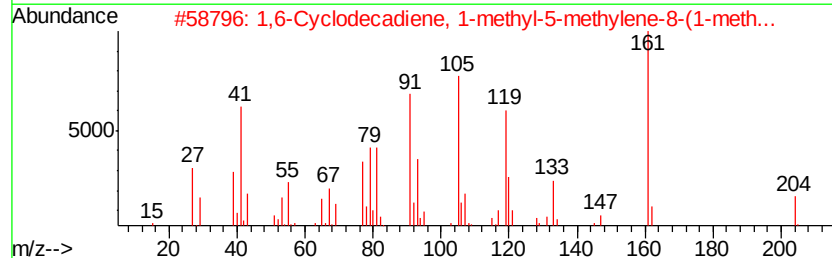
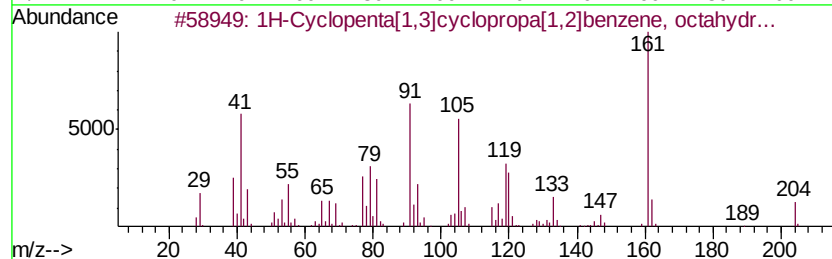
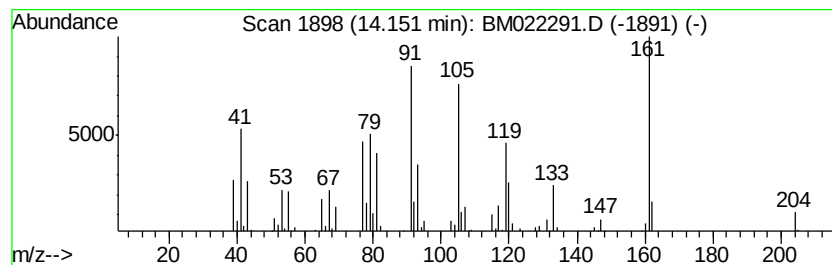
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM082219MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 1H-Cyclopenta[1,3]cycloprop... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.15	7.09 ng/ul	239307	Acenaphthene-d10	14.21
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		1H-Cyclopenta[1,3]cyclopropa[1,2]...	204 C15H24	013744-15-5 94
2		1,6-Cyclodecadiene, 1-methyl-5-m...	204 C15H24	023986-74-5 91
3		1H-Cyclopenta[1,3]cyclopropa[1,2]...	204 C15H24	013744-15-5 90
4		(+)-Epi-bicyclosesquiphellandrene	204 C15H24	054324-03-7 78
5		Naphthalene, 1,2,3,4,4a,5,6,8a-o...	204 C15H24	039029-41-9 68



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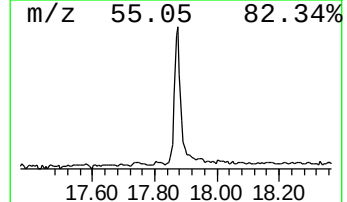
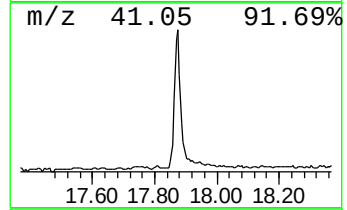
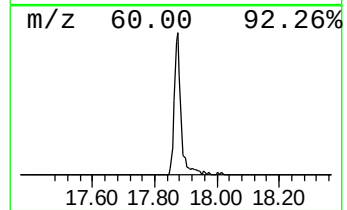
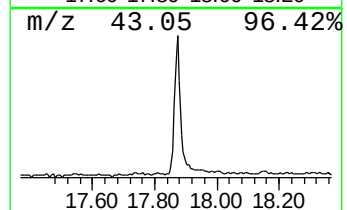
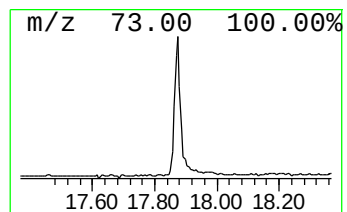
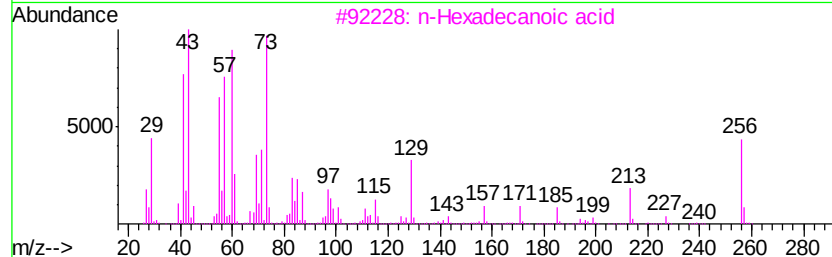
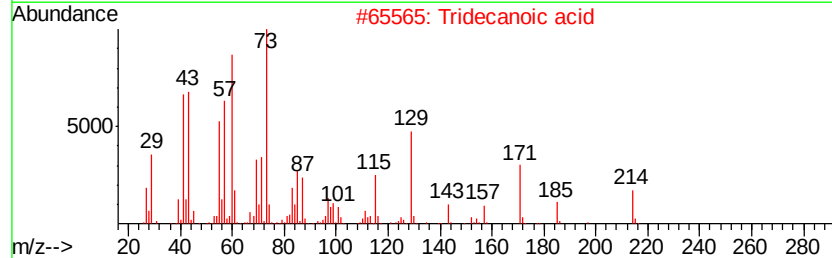
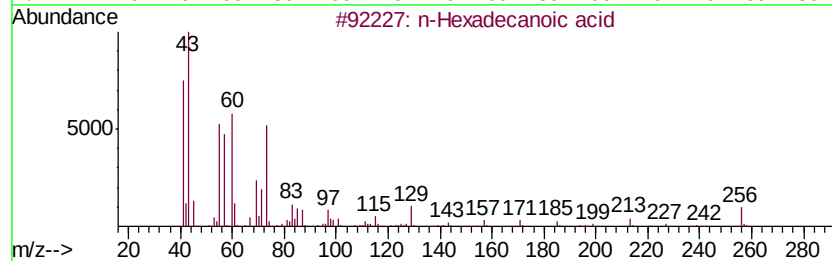
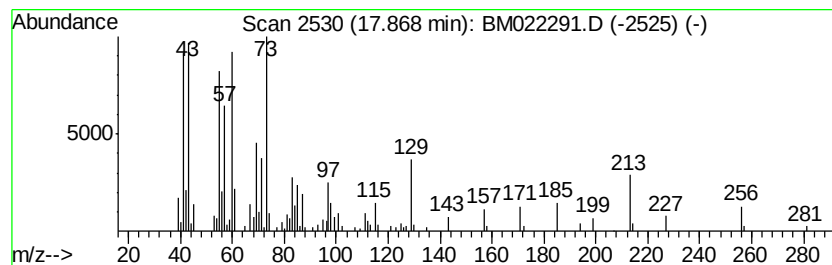
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM082219MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD		R.T.	
17.87	7.95 ng/ul	340991	Phenanthrene-d10		16.97	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98	
2	Tridecanoic acid	214	C13H26O2	000638-53-9	95	
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93	
4	Tridecanoic acid	214	C13H26O2	000638-53-9	89	
5	Tetradecanoic acid	228	C14H28O2	000544-63-8	72	



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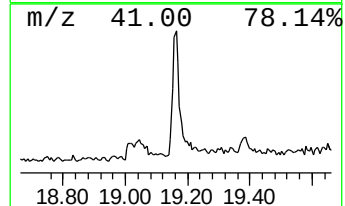
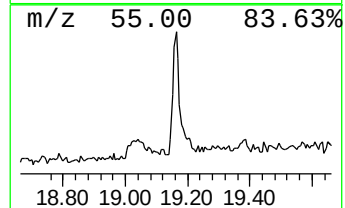
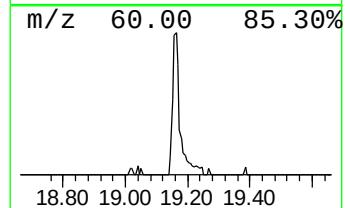
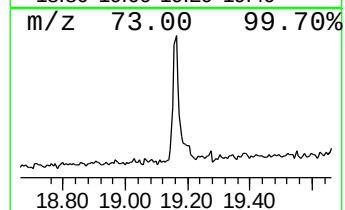
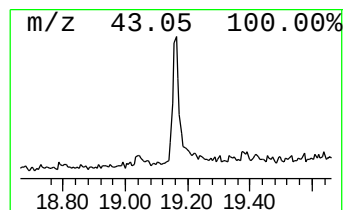
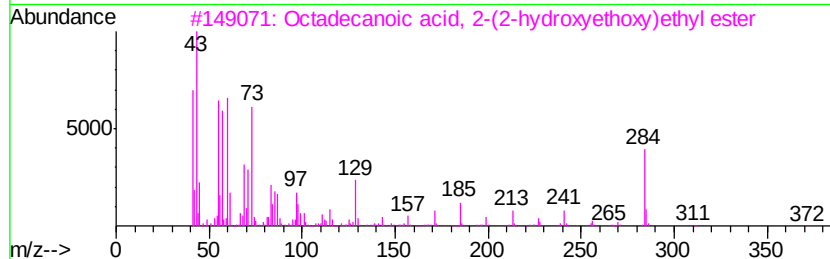
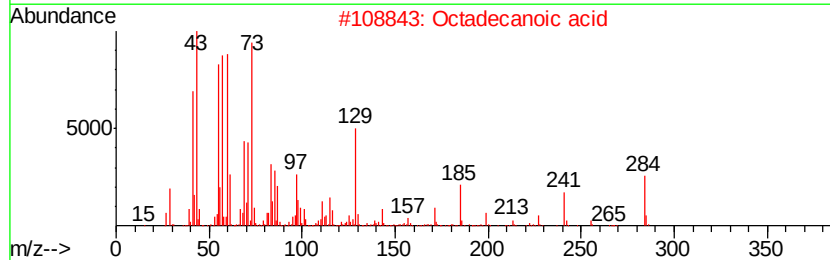
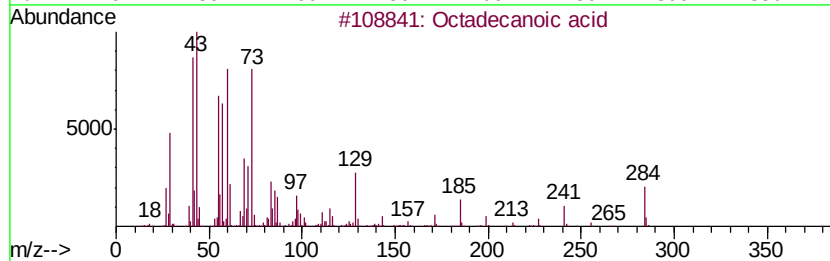
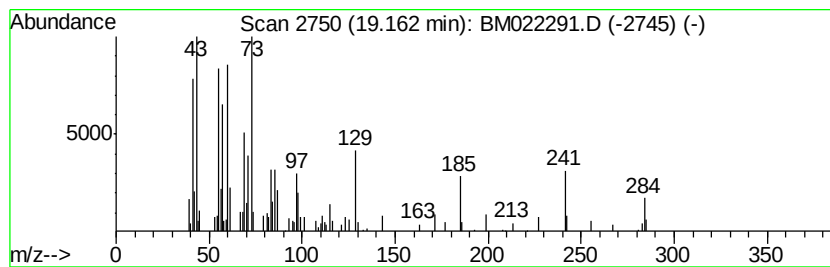
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Octadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.16	2.80 ng/ul	162461	Chrysene-d12	21.19

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	98
2		Octadecanoic acid	284	C18H36O2	000057-11-4	95
3		Octadecanoic acid, 2-(2-hydroxyve...	372	C22H44O4	000106-11-6	91
4		Octadecanoic acid	284	C18H36O2	000057-11-4	90
5		Pentadecanoic acid	242	C15H30O2	001002-84-2	89



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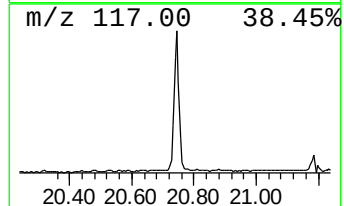
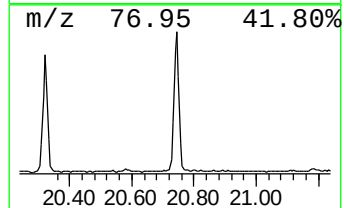
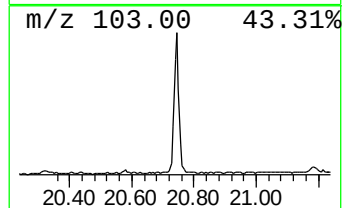
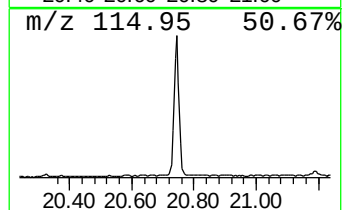
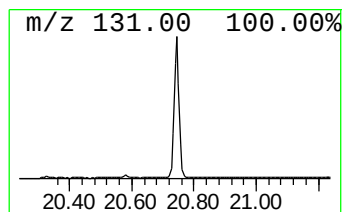
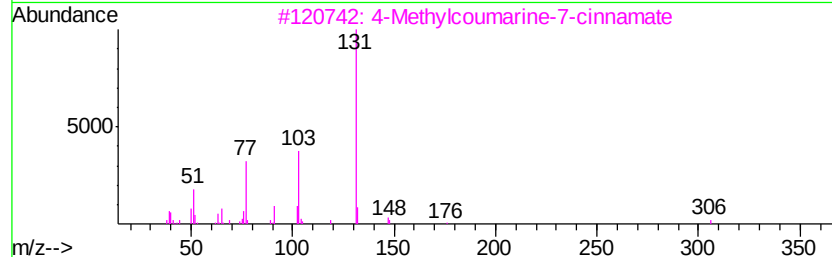
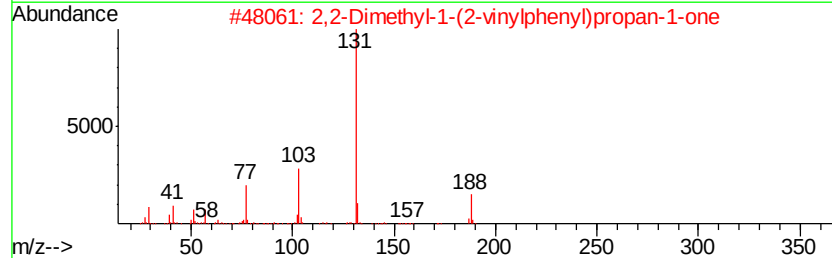
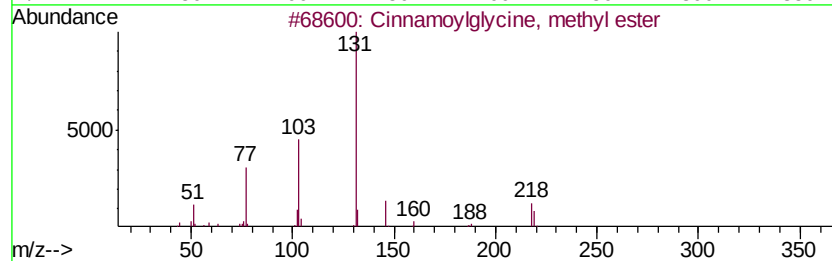
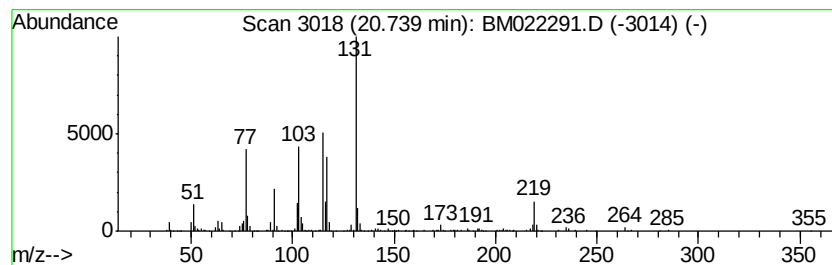
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Cinnamoylglycine, methyl ester Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.		
20.74	8.22 ng/ul	476377	Chrysene-d12	21.19		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cinnamoylglycine, methyl ester	219	C12H13NO3	040778-04-9	50
2		2,2-Dimethyl-1-(2-vinylphenyl)pr...	188	C13H16O	1000210-99-9	47
3		4-Methylcoumarine-7-cinnamate	306	C19H14O4	300829-05-4	47
4		1-Penten-3-one, 1-phenyl-	160	C11H12O	003152-68-9	47
5		N-(2-Methyl-1,3-dioxo-5-isoindol...	306	C18H14N2O3	1000224-44-2	47



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082319\
Data File : BM022291.D
Acq On : 24 Aug 2019 13:22
Operator : HP/JU
Sample : K4242-12
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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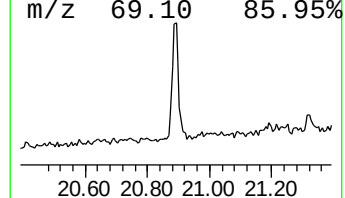
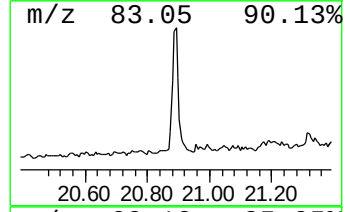
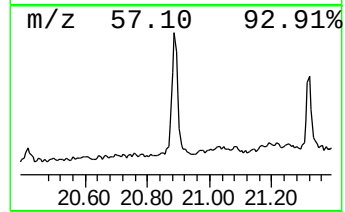
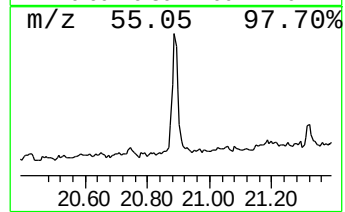
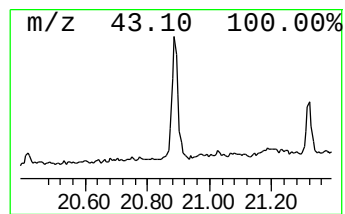
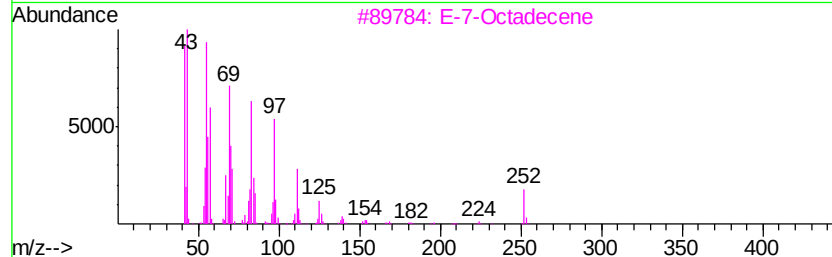
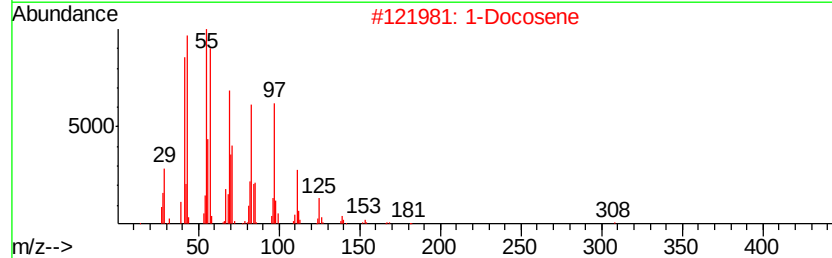
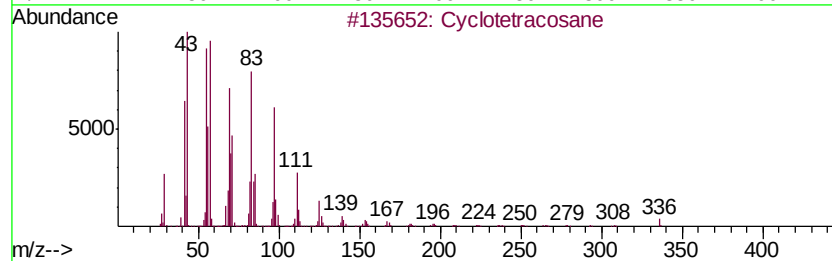
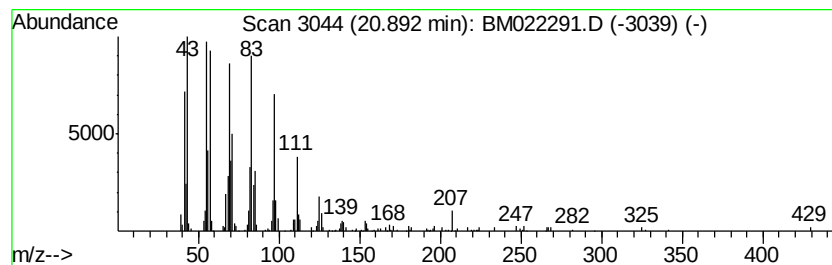
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 (DEL) Alkane: Cyclic20.89 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.89	5.19 ng/ul	300737	Chrysene-d12	21.19

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclotetracosane	336	C24H48	000297-03-0	94
2		1-Docosene	308	C22H44	001599-67-3	93
3		E-7-Octadecene	252	C18H36	1000130-92-0	93
4		Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	91
5		1-Tricosene	322	C23H46	018835-32-0	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082319\
Data File : BM022291.D
Acq On : 24 Aug 2019 13:22
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Sample : K4242-12
Misc :
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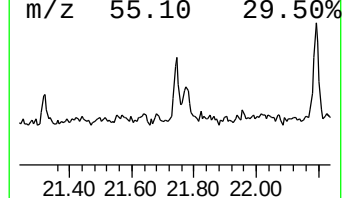
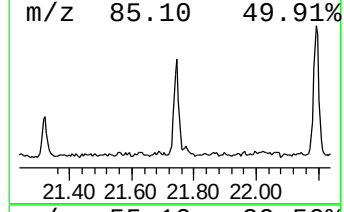
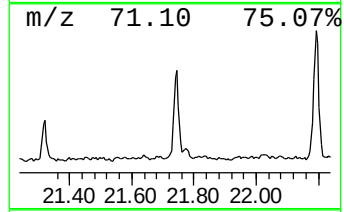
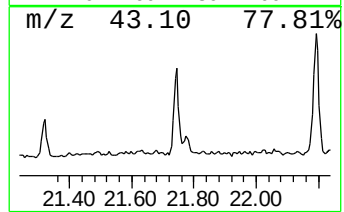
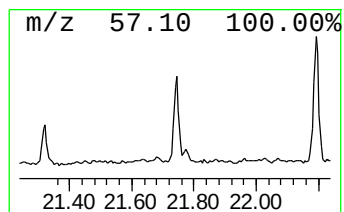
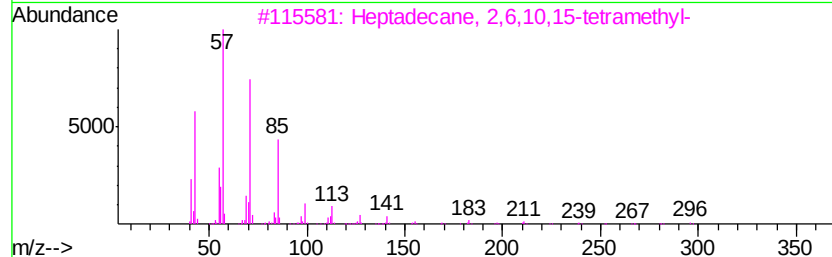
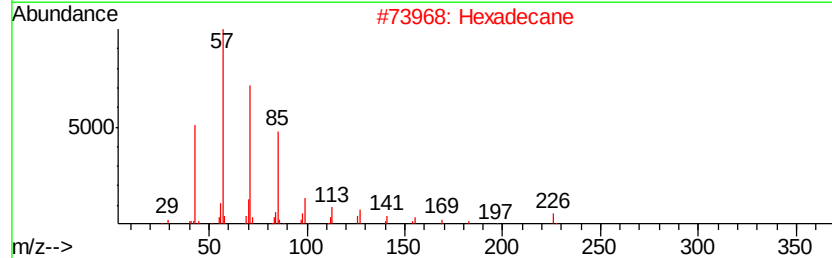
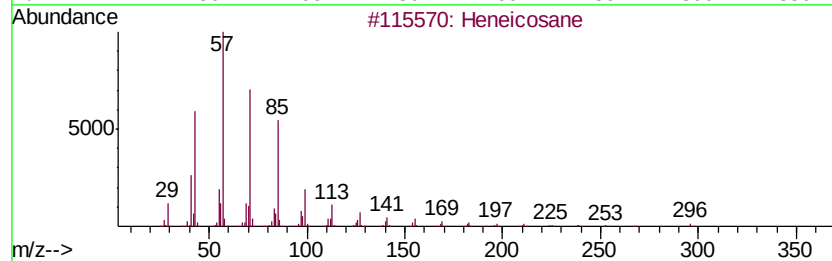
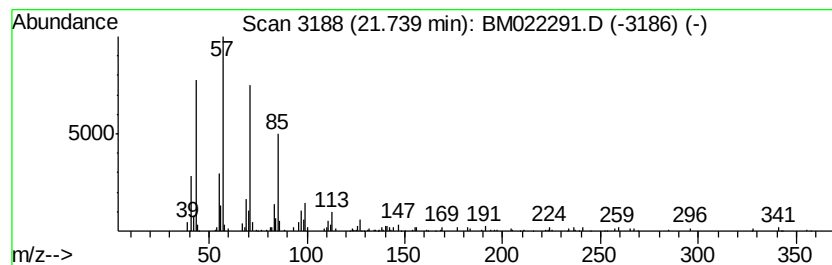
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.74	2.27 ng/ul	131469	Chrysene-d12	21.19

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	98
2	Hexadecane		226	C16H34	000544-76-3	93
3	Heptadecane, 2,6,10,15-tetramethyl-		296	C21H44	054833-48-6	93
4	Octacosane		394	C28H58	000630-02-4	90
5	Hexadecane		226	C16H34	000544-76-3	89



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082319\
Data File : BM022291.D
Acq On : 24 Aug 2019 13:22
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Sample : K4242-12
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument :
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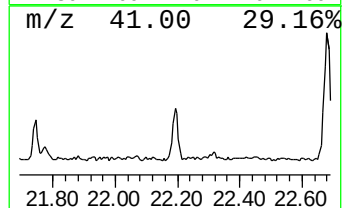
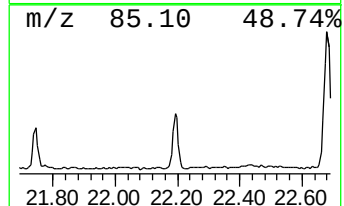
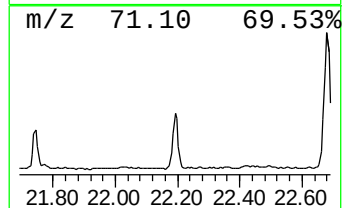
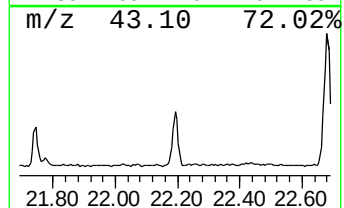
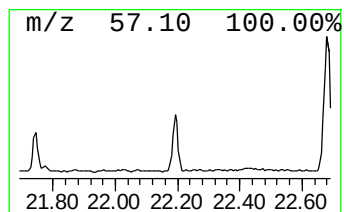
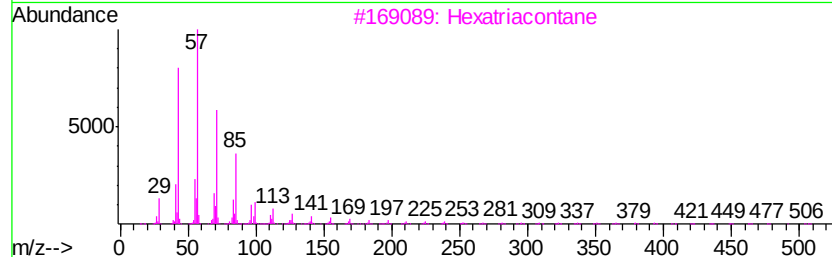
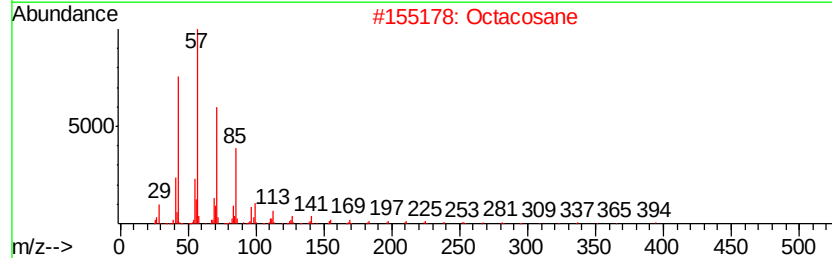
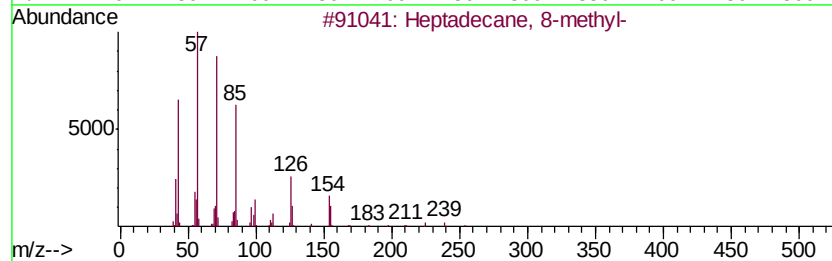
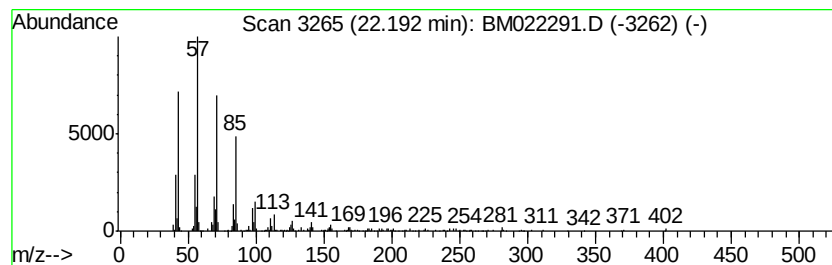
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.19	4.60 ng/ul	266740	Chrysene-d12	21.19

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 8-methyl-	254	C18H38	013287-23-5	93
2		Octacosane	394	C28H58	000630-02-4	90
3		Hexatriacontane	507	C36H74	000630-06-8	90
4		Tetracosane	338	C24H50	000646-31-1	87
5		Heneicosane, 11-decyl-	437	C31H64	055320-06-4	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082319\
Data File : BM022291.D
Acq On : 24 Aug 2019 13:22
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Sample : K4242-12
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument :
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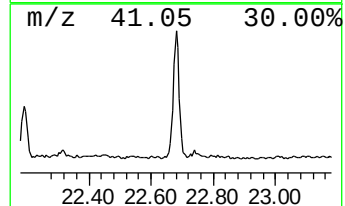
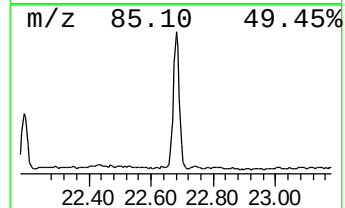
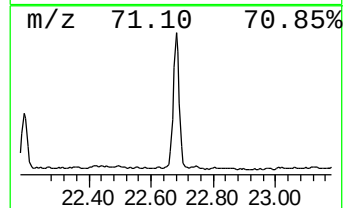
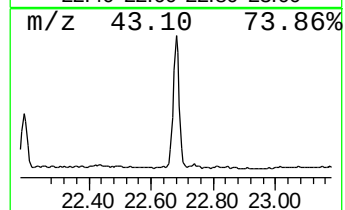
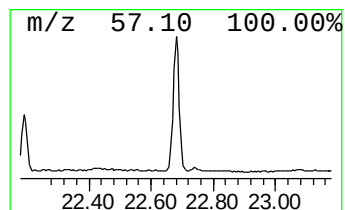
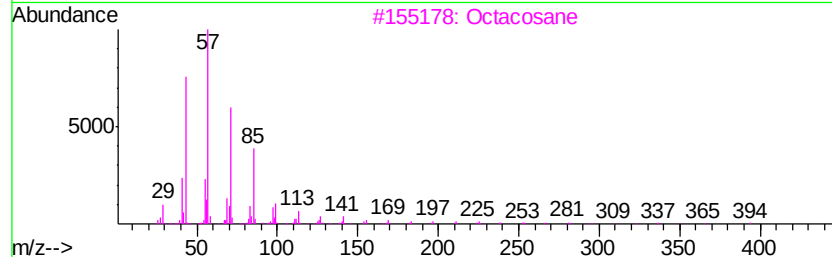
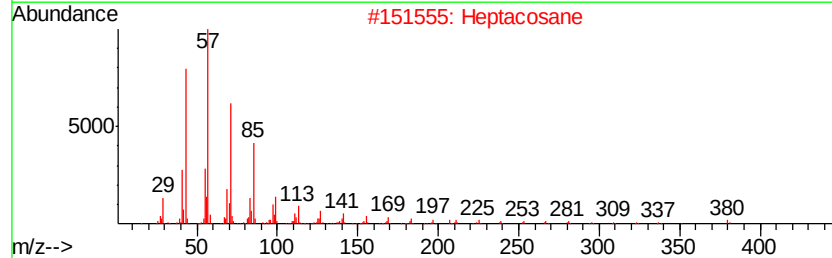
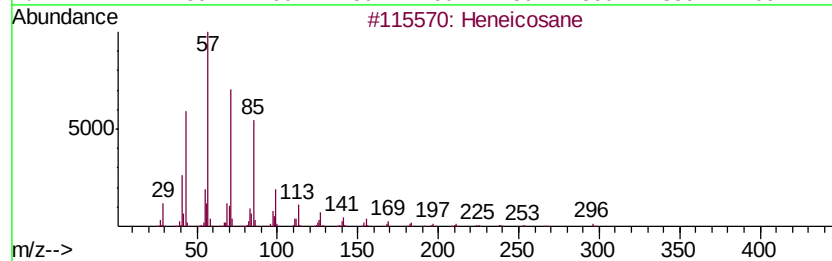
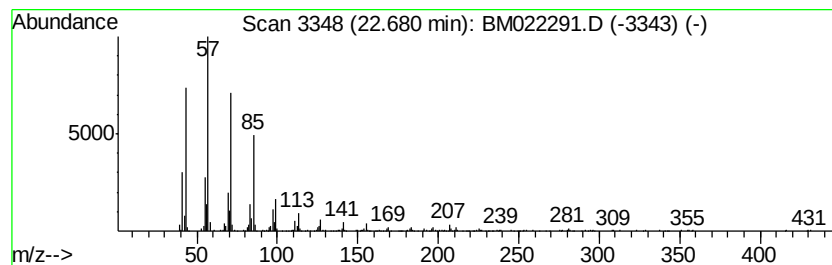
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.68	16.98 ng/ul	727996	Perylene-d12	23.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	97
2		Heptacosane	380	C27H56	000593-49-7	91
3		Octacosane	394	C28H58	000630-02-4	90
4		Tetracosane, 11-decyl-	479	C34H70	055429-84-0	90
5		Tetratetracontane	619	C44H90	007098-22-8	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082319\
Data File : BM022291.D
Acq On : 24 Aug 2019 13:22
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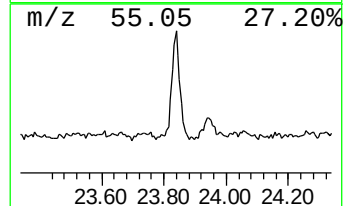
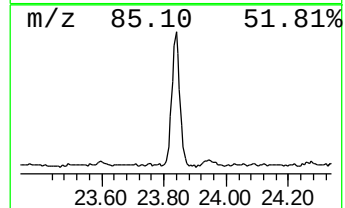
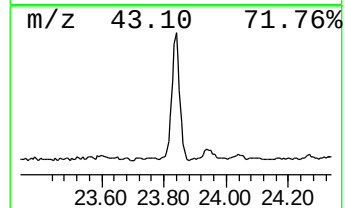
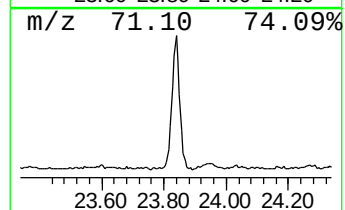
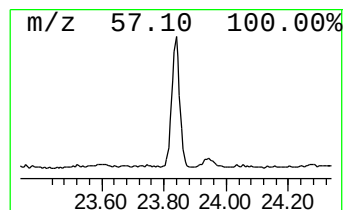
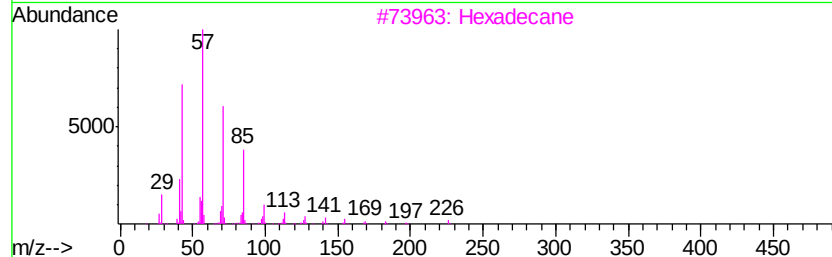
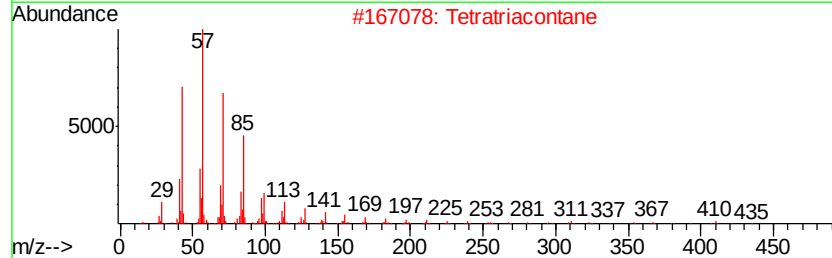
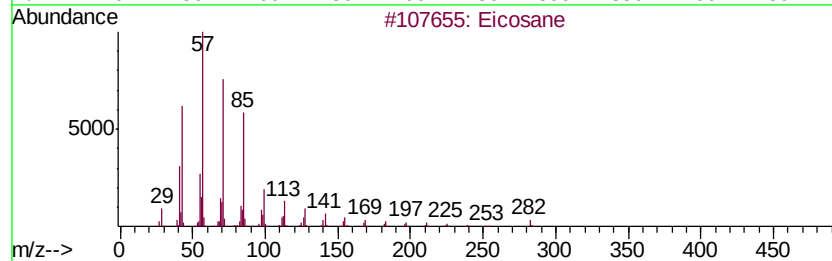
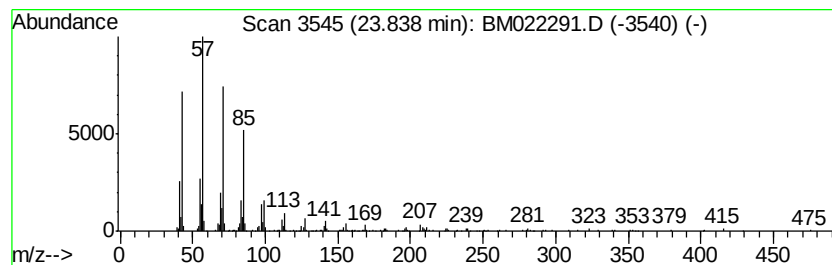
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.84	12.93 ng/ul	554051	Perylene-d12	23.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	95
2		Tetratriacontane	479	C34H70	014167-59-0	94
3		Hexadecane	226	C16H34	000544-76-3	91
4		Heptacosane	380	C27H56	000593-49-7	91
5		Hexadecane	226	C16H34	000544-76-3	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM082319\
Data File : BM022291.D
Acq On : 24 Aug 2019 13:22
Operator : HP/JU
Sample : K4242-12
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1H-Cyclopenta[1,3...	14.15	7.1	ng/ul	239307	3	14.21	675518	20.0
n-Hexadecanoic acid	17.87	8.0	ng/ul	340991	4	16.97	858181	20.0
Octadecanoic acid	19.16	2.8	ng/ul	162461	5	21.19	1158580	20.0
Cinnamoylglycine,...	20.74	8.2	ng/ul	476377	5	21.19	1158580	20.0
(DEL) Alkane: Cyc...	20.89	5.2	ng/ul	300737	5	21.19	1158580	20.0
(DEL) Alkane: Str...	21.74	2.3	ng/ul	131469	5	21.19	1158580	20.0
(DEL) Alkane: Str...	22.19	4.6	ng/ul	266740	5	21.19	1158580	20.0
(DEL) Alkane: Str...	22.68	17.0	ng/ul	727996	6	23.38	857308	20.0
(DEL) Alkane: Str...	23.84	12.9	ng/ul	554051	6	23.38	857308	20.0