

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102219\
 Data File : BM023335.D
 Acq On : 23 Oct 2019 00:55
 Operator : JU
 Sample : K5354-14
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BFB87

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-BM101419MA.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	3.169	28	31	38	rVB	64935	97207	0.99%	0.067%
2	3.675	114	117	121	rVB2	39491	47393	0.48%	0.033%
3	3.804	135	139	144	rVB	58580	88884	0.91%	0.061%
4	3.887	149	153	158	rBV	72110	99636	1.02%	0.068%
5	6.669	621	626	631	rVV3	26690	47129	0.48%	0.032%
6	6.828	643	653	667	rVB	1028126	1877698	19.20%	1.288%
7	6.987	673	680	688	rBV	759249	1202939	12.30%	0.825%
8	7.181	707	713	723	rVB	1077302	1712596	17.51%	1.175%
9	7.292	729	732	738	rBV6	22698	43533	0.45%	0.030%
10	7.645	785	792	801	rBV	1472644	2287807	23.39%	1.569%
11	8.357	905	913	919	rVV	1228555	2007722	20.53%	1.377%
12	8.410	919	922	928	rVV	31546	65130	0.67%	0.045%
13	8.469	928	932	937	rVB4	29786	51550	0.53%	0.035%
14	8.792	973	987	995	rBV	962138	1616487	16.53%	1.109%
15	9.269	1063	1068	1072	rVV3	28413	47181	0.48%	0.032%
16	9.510	1100	1109	1117	rBV	780523	1316811	13.46%	0.903%
17	10.051	1193	1201	1214	rVB	1335029	2184065	22.33%	1.498%
18	10.428	1257	1265	1271	rBV	1945085	3235855	33.09%	2.219%
19	10.475	1271	1273	1279	rVB	82497	119456	1.22%	0.082%
20	10.563	1281	1288	1296	rBV	270638	530060	5.42%	0.364%
21	12.022	1530	1536	1542	rVB3	25613	53627	0.55%	0.037%
22	12.098	1542	1549	1557	rVB	65999	119730	1.22%	0.082%
23	12.322	1580	1587	1593	rVB	47590	81425	0.83%	0.056%
24	13.139	1719	1726	1730	rBV3	29075	69480	0.71%	0.048%
25	13.616	1803	1807	1811	rVV2	69363	113232	1.16%	0.078%
26	13.674	1811	1817	1818	rVV3	36748	57070	0.58%	0.039%
27	13.710	1818	1823	1827	rVV	2286074	3180423	32.52%	2.181%
28	13.757	1827	1831	1837	rVB	976883	1333950	13.64%	0.915%
29	13.851	1843	1847	1851	rBV3	33247	42913	0.44%	0.029%
30	13.980	1862	1869	1882	rBV	2700578	4322705	44.20%	2.965%
31	14.221	1906	1910	1915	rBV2	32219	44143	0.45%	0.030%
32	14.292	1915	1922	1929	rBV2	3005534	4393888	44.93%	3.014%
33	14.357	1929	1933	1941	rVB	135729	229817	2.35%	0.158%
34	14.498	1951	1957	1966	rBV	801407	1298754	13.28%	0.891%

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Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM101419MA.M
 Title : SVOA CALIBRATION

35	14.698	1985	1991	2000	rVV5	114133	336694	3.44%	0.231%
36	14.774	2002	2004	2009	rVB2	50592	63430	0.65%	0.044%
37	14.921	2021	2029	2033	rBV3	43420	108805	1.11%	0.075%
38	14.968	2034	2037	2041	rVB2	39887	55143	0.56%	0.038%
39	15.115	2060	2062	2067	rVB5	26259	38797	0.40%	0.027%
40	15.174	2067	2072	2076	rVB2	48858	77282	0.79%	0.053%
41	15.233	2079	2082	2085	rBV2	23799	38932	0.40%	0.027%
42	15.286	2085	2091	2097	rBV	3406287	5029431	51.43%	3.449%
43	15.345	2097	2101	2106	rVB	179348	248088	2.54%	0.170%
44	15.410	2106	2112	2120	rBV	698285	1015680	10.39%	0.697%
45	15.533	2126	2133	2136	rBV3	41043	77290	0.79%	0.053%
46	15.645	2147	2152	2155	rBV	64734	93900	0.96%	0.064%
47	15.792	2172	2177	2184	rVB	74774	150254	1.54%	0.103%
48	15.874	2185	2191	2195	rVV7	41465	74816	0.76%	0.051%
49	16.039	2212	2219	2222	rBV5	61371	128178	1.31%	0.088%
50	16.068	2222	2224	2230	rVB6	44449	58160	0.59%	0.040%
51	16.351	2262	2272	2277	rBV4	73822	215657	2.21%	0.148%
52	16.398	2277	2280	2286	rVB2	59292	91750	0.94%	0.063%
53	16.621	2315	2318	2321	rVB3	52444	59405	0.61%	0.041%
54	16.692	2326	2330	2339	rVB4	127925	235163	2.40%	0.161%
55	16.780	2341	2345	2346	rBV3	47774	58659	0.60%	0.040%
56	16.857	2351	2358	2362	rBV	108893	195612	2.00%	0.134%
57	17.039	2383	2389	2393	rBV2	3683568	5317511	54.37%	3.647%
58	17.080	2393	2396	2401	rVB	2447937	3138195	32.09%	2.152%
59	17.139	2401	2406	2410	rBV	3737151	5440796	55.63%	3.732%
60	17.233	2418	2422	2428	rBV2	99638	180997	1.85%	0.124%
61	17.439	2453	2457	2463	rVB	197558	303349	3.10%	0.208%
62	17.568	2476	2479	2482	rBV3	81078	94662	0.97%	0.065%
63	17.915	2534	2538	2542	rBV	399768	542281	5.54%	0.372%
64	17.962	2542	2546	2553	rVB	971729	1332094	13.62%	0.914%
65	18.109	2565	2571	2575	rBV2	525228	993217	10.16%	0.681%
66	18.145	2575	2577	2585	rVB	311669	416983	4.26%	0.286%
67	18.262	2595	2597	2602	rVB3	111747	130895	1.34%	0.090%
68	18.421	2620	2624	2627	rBV	273031	347941	3.56%	0.239%
69	18.456	2627	2630	2637	rVB2	213893	309670	3.17%	0.212%
70	18.533	2640	2643	2652	rVB4	125734	207148	2.12%	0.142%
71	18.674	2663	2667	2672	rVB2	257326	320093	3.27%	0.220%

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72	18.721	2672	2675	2679	rBV2	135529	201056	2.06%	0.138%
73	18.839	2691	2695	2700	rBV	360820	551251	5.64%	0.378%
74	18.903	2703	2706	2709	rVB3	192261	235235	2.41%	0.161%
75	18.974	2715	2718	2728	rBV3	230439	454105	4.64%	0.311%
76	19.103	2735	2740	2745	rVB	4199375	5490590	56.14%	3.766%
77	19.168	2747	2751	2755	rBV	1003086	1337986	13.68%	0.918%
78	19.250	2762	2765	2769	rBV	289041	355400	3.63%	0.244%
79	19.439	2791	2797	2799	rBV	5090556	7381797	75.48%	5.063%
80	19.462	2799	2801	2809	rVB	3718427	4719706	48.26%	3.237%
81	19.892	2871	2874	2877	rBV2	312326	388527	3.97%	0.266%
82	19.962	2884	2886	2893	rVB2	549933	775817	7.93%	0.532%
83	20.021	2893	2896	2899	rBV	254825	289386	2.96%	0.198%
84	20.068	2900	2904	2908	rBV	340625	517521	5.29%	0.355%
85	20.121	2910	2913	2917	rVB2	420521	576985	5.90%	0.396%
86	20.515	2978	2980	2984	rBV2	283058	306321	3.13%	0.210%
87	20.968	3053	3057	3062	rBV2	4046145	4993387	51.06%	3.425%
88	21.227	3095	3101	3105	rBV2	5640402	9780072	100.00%	6.708%
89	21.262	3105	3107	3118	rVB	2137377	2724401	27.86%	1.869%
90	21.415	3130	3133	3142	rBV4	786539	1259221	12.88%	0.864%
91	21.856	3204	3208	3218	rVB2	4384915	8365947	85.54%	5.738%
92	22.309	3282	3285	3288	rBV	705309	796297	8.14%	0.546%
93	22.780	3360	3365	3367	rBV	1960899	3525484	36.05%	2.418%
94	22.850	3374	3377	3388	rVB3	2024490	3498178	35.77%	2.399%
95	23.244	3441	3444	3447	rBV	825106	1081734	11.06%	0.742%
96	23.291	3448	3452	3457	rVV	3222677	5343267	54.63%	3.665%
97	23.338	3457	3460	3464	rVB	1220211	1593654	16.29%	1.093%
98	23.433	3471	3476	3482	rBV	3374579	5948613	60.82%	4.080%
99	24.015	3571	3575	3582	rVB	1450917	2669148	27.29%	1.831%
100	25.903	3890	3896	3906	rVB	3542599	9093779	92.98%	6.237%

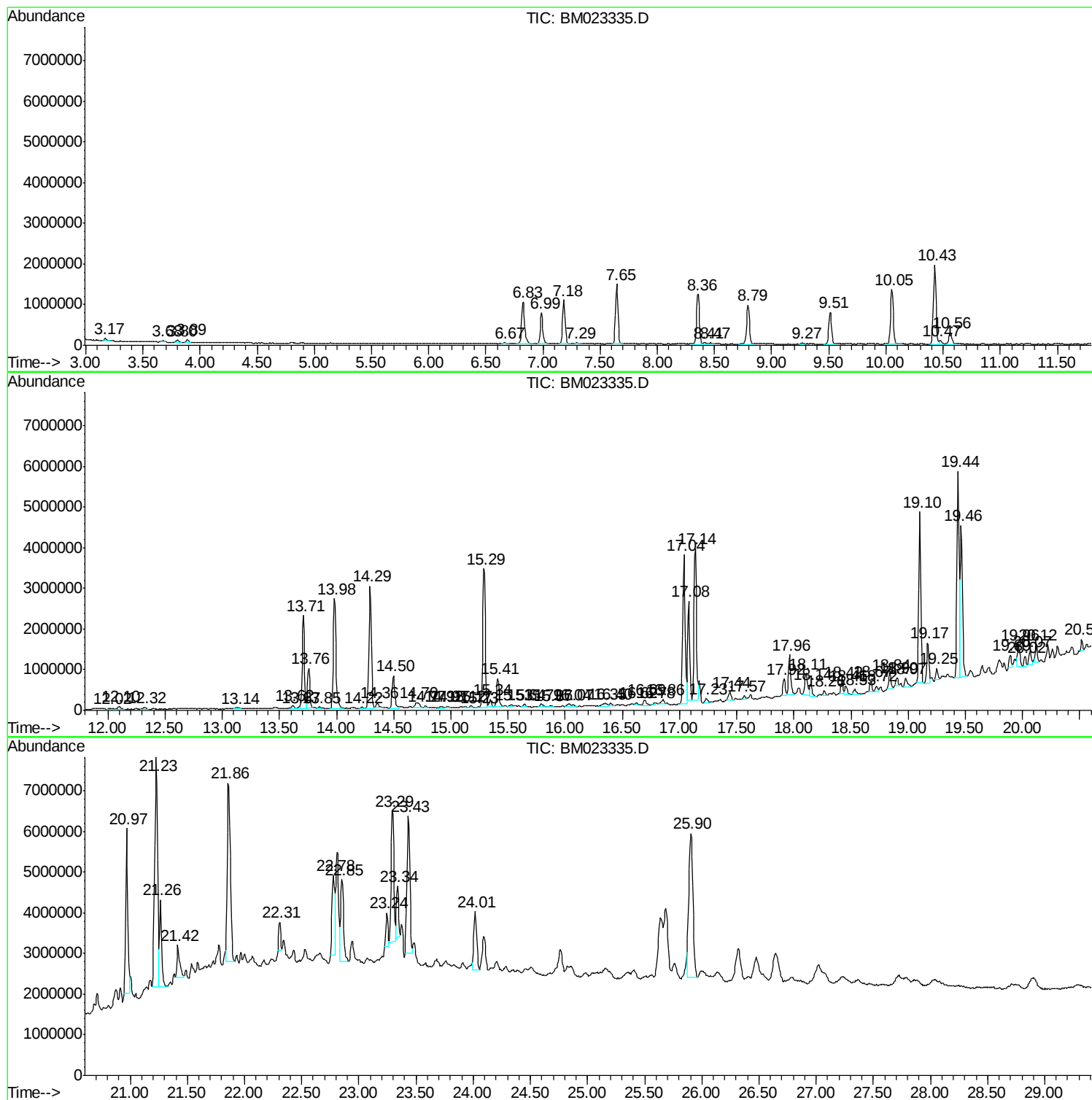
Sum of corrected areas: 145802119

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM101419MA.M
Quant Title : SVOA CALIBRATION

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



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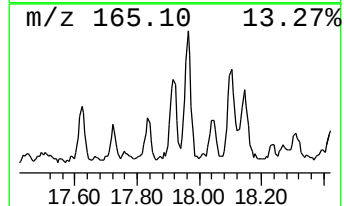
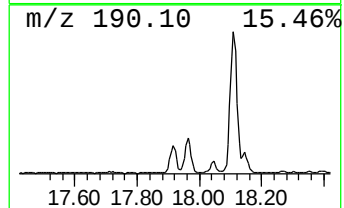
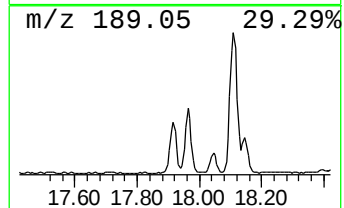
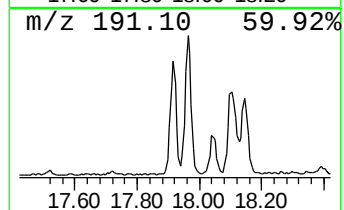
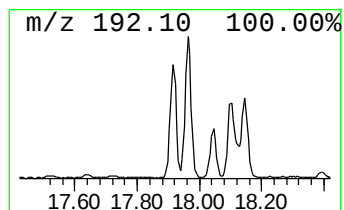
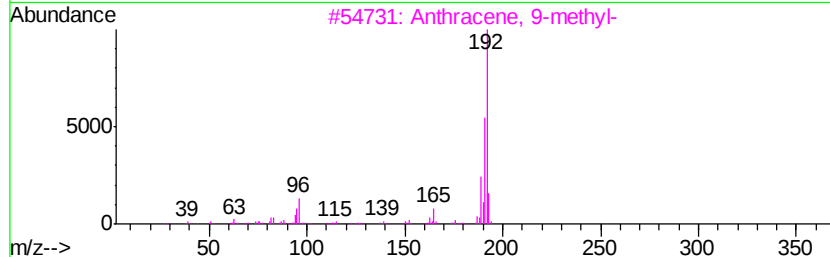
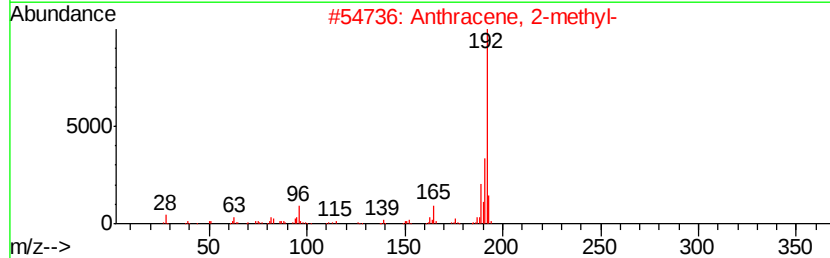
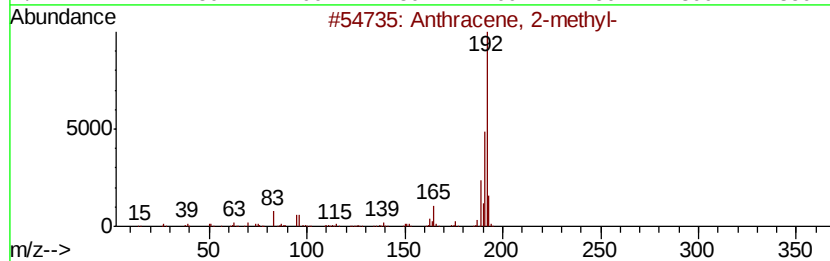
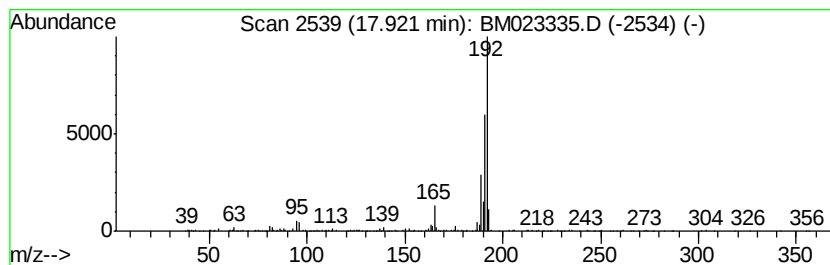
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM101419MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Anthracene, 2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.92	2.04 ng/ul	542281	Phenanthrene-d10	17.04

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
2		Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
3		Anthracene, 9-methyl-	192	C15H12	000779-02-2	93
4		5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	87
5		Anthracene, 9-methyl-	192	C15H12	000779-02-2	87



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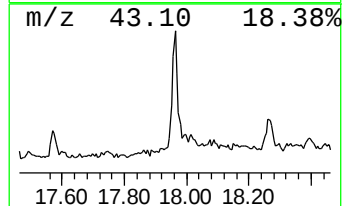
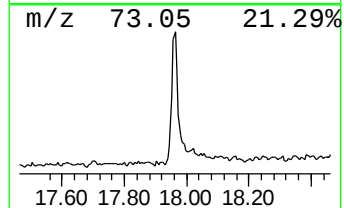
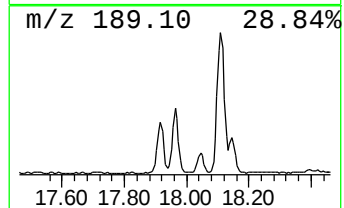
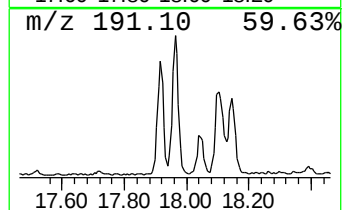
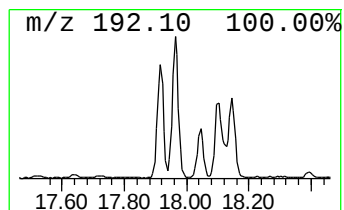
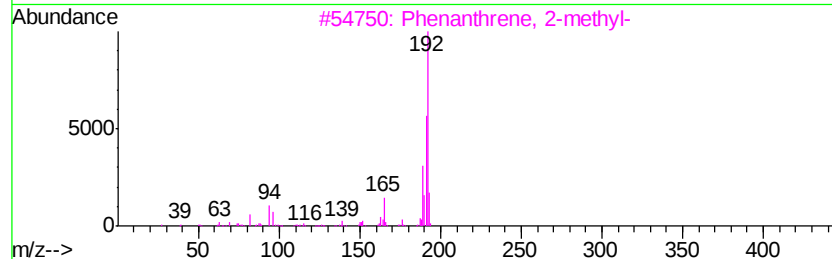
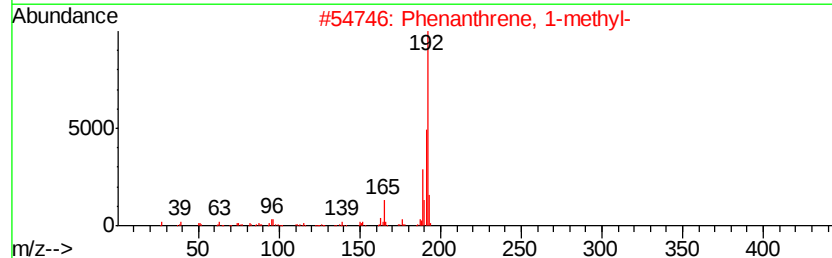
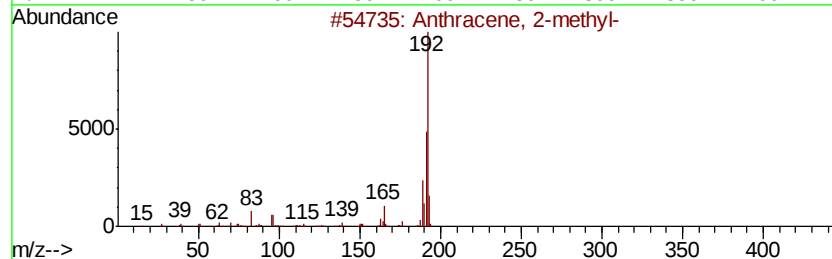
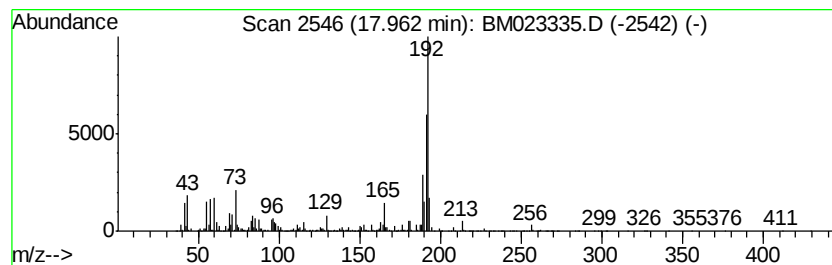
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 1H-Cyclopropa[l]phenanthren... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.96	5.01 ng/ul	1332090	Phenanthrene-d10	17.04

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Anthracene, 2-methyl-	192	C15H12	000613-12-7	97
2		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
3		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
4		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
5		1H-Cyclopropa[l]phenanthrene, 1a, ...	192	C15H12	000949-41-7	96



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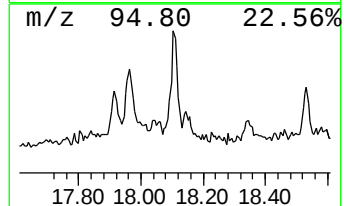
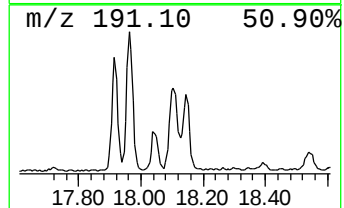
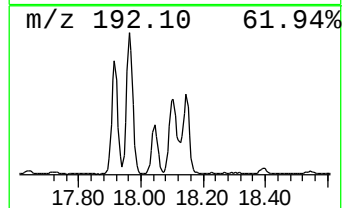
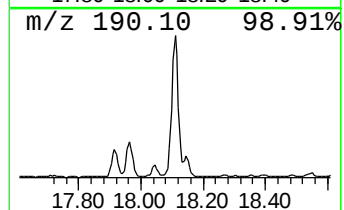
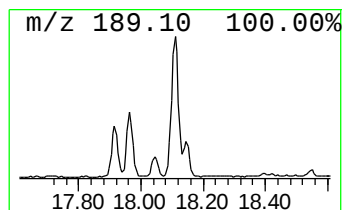
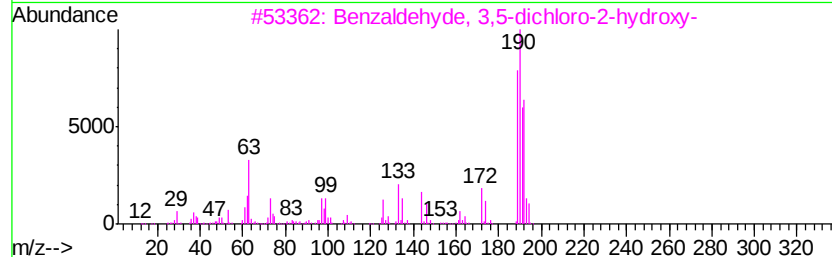
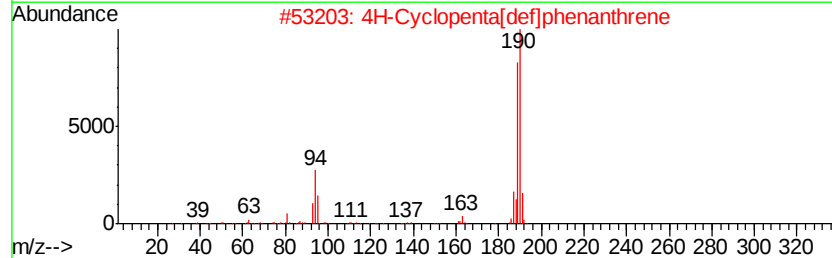
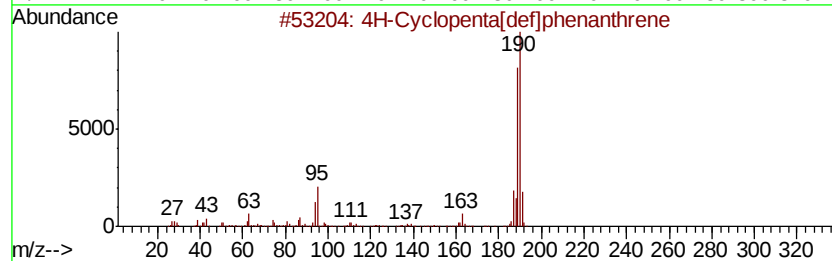
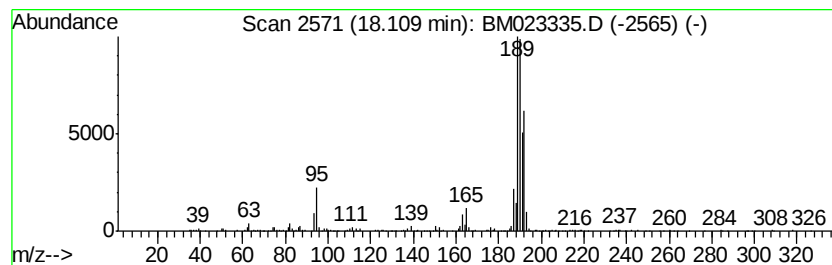
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TIC Integration Parameters: LSCINT.P

Peak Number 3 4H-Cyclopenta[def]phenanthrene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.11	3.74 ng/ul	993217	Phenanthrene-d10	17.04

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	64
2		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	52
3		Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	50
4		Pyrazole-4-carboxaldehyde, 3-(4-...	190	C10H7FN2O	306936-57-2	49
5		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	43



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Data File : BM023335.D
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Sample : K5354-14
Misc :
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ClientSampleId :
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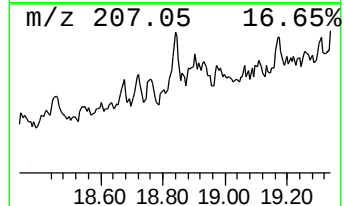
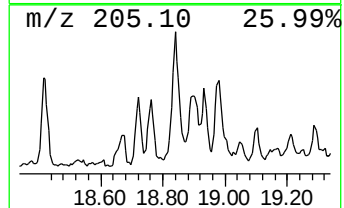
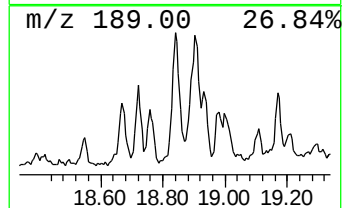
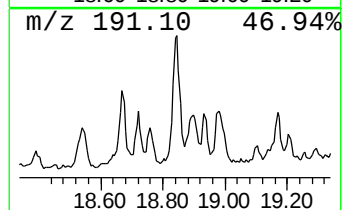
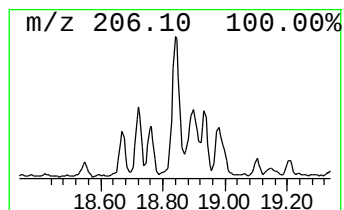
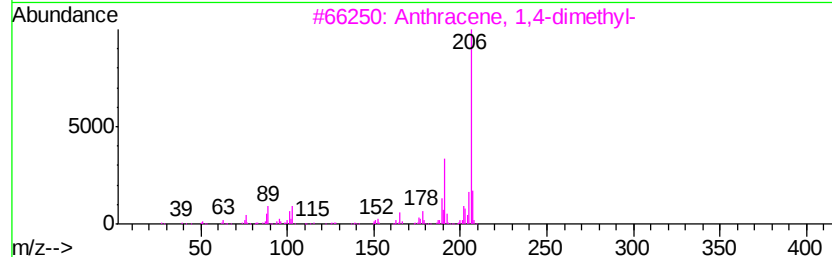
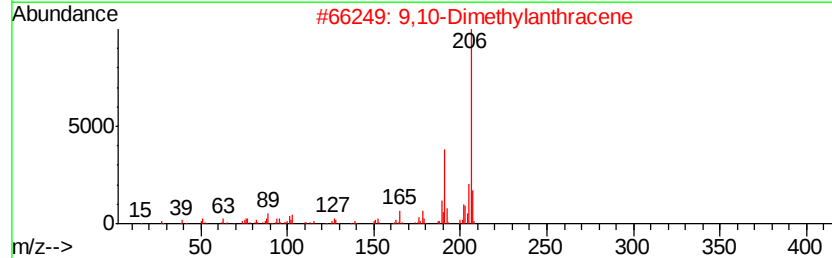
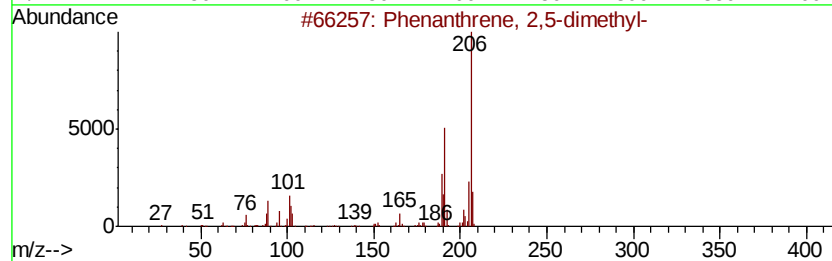
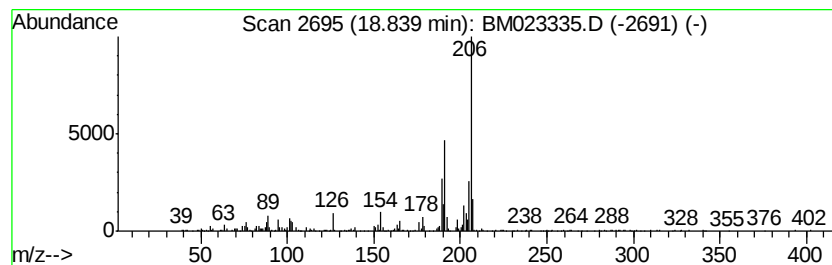
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Phenanthrene, 2,5-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.84	2.07 ng/ul	551251	Phenanthrene-d10	17.04

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	93
2		9,10-Dimethylantracene	206	C16H14	000781-43-1	90
3		Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	86
4		di-p-Tolylacetylene	206	C16H14	002789-88-0	86
5		Phenanthrene, 2,7-dimethyl-	206	C16H14	001576-69-8	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102219\
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ClientSampleId :
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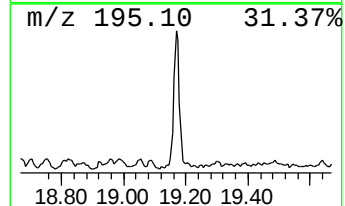
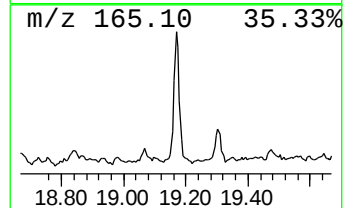
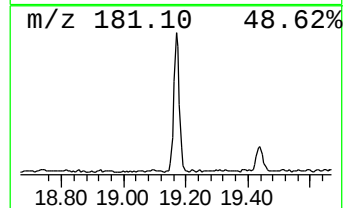
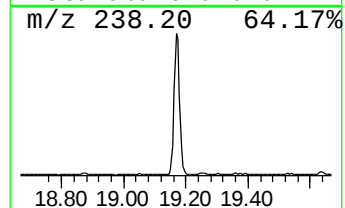
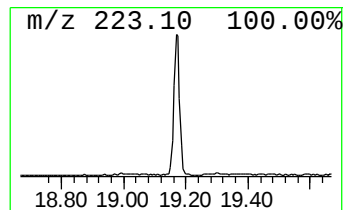
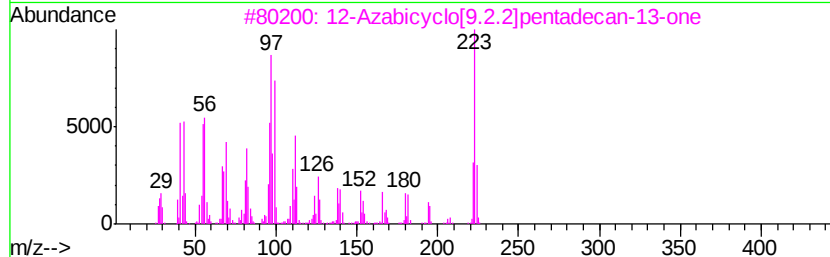
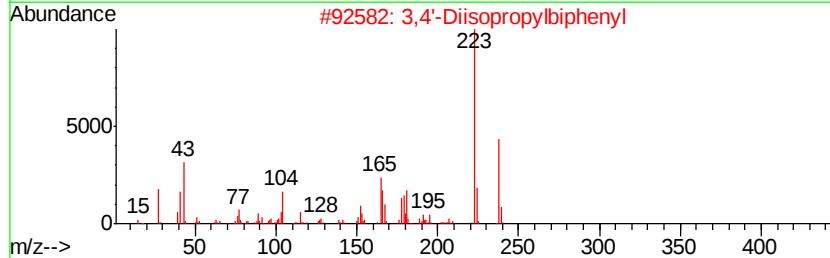
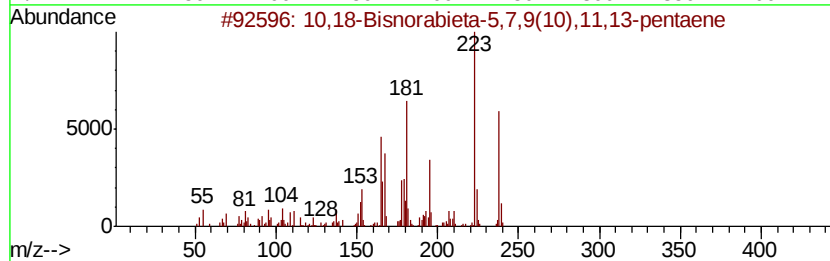
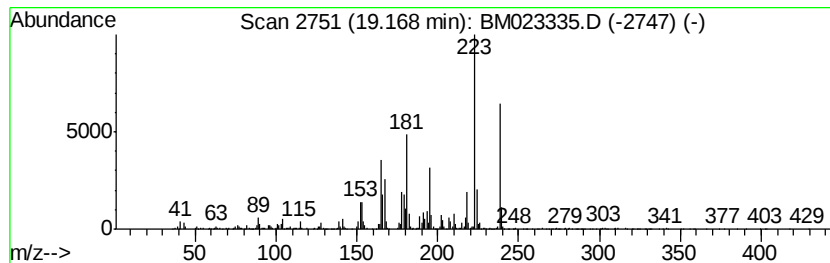
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.17	2.74 ng/ul	1337990	Chrysene-d12	21.23

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		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		12-Azabicyclo[9.2.2]pentadecan-1...	223	C14H25NO	1000191-26-7	70
4		4,4'-Diisopropylbiphenyl	238	C18H22	018970-30-4	60
5		4,4'-Diisopropylbiphenyl	238	C18H22	018970-30-4	52



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102219\
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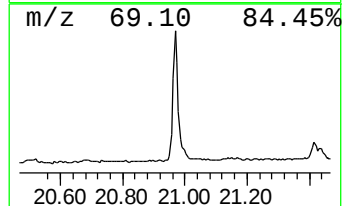
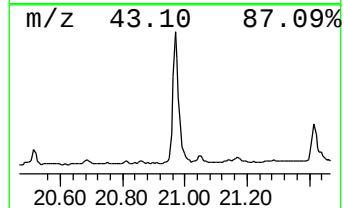
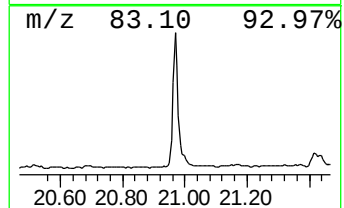
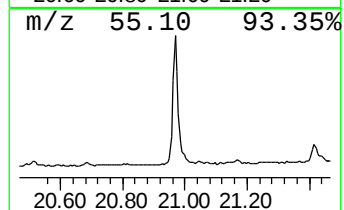
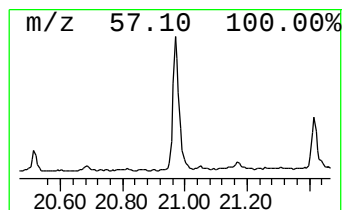
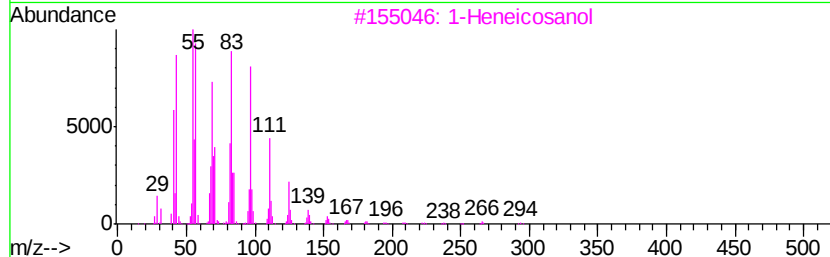
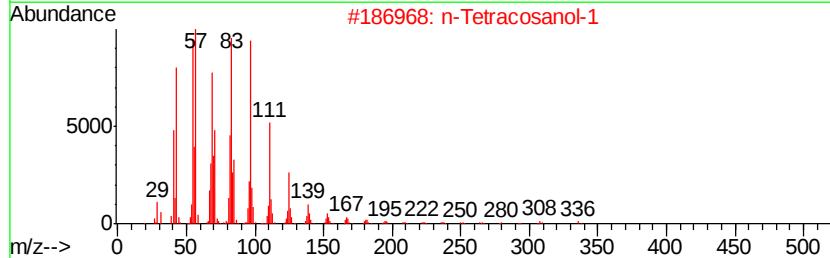
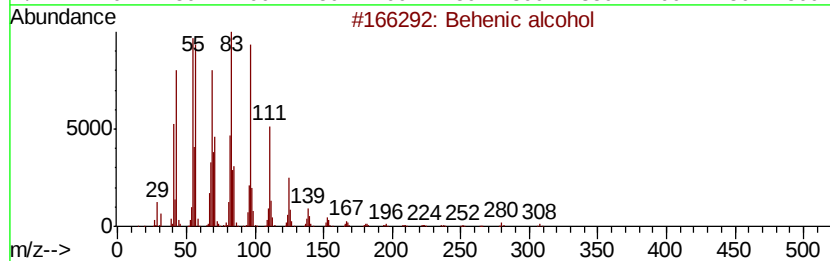
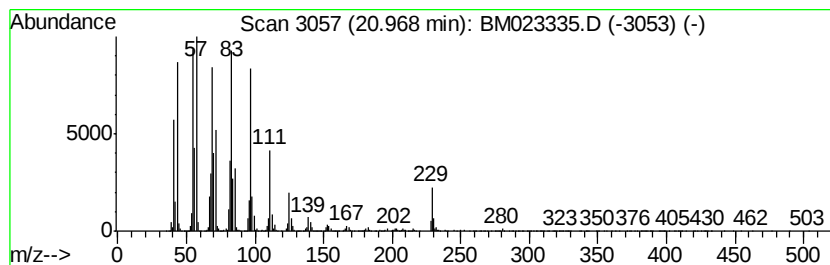
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Behenic alcohol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.97	10.21 ng/ul	4993390	Chrysene-d12	21.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Behenic alcohol	326	C22H46O	000661-19-8	94
2		n-Tetracosanol-1	354	C24H50O	000506-51-4	94
3		1-Heneicosanol	312	C21H44O	015594-90-8	94
4		Cyclopentadecane	210	C15H30	000295-48-7	93
5		5-Eicosene, (E)-	280	C20H40	074685-30-6	92



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102219\
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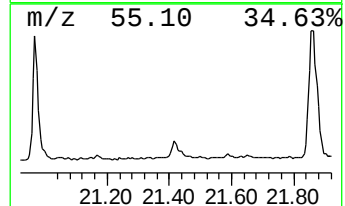
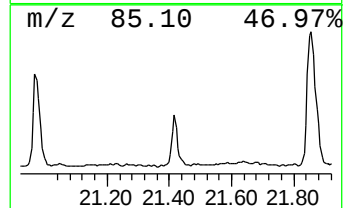
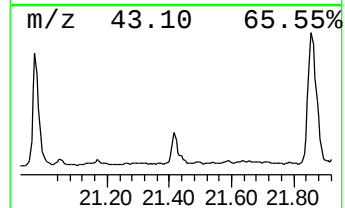
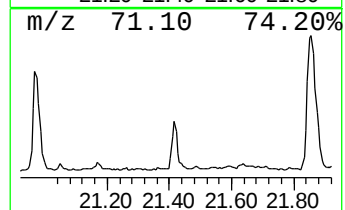
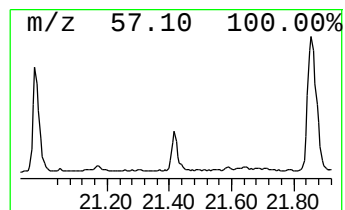
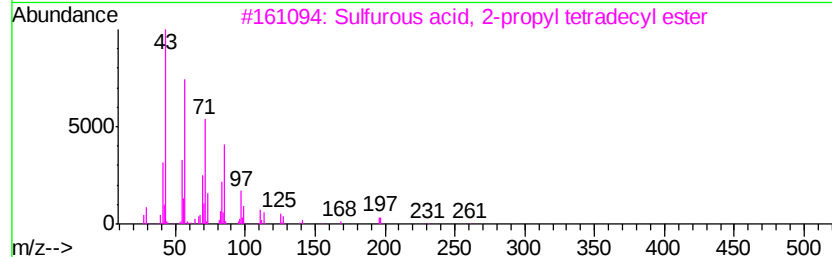
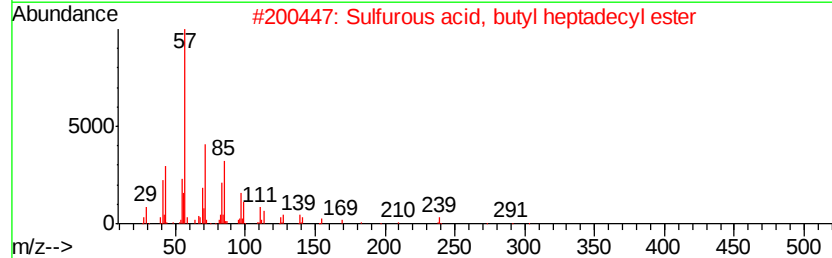
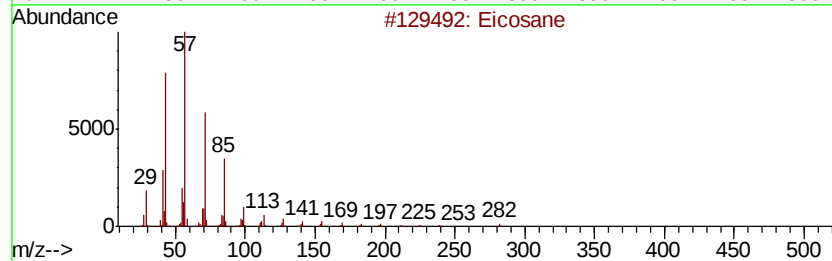
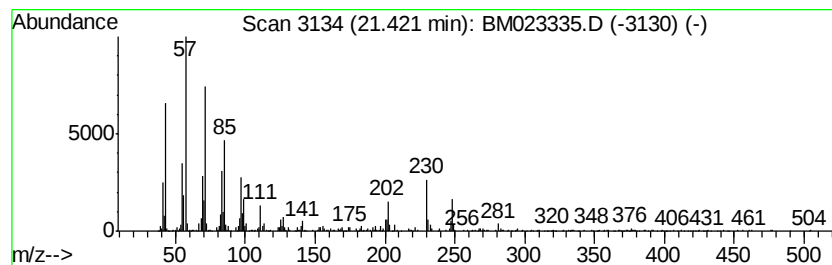
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.42	2.58 ng/ul	1259220	Chrysene-d12	21.23
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Eicosane		282 C20H42	000112-95-8 83
2	Sulfurous acid, butyl heptadecyl...		376 C21H44O3S	1000309-18-4 74
3	Sulfurous acid, 2-propyl tetradec...		320 C17H36O3S	1000309-12-5 74
4	Sulfurous acid, butyl octadecyl ...		390 C22H46O3S	1000309-18-5 74
5	Tetradecane, 1-chloro-		232 C14H29Cl	002425-54-9 64



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102219\
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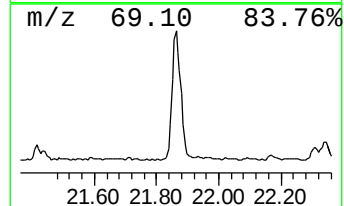
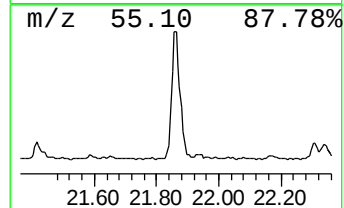
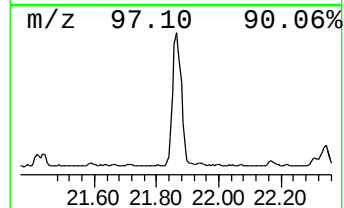
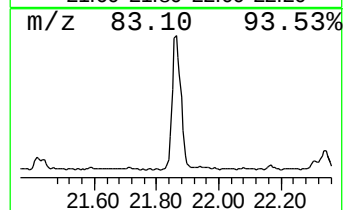
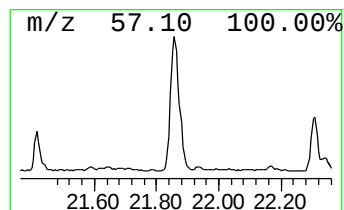
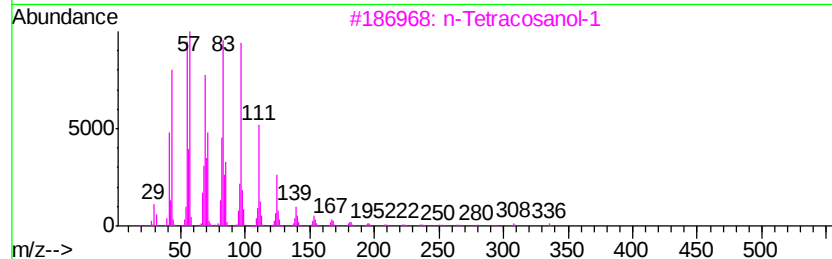
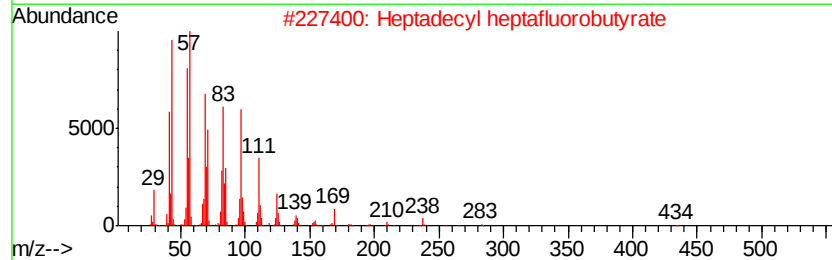
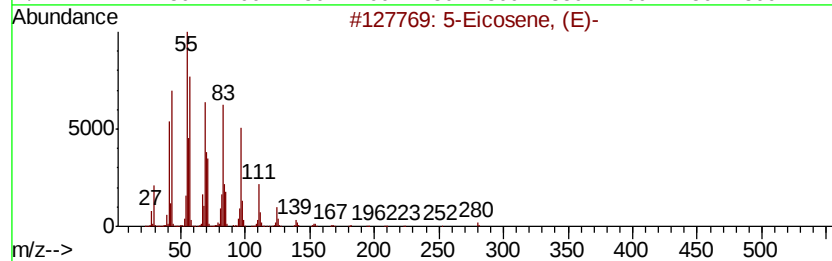
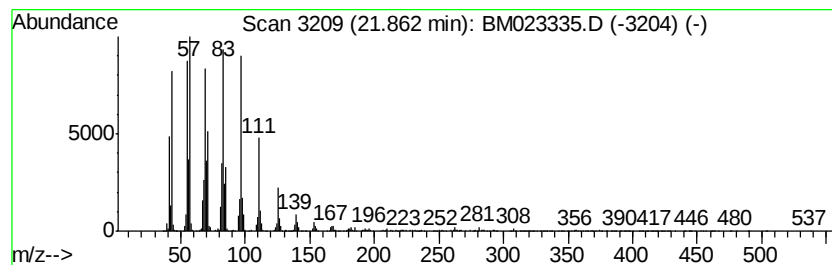
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Cyclic21.86 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.86	17.11 ng/ul	8365950	Chrysene-d12	21.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Eicosene, (E)-	280	C20H40	074685-30-6	97
2		Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
3		n-Tetracosanol-1	354	C24H50O	000506-51-4	93
4		Behenic alcohol	326	C22H46O	000661-19-8	93
5		1-Heneicosanol	312	C21H44O	015594-90-8	93



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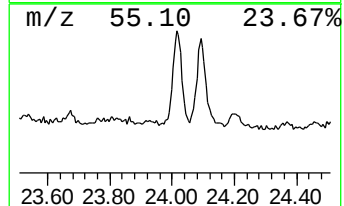
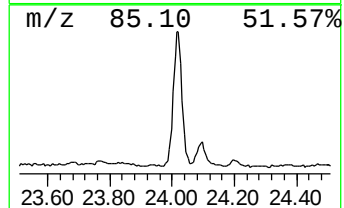
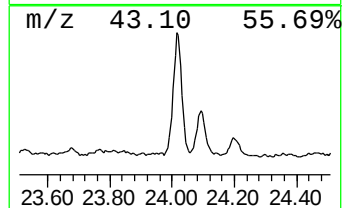
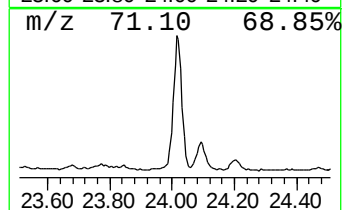
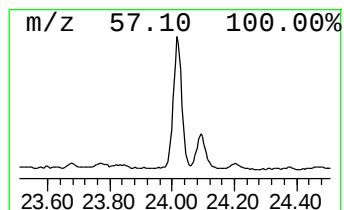
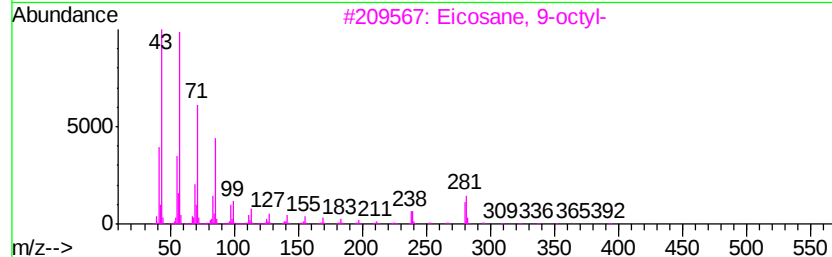
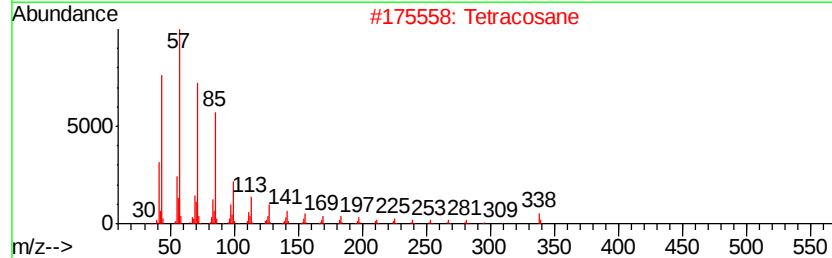
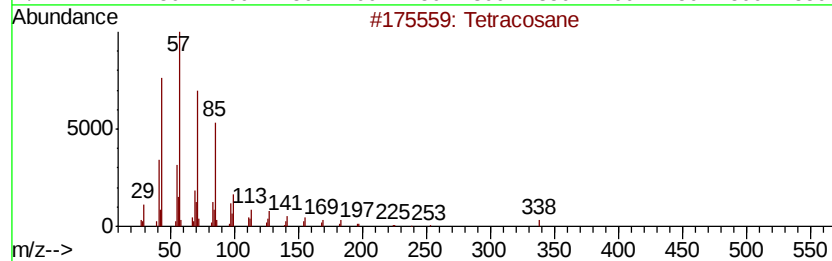
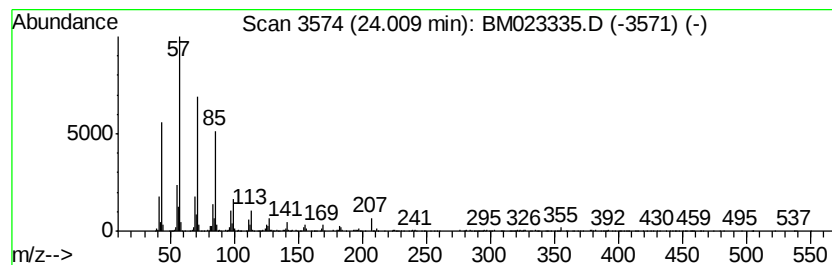
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.01	8.97 ng/ul	2669150	Perylene-d12	23.43

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosane	338	C24H50	000646-31-1	95
2		Tetracosane	338	C24H50	000646-31-1	94
3		Eicosane, 9-octyl-	394	C28H58	013475-77-9	93
4		Octacosane	394	C28H58	000630-02-4	91
5		2-Bromo dodecane	248	C12H25Br	013187-99-0	91



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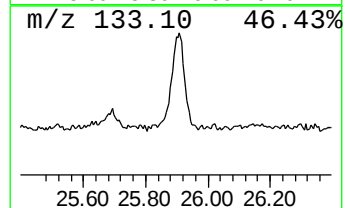
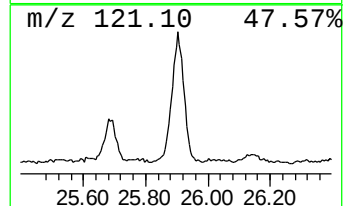
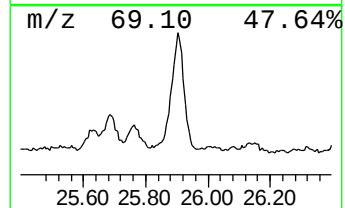
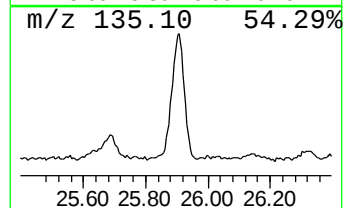
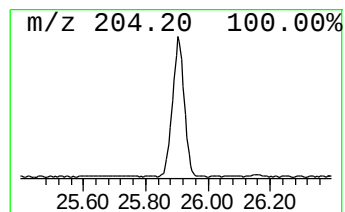
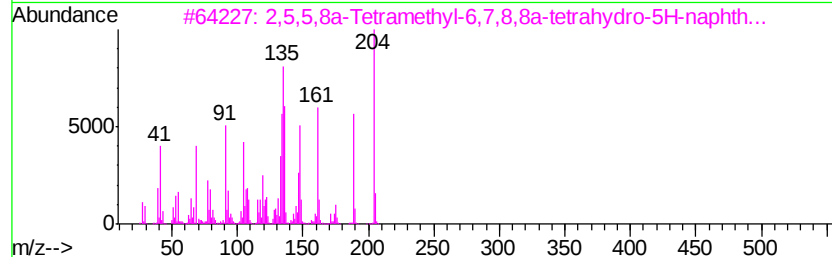
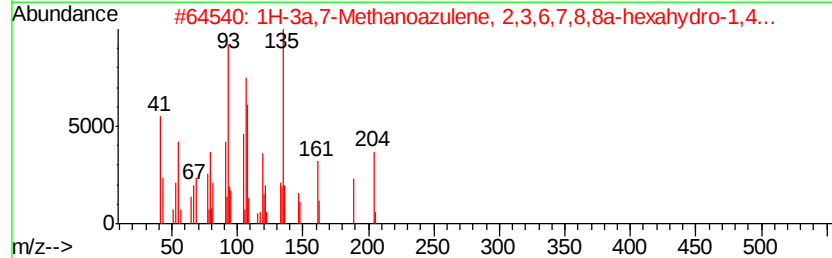
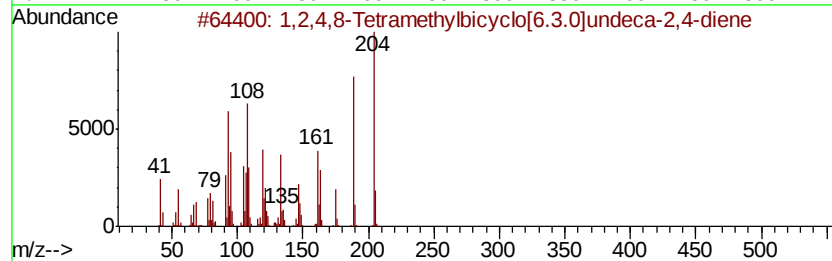
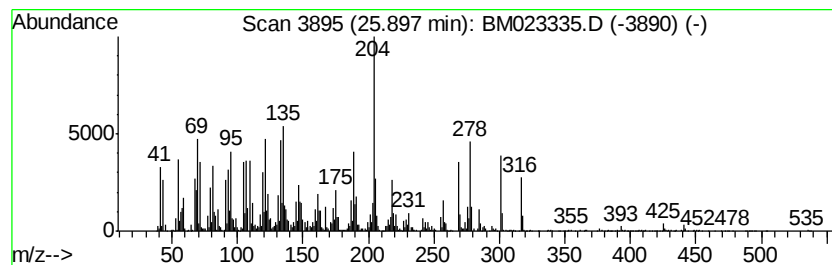
Instrument :
BNA_M
ClientSampleId :
BFB87

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM101419MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 1,2,4,8-Tetramethylbicyclo[...] Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.		
25.90	30.57 ng/ul	9093780	Perylene-d12	23.43		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2,4,8-Tetramethylbicyclo[6.3.0]undeca-2,4-diene	204	C15H24	137235-51-9	56	
2	1H-3a,7-Methanoazulene, 2,3,6,7,8a-hexahydro-1,4-dithia-	204	C15H24	000560-32-7	43	
3	2,5,5,8a-Tetramethyl-6,7,8,8a-tetrahydro-5H-naphtho[2,3-b]pyridine	204	C14H20O	124957-09-1	41	
4	3,5-Dichloro-2,4-dimethyl-1-methoxybenzene	204	C9H10Cl2O	1000250-17-8	38	
5	1H-Cycloprop[e]azulene, 1a,2,3,4,8a-hexahydro-	204	C15H24	000489-40-7	38	



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM102219\
Data File : BM023335.D
Acq On : 23 Oct 2019 00:55
Operator : JU
Sample : K5354-14
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BFB87

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM101419MA.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
Anthracene, 2-met...	17.92	2.0	ng/ul	542281	4	17.04	5317510	20.0
1H-Cyclopropa[l]p...	17.96	5.0	ng/ul	1332090	4	17.04	5317510	20.0
4H-Cyclopenta[def...	18.11	3.7	ng/ul	993217	4	17.04	5317510	20.0
Phenanthrene, 2,5...	18.84	2.1	ng/ul	551251	4	17.04	5317510	20.0
10,18-Bisnorabiet...	19.17	2.7	ng/ul	1337990	5	21.23	9780070	20.0
Behenic alcohol	20.97	10.2	ng/ul	4993390	5	21.23	9780070	20.0
(DEL) Alkane: Str...	21.42	2.6	ng/ul	1259220	5	21.23	9780070	20.0
(DEL) Alkane: Cyc...	21.86	17.1	ng/ul	8365950	5	21.23	9780070	20.0
(DEL) Alkane: Str...	24.01	9.0	ng/ul	2669150	6	23.43	5948610	20.0
1,2,4,8-Tetrameth...	25.90	30.6	ng/ul	9093780	6	23.43	5948610	20.0