

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM111219\
 Data File : BM023668.D
 Acq On : 12 Nov 2019 17:14
 Operator : JU
 Sample : K5565-02
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 JLNA1

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-BM111119MA.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	4.116	206	209	222	rVB	158434	251250	2.35%	0.157%
2	4.704	304	309	322	rBV2	198250	326919	3.06%	0.205%
3	6.728	645	653	666	rBV	1262332	2156733	20.19%	1.349%
4	6.886	674	680	692	rBV	950412	1571755	14.71%	0.983%
5	7.081	706	713	725	rVB	1309299	2050171	19.19%	1.283%
6	7.545	783	792	799	rBV	1747484	2847308	26.65%	1.781%
7	8.257	906	913	928	rBV	1648404	2636346	24.67%	1.649%
8	8.692	979	987	998	rBV	1217514	2008916	18.80%	1.257%
9	9.410	1102	1109	1117	rBV	1018655	1658029	15.52%	1.037%
10	9.945	1193	1200	1212	rBV	1794390	3015710	28.22%	1.887%
11	10.316	1256	1263	1269	rVV	2693402	4585330	42.91%	2.868%
12	10.369	1269	1272	1278	rVB	411245	622805	5.83%	0.390%
13	10.463	1281	1288	1300	rBV	1179448	2168012	20.29%	1.356%
14	11.992	1540	1548	1555	rVB	129896	221160	2.07%	0.138%
15	12.216	1579	1586	1593	rBV	112166	185289	1.73%	0.116%
16	13.039	1719	1726	1733	rVV2	78835	165067	1.54%	0.103%
17	13.374	1775	1783	1792	rBV4	73553	214840	2.01%	0.134%
18	13.521	1802	1808	1813	rVV3	153896	291447	2.73%	0.182%
19	13.621	1819	1825	1829	rVV	3476009	4916788	46.02%	3.076%
20	13.668	1829	1833	1838	rVV	1441482	2001310	18.73%	1.252%
21	13.886	1863	1870	1882	rBV	3906324	5831358	54.58%	3.648%
22	14.127	1907	1911	1915	rBV	122419	164375	1.54%	0.103%
23	14.198	1915	1923	1929	rVV	4682975	6978288	65.31%	4.365%
24	14.262	1929	1934	1942	rVV	435008	672440	6.29%	0.421%
25	14.415	1954	1960	1971	rVV	974128	1548007	14.49%	0.968%
26	14.604	1987	1992	1994	rVV	329667	547791	5.13%	0.343%
27	14.633	1994	1997	2002	rVV	456640	684135	6.40%	0.428%
28	14.686	2002	2006	2010	rVV	105861	160730	1.50%	0.101%
29	14.827	2025	2030	2035	rVV4	71207	135463	1.27%	0.085%
30	14.998	2052	2059	2062	rBV3	65997	127174	1.19%	0.080%
31	15.086	2070	2074	2079	rVV	165430	207960	1.95%	0.130%
32	15.198	2086	2093	2098	rVV	5329699	7637397	71.48%	4.778%
33	15.251	2098	2102	2108	rVV	729150	1075055	10.06%	0.673%
34	15.327	2108	2115	2121	rVV	1326678	1955351	18.30%	1.223%

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Integrator: RTE
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35	15.374	2121	2123	2127	rVV	105622	168568	1.58%	0.105%
36	15.415	2127	2130	2137	rVV3	151562	346904	3.25%	0.217%
37	15.498	2137	2144	2148	rVV5	138356	300913	2.82%	0.188%
38	15.551	2148	2153	2163	rVB	198278	370781	3.47%	0.232%
39	15.715	2174	2181	2186	rVV2	305395	634809	5.94%	0.397%
40	15.786	2186	2193	2202	rVB5	69664	169115	1.58%	0.106%
41	15.951	2212	2221	2223	rBV3	211283	355469	3.33%	0.222%
42	15.974	2223	2225	2233	rVB2	169150	238770	2.23%	0.149%
43	16.233	2263	2269	2271	rBV2	165906	273577	2.56%	0.171%
44	16.262	2271	2274	2278	rVV2	179394	279228	2.61%	0.175%
45	16.309	2278	2282	2287	rVV2	113191	186365	1.74%	0.117%
46	16.409	2296	2299	2304	rVB	100774	142940	1.34%	0.089%
47	16.533	2316	2320	2323	rVB3	94853	127646	1.19%	0.080%
48	16.609	2330	2333	2342	rVB3	63076	128761	1.21%	0.081%
49	16.692	2342	2347	2352	rBV3	127193	294158	2.75%	0.184%
50	16.745	2352	2356	2357	rBV2	180607	210248	1.97%	0.132%
51	16.768	2357	2360	2363	rVB	154721	192451	1.80%	0.120%
52	16.951	2385	2391	2394	rBV	5618399	8019775	75.06%	5.017%
53	16.992	2394	2398	2402	rVV	4299816	5819553	54.47%	3.641%
54	17.051	2402	2408	2412	rVV	5822003	8545767	79.98%	5.346%
55	17.080	2412	2413	2419	rVB	761734	662734	6.20%	0.415%
56	17.156	2420	2426	2431	rBV5	132151	248615	2.33%	0.156%
57	17.356	2455	2460	2464	rBV	237647	338296	3.17%	0.212%
58	17.398	2464	2467	2474	rVB2	82694	125071	1.17%	0.078%
59	17.480	2476	2481	2486	rBV2	219705	298828	2.80%	0.187%
60	17.827	2535	2540	2543	rVV	494060	728280	6.82%	0.456%
61	17.874	2543	2548	2555	rVV	2065077	2857948	26.75%	1.788%
62	17.950	2557	2561	2566	rVV2	297050	587384	5.50%	0.367%
63	18.015	2567	2572	2576	rVV2	806865	1421759	13.31%	0.889%
64	18.056	2576	2579	2587	rVB	366723	543963	5.09%	0.340%
65	18.174	2596	2599	2604	rVB	197669	253585	2.37%	0.159%
66	18.333	2621	2626	2630	rBV	397070	517467	4.84%	0.324%
67	18.580	2665	2668	2672	rVB2	185562	246754	2.31%	0.154%
68	18.633	2673	2677	2681	rBV2	126938	199652	1.87%	0.125%
69	18.750	2693	2697	2703	rBV2	240929	455403	4.26%	0.285%
70	18.815	2704	2708	2712	rBV4	296199	457437	4.28%	0.286%
71	18.892	2717	2721	2729	rBV4	115258	285177	2.67%	0.178%

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72	19.015	2737	2742	2749	rVB	4289272	5694119	53.29%	3.562%
73	19.086	2749	2754	2758	rBV2	706089	911505	8.53%	0.570%
74	19.168	2763	2768	2773	rBV	717954	1069461	10.01%	0.669%
75	19.350	2793	2799	2802	rBV	7611616	10684788	100.00%	6.684%
76	19.380	2802	2804	2811	rVV	3079003	3668456	34.33%	2.295%
77	19.456	2814	2817	2825	rVB2	200660	387534	3.63%	0.242%
78	19.562	2831	2835	2838	rBV	226323	325660	3.05%	0.204%
79	19.709	2855	2860	2867	rBV3	247521	614287	5.75%	0.384%
80	19.809	2873	2877	2880	rBV	155070	208387	1.95%	0.130%
81	19.850	2880	2884	2885	rVV2	163250	217319	2.03%	0.136%
82	19.880	2885	2889	2895	rVB	692384	1001743	9.38%	0.627%
83	19.939	2896	2899	2902	rBV	162827	176493	1.65%	0.110%
84	19.980	2903	2906	2910	rVV	623944	743645	6.96%	0.465%
85	20.039	2913	2916	2920	rVB2	283764	395195	3.70%	0.247%
86	20.180	2937	2940	2943	rVB	148124	173692	1.63%	0.109%
87	20.227	2945	2948	2950	rBV	180595	251220	2.35%	0.157%
88	20.433	2981	2983	2986	rBV2	156011	152582	1.43%	0.095%
89	20.833	3049	3051	3054	rBV	183410	187675	1.76%	0.117%
90	20.891	3057	3061	3079	rVB3	1452497	2634251	24.65%	1.648%
91	21.091	3092	3095	3098	rBV	529067	541657	5.07%	0.339%
92	21.150	3098	3105	3109	rVV2	6075382	9097487	85.14%	5.691%
93	21.186	3109	3111	3124	rVB	1571345	1850471	17.32%	1.158%
94	21.333	3133	3136	3140	rBV	522390	567399	5.31%	0.355%
95	21.762	3205	3209	3220	rVB2	1002432	1744851	16.33%	1.092%
96	22.209	3281	3285	3289	rBV	1144279	1382399	12.94%	0.865%
97	22.697	3360	3368	3373	rBV2	1720278	3500182	32.76%	2.190%
98	23.174	3444	3449	3454	rBV	3502386	5831721	54.58%	3.648%
99	23.315	3467	3473	3478	rBV	3157546	5596385	52.38%	3.501%
100	23.862	3562	3566	3573	rVB	886523	1585890	14.84%	0.992%

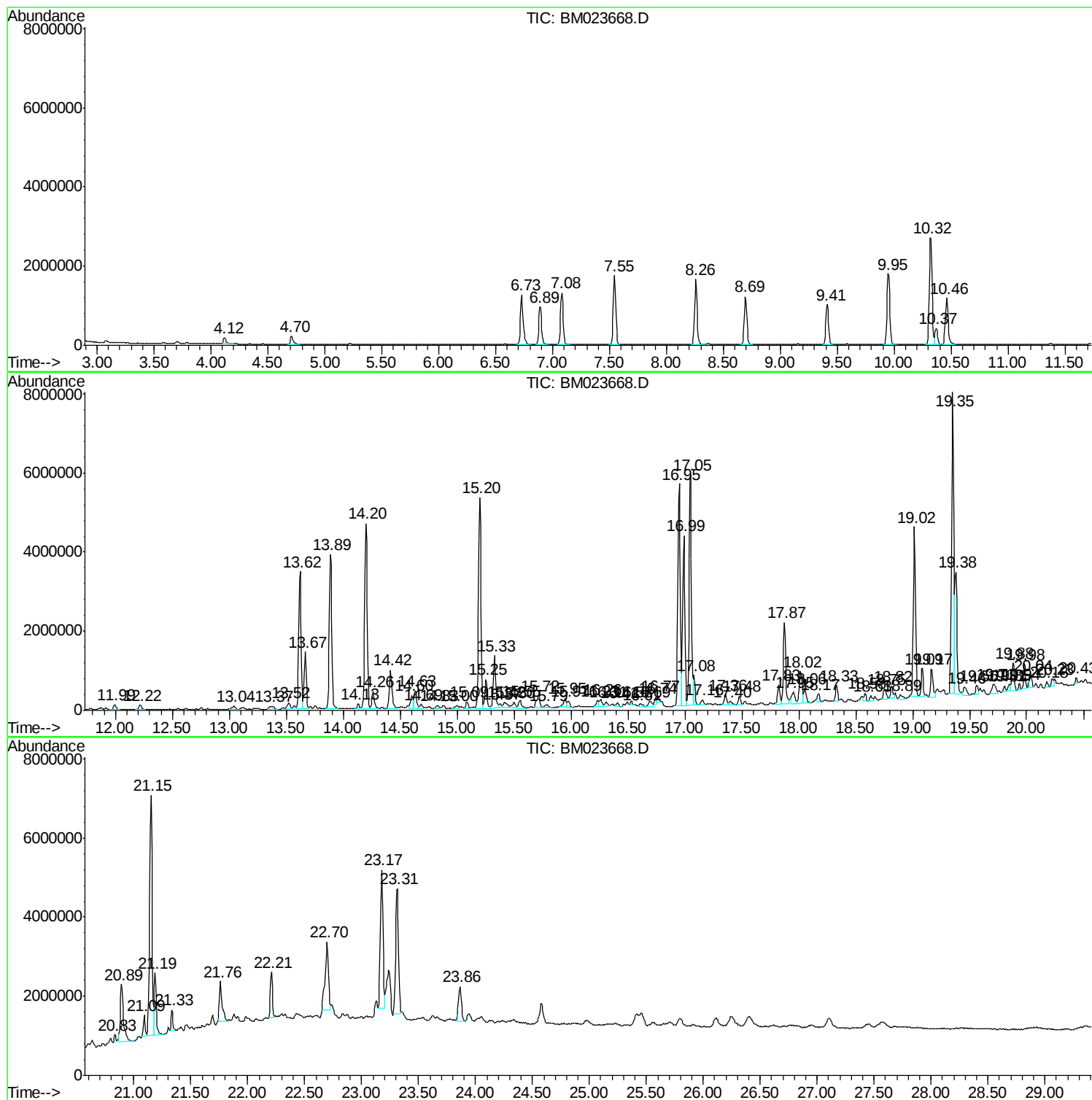
Sum of corrected areas: 159851314

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM11119MA.M
Quant Title : SVOA CALIBRATION

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



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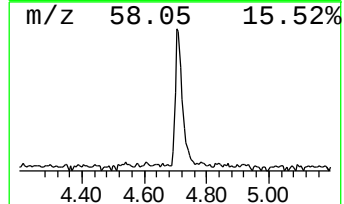
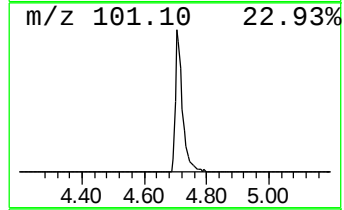
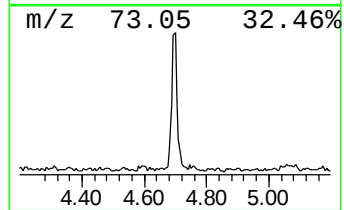
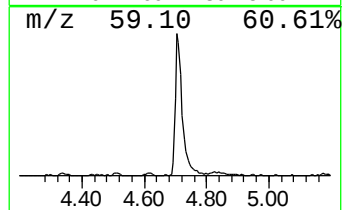
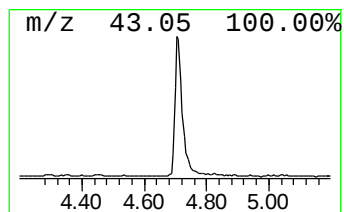
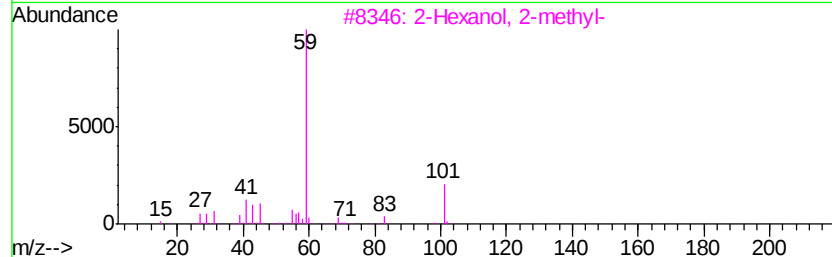
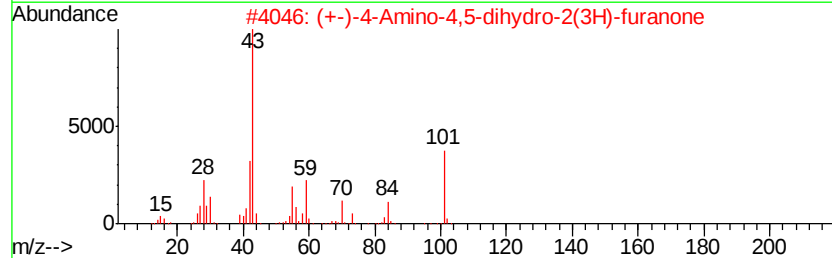
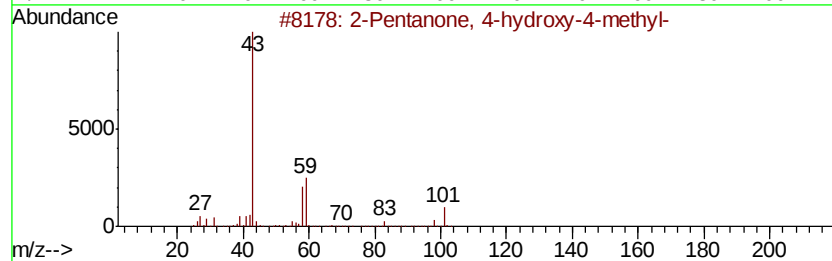
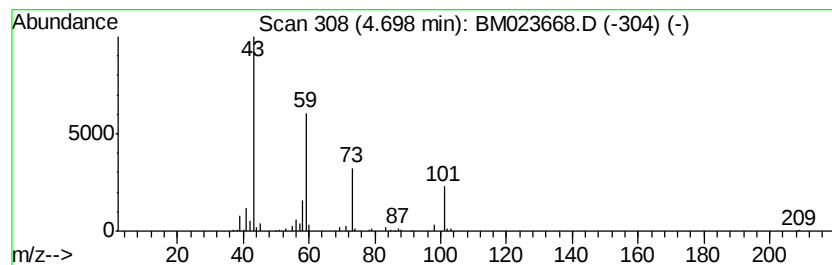
Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM11119MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.70	2.30 ng/ul	326919	1,4-Dichlorobenzene-d4	7.55
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116 C6H12O2	000123-42-2	53
2	(+)-4-Amino-4,5-dihydro-2(3H)-furanone	101 C4H7NO2	016504-58-8	40
3	2-Hexanol, 2-methyl-	116 C7H16O	000625-23-0	40
4	Methyl 2,5,8,11-tetraoxatridecan-1-yl	236 C10H20O6	1000366-78-2	37
5	2-Pentanone, 4-hydroxy-4-methyl-	116 C6H12O2	000123-42-2	36



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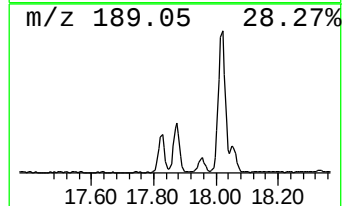
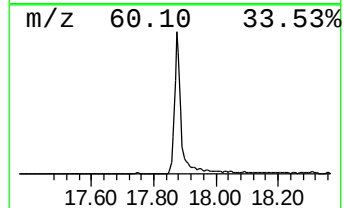
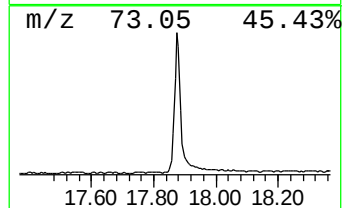
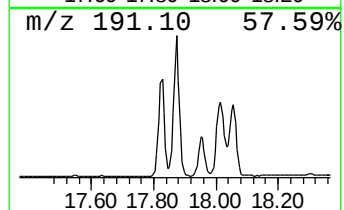
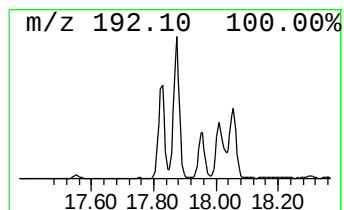
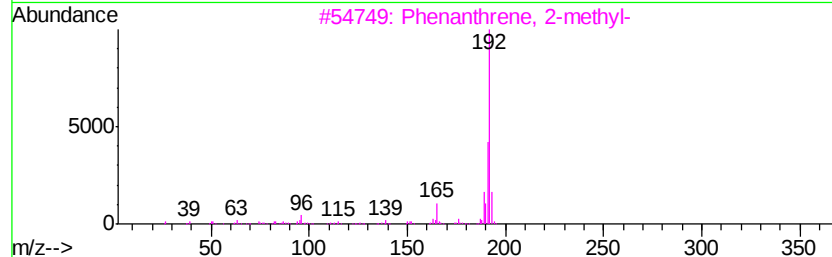
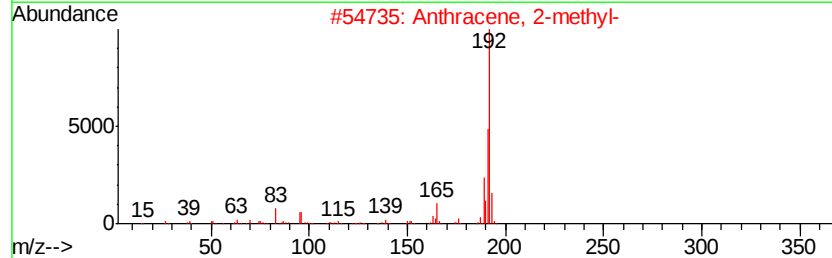
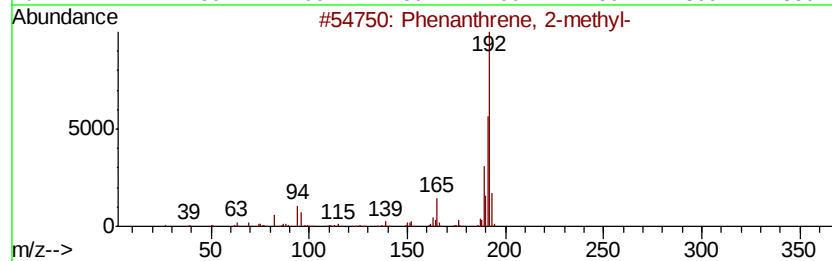
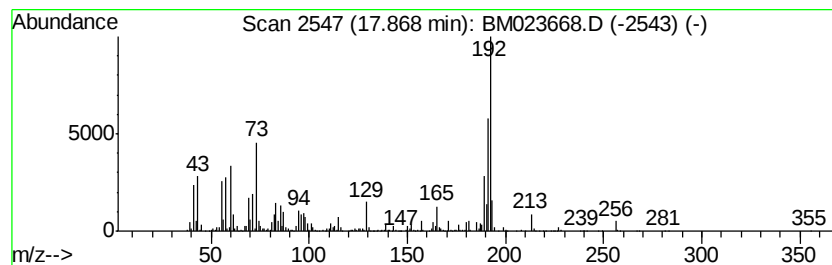
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 Phenanthrene, 2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.87	7.13 ng/ul	2857950	Phenanthrene-d10	16.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
2		Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
3		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
4		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
5		Anthracene, 9-methyl-	192	C15H12	000779-02-2	95



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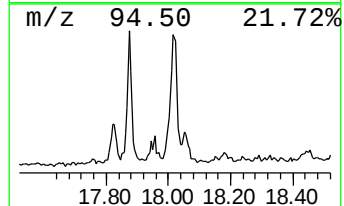
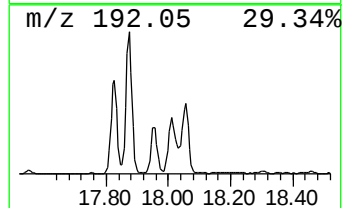
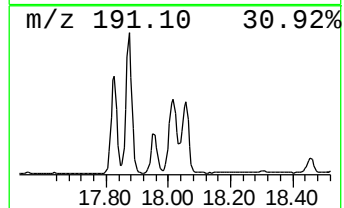
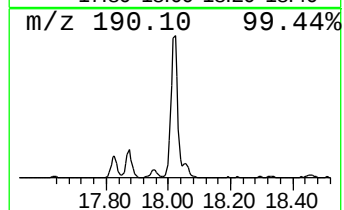
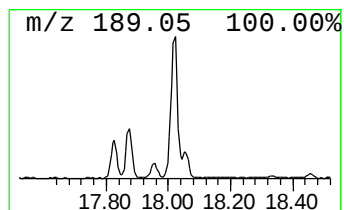
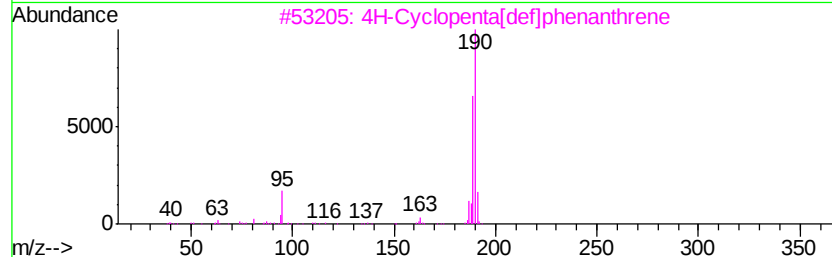
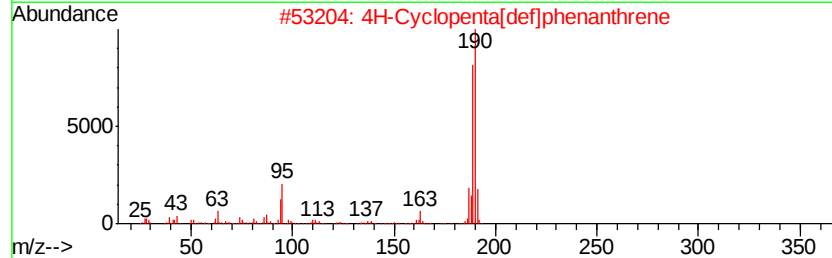
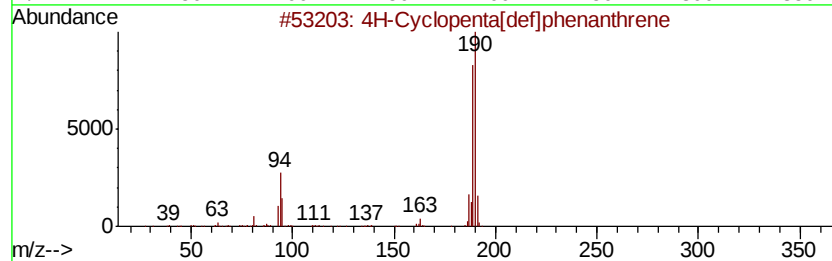
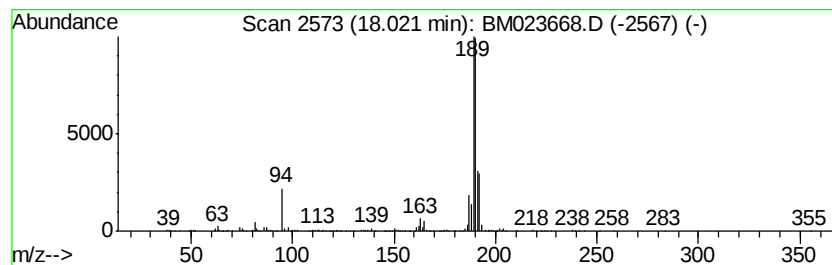
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 4H-Cyclopenta[def]phenanthrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.02	3.55 ng/ul	1421760	Phenanthrene-d10	16.95

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	93	
2	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	64	
3	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	58	
4	6H-Cyclobuta[flk]phenanthrene	190	C15H10	083469-43-6	53	
5	Methyl diselenide	190	C2H6Se2	007101-31-7	43	



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM111219\
Data File : BM023668.D
Acq On : 12 Nov 2019 17:14
Operator : JU
Sample : K5565-02
Misc :
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ClientSampleId :
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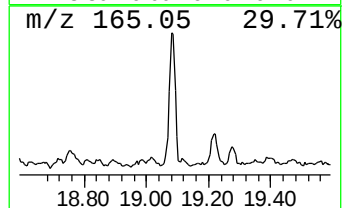
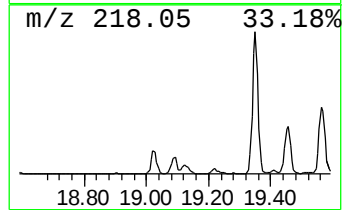
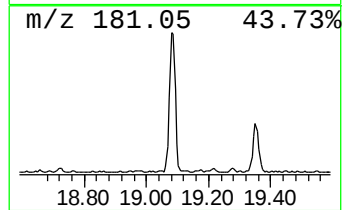
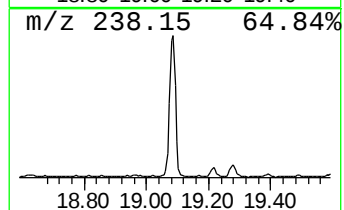
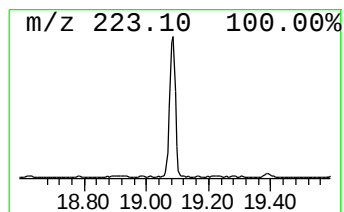
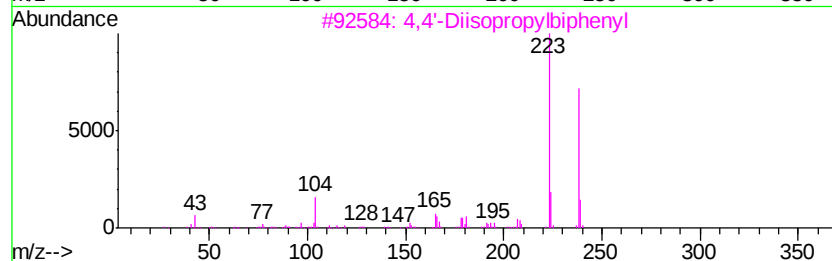
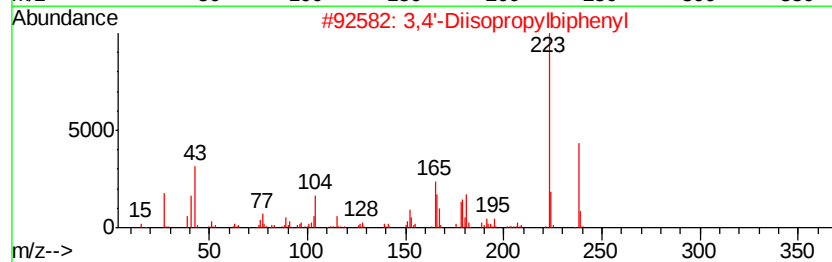
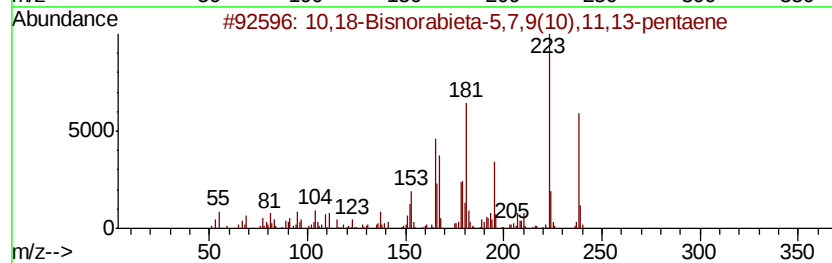
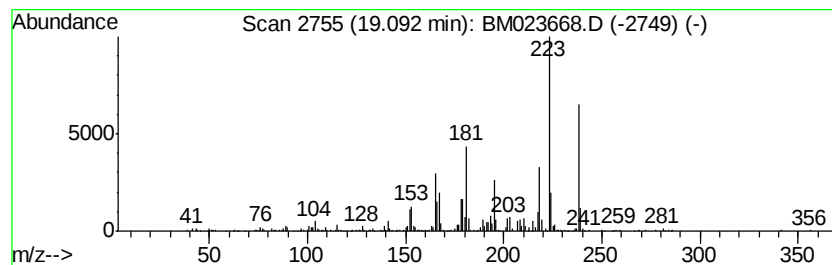
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 10,18-Bisnorabieta-5,7,9(10... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.09	2.00 ng/ul	911505	Chrysene-d12	21.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	10,18-Bisnorabieta-5,7,9(10),11,...	238	C18H22	006566-19-4	99
2		3,4'-Diisopropylbiphenyl	238	C18H22	061434-46-6	86
3		4,4'-Diisopropylbiphenyl	238	C18H22	018970-30-4	60
4		4,4'-Diisopropylbiphenyl	238	C18H22	018970-30-4	49
5		2-Propenoic acid, 3-(3,4,5-trime...	238	C12H14O5	000090-50-6	49



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM111219\
Data File : BM023668.D
Acq On : 12 Nov 2019 17:14
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ClientSampleId :
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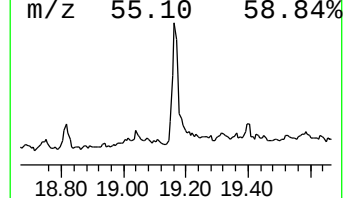
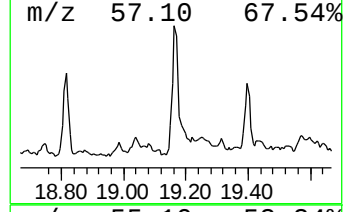
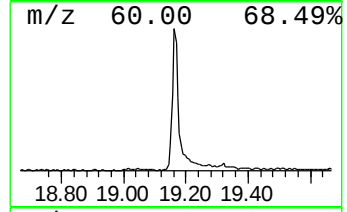
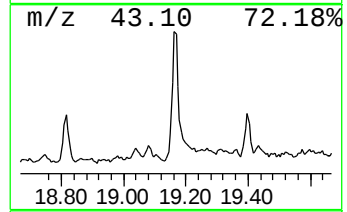
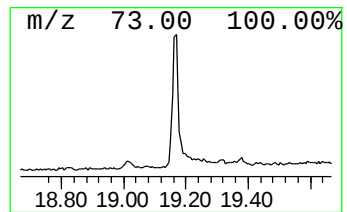
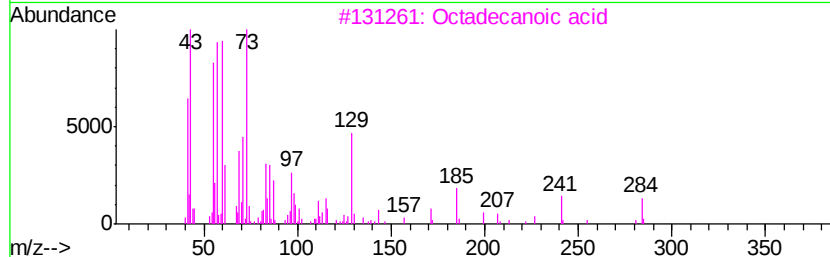
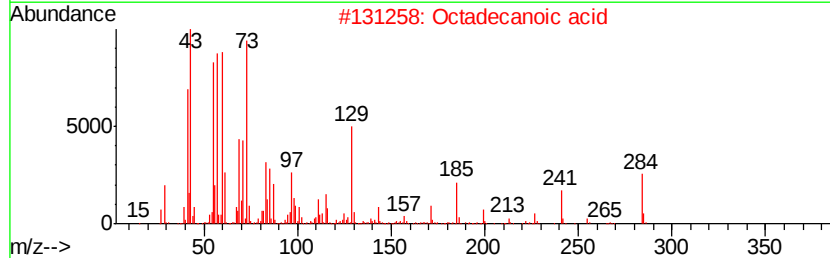
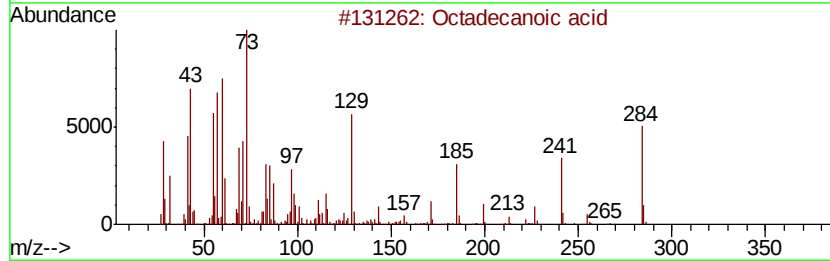
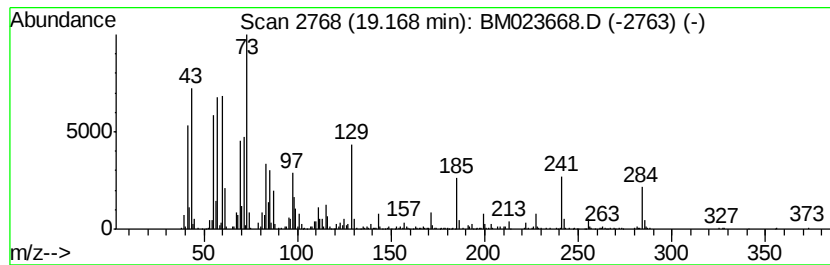
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Octadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.17	2.35 ng/ul	1069460	Chrysene-d12	21.15

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Octadecanoic acid	284	C18H36O2	000057-11-4	98
3			Octadecanoic acid	284	C18H36O2	000057-11-4	95
4			Octadecanoic acid	284	C18H36O2	000057-11-4	94
5			Octadecanoic acid	284	C18H36O2	000057-11-4	89



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM111219\
Data File : BM023668.D
Acq On : 12 Nov 2019 17:14
Operator : JU
Sample : K5565-02
Misc :
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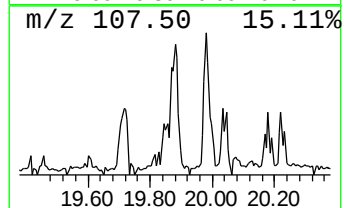
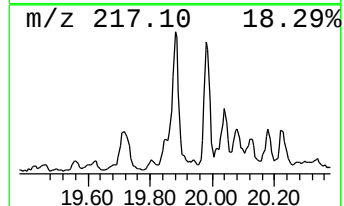
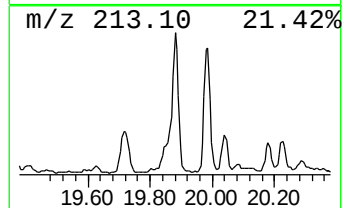
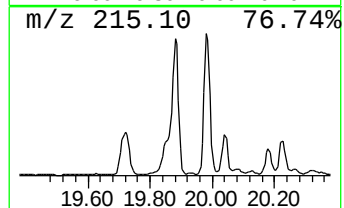
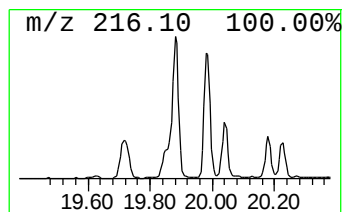
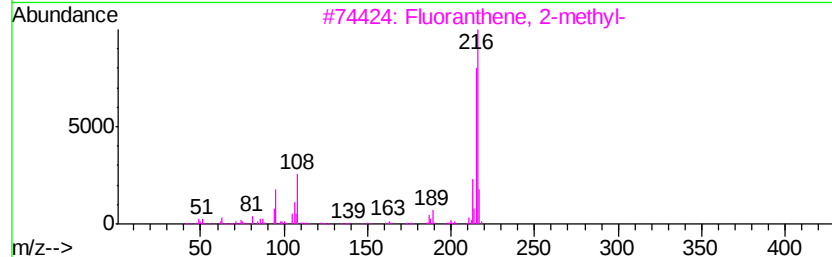
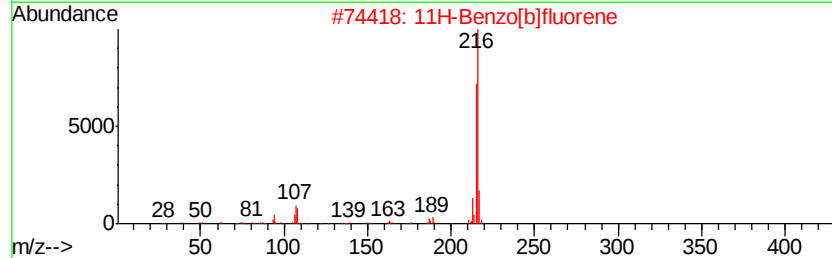
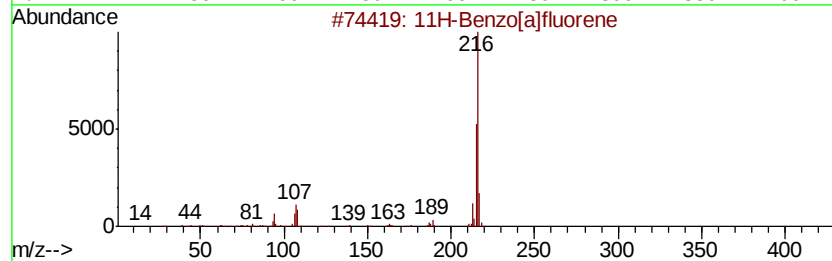
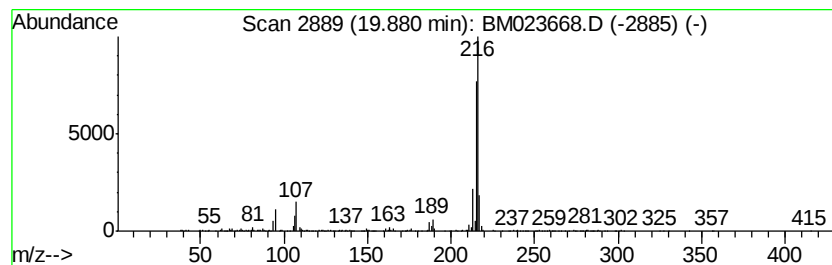
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 11H-Benzo[a]fluorene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.88	2.20 ng/ul	1001740	Chrysene-d12	21.15

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	95
2		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	93
3		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	87
4		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	87
5		Pyrene, 1-methyl-	216	C17H12	002381-21-7	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM111219\
Data File : BM023668.D
Acq On : 12 Nov 2019 17:14
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Misc :
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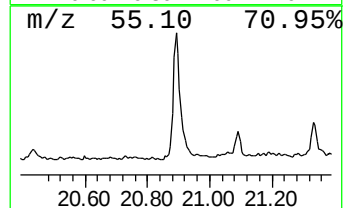
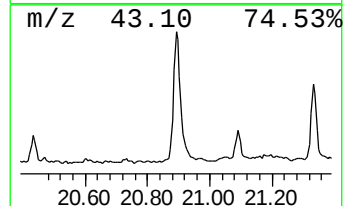
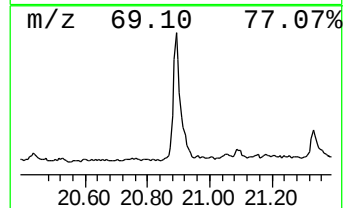
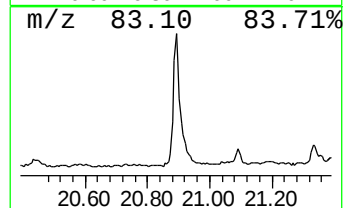
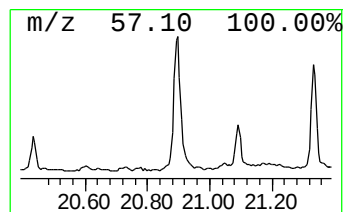
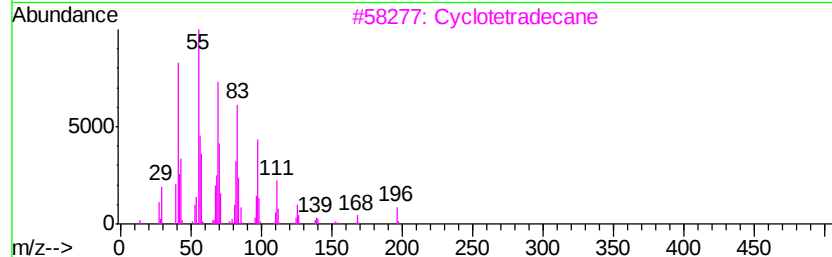
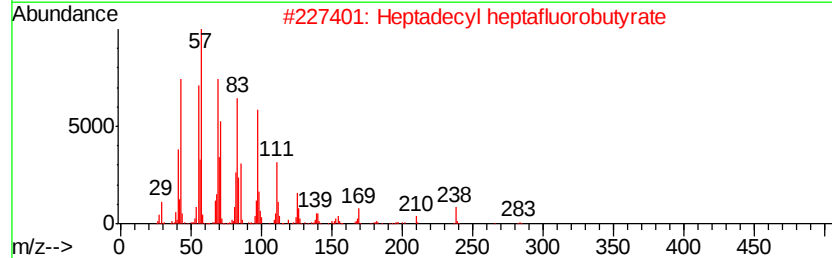
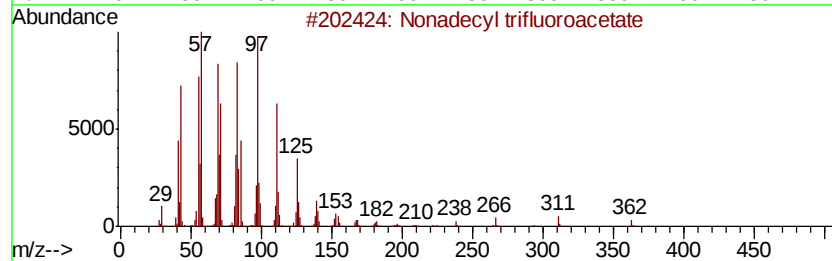
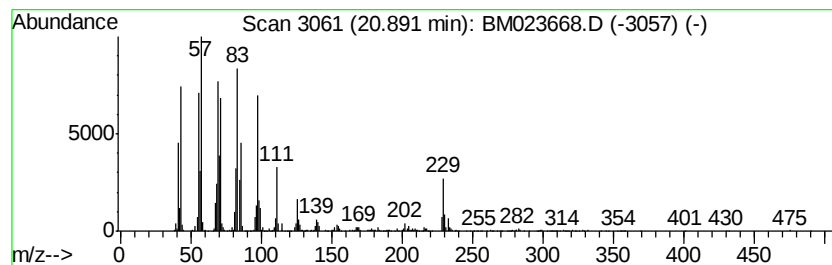
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Nonadecyl trifluoroacetate Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.89	5.79 ng/ul	2634250	Chrysene-d12	21.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecyl trifluoroacetate	380	C21H39F3O2	1000351-76-3	87
2		Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	87
3		Cyclotetradecane	196	C14H28	000295-17-0	86
4		n-Nonadecanol-1	284	C19H40O	001454-84-8	81
5		1-Heneicosanol	312	C21H44O	015594-90-8	76



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM111219\
Data File : BM023668.D
Acq On : 12 Nov 2019 17:14
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Misc :
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ClientSampleId :
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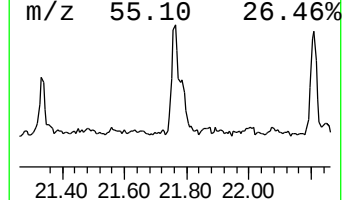
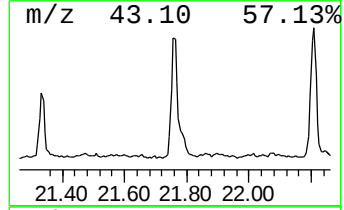
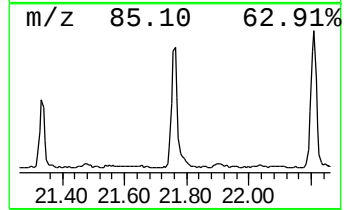
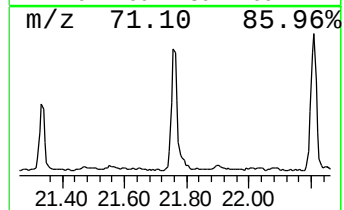
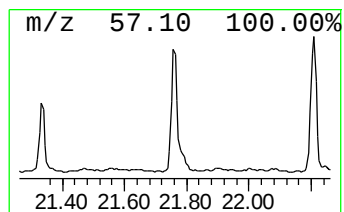
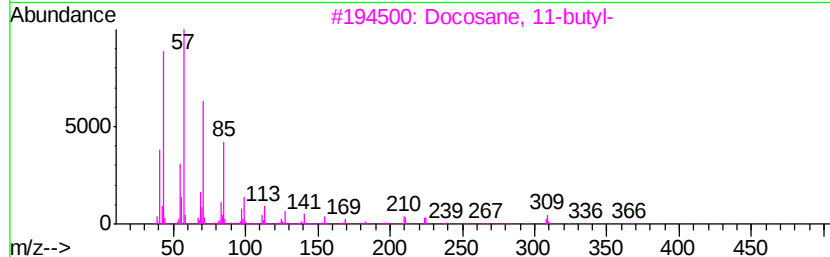
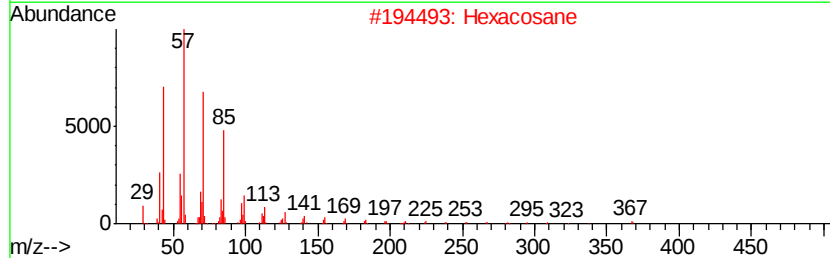
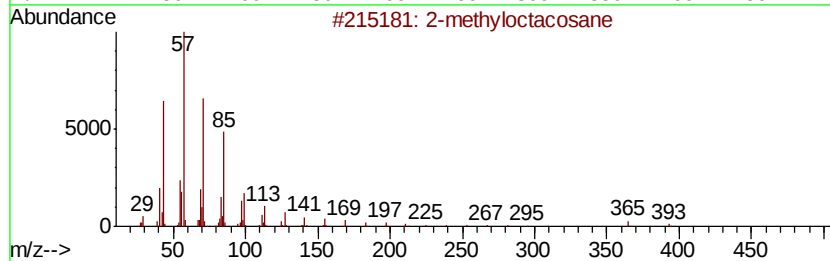
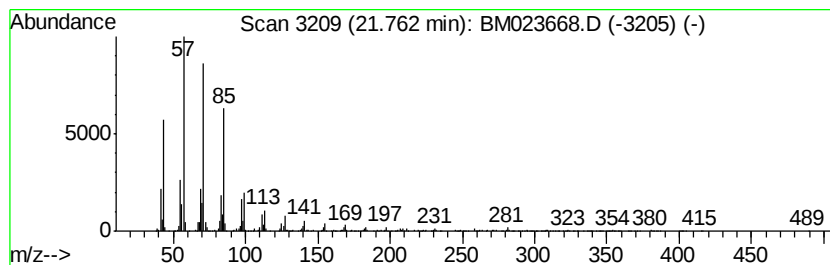
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.76	3.84 ng/ul	1744850	Chrysene-d12	21.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-methyloctacosane	408	C29H60	1000376-72-8	91
2		Hexacosane	366	C26H54	000630-01-3	91
3		Docosane, 11-butyl-	366	C26H54	013475-76-8	90
4		10-Methylnonadecane	282	C20H42	056862-62-5	90
5		Heneicosane	296	C21H44	000629-94-7	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM111219\
Data File : BM023668.D
Acq On : 12 Nov 2019 17:14
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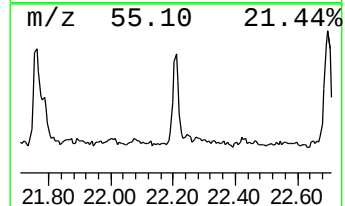
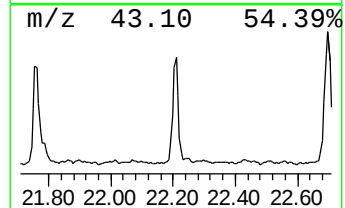
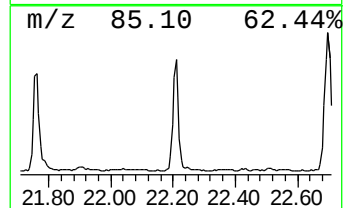
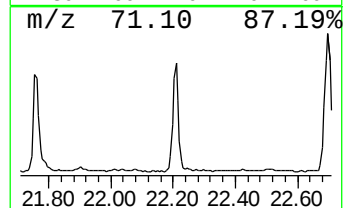
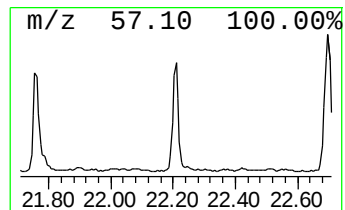
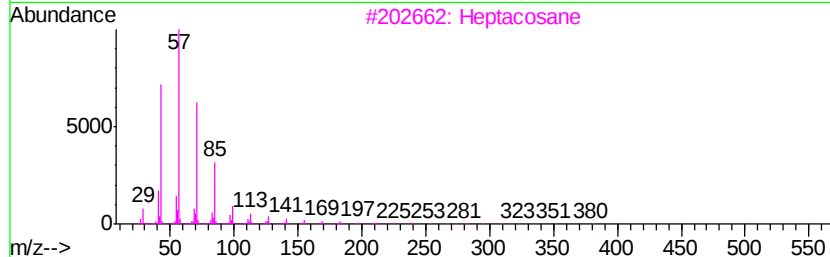
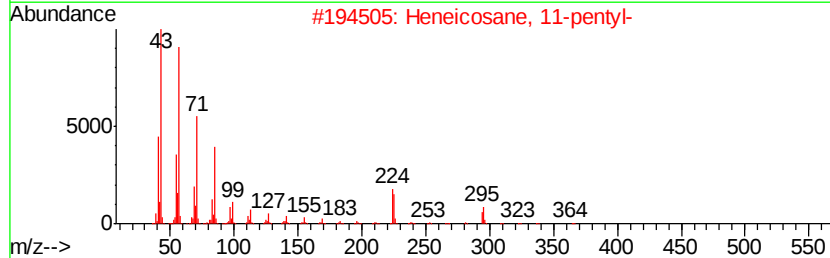
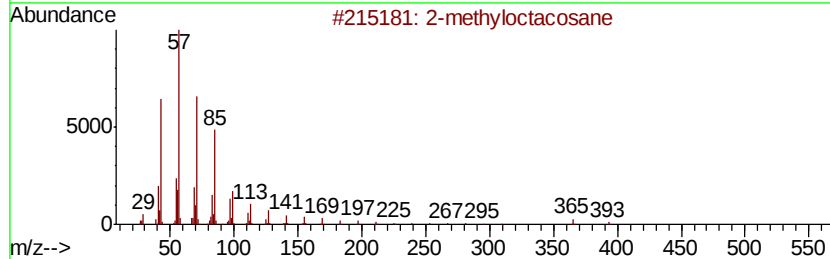
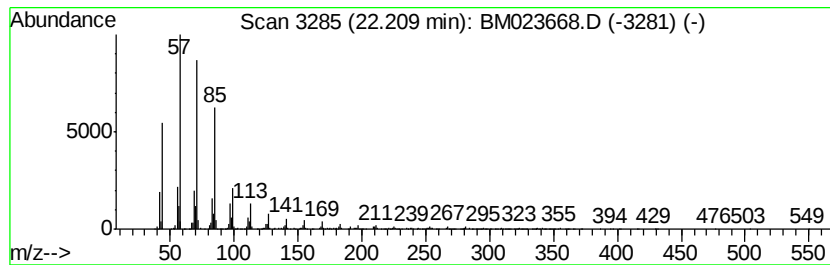
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.21	3.04 ng/ul	1382400	Chrysene-d12	21.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-methyloctacosane	408	C29H60	1000376-72-8	91
2		Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	91
3		Heptacosane	380	C27H56	000593-49-7	91
4		Docosane, 5-butyl-	366	C26H54	055282-16-1	91
5		Octadecane, 2-methyl-	268	C19H40	001560-88-9	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM111219\
Data File : BM023668.D
Acq On : 12 Nov 2019 17:14
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ClientSampleId :
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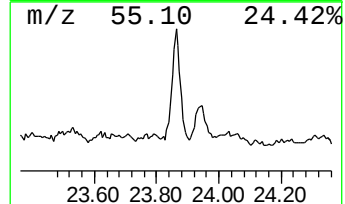
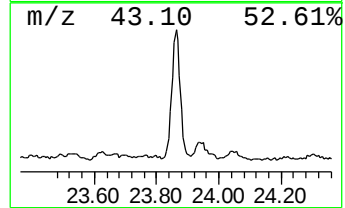
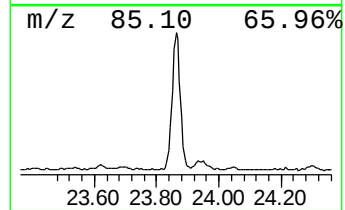
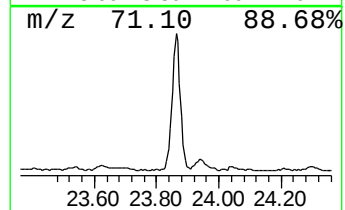
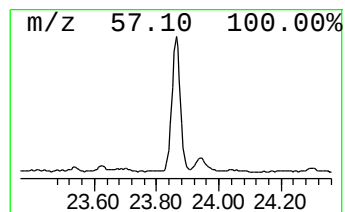
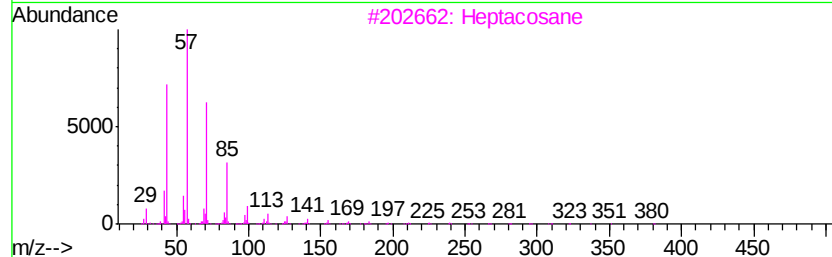
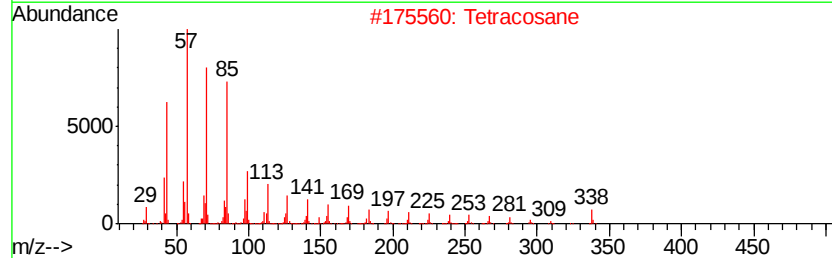
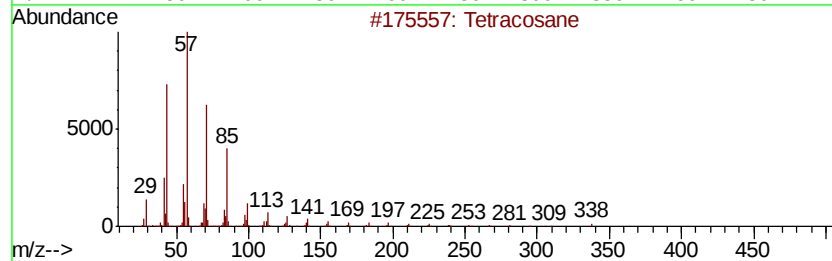
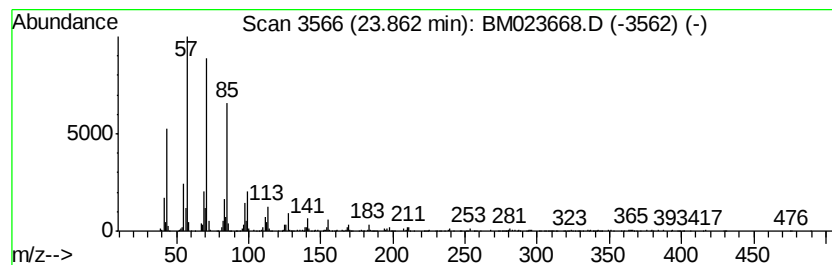
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM11119MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.86	5.67 ng/ul	1585890	Perylene-d12	23.31

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	94
2			Tetracosane	338	C24H50	000646-31-1	94
3			Heptacosane	380	C27H56	000593-49-7	93
4			Tetracosane	338	C24H50	000646-31-1	93
5			Eicosane	282	C20H42	000112-95-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM111219\
Data File : BM023668.D
Acq On : 12 Nov 2019 17:14
Operator : JU
Sample : K5565-02
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
JLNA1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111119MA.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
2-Pentanone, 4-hy...	4.70	2.3	ng/ul	326919	1	7.55	2847310	20.0
Phenanthrene, 2-m...	17.87	7.1	ng/ul	2857950	4	16.95	8019780	20.0
4H-Cyclopenta[def...	18.02	3.5	ng/ul	1421760	4	16.95	8019780	20.0
10,18-Bisnorabiet...	19.09	2.0	ng/ul	911505	5	21.15	9097490	20.0
Octadecanoic acid	19.17	2.4	ng/ul	1069460	5	21.15	9097490	20.0
11H-Benzo[a]fluorene	19.88	2.2	ng/ul	1001740	5	21.15	9097490	20.0
Nonadecyl trifluo...	20.89	5.8	ng/ul	2634250	5	21.15	9097490	20.0
(DEL) Alkane: Str...	21.76	3.8	ng/ul	1744850	5	21.15	9097490	20.0
(DEL) Alkane: Str...	22.21	3.0	ng/ul	1382400	5	21.15	9097490	20.0
(DEL) Alkane: Str...	23.86	5.7	ng/ul	1585890	6	23.31	5596390	20.0