LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM011516\
Data File : BM004029.D
Acq On : 15 Jan 2016 01:54
Operator : SJ/UM
Sample : H1099-09
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BC9A2

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM010116.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	3.646	92	95	104	rVB	28383	39084	3.23%	0.256%
2	6.840	633	638	649	rBV	141814	226445	18.73%	1.484%
3	6.998	659	665	674	rBV	124276	188275	15.57%	1.234%
4	7.198	693	699	708	rBB	157125	239636	19.82%	1.571%
5	7.663	772	778	785	rBB	151789	230387	19.05%	1.510%
6	8.375	894	899	908	rBB	92949	147953	12.24%	0.970%
7	8.486	911	918	924	rBB	14223	21540	1.78%	0.141%
8	8.822	969	975	983	rBV	158691	243129	20.11%	1.594%
9	9.539	1092	1097	1106	rBB	125541	205235	16.97%	1.345%
10	10.080	1183	1189	1200	rBB	188770	306919	25.38%	2.012%
11	10.451	1245	1252	1259	rBV	234219	386819	31.99%	2.536%
12	10.598	1270	1277	1294	rBB	117594	206796	17.10%	1.356%
13	13.733	1804	1810	1814	rVV	393133	523935	43.33%	3.434%
14	13.780	1814	1818	1824	rVB	215182	279094	23.08%	1.829%
15	14.010	1851	1857	1872	rVB	439257	658797	54.48%	4.318%
16	14.221	1889	1893	1897	rBB	12106	14107	1.17%	0.092%
17	14.321	1904	1910	1917	rBB2	354868	530515	43.87%	3.477%
18	14.539	1942	1947	1962	rBV	124907	196707	16.27%	1.289%
19	15.174	2051	2055	2061	rVB	21694	26537	2.19%	0.174%
20	15.315	2073	2079	2085	rBV	566775	829815	68.62%	5.439%
21	15.374	2085	2089	2097	rVB3	10060	14880	1.23%	0.098%
22	15.451	2097	2102	2108	rBV	134813	172706	14.28%	1.132%
23	16.039	2197	2202	2214	rBV2	26525	49694	4.11%	0.326%
24	16.615	2296	2300	2303	rBV2	13808	17049	1.41%	0.112%
25	16.833	2329	2337	2340	rBV2	13180	19573	1.62%	0.128%
26	17.074	2372	2378	2382	rBV	451511	649602	53.72%	4.258%
27	17.115	2382	2385	2389	rVB	167274	207676	17.17%	1.361%
28	17.168	2389	2394	2400	rBV2	601500	882360	72.97%	5.784%
29	17.945	2522	2526	2531	rVV	53061	84499	6.99%	0.554%
30	17.998	2531	2535	2540	rVB	64238	89468	7.40%	0.586%
31	18.074	2544	2548	2553	rVV	11214	16530	1.37%	0.108%
32	18.139	2553	2559	2563	rVV2	69915	122175	10.10%	0.801%
33	18.180	2563	2566	2574	rVB	46151	61666	5.10%	0.404%
34	18.451	2607	2612	2616	rVV	28951	39367	3.26%	0.258%

LSC Area Percent Report

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Acq On : 15 Jan 2016 01:54
Operator : SJ/UM
Sample : H1099-09
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BC9A2

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM010116.M
Title : SVOA CALIBRATION

35	18.498	2616	2620	2626	rVV2	32122	45245	3.74%	0.297%
36	18.698	2651	2654	2659	rVV	14395	19888	1.64%	0.130%
37	18.751	2659	2663	2667	rVV	17493	23174	1.92%	0.152%
38	18.792	2667	2670	2674	rVB2	11486	13457	1.11%	0.088%
39	18.874	2679	2684	2690	rBV	52114	86724	7.17%	0.568%
40	18.933	2690	2694	2698	rVV3	21166	39762	3.29%	0.261%
41	18.962	2698	2699	2703	rVB	14666	14657	1.21%	0.096%
42	19.015	2703	2708	2716	rBV3	26783	57470	4.75%	0.377%
43	19.139	2724	2729	2734	rVB	307142	397173	32.85%	2.603%
44	19.239	2743	2746	2751	rVB3	14785	19863	1.64%	0.130%
45	19.345	2758	2764	2767	rBV4	10658	17543	1.45%	0.115%
46	19.386	2768	2771	2775	rVV3	11761	16550	1.37%	0.108%
47	19.480	2780	2787	2790	rBV	769417	1209233	100.00%	7.926%
48	19.503	2790	2791	2800	rVV	338960	348746	28.84%	2.286%
49	19.580	2800	2804	2809	rVB2	21306	32434	2.68%	0.213%
50	19.686	2817	2822	2828	rVV2	10101	19296	1.60%	0.126%
51	19.745	2828	2832	2838	rVB3	13231	20192	1.67%	0.132%
52	19.839	2842	2848	2858	rBV5	33889	83625	6.92%	0.548%
53	19.968	2865	2870	2872	rVV4	28062	43840	3.63%	0.287%
54	20.003	2872	2876	2882	rVB2	62468	99175	8.20%	0.650%
55	20.103	2890	2893	2897	rVV	36989	50797	4.20%	0.333%
56	20.162	2897	2903	2908	rVV2	56518	91583	7.57%	0.600%
57	20.203	2908	2910	2914	rVV2	12943	16605	1.37%	0.109%
58	20.245	2914	2917	2923	rVV3	9337	17040	1.41%	0.112%
59	20.303	2923	2927	2931	rVV	47588	61748	5.11%	0.405%
60	20.350	2931	2935	2939	rVV	46202	61665	5.10%	0.404%
61	20.474	2950	2956	2959	rVB6	8706	16405	1.36%	0.108%
62	20.568	2967	2972	2974	rBV3	10303	15880	1.31%	0.104%
63	20.633	2980	2983	2987	rVB	11620	15449	1.28%	0.101%
64	20.721	2992	2998	3001	rBV6	10347	21633	1.79%	0.142%
65	20.756	3001	3004	3011	rVB	33092	51177	4.23%	0.335%
66	20.921	3024	3032	3036	rBV3	31364	64289	5.32%	0.421%
67	20.962	3036	3039	3040	rBV	34750	36711	3.04%	0.241%
68	20.986	3040	3043	3050	rVV3	65160	113967	9.42%	0.747%
69	21.056	3051	3055	3062	rVB2	29456	46587	3.85%	0.305%
70	21.162	3066	3073	3075	rBV	9546	17672	1.46%	0.116%
71	21.274	3085	3092	3096	rBV2	569735	945749	78.21%	6.199%

LSC Area Percent Report

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Sample : H1099-09
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BC9A2

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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Peak separation: 5

Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM010116.M
Title : SVOA CALIBRATION

72	21.315	3096	3099	3108	rVB	178693	231609	19.15%	1.518%
73	21.392	3108	3112	3115	rBV2	15770	20208	1.67%	0.132%
74	21.433	3115	3119	3124	rBV3	22726	31390	2.60%	0.206%
75	21.486	3124	3128	3131	rBV4	10578	15141	1.25%	0.099%
76	21.597	3141	3147	3151	rBV4	15385	26133	2.16%	0.171%
77	21.768	3173	3176	3179	rVV	12245	16802	1.39%	0.110%
78	21.827	3179	3186	3189	rVV	53806	95515	7.90%	0.626%
79	21.874	3189	3194	3201	rVV2	45309	104343	8.63%	0.684%
80	21.944	3202	3206	3209	rVV2	20751	38385	3.17%	0.252%
81	21.980	3210	3212	3215	rVV2	20210	23508	1.94%	0.154%
82	22.027	3215	3220	3223	rVV2	28916	50947	4.21%	0.334%
83	22.056	3223	3225	3231	rVV3	20892	25826	2.14%	0.169%
84	22.121	3231	3236	3241	rVB2	17435	27284	2.26%	0.179%
85	22.315	3264	3269	3283	rVB6	25229	54992	4.55%	0.360%
86	22.444	3284	3291	3295	rBV5	15280	31754	2.63%	0.208%
87	22.597	3311	3317	3321	rVB5	10533	20687	1.71%	0.136%
88	22.844	3350	3359	3364	rBV2	93494	263041	21.75%	1.724%
89	23.015	3384	3388	3393	rVB	16085	24649	2.04%	0.162%
90	23.321	3434	3440	3443	rBV	56255	98539	8.15%	0.646%
91	23.374	3443	3449	3454	rVV	448402	830762	68.70%	5.445%
92	23.415	3454	3456	3463	rVB	91887	134212	11.10%	0.880%
93	23.515	3465	3473	3479	rBV	294331	547972	45.32%	3.592%
94	24.015	3551	3558	3569	rVB2	36695	83392	6.90%	0.547%
95	24.385	3617	3621	3626	rVB2	7666	14081	1.16%	0.092%
96	24.756	3678	3684	3691	rBV2	19263	41641	3.44%	0.273%
97	25.615	3824	3830	3837	rVB4	13623	33551	2.77%	0.220%
98	25.762	3846	3855	3865	rVB2	36806	103231	8.54%	0.677%
99	26.121	3910	3916	3925	rVB6	6496	16939	1.40%	0.111%
100	26.450	3964	3972	3983	rVB	22359	67442	5.58%	0.442%

Sum of corrected areas: 15255950

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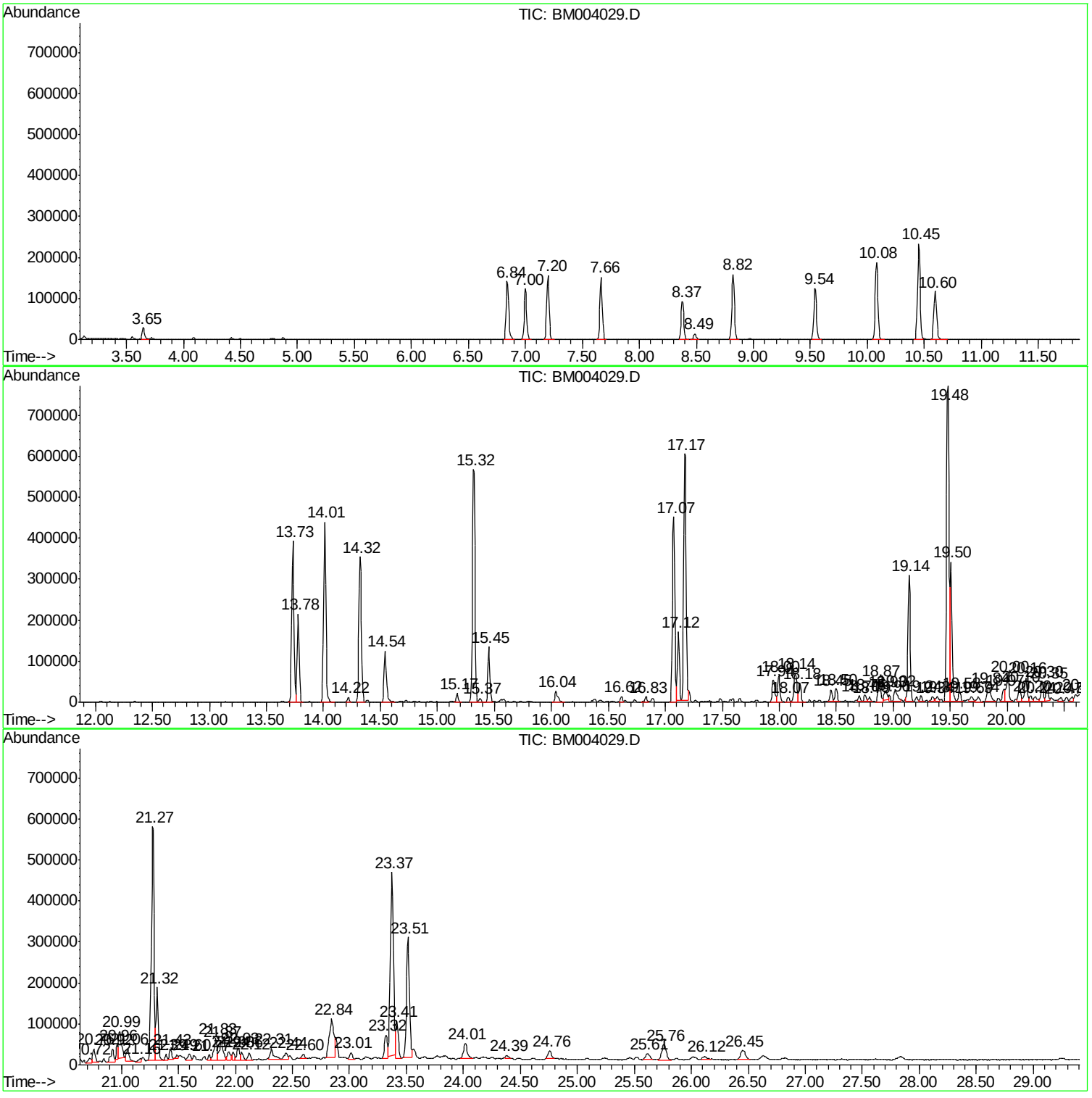
Instrument :
BNA_M
ClientSampleId :
BC9A2

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM010116.M
Quant Title : SVOA CALIBRATION

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



Library Search Compound Report

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Instrument : BNA_M
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Quant Title : SVOA CALIBRATION

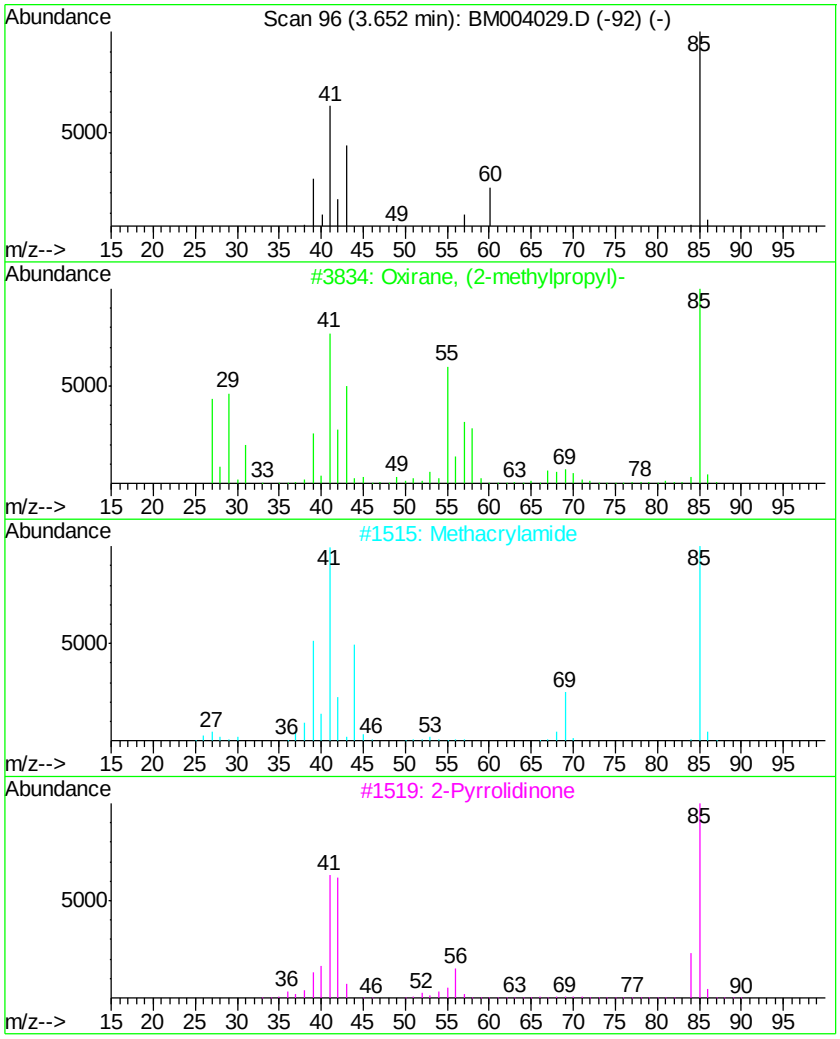
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

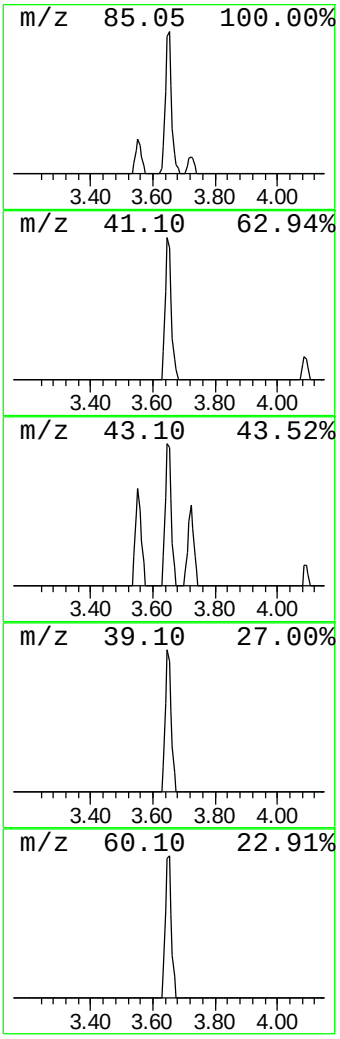
****

Peak Number 1 unknown3.65 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.65	3.39 ng/ul	39084	1,4-Dichlorobenzene-d4	7.66

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxirane, (2-methylpropyl)-	100	C6H12O	023850-78-4	36
2	Methacrylamide	85	C4H7NO	000079-39-0	33
3	2-Pyrrolidinone	85	C4H7NO	000616-45-5	9
4	Methacrylamide	85	C4H7NO	000079-39-0	7
5	2-Pyrrolidinone	85	C4H7NO	000616-45-5	7





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM011516\
Data File : BM004029.D
Acq On : 15 Jan 2016 01:54
Operator : SJ/UM
Sample : H1099-09
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BC9A2

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM010116.M
Quant Title : SVOA CALIBRATION

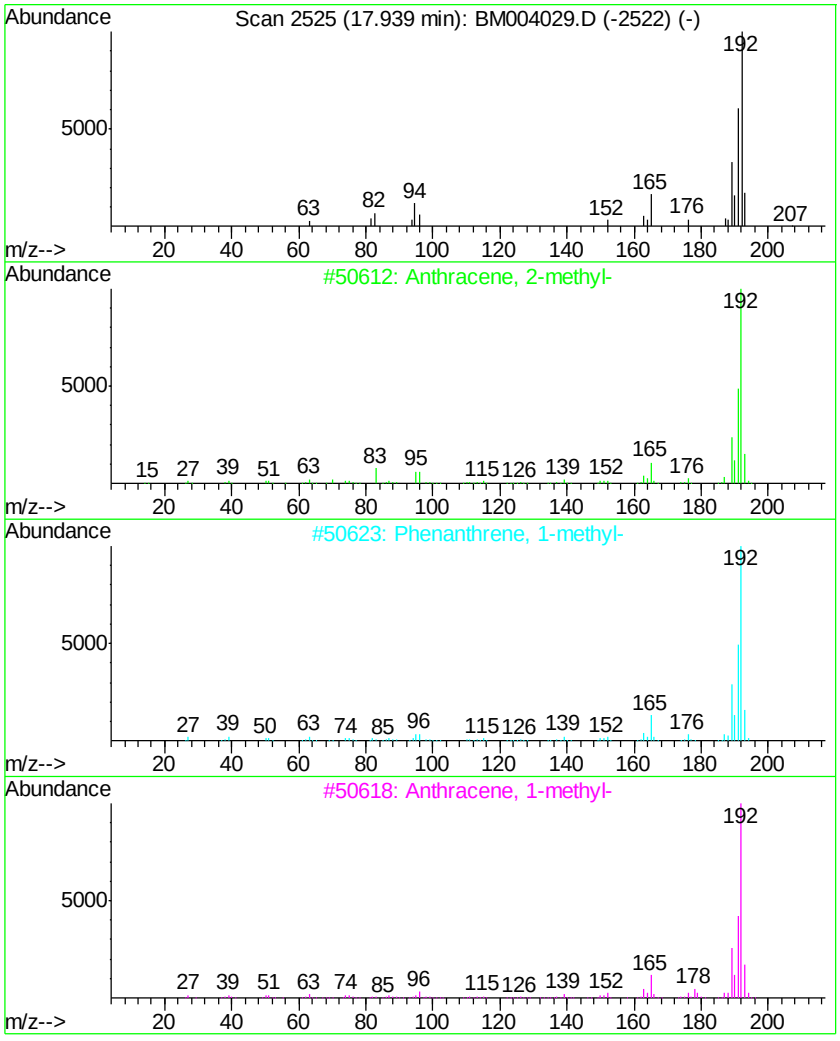
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

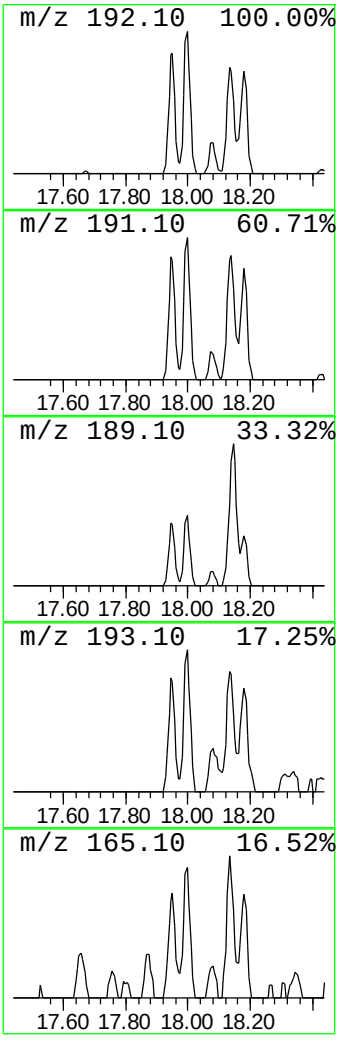
****

Peak Number 2 Anthracene, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.94	2.60 ng/ul	84499	Phenanthrene-d10	17.07

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
3	Anthracene, 1-methyl-	192	C15H12	000610-48-0	96
4	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
5	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	95





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM011516\
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Acq On : 15 Jan 2016 01:54
Operator : SJ/UM
Sample : H1099-09
Misc :
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Instrument : BNA_M
ClientSampleId : BC9A2

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM010116.M
Quant Title : SVOA CALIBRATION

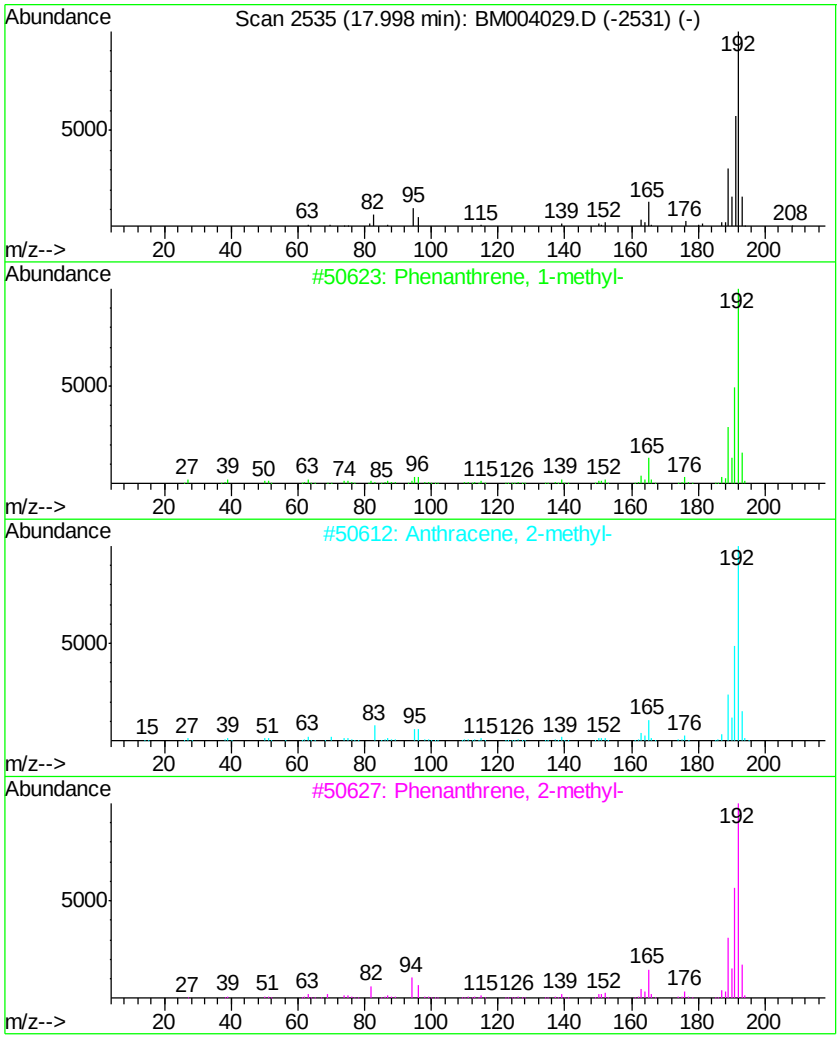
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

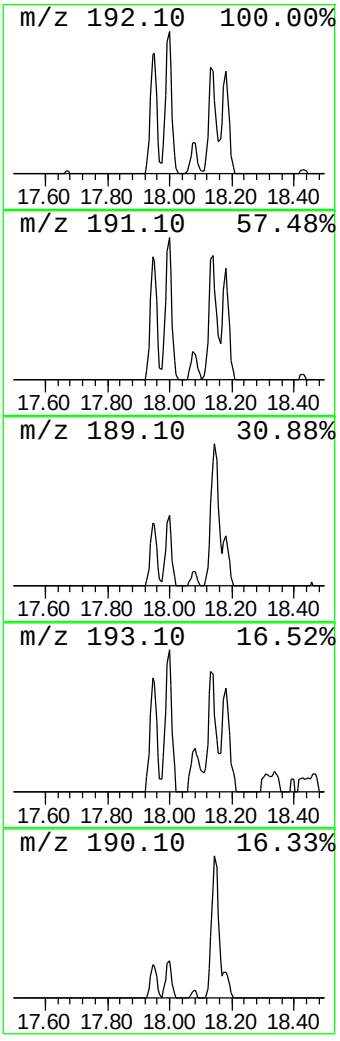
****

Peak Number 3 Phenanthrene, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.00	2.75 ng/ul	89468	Phenanthrene-d10	17.07

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
2	Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
3	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
4	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	96
5	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	95





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM011516\
Data File : BM004029.D
Acq On : 15 Jan 2016 01:54
Operator : SJ/UM
Sample : H1099-09
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BC9A2

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM010116.M
Quant Title : SVOA CALIBRATION

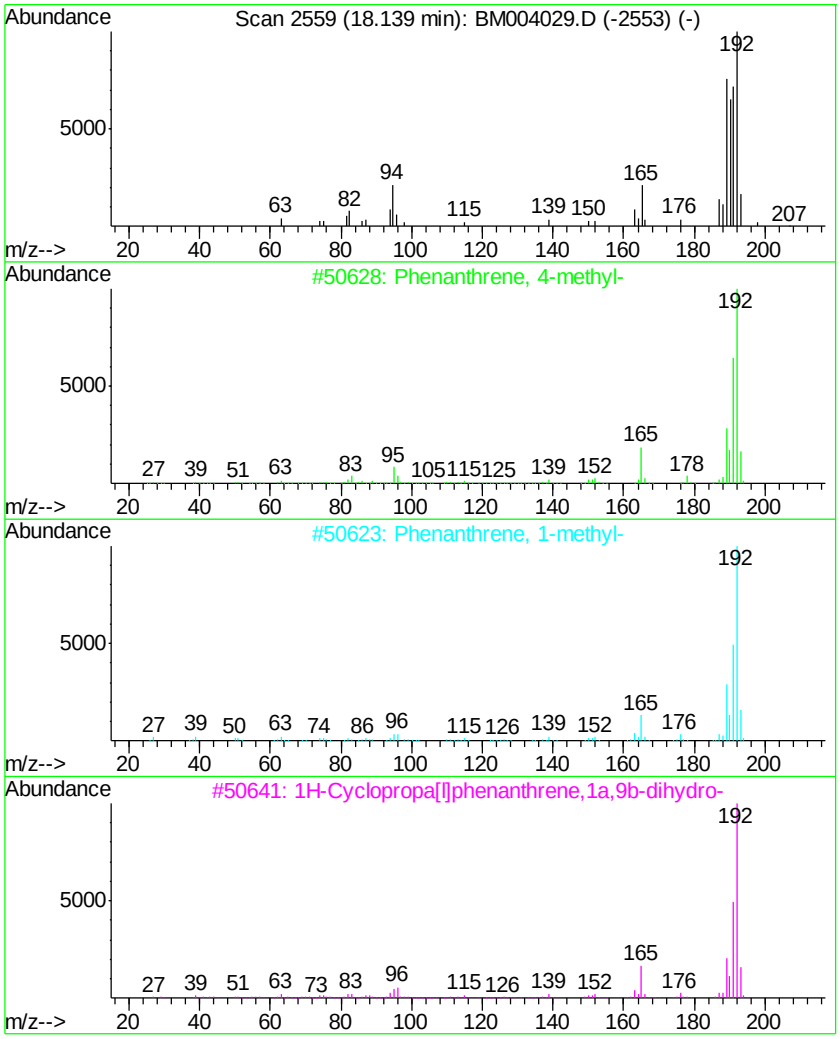
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

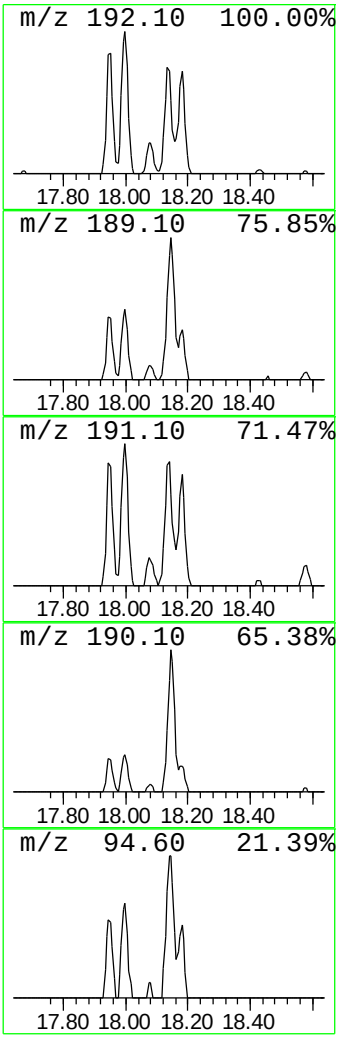
****

Peak Number 4 Phenanthrene, 4-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.14	3.76 ng/ul	122175	Phenanthrene-d10	17.07

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	50
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	50
3	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	50
4	Anthracene, 1-methyl-	192	C15H12	000610-48-0	50
5	1a,9b-Dihydro-1H-cyclopropa[a]an...	192	C15H12	006109-24-6	45





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Data File : BM004029.D
Acq On : 15 Jan 2016 01:54
Operator : SJ/UM
Sample : H1099-09
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BC9A2

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM010116.M
Quant Title : SVOA CALIBRATION

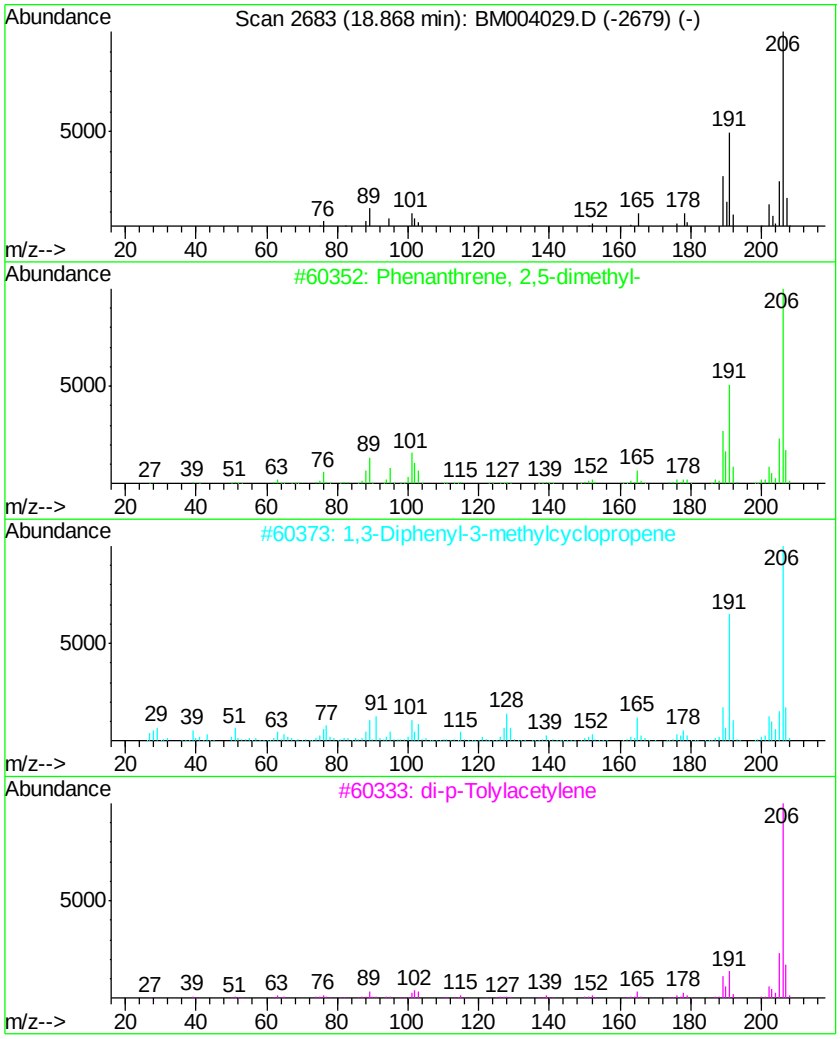
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

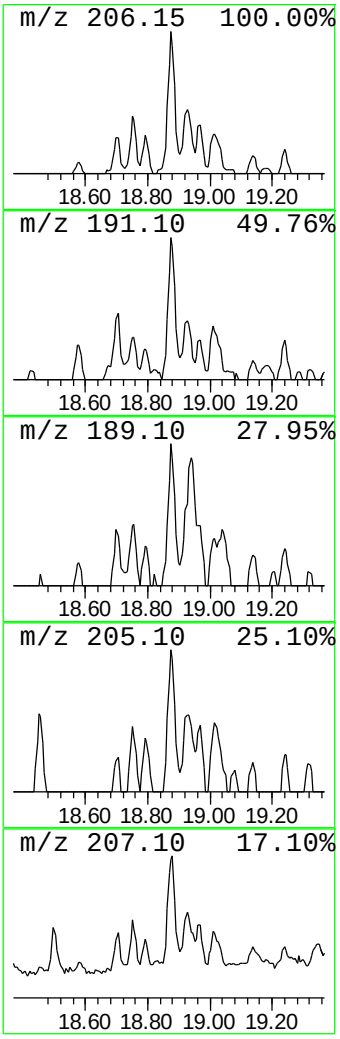
****

Peak Number 5 Phenanthrene, 2,5-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.87	2.67 ng/ul	86724	Phenanthrene-d10	17.07

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	96
2	1,3-Diphenyl-3-methylcyclopropene	206	C16H14	086544-79-8	93
3	di-p-Tolylacetylene	206	C16H14	002789-88-0	93
4	9,10-Dimethylanthracene	206	C16H14	000781-43-1	91
5	Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	87





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Data File : BM004029.D
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Operator : SJ/UM
Sample : H1099-09
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BC9A2

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM010116.M
Quant Title : SVOA CALIBRATION

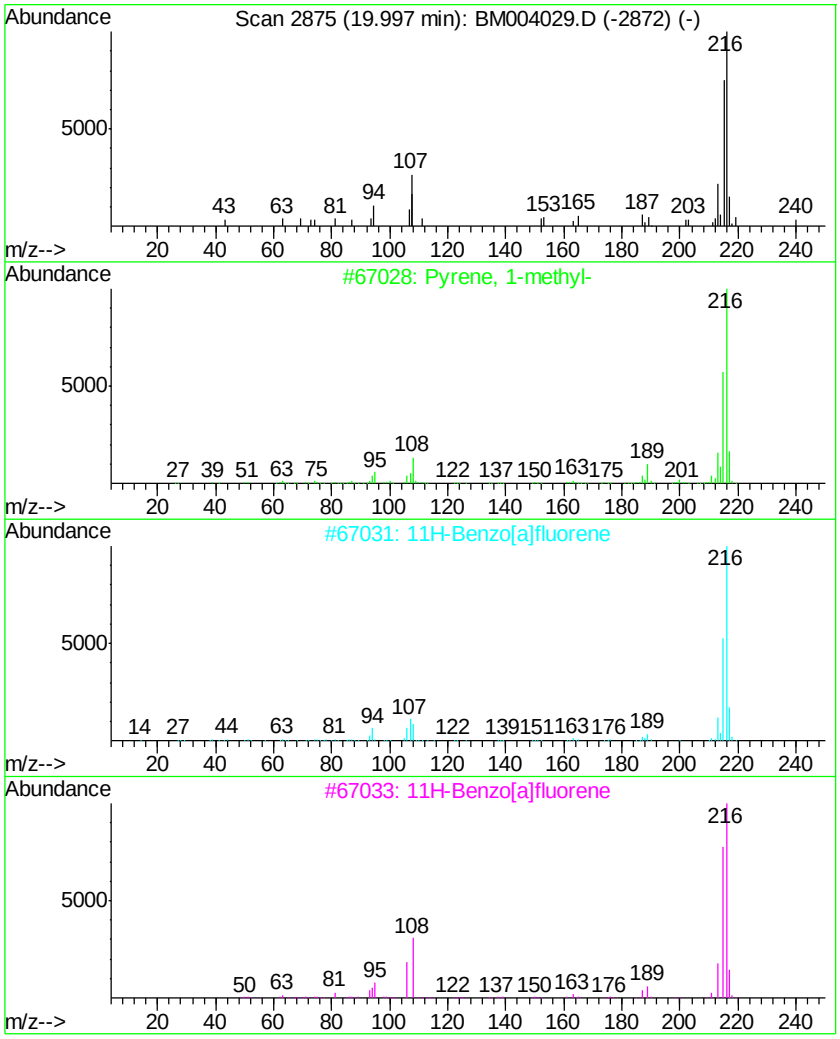
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

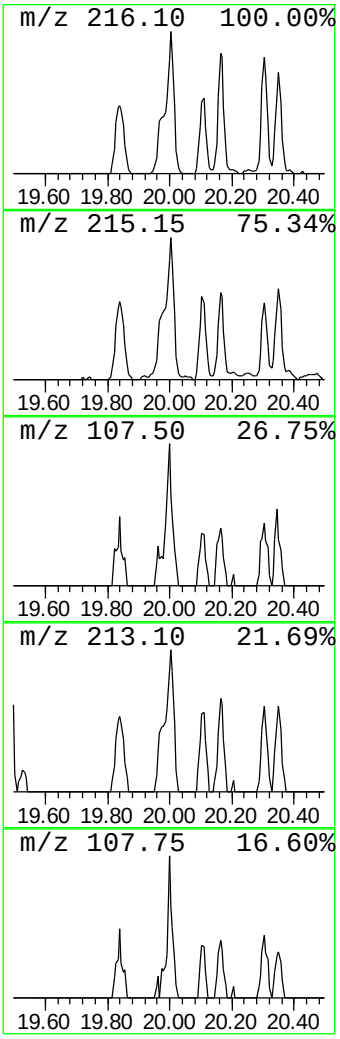
****

Peak Number 6 Pyrene, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.00	2.10 ng/ul	99175	Chrysene-d12	21.28

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	87
2	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	87
3	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	87
4	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	83
5	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	76





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Acq On : 15 Jan 2016 01:54
Operator : SJ/UM
Sample : H1099-09
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BC9A2

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM010116.M
Quant Title : SVOA CALIBRATION

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

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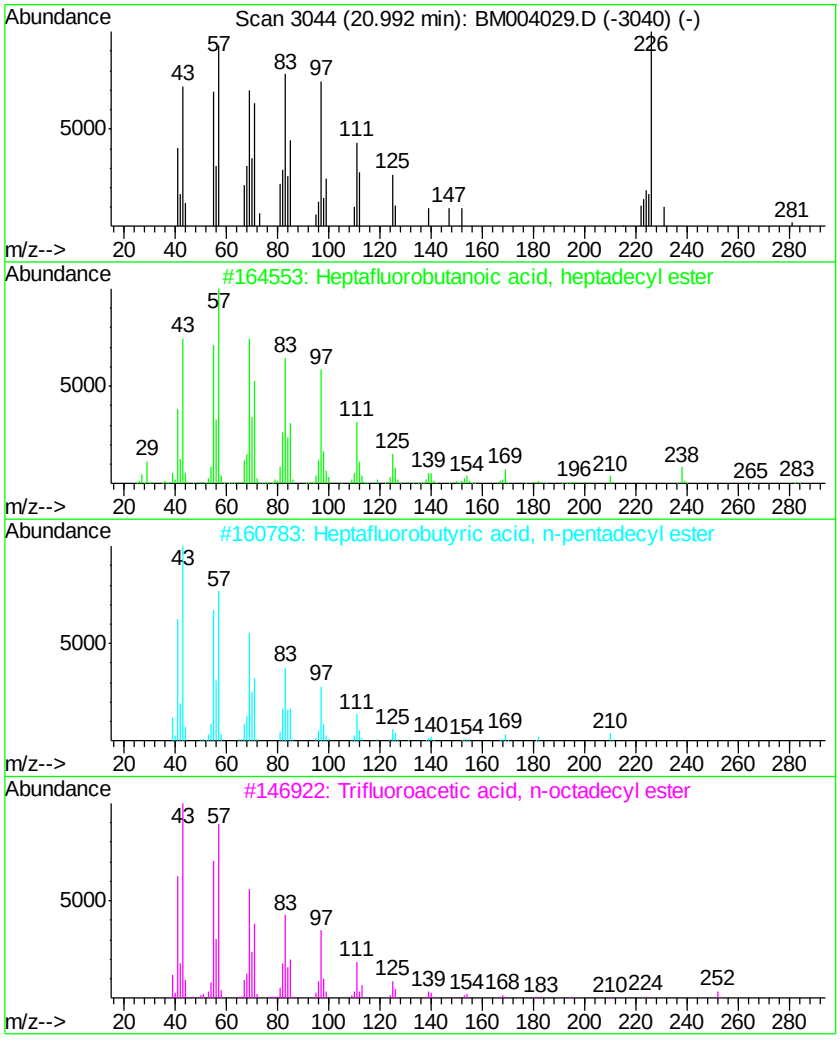
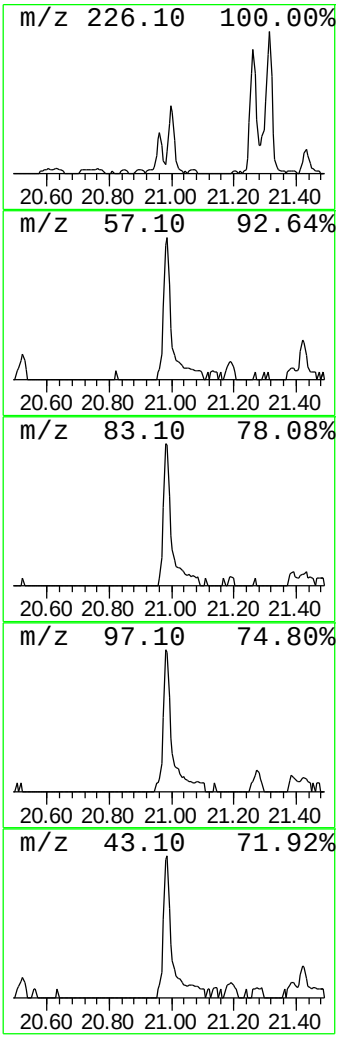
Peak Number 7 Heptafluorobutanoic acid, h... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.99	2.41 ng/ul	113967	Chrysene-d12	21.28

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	91
2	Heptafluorobutyric acid, n-penta...	424	C19H31F7O2	1000216-79-3	91
3	Trifluoroacetic acid, n-octadecy...	366	C20H37F3O2	079392-43-1	91
4	17-Pentatriacontene	491	C35H70	006971-40-0	91
5	1-Hexadecanesulfonyl chloride	324	C16H33ClO2S	038775-38-1	81





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Data File : BM004029.D
Acq On : 15 Jan 2016 01:54
Operator : SJ/UM
Sample : H1099-09
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BC9A2

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM010116.M
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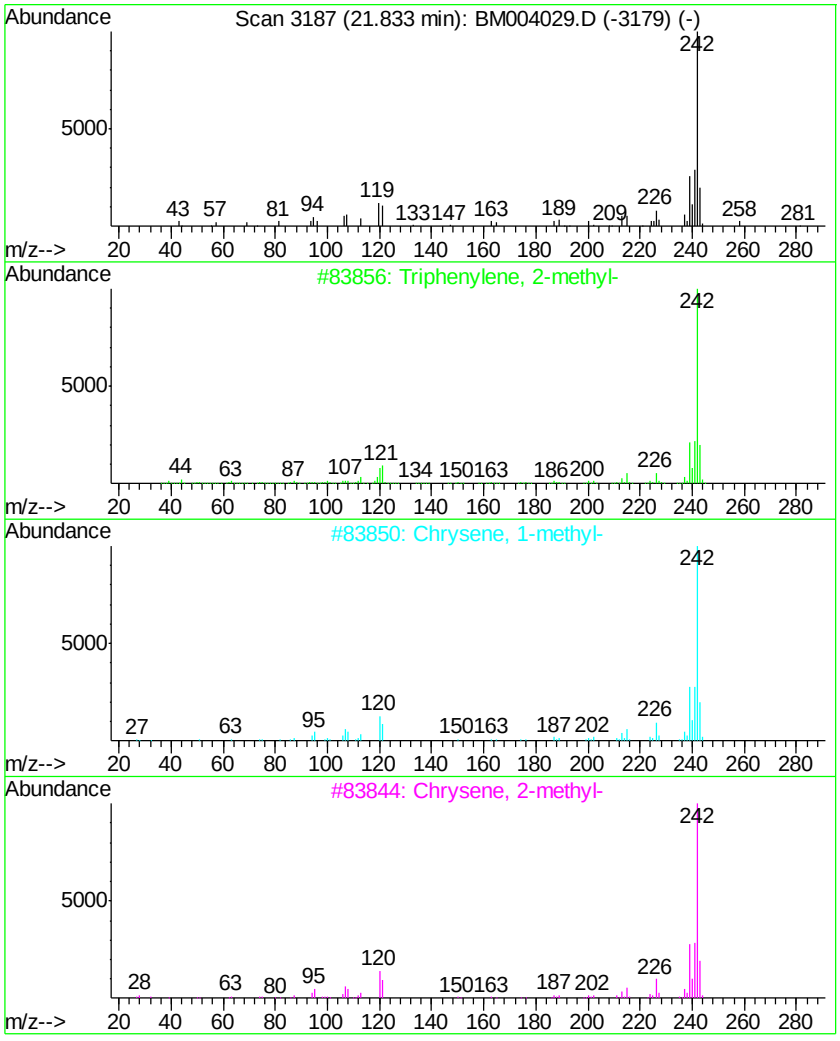
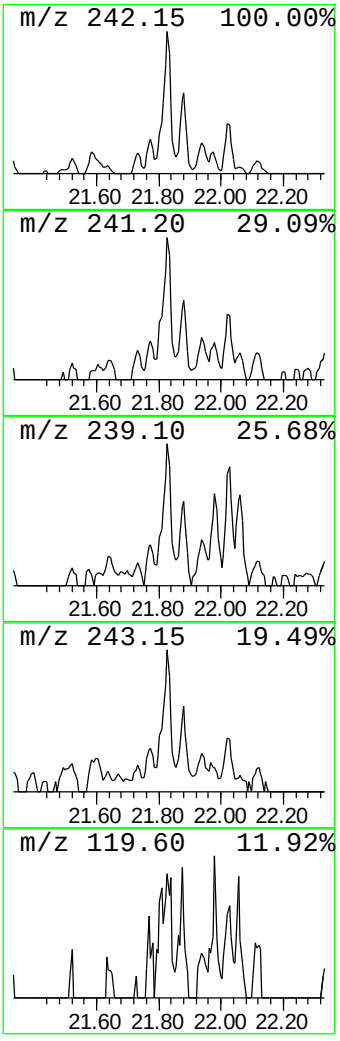
Peak Number 8 Triphenylene, 2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.83	2.02 ng/ul	95515	Chrysene-d12	21.28

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Triphenylene, 2-methyl-	242	C19H14	001705-84-6	91
2	Chrysene, 1-methyl-	242	C19H14	003351-28-8	90
3	Chrysene, 2-methyl-	242	C19H14	003351-32-4	89
4	Benz[a]anthracene, 1-methyl-	242	C19H14	002498-77-3	89
5	Benz[a]anthracene, 8-methyl-	242	C19H14	002381-31-9	86





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Data File : BM004029.D
Acq On : 15 Jan 2016 01:54
Operator : SJ/UM
Sample : H1099-09
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BC9A2

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM010116.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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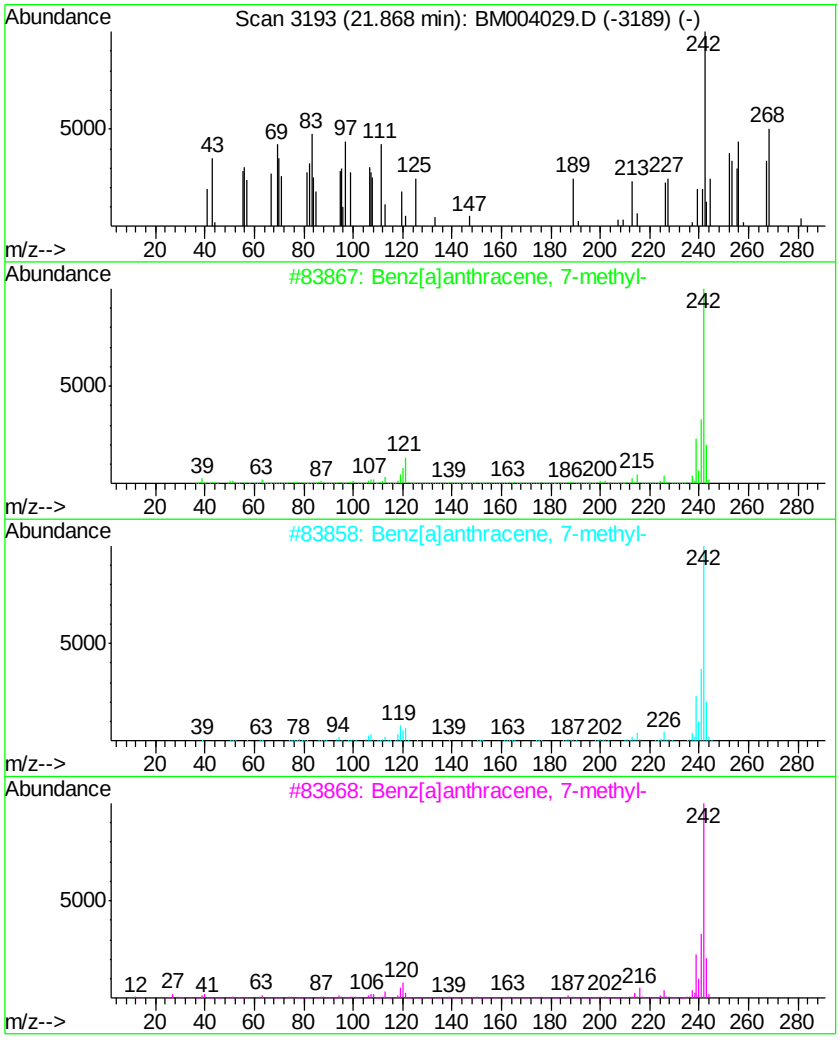
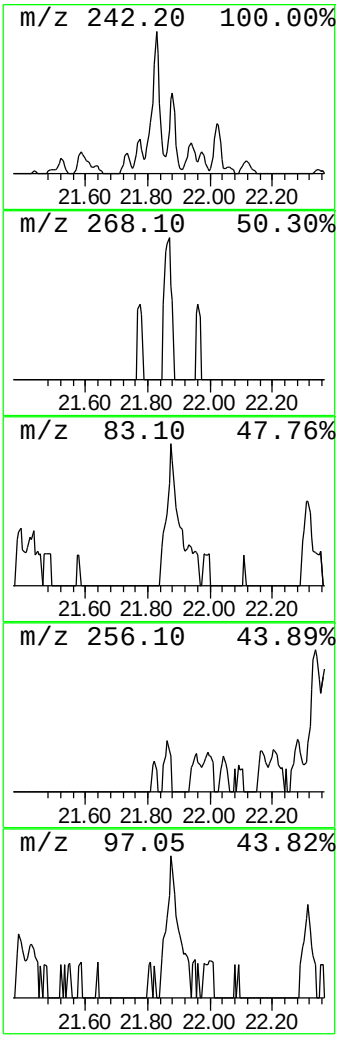
Peak Number 9 Benz[a]anthracene, 7-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.87	2.21 ng/ul	104343	Chrysene-d12	21.28

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benz[a]anthracene, 7-methyl-	242	C19H14	002541-69-7	70
2	Benz[a]anthracene, 7-methyl-	242	C19H14	002541-69-7	70
3	Benz[a]anthracene, 7-methyl-	242	C19H14	002541-69-7	58
4	Chrysene, 6-methyl-	242	C19H14	001705-85-7	58
5	Chrysene, 6-methyl-	242	C19H14	001705-85-7	58





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM011516\
Data File : BM004029.D
Acq On : 15 Jan 2016 01:54
Operator : SJ/UM
Sample : H1099-09
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BC9A2

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM010116.M
Quant Title : SVOA CALIBRATION

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

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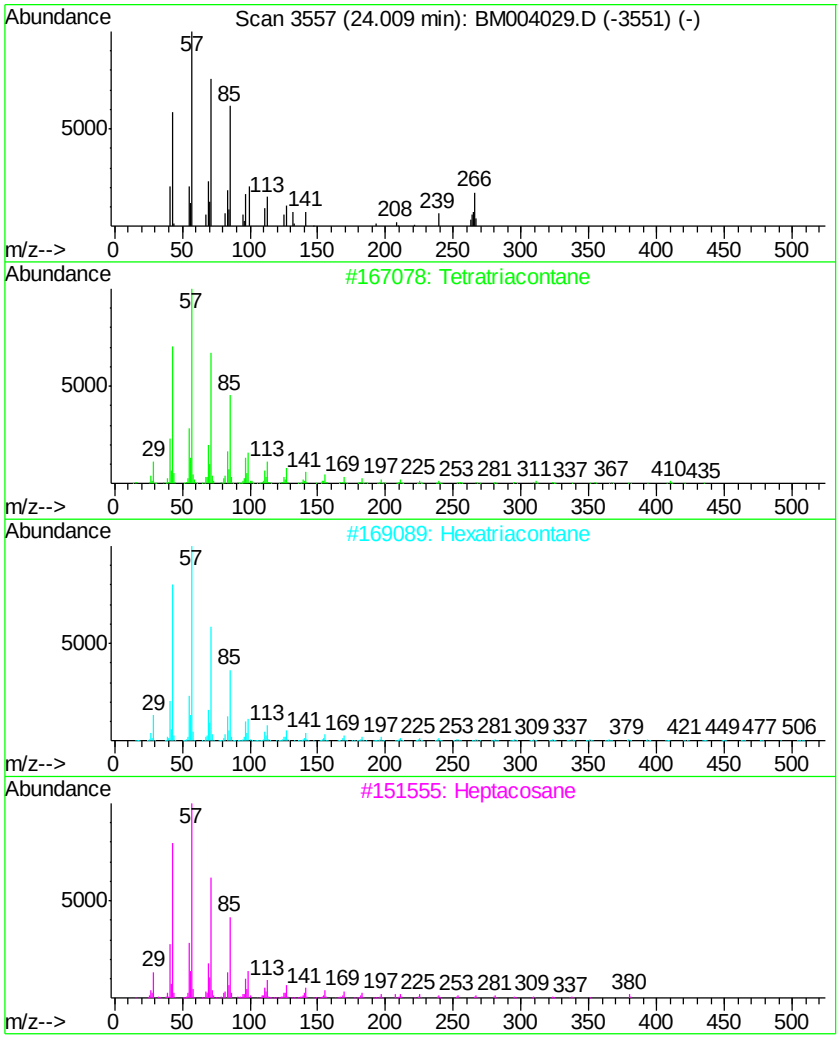
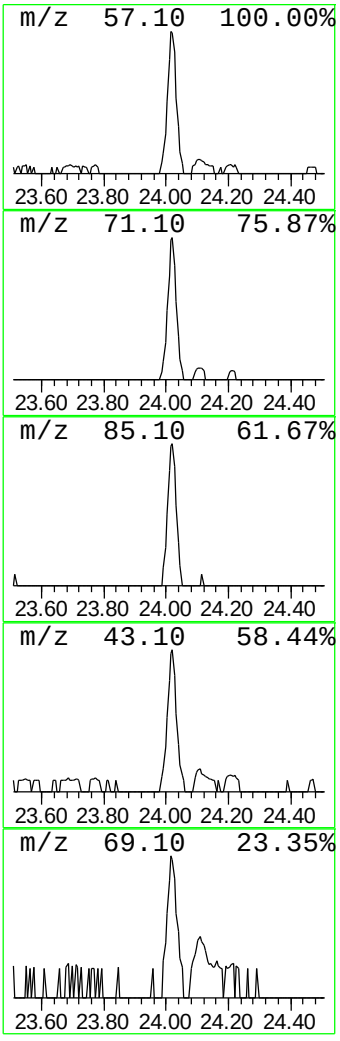
Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.01	3.04 ng/ul	83392	Perylene-d12	23.51

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetratriacontane	479	C34H70	014167-59-0	91
2	Hexatriacontane	507	C36H74	000630-06-8	91
3	Heptacosane	380	C27H56	000593-49-7	91
4	Octacosane	394	C28H58	000630-02-4	90
5	Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	90





Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM011516\
Data File : BM004029.D
Acq On : 15 Jan 2016 01:54
Operator : SJ/UM
Sample : H1099-09
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BC9A2

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM010116.M
Quant Title : SVOA CALIBRATION

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

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
unknown3.65	3.65	3.4	ng/ul	39084	1	7.66	230387	20.0
Anthracene, 2-met...	17.94	2.6	ng/ul	84499	4	17.07	649602	20.0
Phenanthrene, 1-m...	18.00	2.8	ng/ul	89468	4	17.07	649602	20.0
Phenanthrene, 4-m...	18.14	3.8	ng/ul	122175	4	17.07	649602	20.0
Phenanthrene, 2,5...	18.87	2.7	ng/ul	86724	4	17.07	649602	20.0
Pyrene, 1-methyl-	20.00	2.1	ng/ul	99175	5	21.28	945749	20.0
Heptafluorobutano...	20.99	2.4	ng/ul	113967	5	21.28	945749	20.0
Triphenylene, 2-m...	21.83	2.0	ng/ul	95515	5	21.28	945749	20.0
Benz[a]anthracene...	21.87	2.2	ng/ul	104343	5	21.28	945749	20.0
(DEL) Alkane: Str...	24.01	3.0	ng/ul	83392	6	23.51	547972	20.0