LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM042018\
Data File : BM014768.D
Acq On : 20 Apr 2018 18:06
Operator : SJ/JU
Sample : J2434-16
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AD1

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 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\SOM-EPA-BM041918.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	3.316	3	5	10	rVB	415507	498875	3.81%	0.309%
2	3.369	10	14	20	rVV	1029168	1337235	10.21%	0.827%
3	3.422	20	23	38	rVB	140548	220763	1.69%	0.137%
4	4.263	161	166	176	rVB	382550	531228	4.06%	0.329%
5	5.516	373	379	388	rVB4	71541	145273	1.11%	0.090%
6	7.663	737	744	760	rVB2	926089	1806859	13.80%	1.118%
7	7.846	765	775	784	rBV	785527	1281444	9.79%	0.793%
8	8.063	804	812	817	rBV	953894	1581847	12.08%	0.979%
9	8.110	817	820	829	rVB	92836	153192	1.17%	0.095%
10	8.546	887	894	905	rVB	1194351	1983578	15.15%	1.227%
11	9.240	1005	1012	1020	rBV	807594	1323714	10.11%	0.819%
12	9.722	1087	1094	1103	rBV	926781	1628870	12.44%	1.008%
13	10.457	1211	1219	1230	rVB	729845	1247856	9.53%	0.772%
14	10.992	1303	1310	1318	rBV	1165567	1937274	14.79%	1.198%
15	11.392	1370	1378	1384	rBV	1551749	2715845	20.74%	1.680%
16	11.510	1391	1398	1418	rBV	831760	1551121	11.84%	0.960%
17	14.533	1905	1912	1916	rBV	1977029	2761430	21.09%	1.708%
18	14.580	1916	1920	1925	rVB	587843	854013	6.52%	0.528%
19	14.845	1958	1965	1978	rBV	2364013	3512505	26.82%	2.173%
20	14.992	1982	1990	1996	rVB	154474	208877	1.60%	0.129%
21	15.145	2005	2016	2022	rBV2	2148000	3494275	26.68%	2.162%
22	15.204	2022	2026	2035	rVB	125748	203170	1.55%	0.126%
23	15.292	2035	2041	2050	rBV	578078	899080	6.87%	0.556%
24	15.933	2145	2150	2155	rVB2	183247	262293	2.00%	0.162%
25	16.122	2175	2182	2188	rBV	2803456	4205949	32.12%	2.602%
26	16.216	2194	2198	2204	rVB	390967	507892	3.88%	0.314%
27	16.333	2213	2218	2230	rVB2	65577	177109	1.35%	0.110%
28	16.780	2290	2294	2309	rVB	160292	332174	2.54%	0.205%
29	17.857	2469	2477	2481	rBV	2643141	3848142	29.39%	2.381%
30	17.898	2481	2484	2488	rVV	818562	1098550	8.39%	0.680%
31	17.951	2488	2493	2504	rVB	3040139	4545463	34.71%	2.812%
32	18.245	2536	2543	2548	rBV2	164951	290399	2.22%	0.180%
33	18.568	2593	2598	2603	rBV3	88084	143057	1.09%	0.088%
34	18.680	2612	2617	2620	rBV	486469	667235	5.10%	0.413%

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Misc :
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Instrument : BNA_M
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Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
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Start Thrs: 0.2 Max Peaks: 100
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Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM041918.M
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35	18.757	2626	2630	2635	rVB2	224745	332781	2.54%	0.206%
36	18.910	2649	2656	2659	rBV3	230338	447755	3.42%	0.277%
37	19.598	2767	2773	2778	rBV3	74351	165090	1.26%	0.102%
38	19.862	2812	2818	2823	rVB	4289941	5808944	44.36%	3.594%
39	19.915	2823	2827	2830	rBV2	147871	192572	1.47%	0.119%
40	20.186	2867	2873	2876	rBV2	4020633	6458198	49.32%	3.995%
41	20.215	2876	2878	2885	rVV	4776597	5674278	43.33%	3.510%
42	20.274	2885	2888	2893	rVB	183807	230183	1.76%	0.142%
43	20.386	2903	2907	2912	rVB	234438	302600	2.31%	0.187%
44	20.451	2913	2918	2924	rVB	340436	476242	3.64%	0.295%
45	20.521	2925	2930	2936	rBV2	336816	721571	5.51%	0.446%
46	20.574	2936	2939	2944	rVB	158387	204046	1.56%	0.126%
47	20.686	2946	2958	2963	rVV	1074585	2166569	16.54%	1.340%
48	20.786	2970	2975	2979	rBV2	857144	1207131	9.22%	0.747%
49	20.845	2982	2985	2989	rVV	424019	601490	4.59%	0.372%
50	20.880	2989	2991	2995	rVB2	174413	182199	1.39%	0.113%
51	20.986	3005	3009	3012	rBV	461532	617297	4.71%	0.382%
52	21.027	3013	3016	3023	rVB	471939	642801	4.91%	0.398%
53	21.374	3072	3075	3079	rBV2	179004	310476	2.37%	0.192%
54	21.415	3080	3082	3089	rVB3	249320	404340	3.09%	0.250%
55	21.592	3107	3112	3114	rBV2	1119582	1500024	11.45%	0.928%
56	21.662	3120	3124	3128	rBV2	831641	1209389	9.24%	0.748%
57	21.803	3145	3148	3155	rBV2	239847	471592	3.60%	0.292%
58	21.927	3164	3169	3175	rBV2	5550619	11446914	87.41%	7.081%
59	21.980	3175	3178	3188	rVB	4968483	6442467	49.20%	3.985%
60	22.109	3196	3200	3206	rBV	645209	965359	7.37%	0.597%
61	22.215	3216	3218	3223	rVB	498861	664188	5.07%	0.411%
62	22.274	3224	3228	3234	rBV	962075	1423938	10.87%	0.881%
63	22.592	3279	3282	3287	rVB	467060	645806	4.93%	0.400%
64	22.756	3307	3310	3314	rBV3	372666	605621	4.62%	0.375%
65	23.680	3460	3467	3472	rBV	5306705	13095323	100.00%	8.101%
66	23.868	3495	3499	3505	rVB	704867	1220110	9.32%	0.755%
67	24.227	3553	3560	3565	rBV	3361884	7309538	55.82%	4.522%
68	24.280	3565	3569	3573	rVV2	2188186	4544497	34.70%	2.811%
69	24.333	3573	3578	3586	rVB	5147743	10043276	76.69%	6.213%
70	24.444	3591	3597	3601	rVV	1912607	4106639	31.36%	2.540%
71	24.497	3602	3606	3615	rVB2	1507683	3022905	23.08%	1.870%

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72	27.033	4028	4037	4048	rVB	2924758	9361664	71.49%	5.791%
73	27.838	4165	4174	4195	rVB	2369001	8467848	64.66%	5.238%

Sum of corrected areas: 161650248

LSC Report - Integrated Chromatogram

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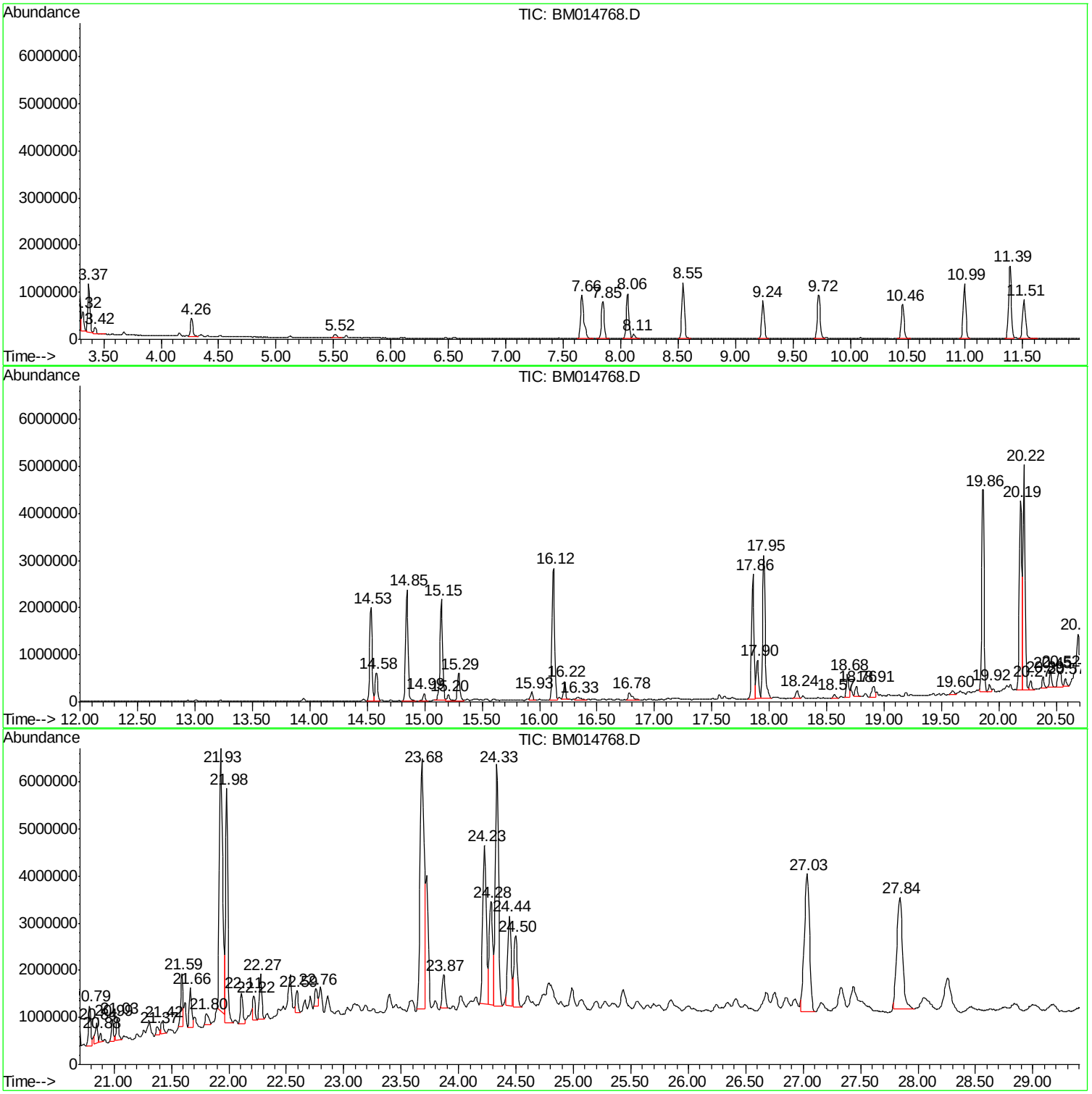
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AD1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM041918.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA M\DATA\BM042018\
Data File : BM014768.D
Acq On : 20 Apr 2018 18:06
Operator : SJ/JU
Sample : J2434-16
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AD1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM041918.M
Quant Title : SVOA CALIBRATION

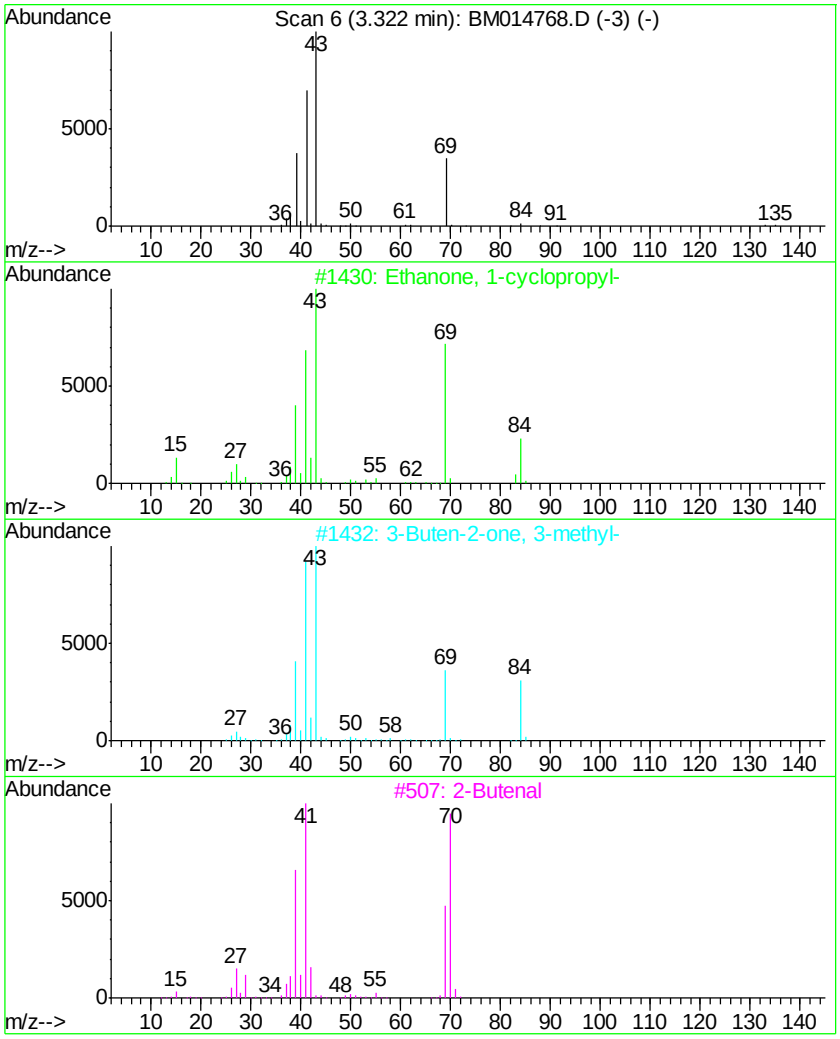
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

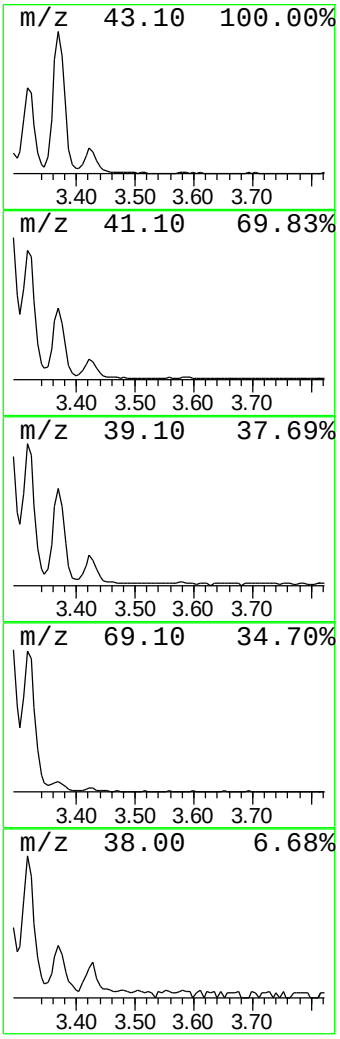
****

Peak Number 1 Ethanone, 1-cyclopropyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.32	5.03 ng/ul	498875	1,4-Dichlorobenzene-d4	8.55

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanone, 1-cyclopropyl-	84	C5H8O	000765-43-5	72
2	3-Buten-2-one, 3-methyl-	84	C5H8O	000814-78-8	56
3	2-Butenal	70	C4H6O	004170-30-3	28
4	Furan, 2,3-dihydro-	70	C4H6O	001191-99-7	28
5	4-Penten-2-one	84	C5H8O	013891-87-7	9





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Sample : J2434-16
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AD1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM041918.M
Quant Title : SVOA CALIBRATION

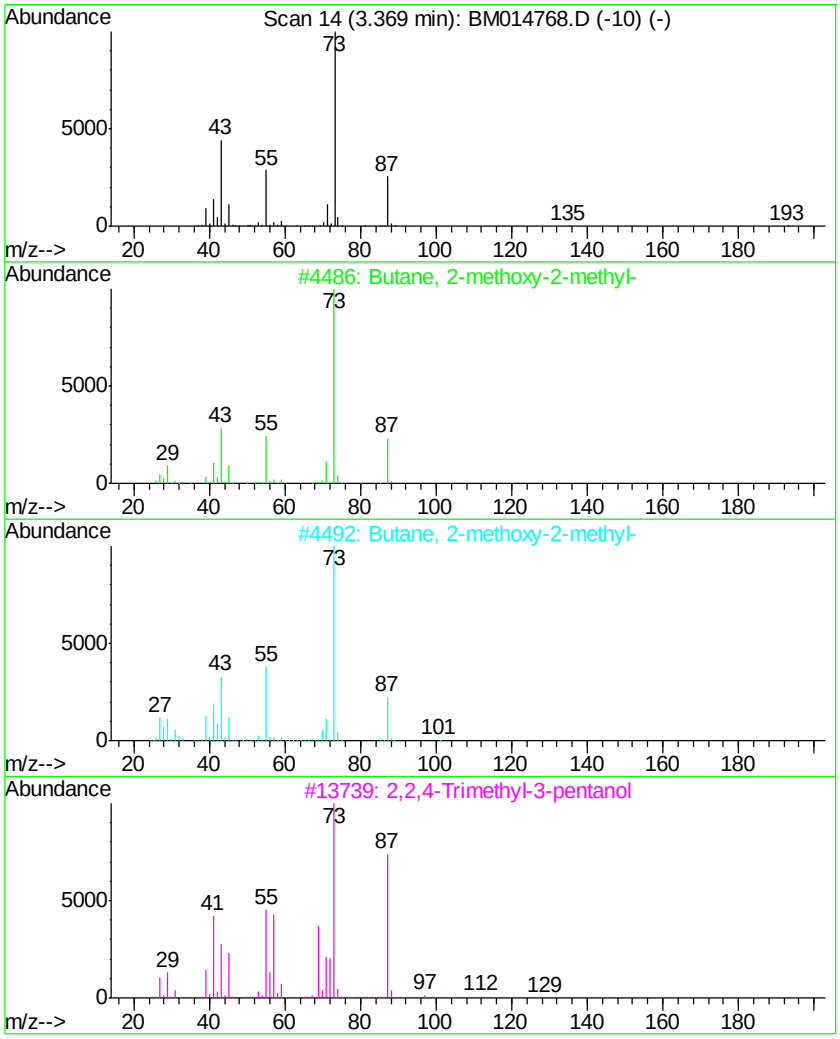
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

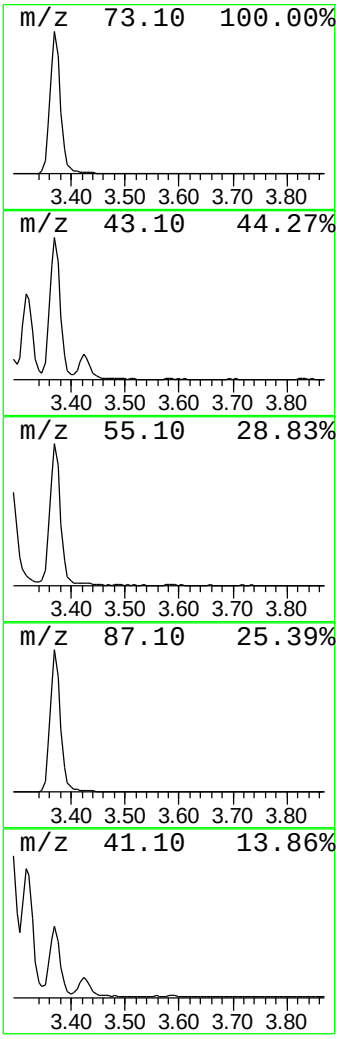
****

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.37	13.48 ng/ul	1337240	1,4-Dichlorobenzene-d4	8.55

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
3	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
4	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
5	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	23





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM042018\
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Acq On : 20 Apr 2018 18:06
Operator : SJ/JU
Sample : J2434-16
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AD1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM041918.M
Quant Title : SVOA CALIBRATION

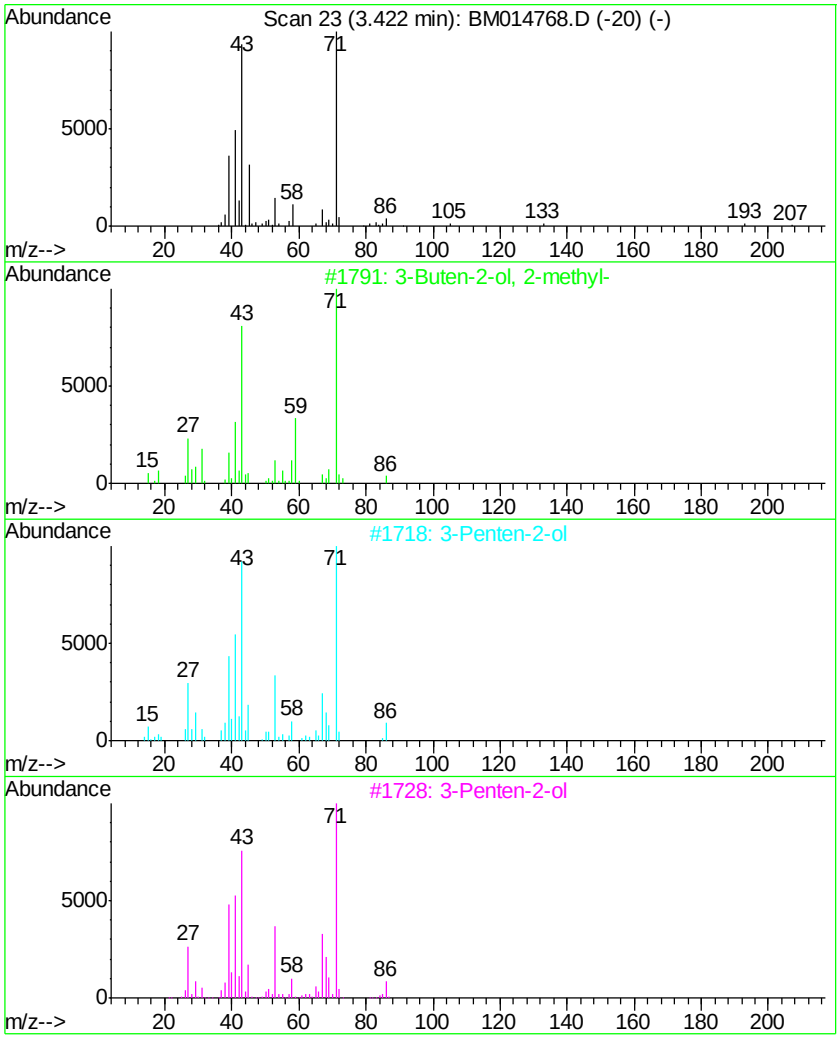
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

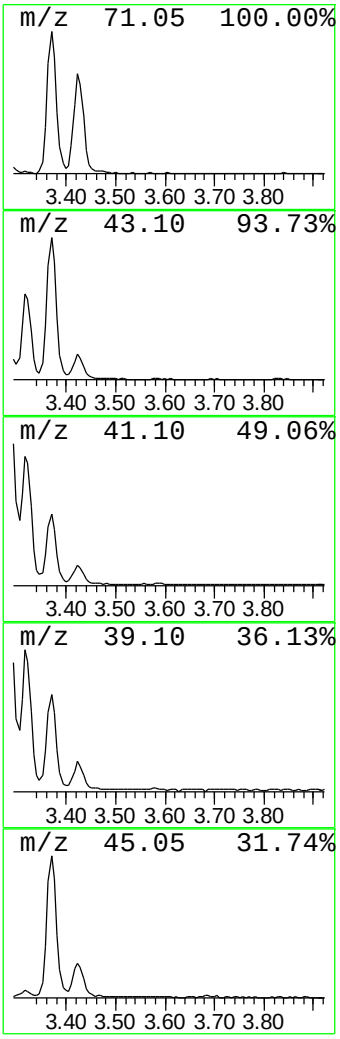
****

Peak Number 3 3-Buten-2-ol, 2-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.42	2.23 ng/ul	220763	1,4-Dichlorobenzene-d4	8.55

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Buten-2-ol, 2-methyl-	86	C5H10O	000115-18-4	80
2	3-Penten-2-ol	86	C5H10O	001569-50-2	64
3	3-Penten-2-ol	86	C5H10O	001569-50-2	64
4	3-Penten-2-ol	86	C5H10O	001569-50-2	56
5	Furan, tetrahydro-2-(methoxymeth...	116	C6H12O2	019354-27-9	56





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Instrument : BNA_M
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM041918.M
Quant Title : SVOA CALIBRATION

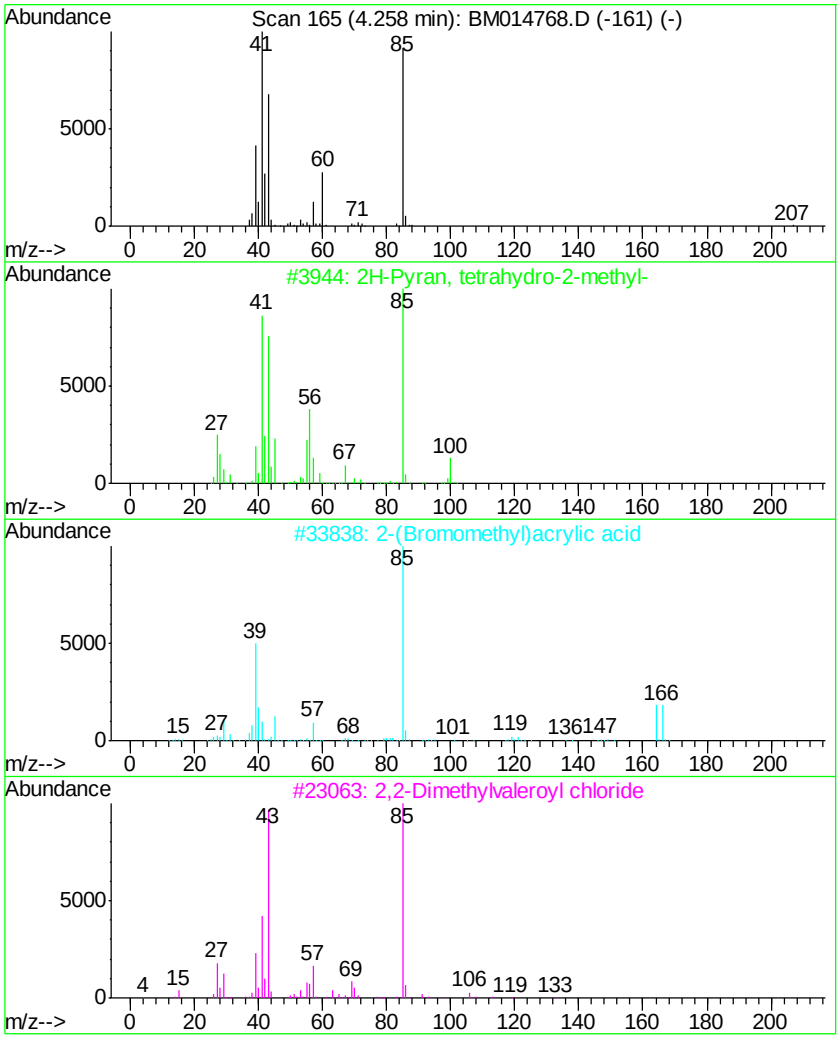
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

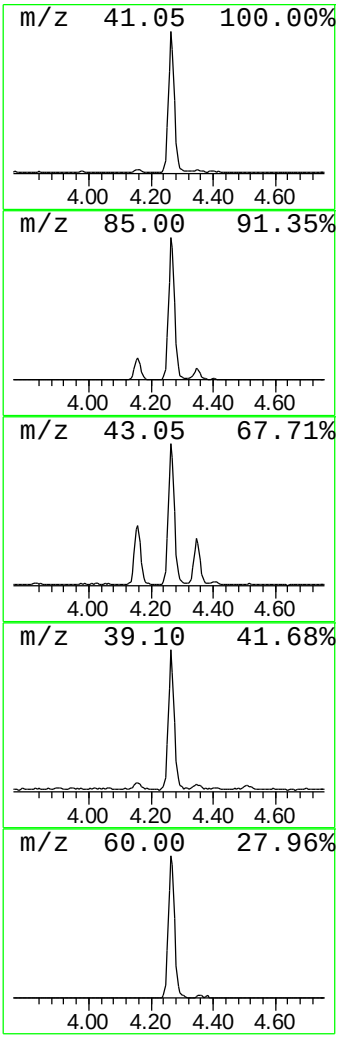
****

Peak Number 4 2H-Pyran, tetrahydro-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.26	5.36 ng/ul	531228	1,4-Dichlorobenzene-d4	8.55

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2H-Pyran, tetrahydro-2-methyl-	100	C6H12O	010141-72-7	50
2	2-(Bromomethyl)acrylic acid	164	C4H5BrO2	072707-66-5	40
3	2,2-Dimethylvaleroyl chloride	148	C7H13ClO	015721-22-9	25
4	Methacrylamide	85	C4H7NO	000079-39-0	25
5	2(3H)-Furanone, 5-acetyldihydro-	128	C6H8O3	029393-32-6	25





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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
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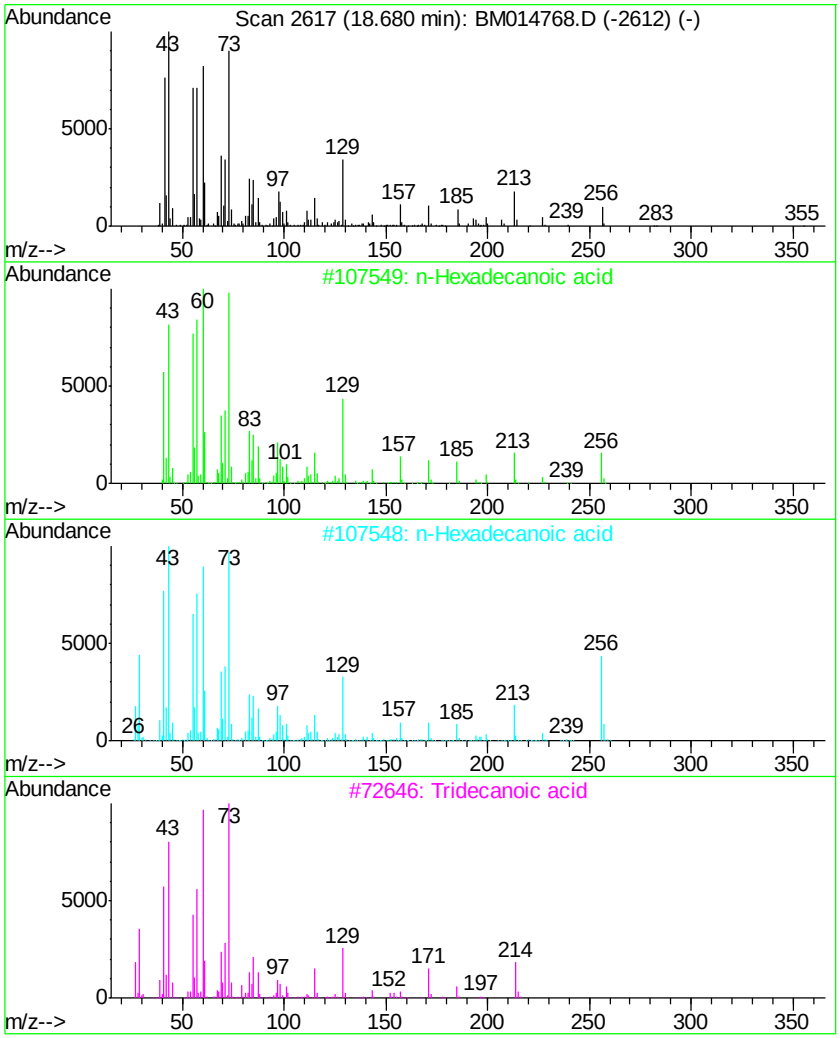
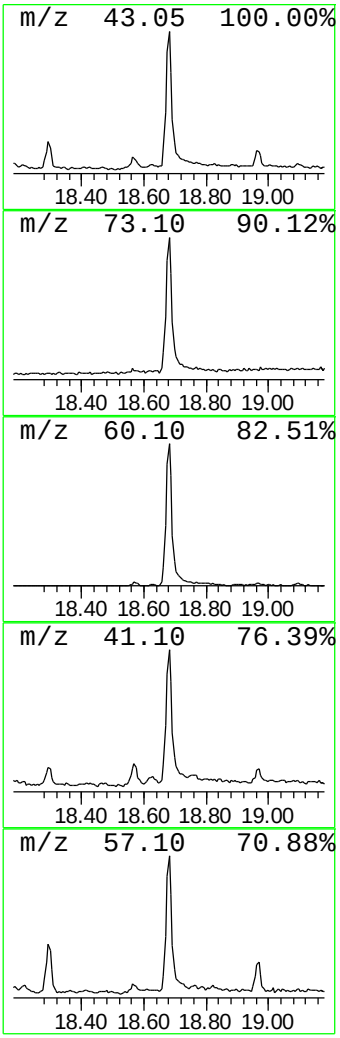
****

Peak Number 5 n-Hexadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.68	3.47 ng/ul	667235	Phenanthrene-d10	17.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94
3	Tridecanoic acid	214	C13H26O2	000638-53-9	93
4	Tridecanoic acid	214	C13H26O2	000638-53-9	93
5	Tetradecanoic acid	228	C14H28O2	000544-63-8	86





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Instrument : BNA_M
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM041918.M
Quant Title : SVOA CALIBRATION

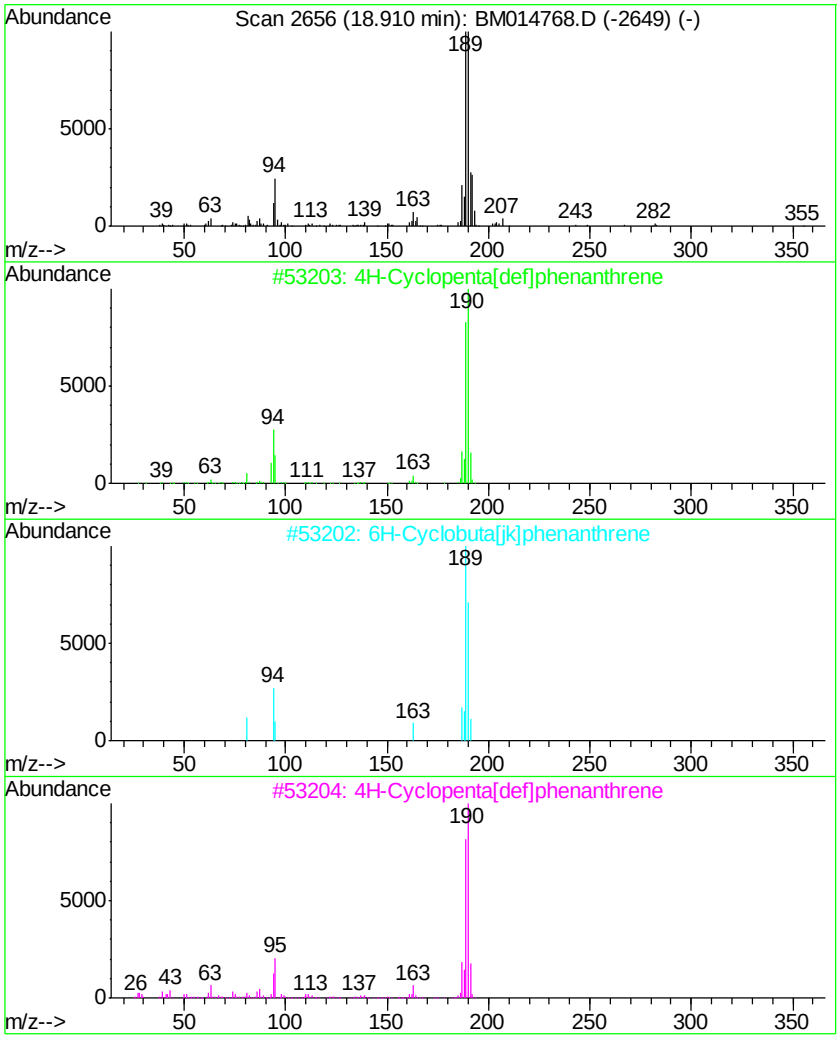
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

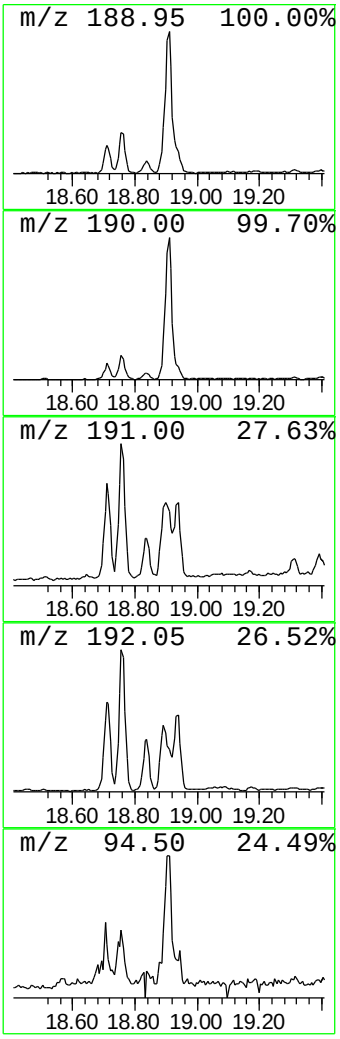
****

Peak Number 6 4H-Cyclopenta[def]phenanthrene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.91	2.33 ng/ul	447755	Phenanthrene-d10	17.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	87
2	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	64
3	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	62
4	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	59
5	Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	52





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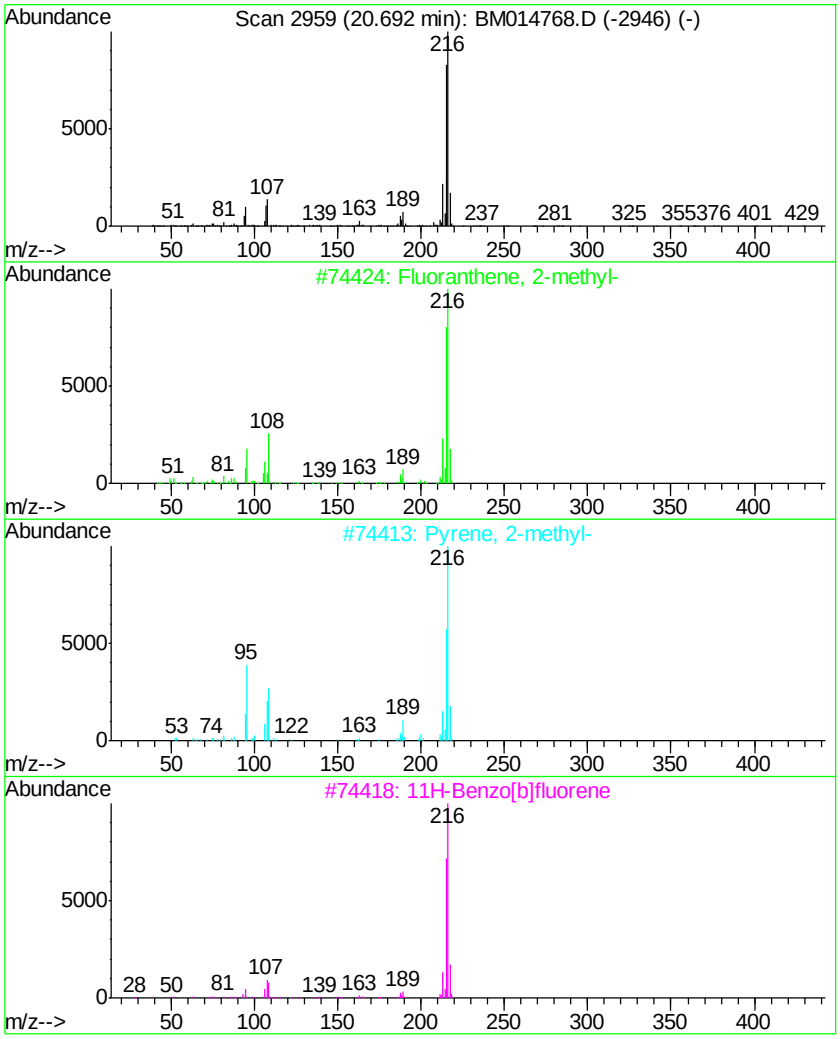
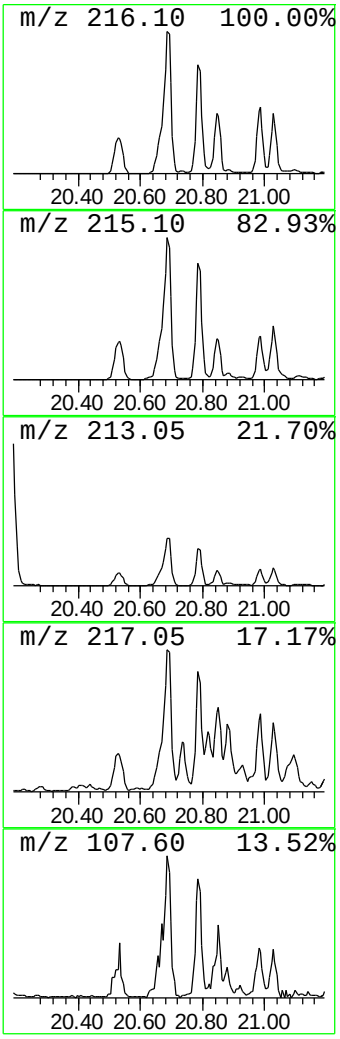
Peak Number 7 Fluoranthene, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.69	3.79 ng/ul	2166570	Chrysene-d12	21.94

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	95
2	Pyrene, 2-methyl-	216	C17H12	003442-78-2	94
3	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	94
4	Pyrene, 1-methyl-	216	C17H12	002381-21-7	94
5	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	91





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Data File : BM014768.D
Acq On : 20 Apr 2018 18:06
Operator : SJ/JU
Sample : J2434-16
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AD1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM041918.M
Quant Title : SVOA CALIBRATION

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

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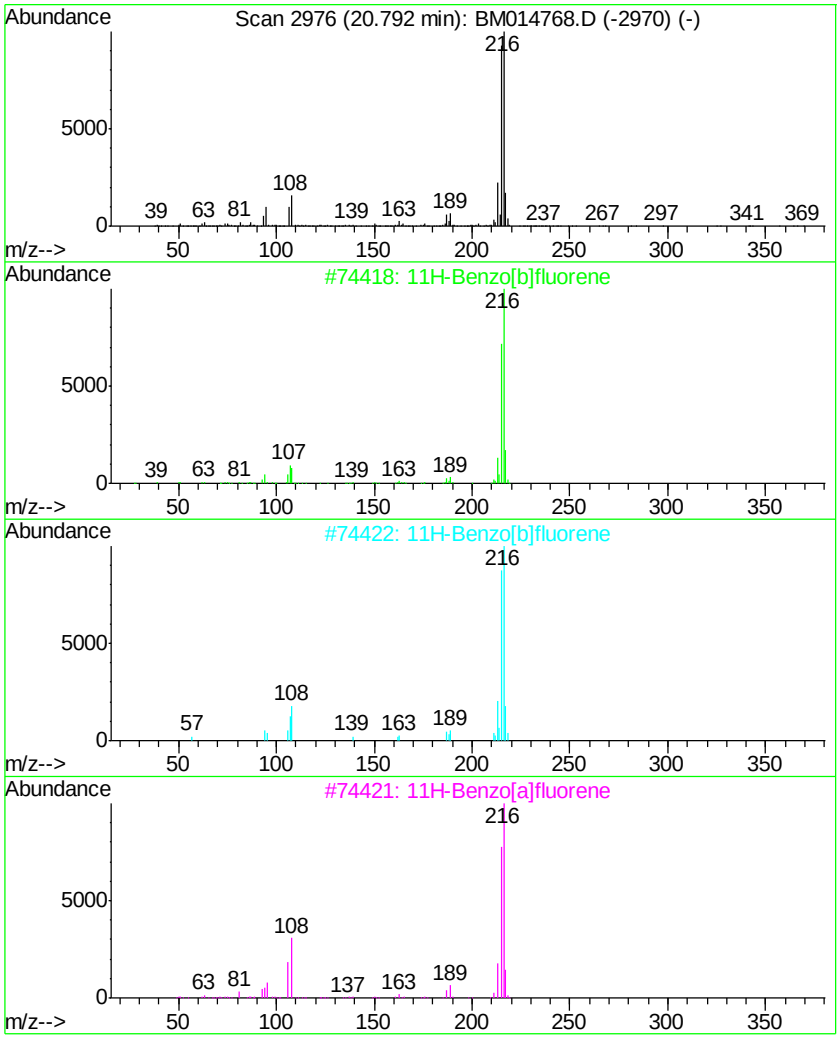
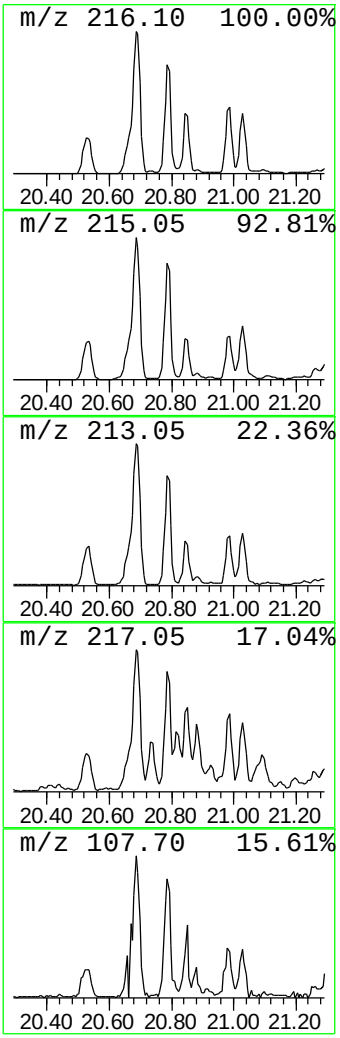
Peak Number 8 11H-Benzo[b]fluorene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.79	2.11 ng/ul	1207130	Chrysene-d12	21.94

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	95
2	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	93
3	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	91
4	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	91
5	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	91





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Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AD1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM041918.M
Quant Title : SVOA CALIBRATION

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

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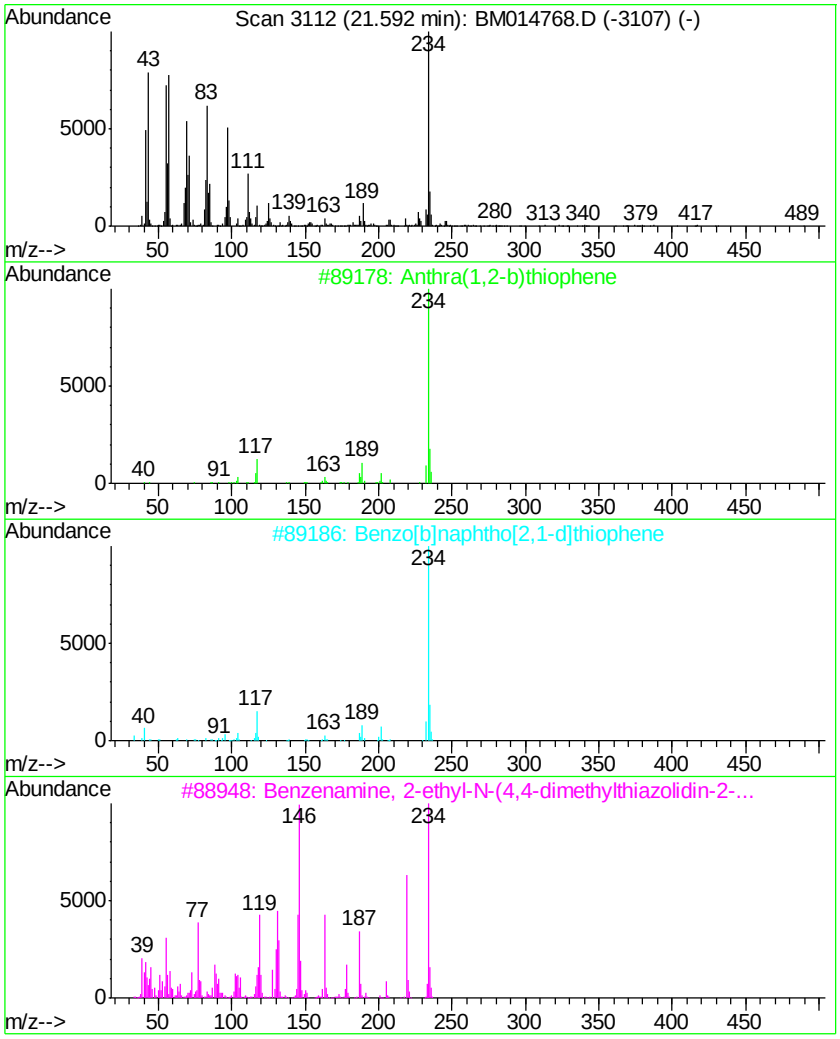
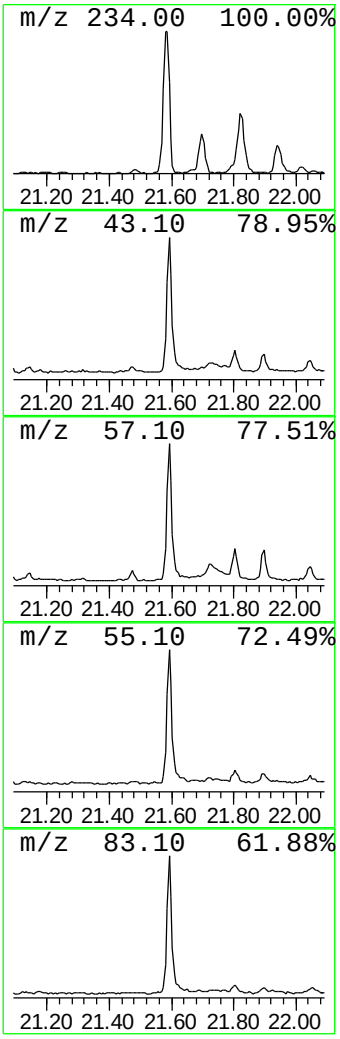
Peak Number 9 Anthra(1,2-b)thiophene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.59	2.62 ng/ul	1500020	Chrysene-d12	21.94

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthra(1,2-b)thiophene	234	C16H10S	000227-86-1	84
2	Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	83
3	Benzenamine, 2-ethyl-N-(4,4-dime...	234	C13H18N2S	1000272-26-2	83
4	1-Heneicosanol	312	C21H44O	015594-90-8	83
5	Pentafluoropropionic acid, hexad...	388	C19H33F5O2	006222-07-7	64





Library Search Compound Report

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Acq On : 20 Apr 2018 18:06
Operator : SJ/JU
Sample : J2434-16
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AD1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM041918.M
Quant Title : SVOA CALIBRATION

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

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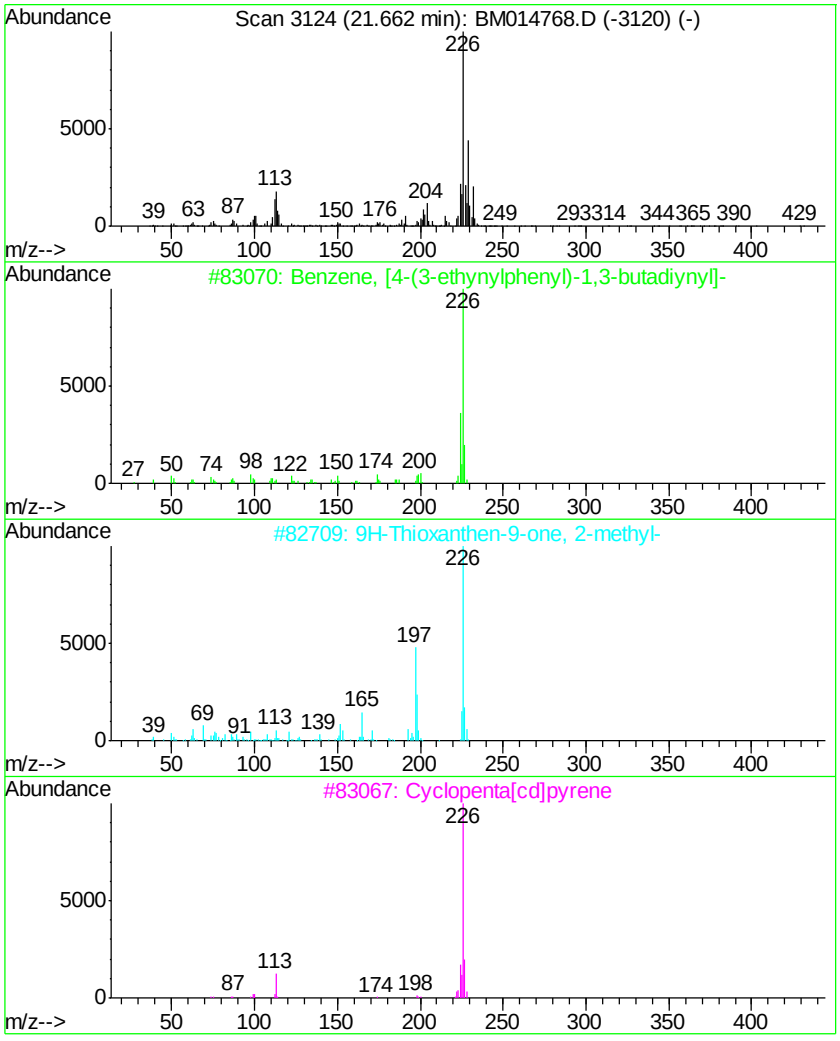
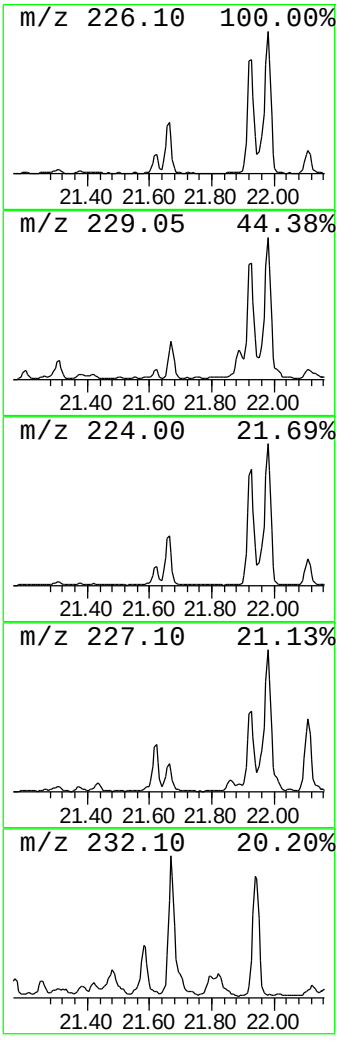
Peak Number 10 Benzene, [4-(3-ethynylpheny... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.66	2.11 ng/ul	1209390	Chrysene-d12	21.94

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, [4-(3-ethynylphenyl)-1,...	226	C18H10	1000115-87-3	50
2	9H-Thioxanthen-9-one, 2-methyl-	226	C14H10OS	015774-82-0	46
3	Cyclopenta[cd]pyrene	226	C18H10	027208-37-3	45
4	4-Naphthalen-2-yl-thiazol-2-ylamine	226	C13H10N2S	1000300-53-4	43
5	.beta.-Carboline, 7-methoxy-1,2-...	226	C14H14N2O	006519-18-2	38





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Acq On : 20 Apr 2018 18:06
Operator : SJ/JU
Sample : J2434-16
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AD1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM041918.M
Quant Title : SVOA CALIBRATION

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

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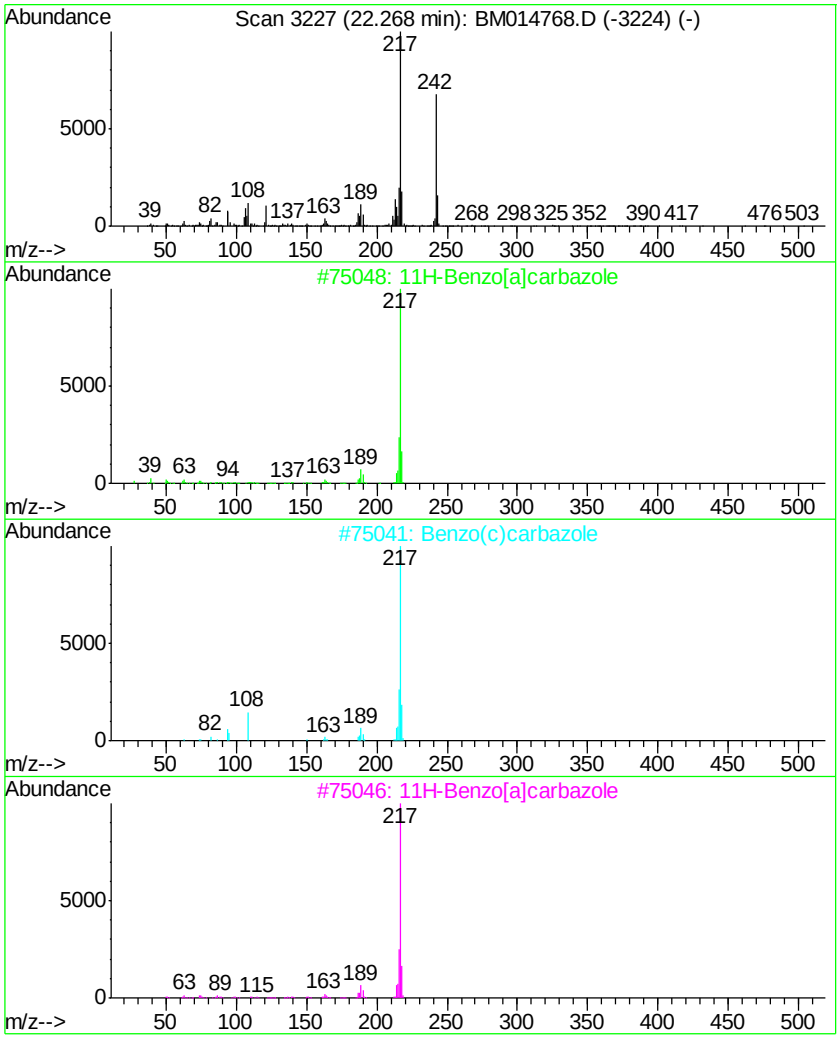
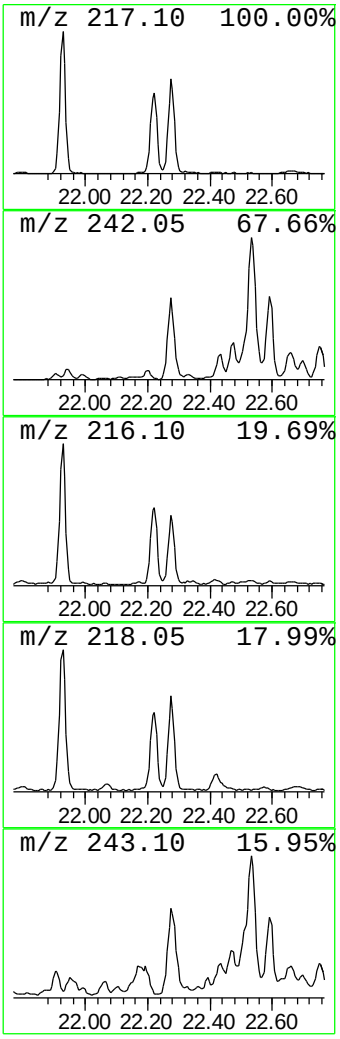
Peak Number 11 11H-Benzo[a]carbazole Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.27	2.49 ng/ul	1423940	Chrysene-d12	21.94

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[a]carbazole	217	C16H11N	000239-01-0	83
2	Benzo(c)carbazole	217	C16H11N	034777-33-8	70
3	11H-Benzo[a]carbazole	217	C16H11N	000239-01-0	64
4	7H-Benzo[c]carbazole	217	C16H11N	000205-25-4	64
5	Benzo(b)carbazole	217	C16H11N	000243-28-7	64





Library Search Compound Report

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Data File : BM014768.D
Acq On : 20 Apr 2018 18:06
Operator : SJ/JU
Sample : J2434-16
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AD1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM041918.M
Quant Title : SVOA CALIBRATION

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

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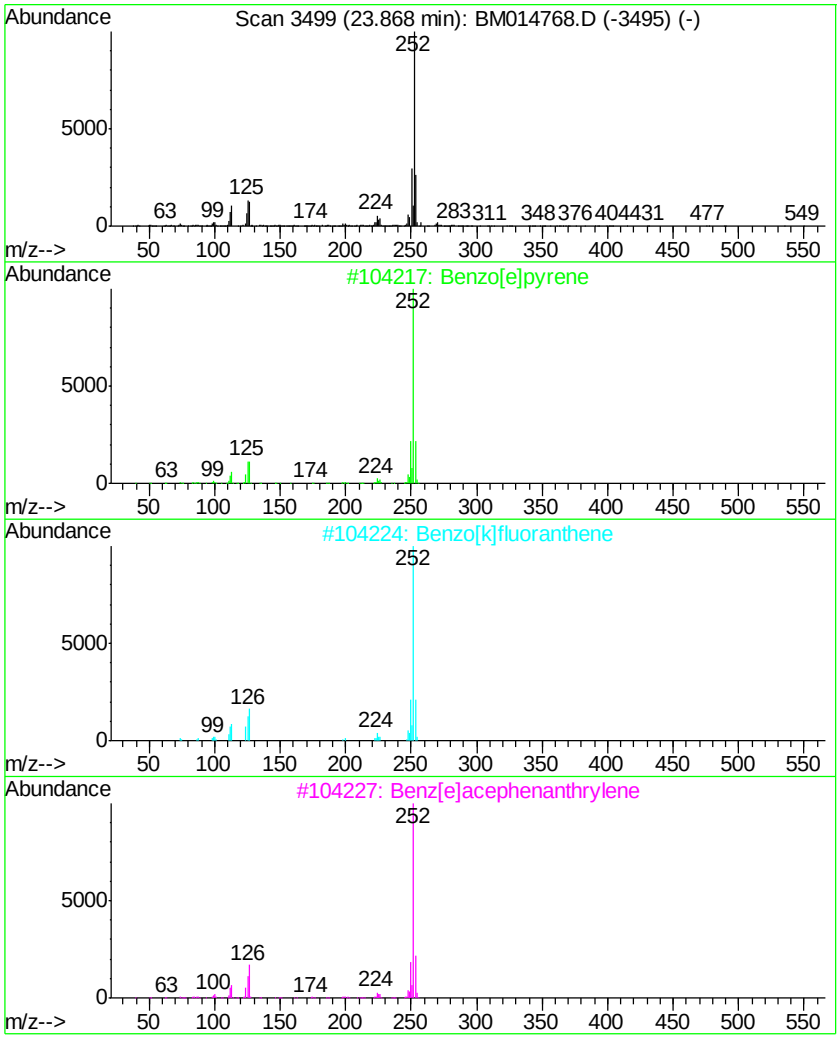
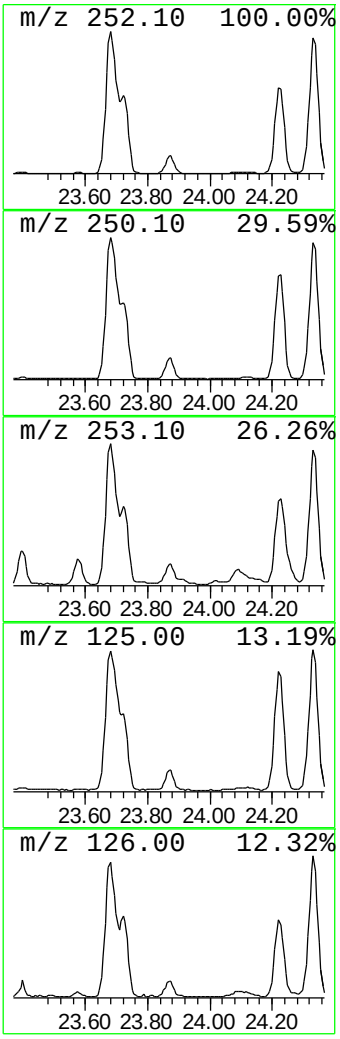
Peak Number 12 Benzo[e]pyrene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.87	5.94 ng/ul	1220110	Perylene-d12	24.44

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]pyrene	252	C20H12	000192-97-2	98
2	Benzo[k]fluoranthene	252	C20H12	000207-08-9	98
3	Benz[e]acephenanthrylene	252	C20H12	000205-99-2	93
4	Perylene	252	C20H12	000198-55-0	93
5	Benzo[k]fluoranthene	252	C20H12	000207-08-9	93





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Operator : SJ/JU
Sample : J2434-16
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AD1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM041918.M
Quant Title : SVOA CALIBRATION

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

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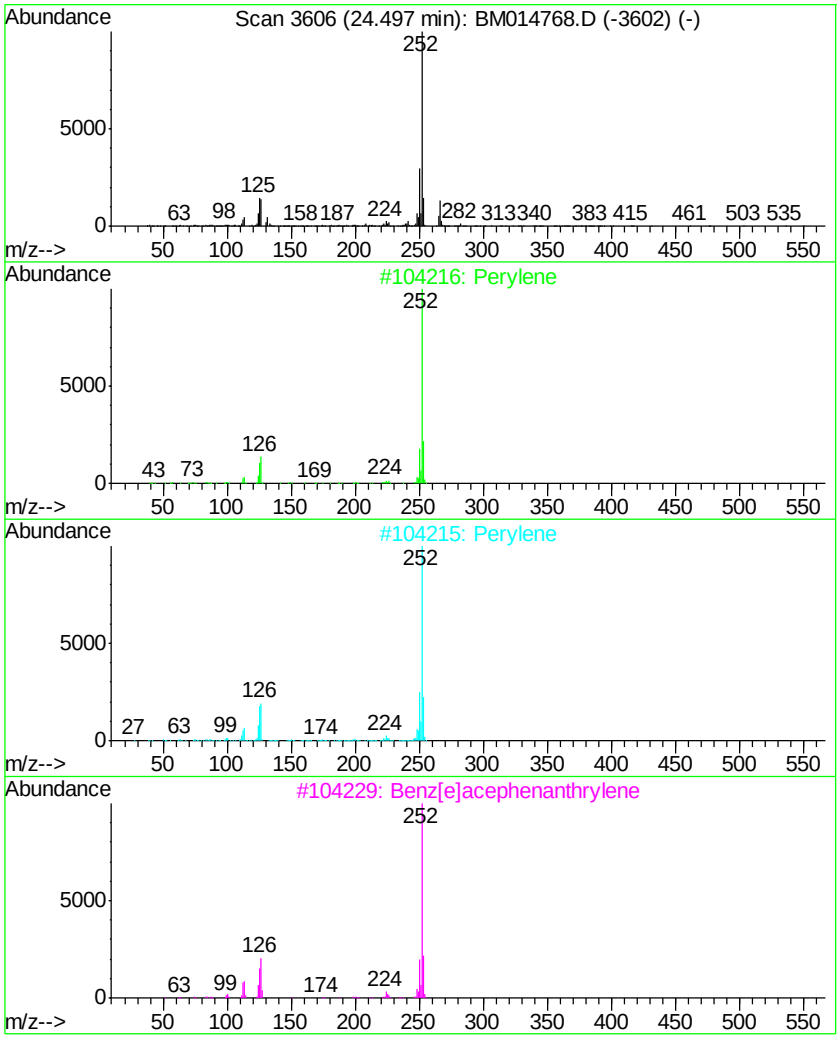
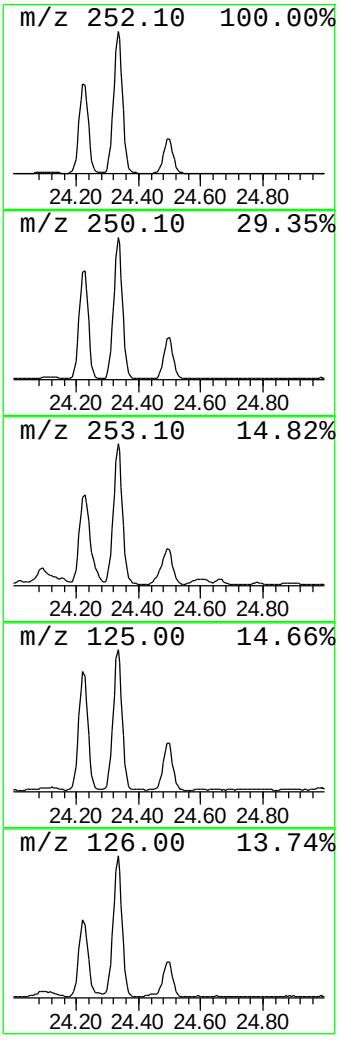
Peak Number 14 Perylene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.50	14.72 ng/ul	3022910	Perylene-d12	24.44

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Perylene	252	C20H12	000198-55-0	93
2	Perylene	252	C20H12	000198-55-0	89
3	Benz[e]acephenanthrylene	252	C20H12	000205-99-2	89
4	Benzo[a]pyrene	252	C20H12	000050-32-8	86
5	Benzo[e]pyrene	252	C20H12	000192-97-2	76





Tentatively Identified Compound (LSC) summary

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Data File : BM014768.D
Acq On : 20 Apr 2018 18:06
Operator : SJ/JU
Sample : J2434-16
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AD1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM041918.M
Quant Title : SVOA CALIBRATION

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

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Ethanone, 1-cyclo...	3.32	5.0	ng/ul	498875	1	8.55	1983580	20.0
Butane, 2-methoxy...	3.37	13.5	ng/ul	1337240	1	8.55	1983580	20.0
3-Buten-2-ol, 2-m...	3.42	2.2	ng/ul	220763	1	8.55	1983580	20.0
2H-Pyran, tetrahy...	4.26	5.4	ng/ul	531228	1	8.55	1983580	20.0
n-Hexadecanoic acid	18.68	3.5	ng/ul	667235	4	17.86	3848140	20.0
4H-Cyclopenta[def...	18.91	2.3	ng/ul	447755	4	17.86	3848140	20.0
Fluoranthene, 2-m...	20.69	3.8	ng/ul	2166570	5	21.94	11446900	20.0
11H-Benzo[b]fluorene	20.79	2.1	ng/ul	1207130	5	21.94	11446900	20.0
Anthra(1,2-b)thio...	21.59	2.6	ng/ul	1500020	5	21.94	11446900	20.0
Benzene, [4-(3-et...	21.66	2.1	ng/ul	1209390	5	21.94	11446900	20.0
11H-Benzo[a]carba...	22.27	2.5	ng/ul	1423940	5	21.94	11446900	20.0
Benzo[e]pyrene	23.87	5.9	ng/ul	1220110	6	24.44	4106640	20.0
Perylene	24.50	14.7	ng/ul	3022910	6	24.44	4106640	20.0