LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM011516\
Data File : BM004035.D
Acq On : 15 Jan 2016 05:29
Operator : SJ/UM
Sample : H1099-12
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BC9D8

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	3.652	92	96	102	rBV	24406	30573	3.41%	0.234%
2	4.787	284	289	295	rVB2	16032	22949	2.56%	0.175%
3	6.845	632	639	649	rVB	121141	196574	21.90%	1.502%
4	6.998	660	665	674	rBB	101259	154833	17.25%	1.183%
5	7.198	693	699	707	rBV	131048	199872	22.27%	1.527%
6	7.663	772	778	785	rBB	151391	230627	25.69%	1.762%
7	8.375	894	899	908	rBB	78592	126581	14.10%	0.967%
8	8.486	913	918	924	rBB	13320	20601	2.30%	0.157%
9	8.822	969	975	985	rBB	128066	199875	22.27%	1.527%
10	9.539	1092	1097	1104	rBB	106318	167268	18.64%	1.278%
11	10.081	1183	1189	1200	rBB	158409	256627	28.59%	1.961%
12	10.451	1245	1252	1264	rBB	240136	388793	43.32%	2.970%
13	10.598	1271	1277	1292	rBV	72637	126798	14.13%	0.969%
14	11.980	1508	1512	1519	rBB	15582	23563	2.63%	0.180%
15	13.733	1804	1810	1814	rBV	314300	442162	49.26%	3.378%
16	13.774	1814	1817	1825	rVB	144238	195106	21.74%	1.491%
17	14.010	1851	1857	1869	rBB	377738	549404	61.21%	4.198%
18	14.216	1888	1892	1897	rBB	19100	24792	2.76%	0.189%
19	14.321	1903	1910	1917	rBB2	370670	538869	60.04%	4.117%
20	14.539	1942	1947	1959	rBV	101724	165600	18.45%	1.265%
21	14.686	1965	1972	1975	rBV4	5322	10320	1.15%	0.079%
22	15.174	2050	2055	2061	rVB	38086	47600	5.30%	0.364%
23	15.315	2073	2079	2085	rVV	485745	691969	77.09%	5.287%
24	15.374	2085	2089	2094	rVB3	10596	15247	1.70%	0.116%
25	15.451	2098	2102	2109	rBV	90701	118095	13.16%	0.902%
26	15.557	2115	2120	2122	rBV2	9524	13612	1.52%	0.104%
27	15.580	2122	2124	2128	rVV	11675	13378	1.49%	0.102%
28	15.792	2155	2160	2165	rBV4	5818	10229	1.14%	0.078%
29	16.039	2196	2202	2215	rBV	42547	78694	8.77%	0.601%
30	16.374	2252	2259	2263	rBV2	7903	16243	1.81%	0.124%
31	16.609	2296	2299	2302	rVV3	9259	11186	1.25%	0.085%
32	16.827	2332	2336	2340	rVV	19378	24740	2.76%	0.189%
33	16.886	2341	2346	2353	rVB3	8088	13553	1.51%	0.104%
34	17.068	2371	2377	2381	rBV	440375	631090	70.31%	4.822%

LSC Area Percent Report

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Data File : BM004035.D
Acq On : 15 Jan 2016 05:29
Operator : SJ/UM
Sample : H1099-12
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BC9D8

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	17.109	2381	2384	2389	rVB	134888	177056	19.73%	1.353%
36	17.168	2389	2394	2399	rBV2	510272	725903	80.87%	5.546%
37	17.945	2521	2526	2528	rVV	35490	47553	5.30%	0.363%
38	17.992	2532	2534	2539	rVV	46494	57151	6.37%	0.437%
39	18.074	2544	2548	2553	rVV2	9687	13628	1.52%	0.104%
40	18.139	2553	2559	2563	rVV2	48486	86233	9.61%	0.659%
41	18.180	2563	2566	2571	rVB	28101	36688	4.09%	0.280%
42	18.451	2606	2612	2616	rBV	21452	27659	3.08%	0.211%
43	18.498	2616	2620	2625	rVB2	19608	27157	3.03%	0.207%
44	18.698	2651	2654	2659	rVV	8804	11136	1.24%	0.085%
45	18.751	2659	2663	2666	rVV	11731	13790	1.54%	0.105%
46	18.792	2666	2670	2673	rVB2	6431	9907	1.10%	0.076%
47	18.874	2678	2684	2689	rBV	33033	55848	6.22%	0.427%
48	18.933	2689	2694	2698	rVV3	17319	35923	4.00%	0.274%
49	18.968	2698	2700	2703	rVV	9599	9366	1.04%	0.072%
50	19.015	2703	2708	2720	rVB5	18885	49090	5.47%	0.375%
51	19.139	2720	2729	2734	rBV	282525	373833	41.65%	2.856%
52	19.239	2743	2746	2751	rVB3	8771	11290	1.26%	0.086%
53	19.474	2780	2786	2789	rBV2	636294	897579	100.00%	6.858%
54	19.503	2789	2791	2799	rVB	294690	344661	38.40%	2.633%
55	19.580	2799	2804	2811	rVB2	16731	25651	2.86%	0.196%
56	19.686	2817	2822	2827	rBV3	8750	12821	1.43%	0.098%
57	19.745	2827	2832	2837	rVB4	10373	17967	2.00%	0.137%
58	19.833	2840	2847	2854	rBV4	25051	62002	6.91%	0.474%
59	19.968	2865	2870	2871	rBV3	19468	24239	2.70%	0.185%
60	20.003	2871	2876	2883	rVB2	49734	88185	9.82%	0.674%
61	20.103	2889	2893	2897	rVV2	32685	40329	4.49%	0.308%
62	20.162	2897	2903	2907	rVV3	39923	63029	7.02%	0.482%
63	20.203	2907	2910	2914	rVB4	8333	10582	1.18%	0.081%
64	20.303	2922	2927	2931	rBV	30673	41805	4.66%	0.319%
65	20.350	2931	2935	2939	rVV	29290	37771	4.21%	0.289%
66	20.386	2939	2941	2948	rVB5	6548	11699	1.30%	0.089%
67	20.468	2948	2955	2959	rBV8	6574	14199	1.58%	0.108%
68	20.521	2959	2964	2968	rBV5	8103	14580	1.62%	0.111%
69	20.627	2979	2982	2987	rVB3	11540	14881	1.66%	0.114%
70	20.721	2993	2998	3000	rBV4	8409	13508	1.50%	0.103%
71	20.756	3000	3004	3010	rVB	25524	39114	4.36%	0.299%

LSC Area Percent Report

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Operator : SJ/UM
Sample : H1099-12
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BC9D8

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

72	20.921	3024	3032	3035	rBV3	25683	50116	5.58%	0.383%
73	20.980	3035	3042	3050	rVV6	122609	230570	25.69%	1.762%
74	21.050	3051	3054	3062	rVB2	22897	39164	4.36%	0.299%
75	21.186	3074	3077	3082	rVB2	16723	20185	2.25%	0.154%
76	21.274	3085	3092	3096	rBV2	545691	857976	95.59%	6.555%
77	21.309	3096	3098	3108	rVB	135821	177055	19.73%	1.353%
78	21.392	3108	3112	3115	rBV6	14407	20828	2.32%	0.159%
79	21.427	3115	3118	3125	rBV4	21355	31978	3.56%	0.244%
80	21.592	3141	3146	3151	rBV4	14549	25890	2.88%	0.198%
81	21.633	3151	3153	3157	rBV3	8874	10279	1.15%	0.079%
82	21.827	3179	3186	3189	rBV	39099	68654	7.65%	0.525%
83	21.874	3189	3194	3202	rVB6	37417	92552	10.31%	0.707%
84	21.980	3209	3212	3215	rBV2	6858	10149	1.13%	0.078%
85	22.021	3215	3219	3223	rBV2	23389	40629	4.53%	0.310%
86	22.121	3232	3236	3241	rVB3	13776	23190	2.58%	0.177%
87	22.315	3264	3269	3273	rBV4	19657	34887	3.89%	0.267%
88	22.444	3285	3291	3295	rBV	84208	122581	13.66%	0.937%
89	22.815	3349	3354	3356	rBV	76482	114153	12.72%	0.872%
90	23.015	3384	3388	3393	rVB	15614	23815	2.65%	0.182%
91	23.321	3434	3440	3443	rBV	48771	89420	9.96%	0.683%
92	23.374	3443	3449	3453	rVV	352126	646219	72.00%	4.937%
93	23.415	3453	3456	3462	rVB	81211	129254	14.40%	0.988%
94	23.515	3466	3473	3479	rBV2	276747	528672	58.90%	4.039%
95	24.015	3552	3558	3568	rVB	55092	111664	12.44%	0.853%
96	24.756	3679	3684	3691	rBV2	12259	24361	2.71%	0.186%
97	25.609	3825	3829	3837	rVB4	13090	28580	3.18%	0.218%
98	25.756	3846	3854	3865	rBV	35493	99516	11.09%	0.760%
99	26.450	3966	3972	3981	rVB2	19922	52734	5.88%	0.403%
100	27.809	4196	4203	4214	rVB10	24440	82390	9.18%	0.629%

Sum of corrected areas: 13088697

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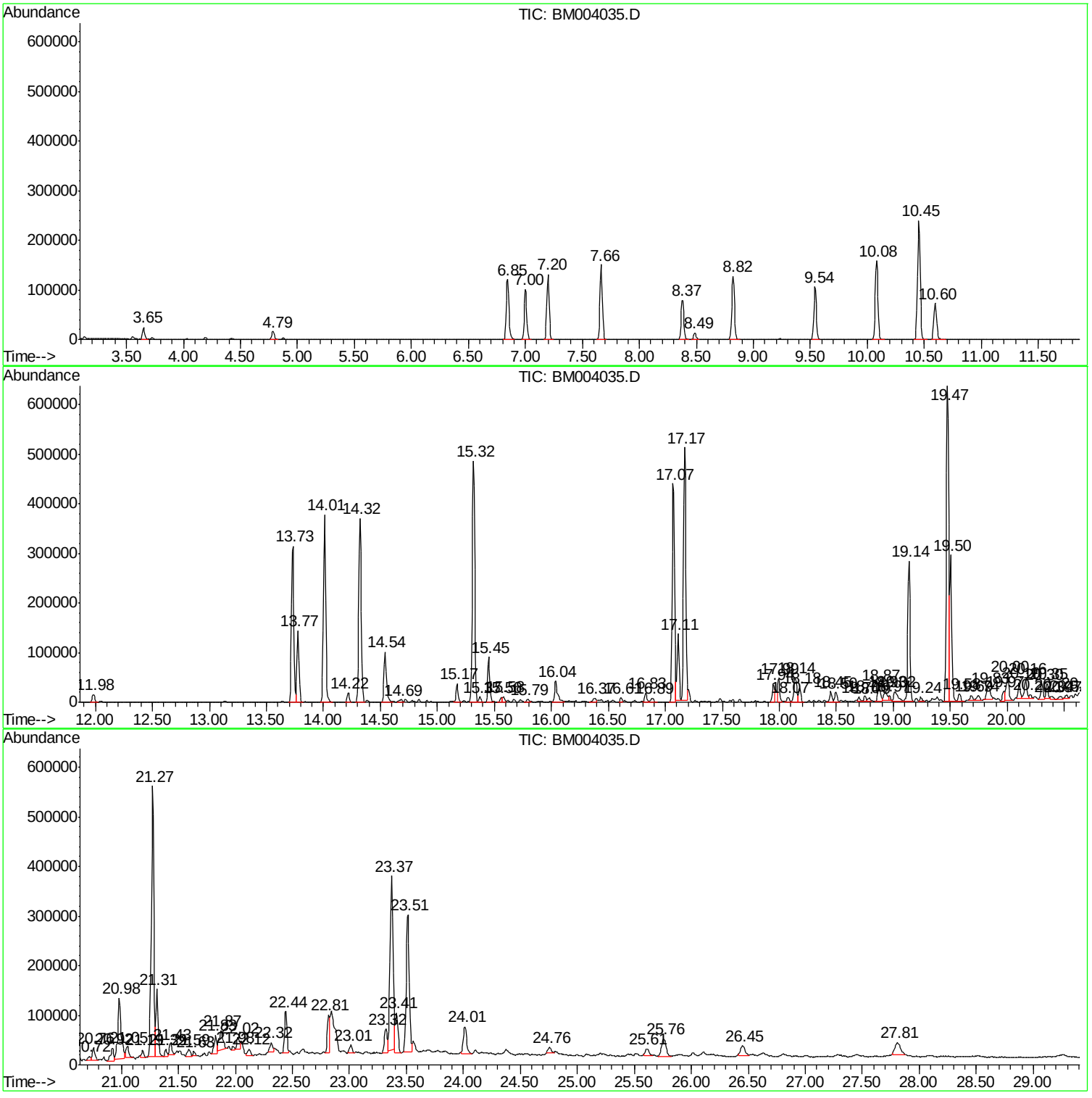
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BC9D8

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

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM011516\
Data File : BM004035.D
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Instrument : BNA_M
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
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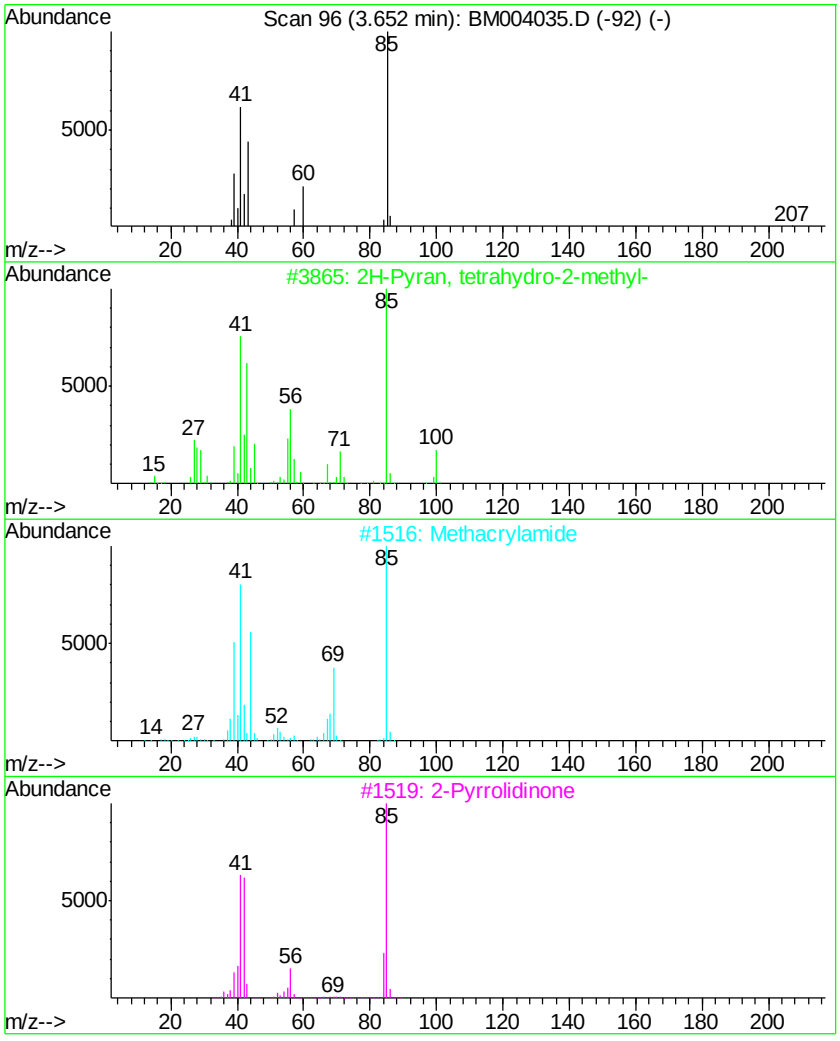
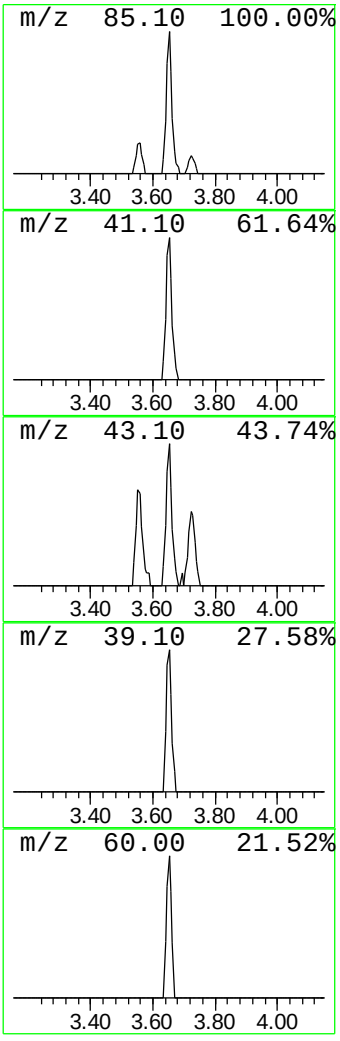
****

Peak Number 1 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.65	2.65 ng/ul	30573	1,4-Dichlorobenzene-d4	7.66

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2H-Pyran, tetrahydro-2-methyl-	100	C6H12O	010141-72-7	36
2	Methacrylamide	85	C4H7NO	000079-39-0	33
3	2-Pyrrolidinone	85	C4H7NO	000616-45-5	9
4	2-Furanmethanol, tetrahydro-5-me...	116	C6H12O2	054774-28-6	9
5	2-Pyrrolidinone	85	C4H7NO	000616-45-5	7





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM011516\
Data File : BM004035.D
Acq On : 15 Jan 2016 05:29
Operator : SJ/UM
Sample : H1099-12
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BC9D8

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

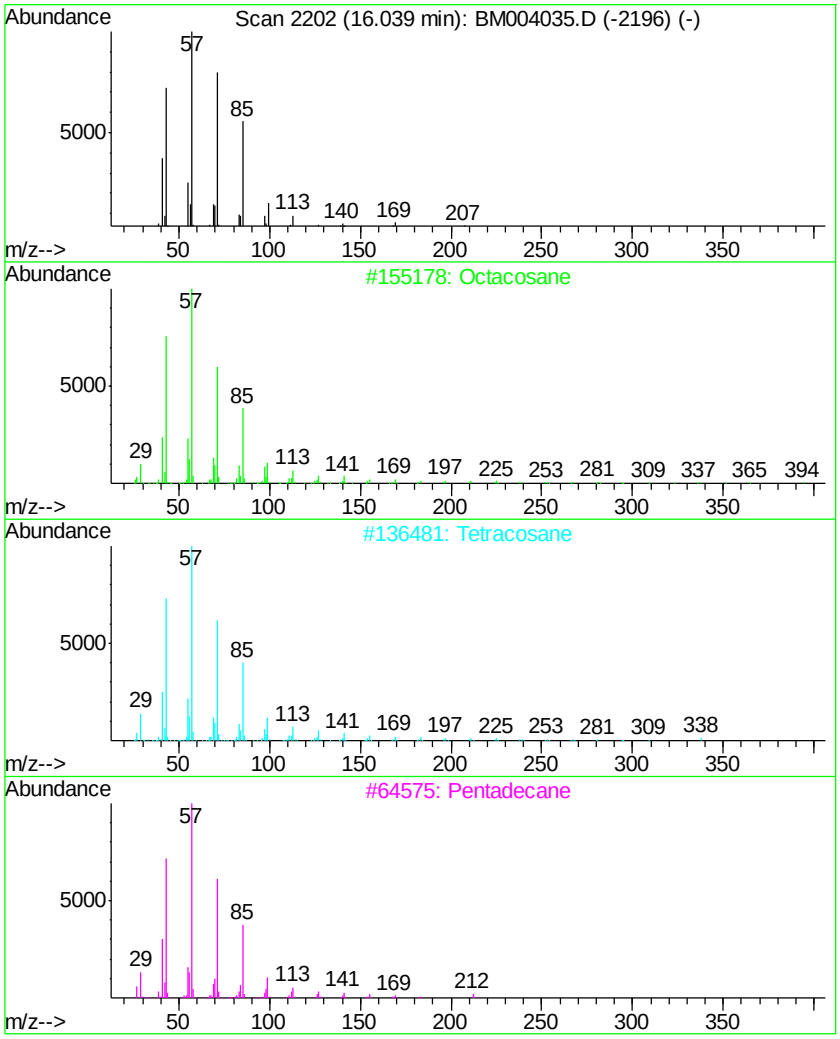
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

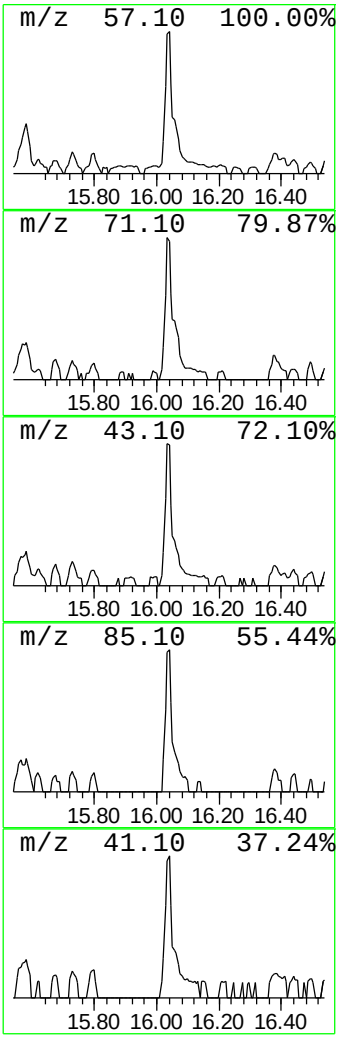
****

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.04	2.49 ng/ul	78694	Phenanthrene-d10	17.07

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octacosane	394	C28H58	000630-02-4	87
2	Tetracosane	338	C24H50	000646-31-1	86
3	Pentadecane	212	C15H32	000629-62-9	86
4	Pentadecane	212	C15H32	000629-62-9	86
5	Tetradecane	198	C14H30	000629-59-4	86





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM011516\
Data File : BM004035.D
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Operator : SJ/UM
Sample : H1099-12
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BC9D8

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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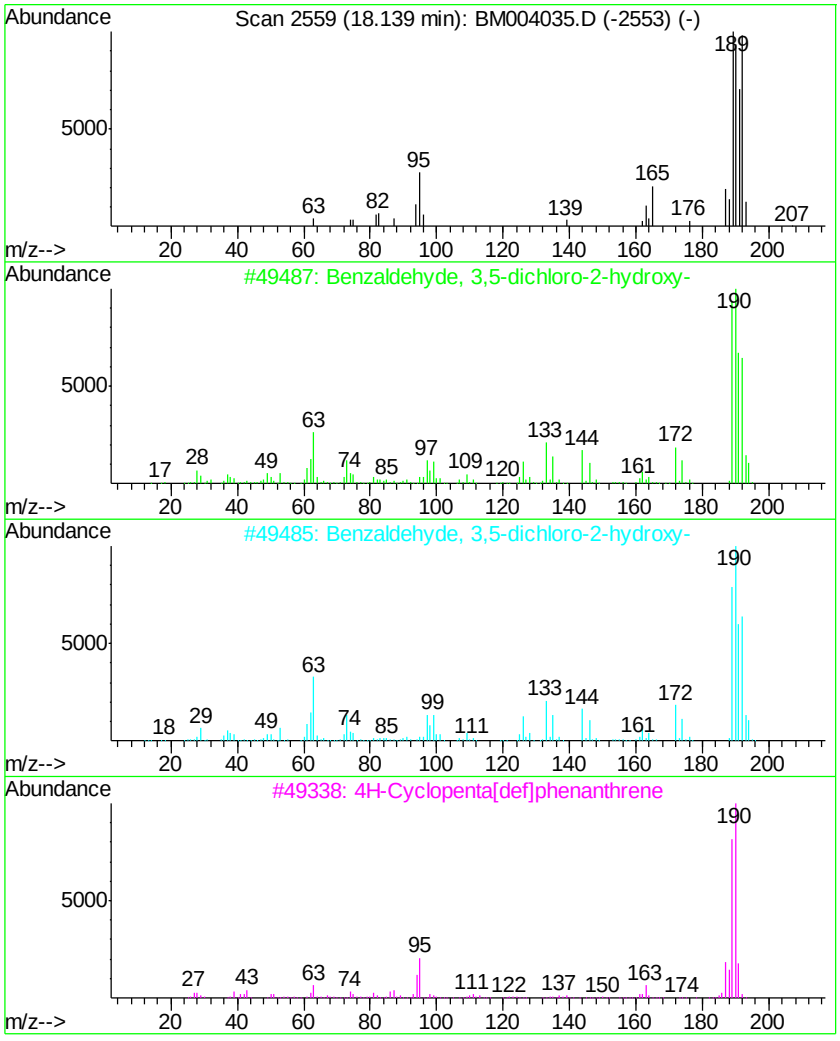
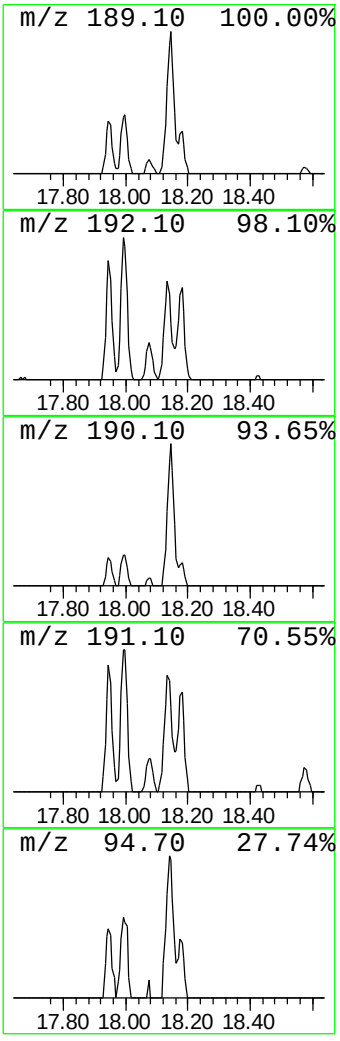
Peak Number 3 Benzaldehyde, 3,5-dichloro-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.14	2.73 ng/ul	86233	Phenanthrene-d10	17.07

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	52
2	Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	52
3	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	49
4	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	43
5	Anthracene, 1-methyl-	192	C15H12	000610-48-0	41





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM011516\
Data File : BM004035.D
Acq On : 15 Jan 2016 05:29
Operator : SJ/UM
Sample : H1099-12
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BC9D8

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

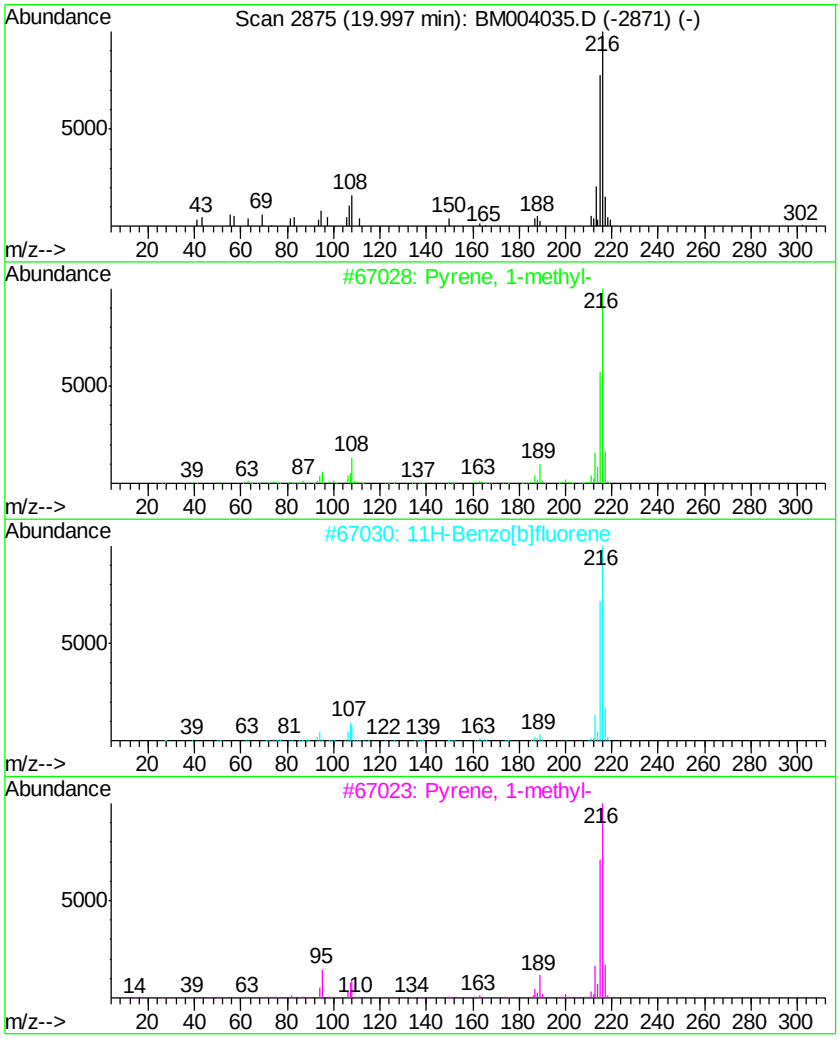
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

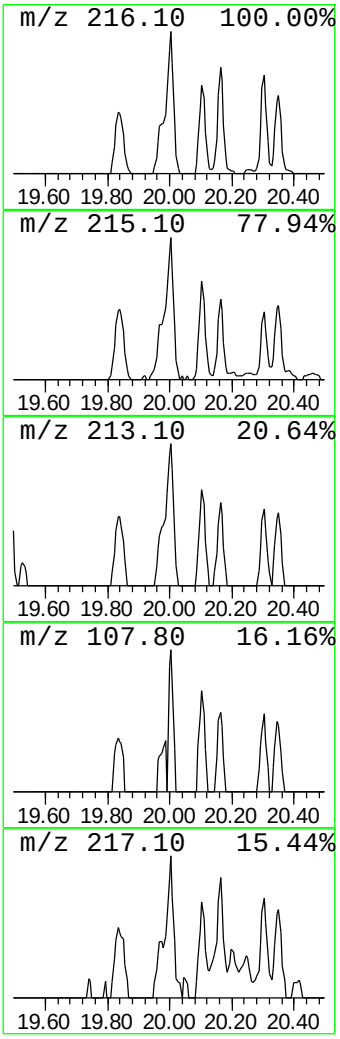
****

Peak Number 4 Pyrene, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.00	2.06 ng/ul	88185	Chrysene-d12	21.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	94
2	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	91
3	Pyrene, 1-methyl-	216	C17H12	002381-21-7	91
4	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	90
5	Pyrene, 1-methyl-	216	C17H12	002381-21-7	90





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM011516\
Data File : BM004035.D
Acq On : 15 Jan 2016 05:29
Operator : SJ/UM
Sample : H1099-12
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BC9D8

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

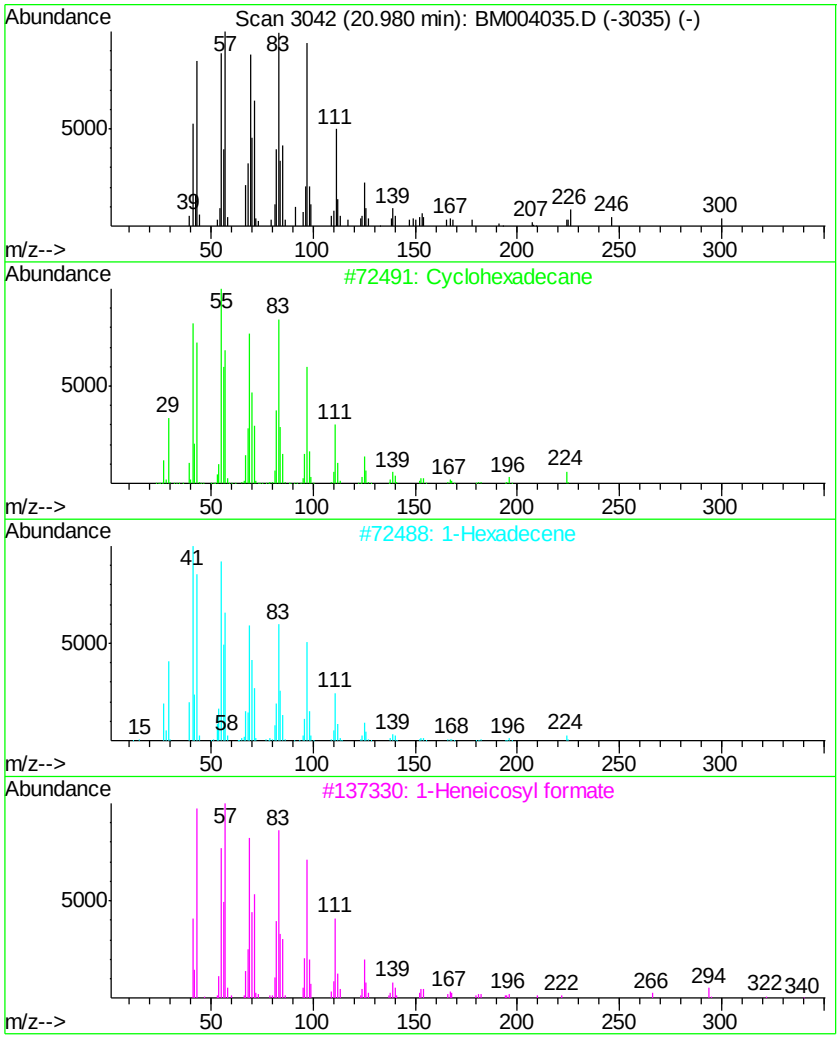
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

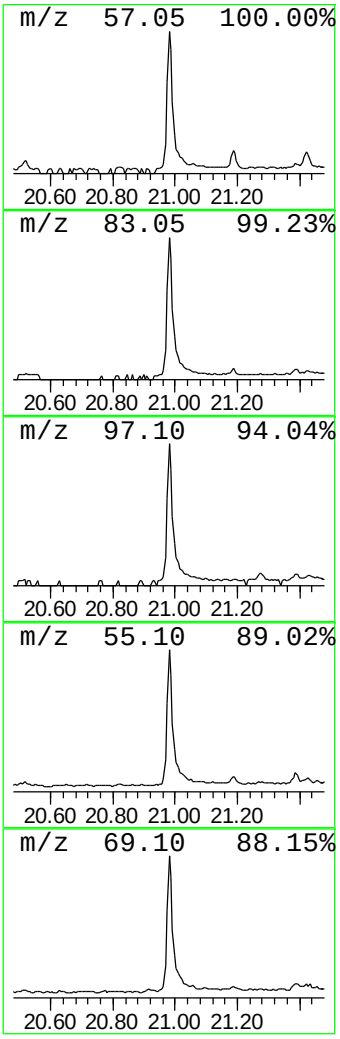
****

Peak Number 5 (DEL) Alkane: Cyclic20.98 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.98	5.37 ng/ul	230570	Chrysene-d12	21.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexadecane	224	C16H32	000295-65-8	98
2	1-Hexadecene	224	C16H32	000629-73-2	94
3	1-Heneicosyl formate	340	C22H44O2	077899-03-7	94
4	Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	94
5	2- Chloropropionic acid, hexadec...	332	C19H37ClO2	086711-81-1	94





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM011516\
Data File : BM004035.D
Acq On : 15 Jan 2016 05:29
Operator : SJ/UM
Sample : H1099-12
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BC9D8

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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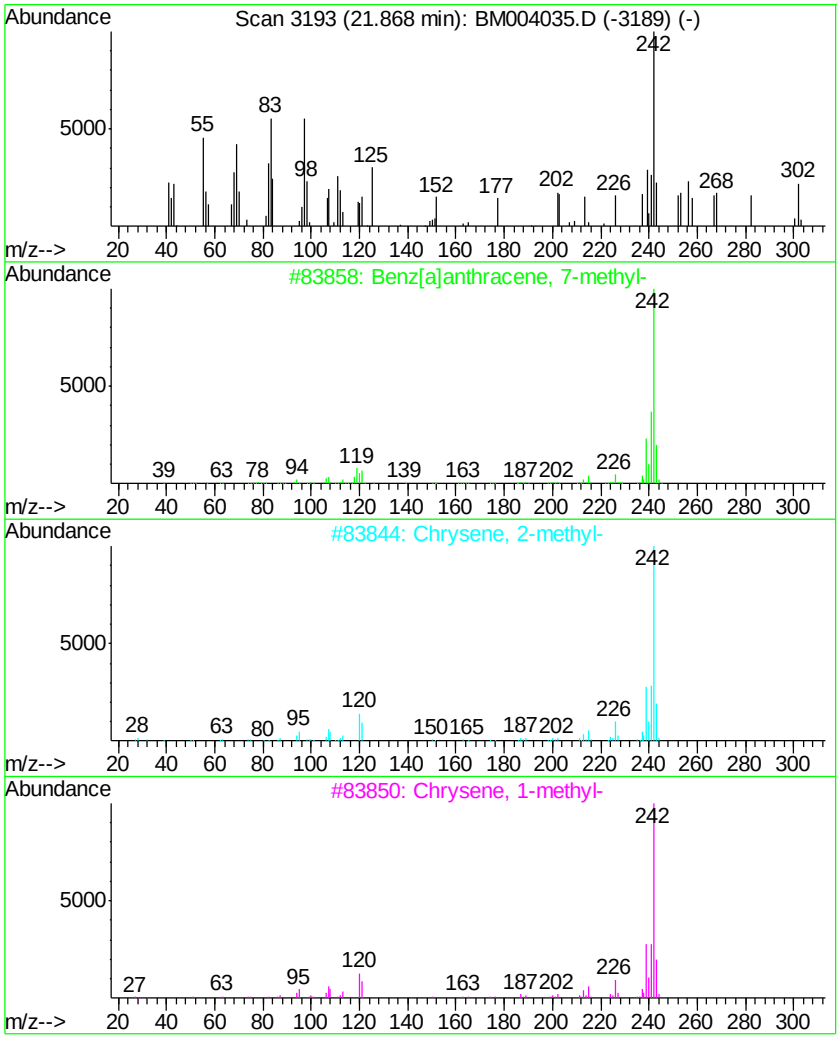
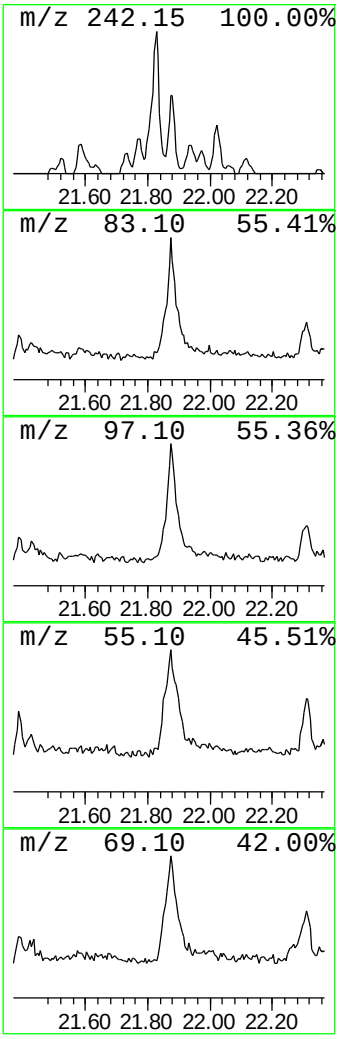
Peak Number 6 Benz[a]anthracene, 7-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.87	2.16 ng/ul	92552	Chrysene-d12	21.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benz[a]anthracene, 7-methyl-	242	C19H14	002541-69-7	60
2	Chrysene, 2-methyl-	242	C19H14	003351-32-4	60
3	Chrysene, 1-methyl-	242	C19H14	003351-28-8	60
4	Chrysene, 6-methyl-	242	C19H14	003697-24-3	58
5	Chrysene, 6-methyl-	242	C19H14	001705-85-7	55





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM011516\
Data File : BM004035.D
Acq On : 15 Jan 2016 05:29
Operator : SJ/UM
Sample : H1099-12
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BC9D8

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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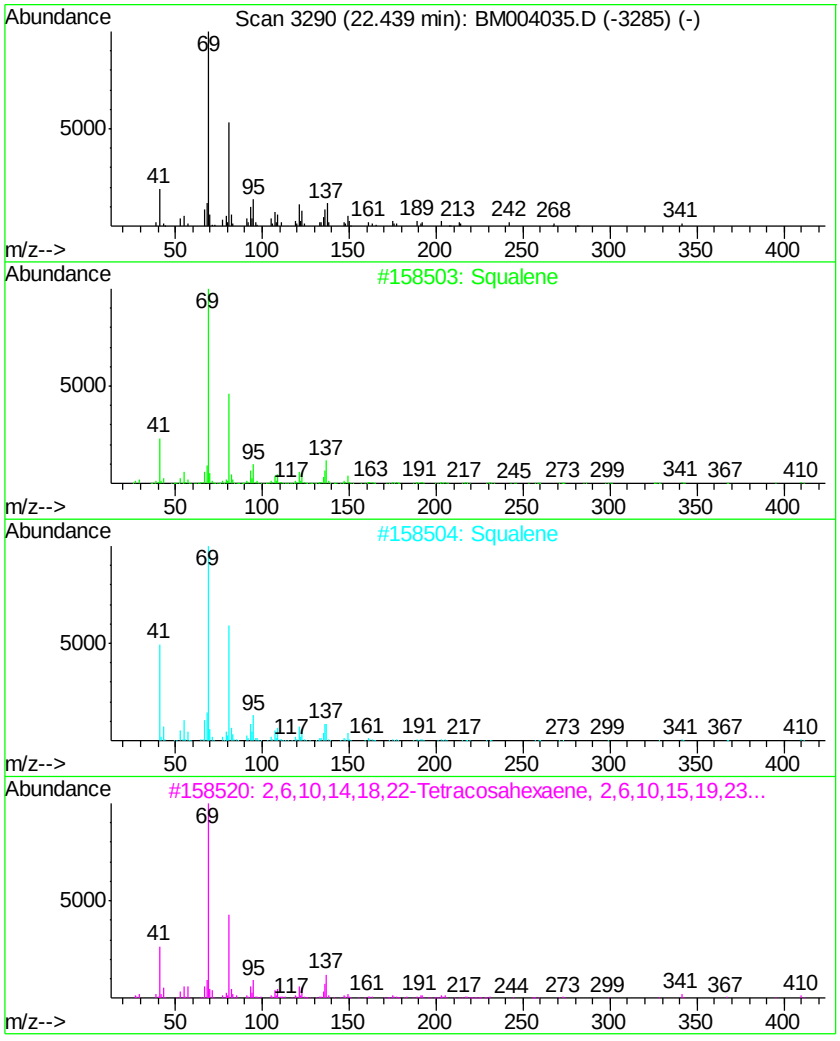
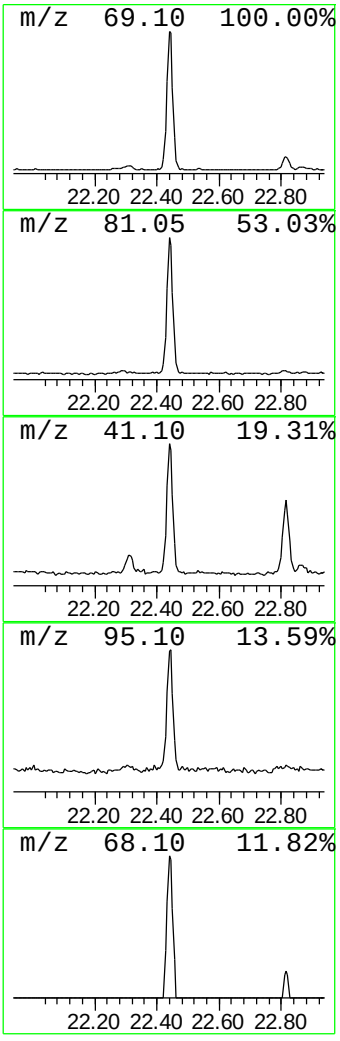
Peak Number 7 Squalene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.44	4.64 ng/ul	122581	Perylene-d12	23.51

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Squalene	410	C30H50	007683-64-9	91
2	Squalene	410	C30H50	007683-64-9	91
3	2,6,10,14,18,22-Tetracosahexaene...	410	C30H50	000111-02-4	91
4	1,5,9-Undecatriene, 2,6,10-trime...	192	C14H24	062951-96-6	87
5	Docosa-2,6,10,14,18-pentaen-22-a...	384	C27H44O	1000163-04-7	83





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM011516\
Data File : BM004035.D
Acq On : 15 Jan 2016 05:29
Operator : SJ/UM
Sample : H1099-12
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BC9D8

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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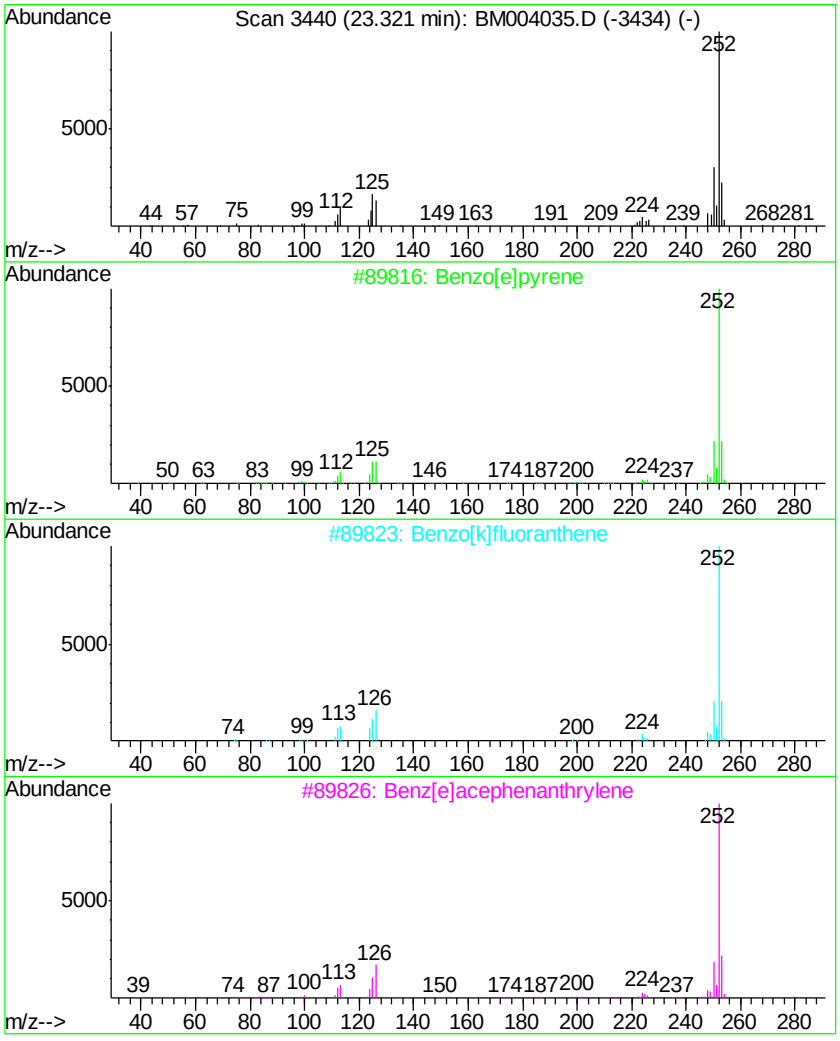
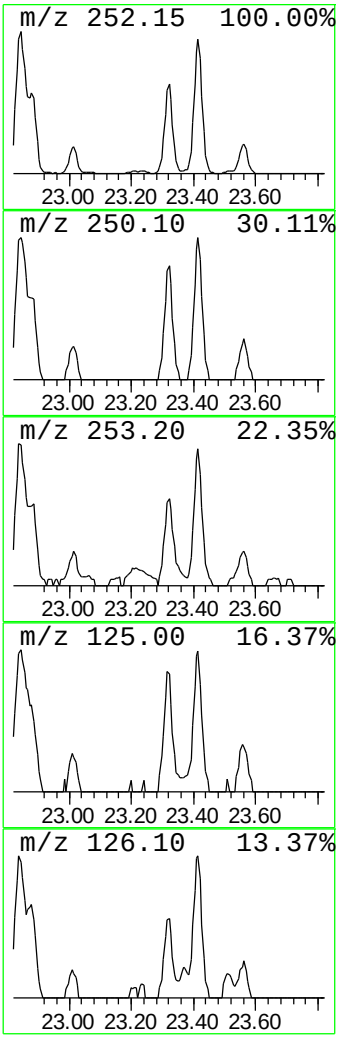
Peak Number 8 Benzo[e]pyrene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.32	3.38 ng/ul	89420	Perylene-d12	23.51

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]pyrene	252	C20H12	000192-97-2	97
2	Benzo[k]fluoranthene	252	C20H12	000207-08-9	96
3	Benz[e]acephenanthrylene	252	C20H12	000205-99-2	95
4	Perylene	252	C20H12	000198-55-0	95
5	Perylene	252	C20H12	000198-55-0	94





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM011516\
Data File : BM004035.D
Acq On : 15 Jan 2016 05:29
Operator : SJ/UM
Sample : H1099-12
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BC9D8

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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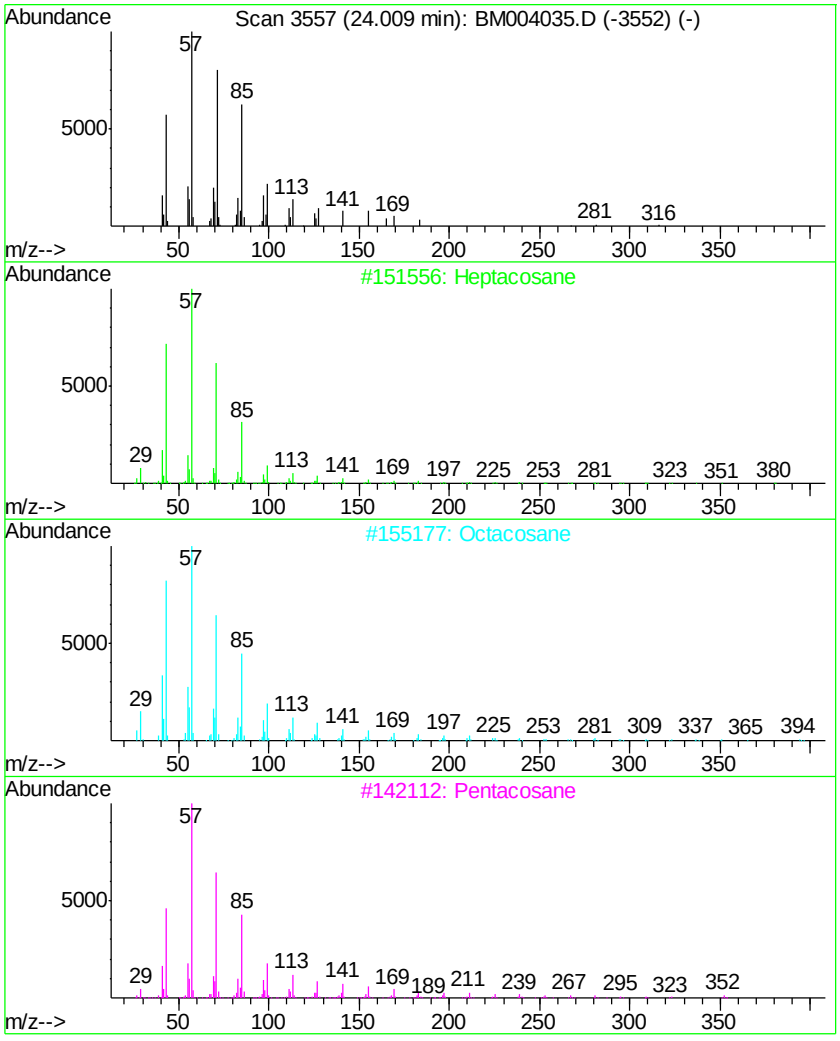
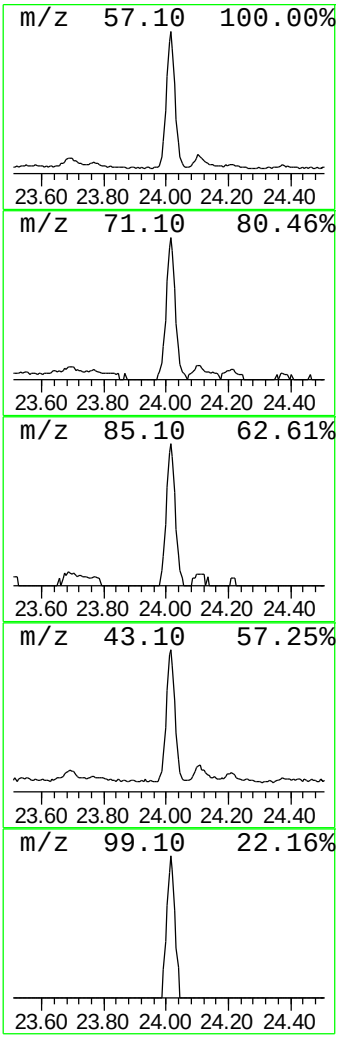
Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.01	4.22 ng/ul	111664	Perylene-d12	23.51

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacosane	380	C27H56	000593-49-7	91
2	Octacosane	394	C28H58	000630-02-4	91
3	Pentacosane	352	C25H52	000629-99-2	91
4	Hexatriacontane	507	C36H74	000630-06-8	91
5	Heptacosane	380	C27H56	000593-49-7	91





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM011516\
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Acq On : 15 Jan 2016 05:29
Operator : SJ/UM
Sample : H1099-12
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BC9D8

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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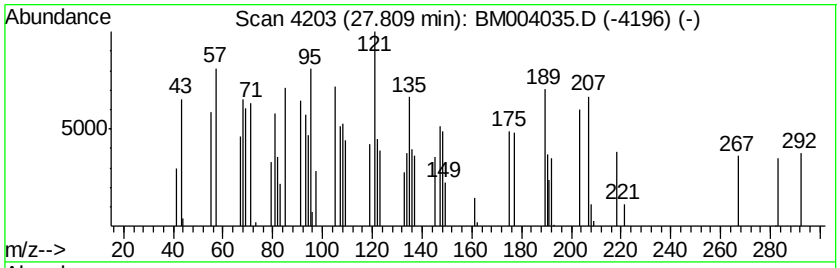
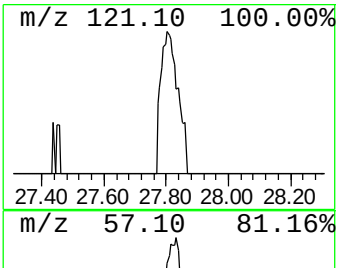
Peak Number 10 7.alpha.-Ethyl-8.beta.-hydr... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.81	3.12 ng/ul	82390	Perylene-d12	23.51

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	7.alpha.-Ethyl-8.beta.-hydroxy-2...	208	C14H24O	1000077-92-4	64
2	2-(Pyridin-2-ylamino)-cyclohexanol	192	C11H16N2O	1000194-39-2	27
3	2H-Cyclopropa[g]benzofuran, 4,5,...	218	C15H22O	102681-49-2	22
4	2-Azidomethyl-1,3,3-trimethyl-cy...	179	C10H17N3	090073-44-2	22
5	Ambrein	428	C30H52O	1000292-95-3	14





Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM011516\
Data File : BM004035.D
Acq On : 15 Jan 2016 05:29
Operator : SJ/UM
Sample : H1099-12
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BC9D8

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
unknown-01	3.65	2.6	ng/ul	30573	1	7.66	230627	20.0
(DEL) Alkane: Str...	16.04	2.5	ng/ul	78694	4	17.07	631090	20.0
Benzaldehyde, 3,5...	18.14	2.7	ng/ul	86233	4	17.07	631090	20.0
Pyrene, 1-methyl-	20.00	2.1	ng/ul	88185	5	21.27	857976	20.0
(DEL) Alkane: Cyc...	20.98	5.4	ng/ul	230570	5	21.27	857976	20.0
Benz[a]anthracene...	21.87	2.2	ng/ul	92552	5	21.27	857976	20.0
Squalene	22.44	4.6	ng/ul	122581	6	23.51	528672	20.0
Benzo[e]pyrene	23.32	3.4	ng/ul	89420	6	23.51	528672	20.0
(DEL) Alkane: Str...	24.01	4.2	ng/ul	111664	6	23.51	528672	20.0
7.alpha.-Ethyl-8....	27.81	3.1	ng/ul	82390	6	23.51	528672	20.0