LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG122115\
Data File : BG020400.D
Acq On : 22 Dec 2015 6:10
Operator : UM/SJ
Sample : G4848-04DL 30X
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** D9N50DL

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_G\METHODS\SOM02.2-EPA-BG121415.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 | 8.091 | 888 | 894 | 903 | rBB | 58172 | 96126 | 3.87% | 0.604% |
| 2 | 8.632 | 979 | 986 | 994 | rBB | 23433 | 38714 | 1.56% | 0.243% |
| 3 | 10.906 | 1365 | 1373 | 1377 | rBV | 83303 | 154628 | 6.23% | 0.971% |
| 4 | 10.959 | 1377 | 1382 | 1393 | rVV | 535070 | 966947 | 38.95% | 6.071% |
| 5 | 11.076 | 1393 | 1402 | 1410 | rVB2 | 15432 | 29212 | 1.18% | 0.183% |
| 6 | 11.734 | 1509 | 1514 | 1521 | rBV2 | 10075 | 17545 | 0.71% | 0.110% |
| 7 | 12.557 | 1647 | 1654 | 1666 | rVB | 284059 | 467574 | 18.83% | 2.936% |
| 8 | 12.774 | 1680 | 1691 | 1703 | rBB | 135099 | 225730 | 9.09% | 1.417% |
| 9 | 13.555 | 1817 | 1824 | 1834 | rBB | 98952 | 150668 | 6.07% | 0.946% |
| 10 | 13.749 | 1852 | 1857 | 1867 | rVB | 32982 | 56858 | 2.29% | 0.357% |
| 11 | 13.884 | 1872 | 1880 | 1881 | rBV | 41241 | 60011 | 2.42% | 0.377% |
| 12 | 14.043 | 1900 | 1907 | 1912 | rBV | 62526 | 112874 | 4.55% | 0.709% |
| 13 | 14.096 | 1912 | 1916 | 1924 | rVV2 | 43303 | 72854 | 2.93% | 0.457% |
| 14 | 14.278 | 1942 | 1947 | 1950 | rBV2 | 24557 | 35925 | 1.45% | 0.226% |
| 15 | 14.443 | 1966 | 1975 | 1982 | rBV2 | 31793 | 55971 | 2.25% | 0.351% |
| 16 | 14.689 | 2011 | 2017 | 2019 | rBV | 41348 | 62492 | 2.52% | 0.392% |
| 17 | 14.725 | 2019 | 2023 | 2028 | rVV2 | 134455 | 222035 | 8.94% | 1.394% |
| 18 | 14.789 | 2028 | 2034 | 2040 | rVV | 569870 | 879100 | 35.41% | 5.520% |
| 19 | 14.854 | 2040 | 2045 | 2051 | rVV | 23288 | 37029 | 1.49% | 0.232% |
| 20 | 14.919 | 2051 | 2056 | 2065 | rVV3 | 9265 | 22338 | 0.90% | 0.140% |
| 21 | 15.030 | 2065 | 2075 | 2080 | rVV3 | 21197 | 39462 | 1.59% | 0.248% |
| 22 | 15.118 | 2083 | 2090 | 2098 | rVV | 379695 | 589896 | 23.76% | 3.704% |
| 23 | 15.189 | 2098 | 2102 | 2111 | rVB2 | 14203 | 21963 | 0.88% | 0.138% |
| 24 | 15.330 | 2120 | 2126 | 2131 | rBV3 | 10508 | 18731 | 0.75% | 0.118% |
| 25 | 15.383 | 2131 | 2135 | 2145 | rVB2 | 11439 | 17978 | 0.72% | 0.113% |
| 26 | 15.500 | 2148 | 2155 | 2159 | rBV3 | 8027 | 16198 | 0.65% | 0.102% |
| 27 | 15.682 | 2182 | 2186 | 2188 | rVV3 | 10472 | 14613 | 0.59% | 0.092% |
| 28 | 15.771 | 2194 | 2201 | 2209 | rVV | 505346 | 780532 | 31.44% | 4.901% |
| 29 | 15.859 | 2211 | 2216 | 2218 | rVV | 22730 | 31649 | 1.27% | 0.199% |
| 30 | 15.888 | 2218 | 2221 | 2224 | rVV2 | 35445 | 56022 | 2.26% | 0.352% |
| 31 | 15.929 | 2224 | 2228 | 2233 | rVV2 | 69424 | 107532 | 4.33% | 0.675% |
| 32 | 15.976 | 2233 | 2236 | 2239 | rVV3 | 14643 | 23190 | 0.93% | 0.146% |
| 33 | 16.017 | 2239 | 2243 | 2246 | rVV2 | 32651 | 51639 | 2.08% | 0.324% |
| 34 | 16.058 | 2246 | 2250 | 2259 | rVB | 74056 | 110275 | 4.44% | 0.692% |

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG122115\
Data File : BG020400.D
Acq On : 22 Dec 2015 6:10
Operator : UM/SJ
Sample : G4848-04DL 30X
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** D9N50DL

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_G\METHODS\SOM02.2-EPA-BG121415.M
Title : SVOA CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 35 | 16.205 | 2265 | 2275 | 2283 | rBV | 78251 | 166043 | 6.69% | 1.043% |
| 36 | 16.293 | 2288 | 2290 | 2296 | rVB2 | 11788 | 15898 | 0.64% | 0.100% |
| 37 | 16.558 | 2326 | 2335 | 2341 | rVB | 43029 | 62337 | 2.51% | 0.391% |
| 38 | 16.769 | 2359 | 2371 | 2375 | rBV2 | 55163 | 123601 | 4.98% | 0.776% |
| 39 | 16.816 | 2375 | 2379 | 2385 | rVB2 | 18272 | 25829 | 1.04% | 0.162% |
| 40 | 16.916 | 2390 | 2396 | 2401 | rVB3 | 21588 | 36712 | 1.48% | 0.231% |
| 41 | 16.993 | 2401 | 2409 | 2412 | rBV4 | 18595 | 37947 | 1.53% | 0.238% |
| 42 | 17.034 | 2412 | 2416 | 2420 | rVB2 | 16960 | 21471 | 0.86% | 0.135% |
| 43 | 17.122 | 2427 | 2431 | 2438 | rVB4 | 10241 | 18450 | 0.74% | 0.116% |
| 44 | 17.192 | 2438 | 2443 | 2454 | rBV3 | 23462 | 64046 | 2.58% | 0.402% |
| 45 | 17.286 | 2454 | 2459 | 2468 | rVB | 116280 | 175935 | 7.09% | 1.105% |
| 46 | 17.469 | 2483 | 2490 | 2493 | rBV | 189275 | 303589 | 12.23% | 1.906% |
| 47 | 17.516 | 2493 | 2498 | 2504 | rVV | 1549818 | 2482558 | 100.00% | 15.587% |
| 48 | 17.574 | 2504 | 2508 | 2509 | rVV2 | 29312 | 44556 | 1.79% | 0.280% |
| 49 | 17.604 | 2509 | 2513 | 2518 | rVV | 106712 | 158630 | 6.39% | 0.996% |
| 50 | 17.651 | 2518 | 2521 | 2526 | rVB4 | 7651 | 15271 | 0.62% | 0.096% |
| 51 | 17.868 | 2550 | 2558 | 2567 | rBV | 76953 | 133379 | 5.37% | 0.837% |
| 52 | 17.986 | 2573 | 2578 | 2583 | rVV3 | 26136 | 39608 | 1.60% | 0.249% |
| 53 | 18.050 | 2583 | 2589 | 2597 | rVB7 | 10787 | 24402 | 0.98% | 0.153% |
| 54 | 18.191 | 2608 | 2613 | 2621 | rVB2 | 10508 | 24274 | 0.98% | 0.152% |
| 55 | 18.344 | 2633 | 2639 | 2642 | rBV | 104977 | 151031 | 6.08% | 0.948% |
| 56 | 18.391 | 2642 | 2647 | 2653 | rVB | 148608 | 222216 | 8.95% | 1.395% |
| 57 | 18.467 | 2653 | 2660 | 2666 | rBV | 49072 | 77307 | 3.11% | 0.485% |
| 58 | 18.538 | 2666 | 2672 | 2676 | rBV2 | 197909 | 351373 | 14.15% | 2.206% |
| 59 | 18.832 | 2717 | 2722 | 2727 | rVV | 89225 | 127316 | 5.13% | 0.799% |
| 60 | 19.078 | 2759 | 2764 | 2768 | rVB4 | 16021 | 20819 | 0.84% | 0.131% |
| 61 | 19.131 | 2768 | 2773 | 2776 | rBV | 13766 | 16964 | 0.68% | 0.107% |
| 62 | 19.167 | 2776 | 2779 | 2783 | rVB2 | 11925 | 15055 | 0.61% | 0.095% |
| 63 | 19.249 | 2787 | 2793 | 2799 | rBV3 | 27087 | 52572 | 2.12% | 0.330% |
| 64 | 19.313 | 2799 | 2804 | 2807 | rBV4 | 14977 | 23657 | 0.95% | 0.149% |
| 65 | 19.384 | 2812 | 2816 | 2826 | rVB5 | 10624 | 27464 | 1.11% | 0.172% |
| 66 | 19.519 | 2832 | 2839 | 2846 | rBV | 1038930 | 1479214 | 59.58% | 9.288% |
| 67 | 19.607 | 2851 | 2854 | 2858 | rVB2 | 16807 | 22061 | 0.89% | 0.139% |
| 68 | 19.760 | 2875 | 2880 | 2888 | rVB | 28214 | 48381 | 1.95% | 0.304% |
| 69 | 19.848 | 2888 | 2895 | 2896 | rBV | 178654 | 256212 | 10.32% | 1.609% |
| 70 | 19.877 | 2896 | 2900 | 2908 | rVV | 638033 | 954246 | 38.44% | 5.991% |
| 71 | 19.942 | 2908 | 2911 | 2918 | rVB2 | 33925 | 53408 | 2.15% | 0.335% |

LSC Area Percent Report

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Acq On : 22 Dec 2015 6:10
Operator : UM/SJ
Sample : G4848-04DL 30X
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** D9N50DL

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_G\METHODS\SOM02.2-EPA-BG121415.M
Title : SVOA CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 72 | 20.054 | 2924 | 2930 | 2935 | rVB | 36757 | 52722 | 2.12% | 0.331% |
| 73 | 20.112 | 2935 | 2940 | 2945 | rVB3 | 9193 | 13722 | 0.55% | 0.086% |
| 74 | 20.201 | 2948 | 2955 | 2960 | rBV3 | 33982 | 85737 | 3.45% | 0.538% |
| 75 | 20.253 | 2960 | 2964 | 2970 | rVB | 16765 | 20534 | 0.83% | 0.129% |
| 76 | 20.365 | 2970 | 2983 | 2989 | rBV | 119828 | 201278 | 8.11% | 1.264% |
| 77 | 20.465 | 2993 | 3000 | 3004 | rBV | 135476 | 192086 | 7.74% | 1.206% |
| 78 | 20.524 | 3004 | 3010 | 3013 | rVV2 | 30676 | 62813 | 2.53% | 0.394% |
| 79 | 20.559 | 3013 | 3016 | 3020 | rVV | 21496 | 23986 | 0.97% | 0.151% |
| 80 | 20.600 | 3020 | 3023 | 3027 | rVB3 | 10787 | 15912 | 0.64% | 0.100% |
| 81 | 20.659 | 3029 | 3033 | 3037 | rBV | 13849 | 18814 | 0.76% | 0.118% |
| 82 | 20.706 | 3037 | 3041 | 3045 | rBV | 15898 | 20068 | 0.81% | 0.126% |
| 83 | 20.894 | 3068 | 3073 | 3076 | rBV | 18218 | 27849 | 1.12% | 0.175% |
| 84 | 20.959 | 3079 | 3084 | 3087 | rVV4 | 8659 | 14344 | 0.58% | 0.090% |
| 85 | 20.994 | 3087 | 3090 | 3095 | rVB3 | 9765 | 15042 | 0.61% | 0.094% |
| 86 | 21.082 | 3100 | 3105 | 3110 | rVV3 | 14762 | 28460 | 1.15% | 0.179% |
| 87 | 21.135 | 3110 | 3114 | 3119 | rVV4 | 9827 | 19865 | 0.80% | 0.125% |
| 88 | 21.317 | 3140 | 3145 | 3149 | rBV | 27989 | 43303 | 1.74% | 0.272% |
| 89 | 21.358 | 3149 | 3152 | 3157 | rVB | 25404 | 33343 | 1.34% | 0.209% |
| 90 | 21.411 | 3157 | 3161 | 3166 | rBV3 | 21772 | 34737 | 1.40% | 0.218% |
| 91 | 21.740 | 3204 | 3217 | 3222 | rVV2 | 290300 | 711969 | 28.68% | 4.470% |
| 92 | 21.787 | 3222 | 3225 | 3237 | rVB | 118385 | 209566 | 8.44% | 1.316% |
| 93 | 21.946 | 3246 | 3252 | 3258 | rBV | 24953 | 40692 | 1.64% | 0.255% |
| 94 | 22.157 | 3282 | 3288 | 3294 | rVV3 | 11049 | 25114 | 1.01% | 0.158% |
| 95 | 22.480 | 3335 | 3343 | 3350 | rVV | 11616 | 30659 | 1.23% | 0.192% |
| 96 | 22.803 | 3396 | 3398 | 3406 | rVB5 | 8753 | 16447 | 0.66% | 0.103% |
| 97 | 23.973 | 3589 | 3597 | 3604 | rBV2 | 25043 | 85448 | 3.44% | 0.537% |
| 98 | 24.707 | 3716 | 3722 | 3731 | rBV2 | 10914 | 30475 | 1.23% | 0.191% |
| 99 | 24.866 | 3741 | 3749 | 3760 | rVB3 | 13194 | 43128 | 1.74% | 0.271% |
| 100 | 25.019 | 3764 | 3775 | 3787 | rBV | 88053 | 290618 | 11.71% | 1.825% |

Sum of corrected areas: 15926874

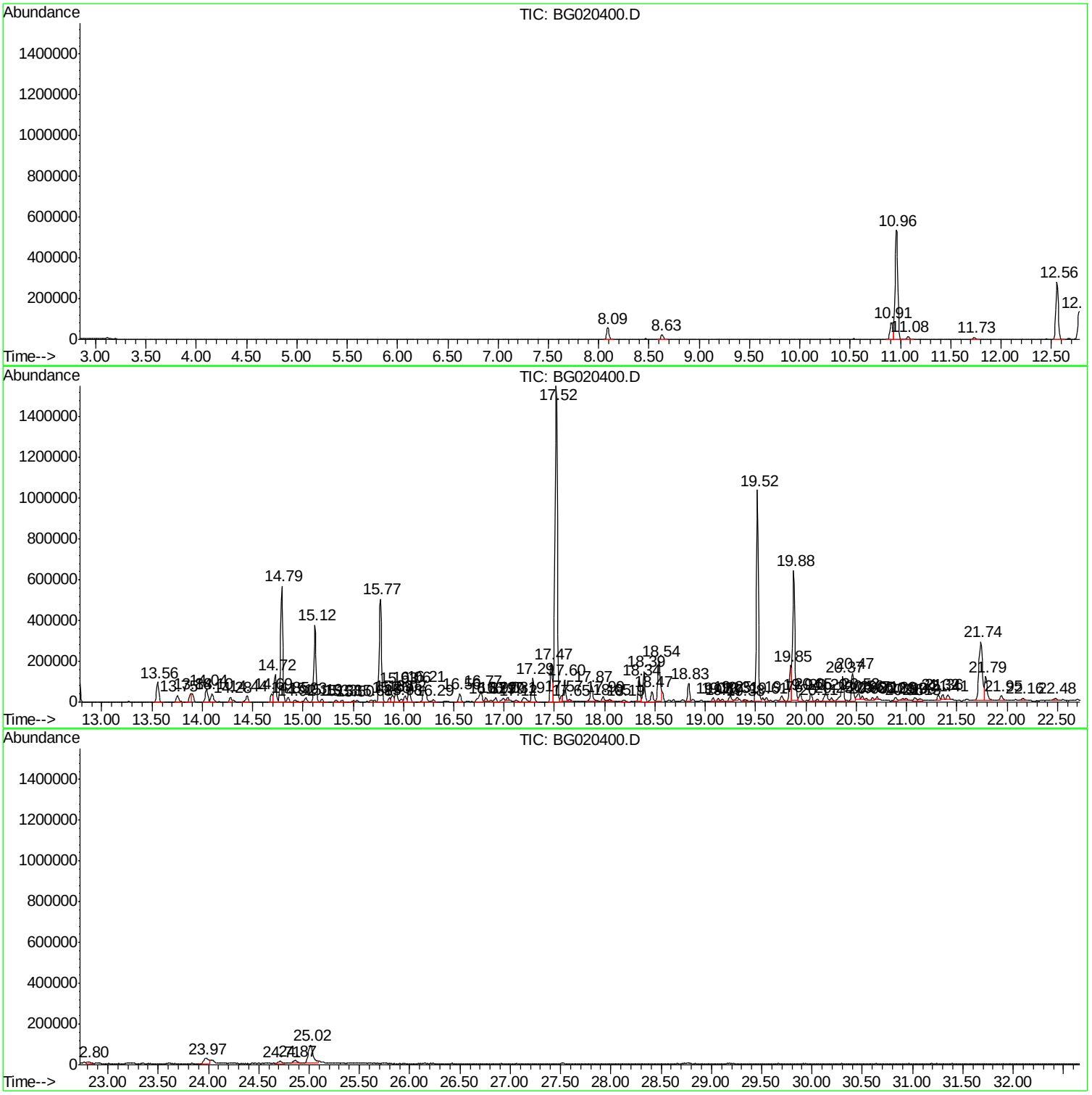
LSC Report - Integrated Chromatogram

Data Path : Z:\HPCHEM1\BNA G\DATA\BG122115\
Data File : BG020400.D
Acq On : 22 Dec 2015 6:10
Operator : UM/SJ
Sample : G4848-04DL 30X
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
D9N50DL

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG121415.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 G\DATA\BG122115\
Data File : BG020400.D
Acq On : 22 Dec 2015 6:10
Operator : UM/SJ
Sample : G4848-04DL 30X
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** D9N50DL

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG121415.M
Quant Title : SVOA CALIBRATION

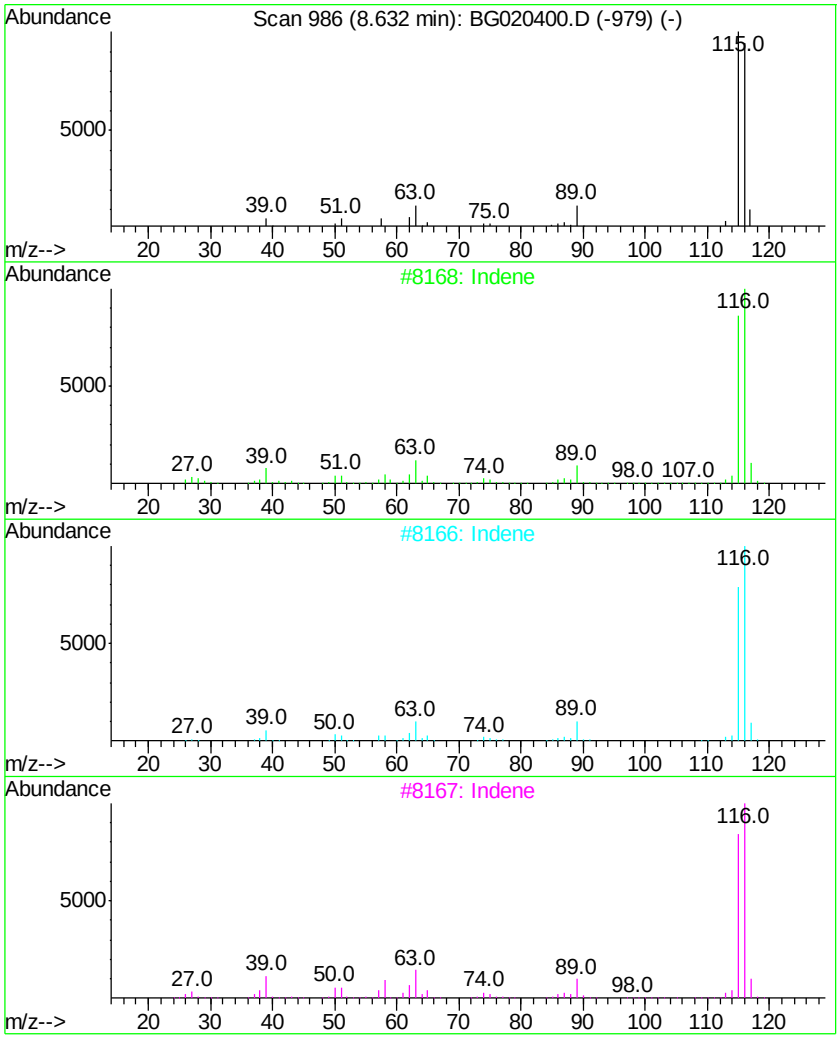
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

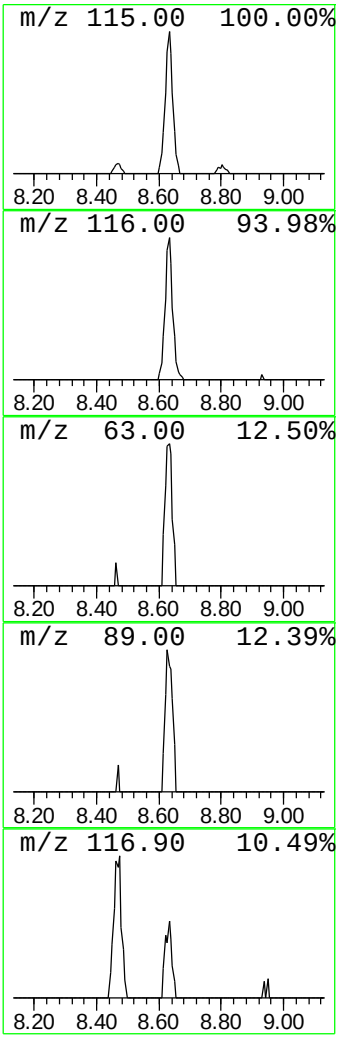
**********

Peak Number 1 Indene Concentration Rank 12

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 8.63 | 8.05 ng/ul | 38714 | 1,4-Dichlorobenzene-d4 | 8.09 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Indene | 116 | C9H8 | 000095-13-6 | 91 |
| 2 | Indene | 116 | C9H8 | 000095-13-6 | 91 |
| 3 | Indene | 116 | C9H8 | 000095-13-6 | 91 |
| 4 | Benzene, 1-ethynyl-4-methyl- | 116 | C9H8 | 000766-97-2 | 91 |
| 5 | Benzene, 1-propynyl- | 116 | C9H8 | 000673-32-5 | 91 |



# Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG122115\
Data File : BG020400.D
Acq On : 22 Dec 2015 6:10
Operator : UM/SJ
Sample : G4848-04DL 30X
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** D9N50DL

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG121415.M
Quant Title : SVOA CALIBRATION

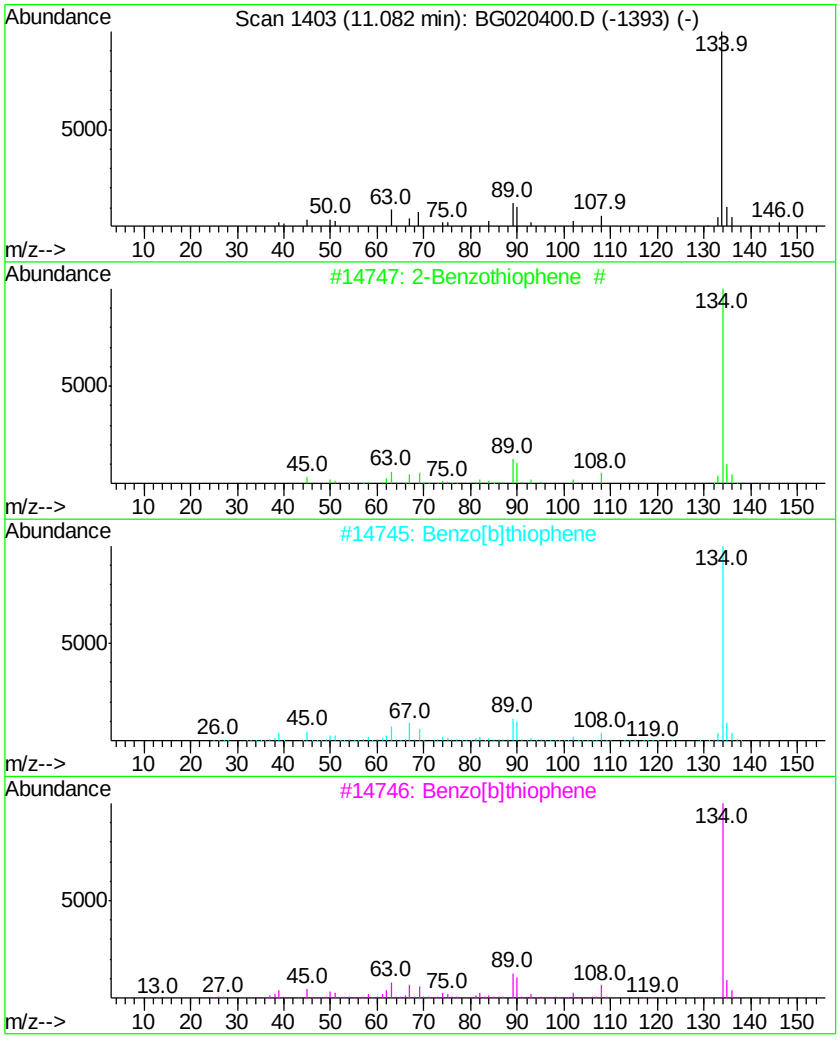
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

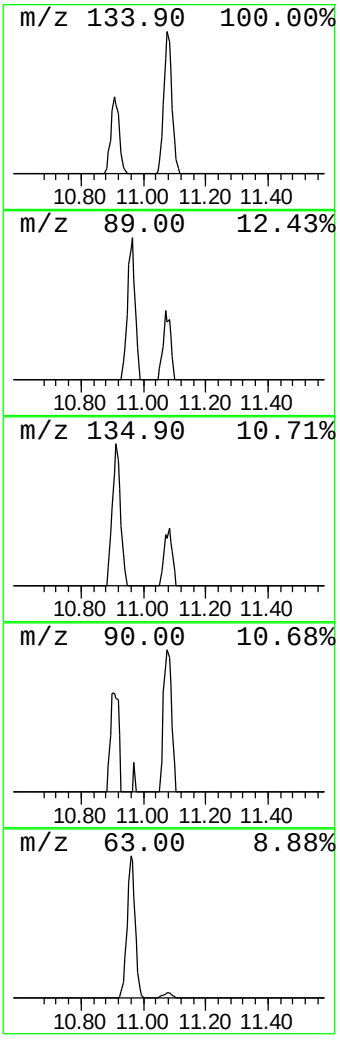
**********

Peak Number 2 2-Benzothiophene # Concentration Rank 23

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.08 | 3.78 ng/ul | 29212 | Naphthalene-d8 | 10.91 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-Benzothiophene # | 134 | C8H6S | 000270-82-6 | 94 |
| 2 | Benzo[b]thiophene | 134 | C8H6S | 000095-15-8 | 94 |
| 3 | Benzo[b]thiophene | 134 | C8H6S | 000095-15-8 | 91 |
| 4 | Cyclopenta[c]thiapyran | 134 | C8H6S | 000270-63-3 | 91 |
| 5 | Benzo[b]thiophene | 134 | C8H6S | 000095-15-8 | 91 |



# Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG122115\
Data File : BG020400.D
Acq On : 22 Dec 2015 6:10
Operator : UM/SJ
Sample : G4848-04DL 30X
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** D9N50DL

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG121415.M
Quant Title : SVOA CALIBRATION

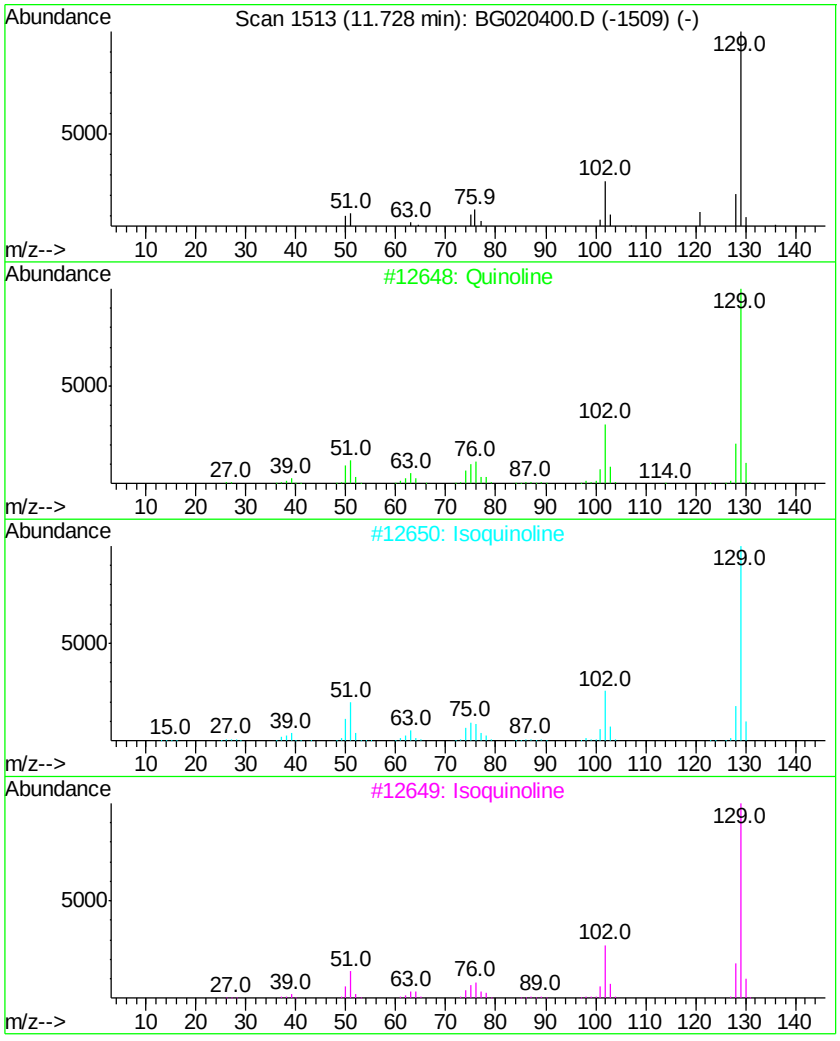
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

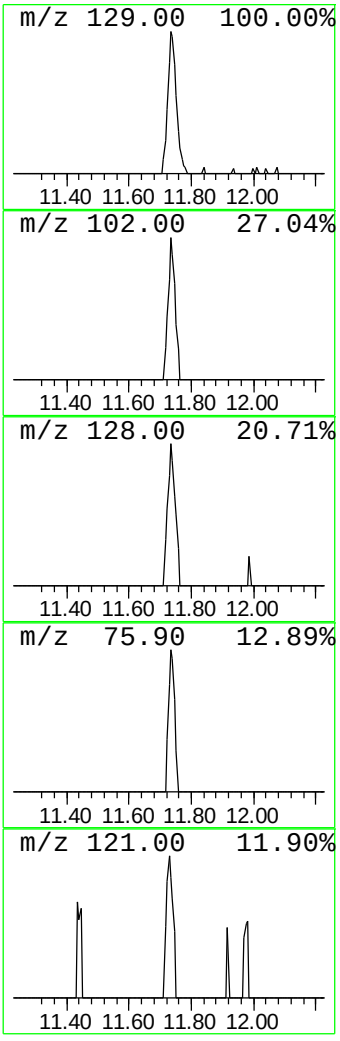
**********

Peak Number 3 Quinoline Concentration Rank 33

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.73 | 2.27 ng/ul | 17545 | Naphthalene-d8 | 10.91 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Quinoline | 129 | C9H7N | 000091-22-5 | 94 |
| 2 | Isoquinoline | 129 | C9H7N | 000119-65-3 | 91 |
| 3 | Isoquinoline | 129 | C9H7N | 000119-65-3 | 91 |
| 4 | Quinoline | 129 | C9H7N | 000091-22-5 | 91 |
| 5 | Quinoline | 129 | C9H7N | 000091-22-5 | 91 |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG122115\
Data File : BG020400.D
Acq On : 22 Dec 2015 6:10
Operator : UM/SJ
Sample : G4848-04DL 30X
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** D9N50DL

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG121415.M
Quant Title : SVOA CALIBRATION

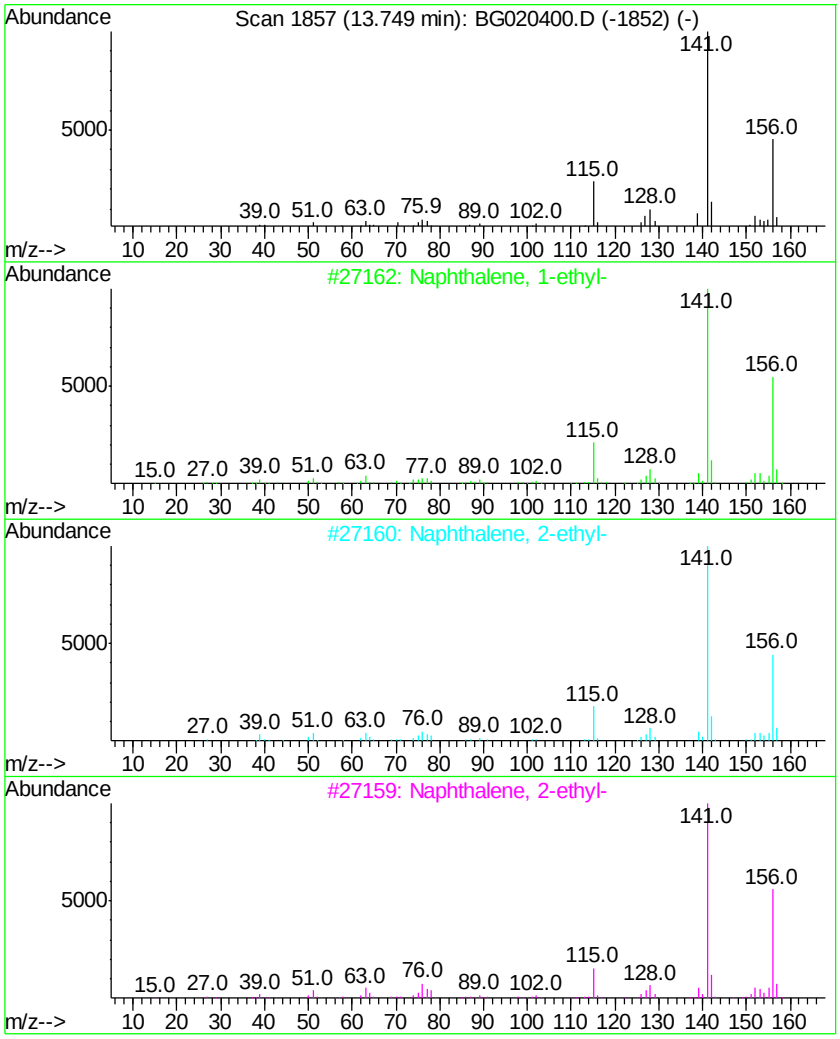
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

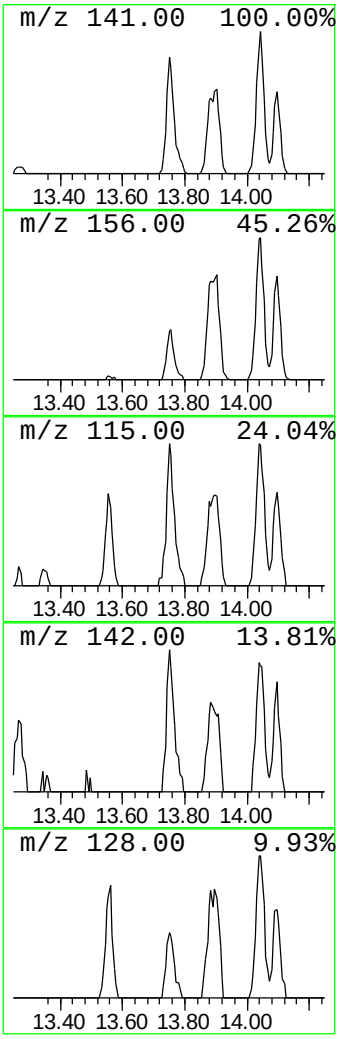
**********

Peak Number 5 Naphthalene, 1-ethyl- Concentration Rank 17

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.75 | 5.12 ng/ul | 56858 | Acenaphthene-d10 | 14.72 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Naphthalene, 1-ethyl- | 156 | C12H12 | 001127-76-0 | 95 |
| 2 | Naphthalene, 2-ethyl- | 156 | C12H12 | 000939-27-5 | 94 |
| 3 | Naphthalene, 2-ethyl- | 156 | C12H12 | 000939-27-5 | 94 |
| 4 | Naphthalene, 2-ethyl- | 156 | C12H12 | 000939-27-5 | 91 |
| 5 | Naphthalene, 1-ethyl- | 156 | C12H12 | 001127-76-0 | 91 |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG122115\
Data File : BG020400.D
Acq On : 22 Dec 2015 6:10
Operator : UM/SJ
Sample : G4848-04DL 30X
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** D9N50DL

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG121415.M
Quant Title : SVOA CALIBRATION

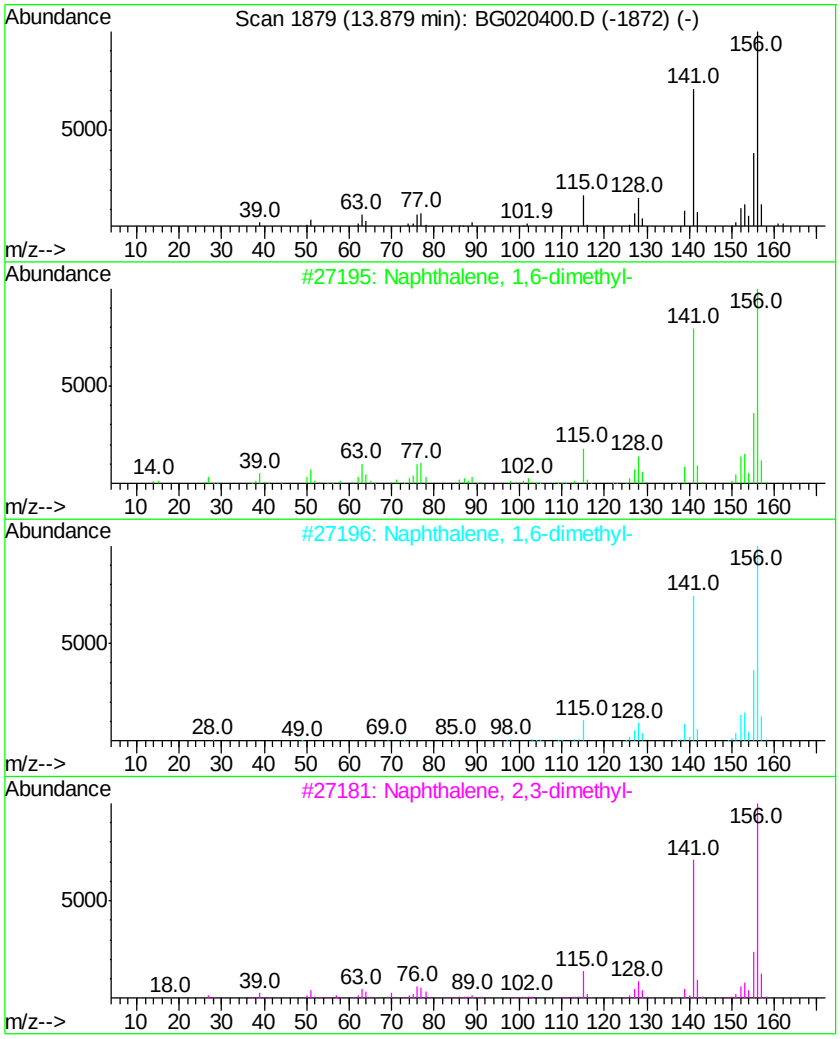
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

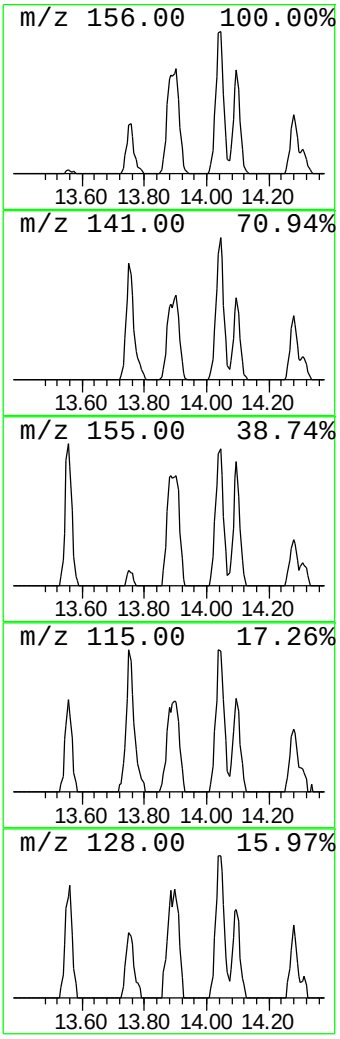
**********

Peak Number 6 Naphthalene, 1,6-dimethyl- Concentration Rank 15

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.88 | 5.41 ng/ul | 60011 | Acenaphthene-d10 | 14.72 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Naphthalene, 1,6-dimethyl- | 156 | C12H12 | 000575-43-9 | 96 |
| 2 | Naphthalene, 1,6-dimethyl- | 156 | C12H12 | 000575-43-9 | 96 |
| 3 | Naphthalene, 2,3-dimethyl- | 156 | C12H12 | 000581-40-8 | 96 |
| 4 | Naphthalene, 2,7-dimethyl- | 156 | C12H12 | 000582-16-1 | 96 |
| 5 | Naphthalene, 2,6-dimethyl- | 156 | C12H12 | 000581-42-0 | 95 |



# Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG122115\
Data File : BG020400.D
Acq On : 22 Dec 2015 6:10
Operator : UM/SJ
Sample : G4848-04DL 30X
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** D9N50DL

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG121415.M
Quant Title : SVOA CALIBRATION

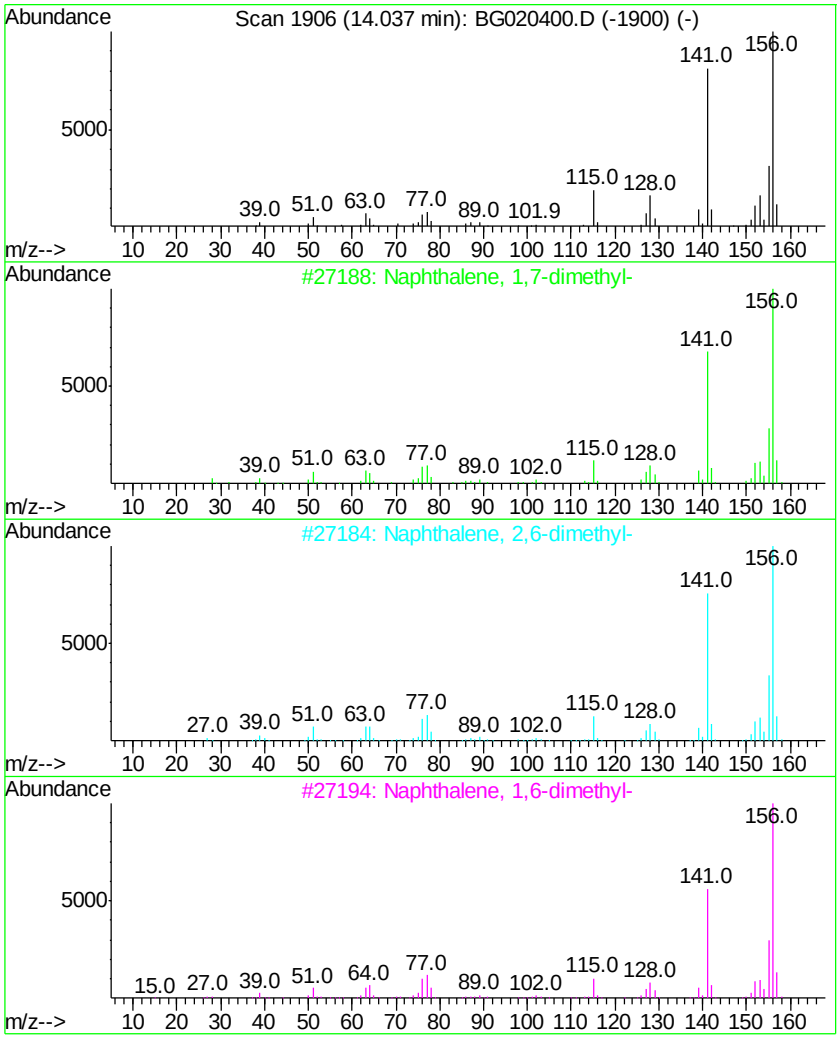
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

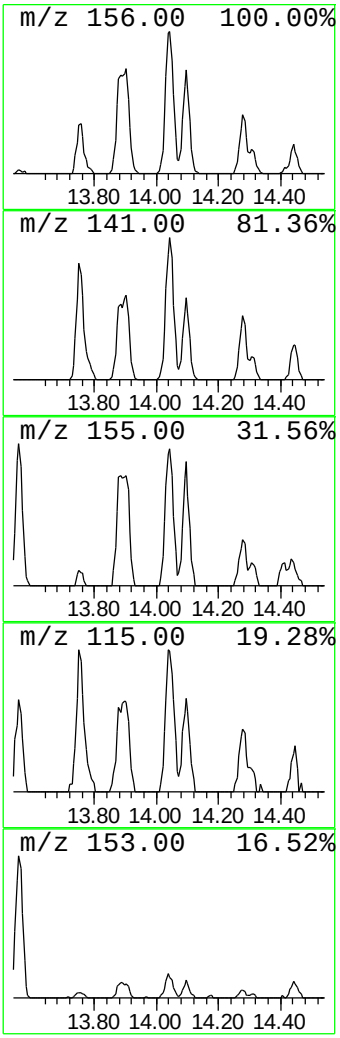
**********

Peak Number 7 Naphthalene, 1,7-dimethyl- Concentration Rank 6

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 14.04 | 10.17 ng/ul | 112874 | Acenaphthene-d10 | 14.72 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Naphthalene, 1,7-dimethyl- | 156 | C12H12 | 000575-37-1 | 97 |
| 2 | Naphthalene, 2,6-dimethyl- | 156 | C12H12 | 000581-42-0 | 97 |
| 3 | Naphthalene, 1,6-dimethyl- | 156 | C12H12 | 000575-43-9 | 96 |
| 4 | Naphthalene, 1,6-dimethyl- | 156 | C12H12 | 000575-43-9 | 96 |
| 5 | Naphthalene, 2,3-dimethyl- | 156 | C12H12 | 000581-40-8 | 95 |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG122115\
Data File : BG020400.D
Acq On : 22 Dec 2015 6:10
Operator : UM/SJ
Sample : G4848-04DL 30X
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** D9N50DL

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG121415.M
Quant Title : SVOA CALIBRATION

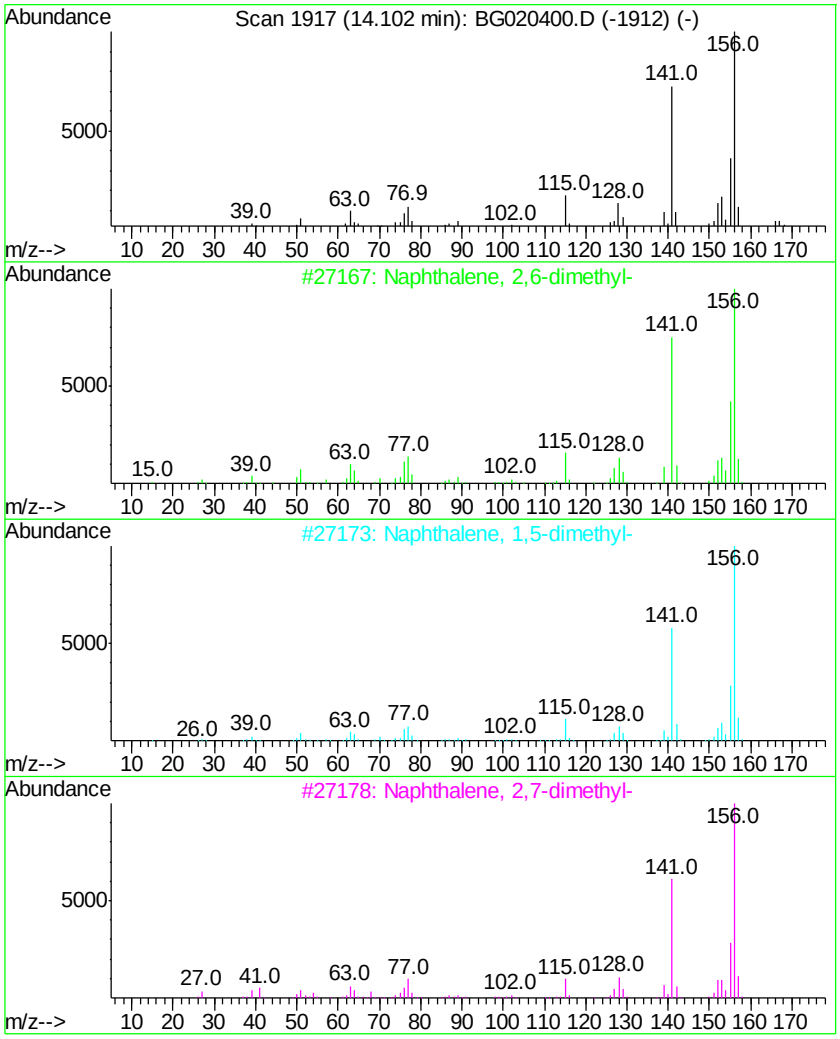
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

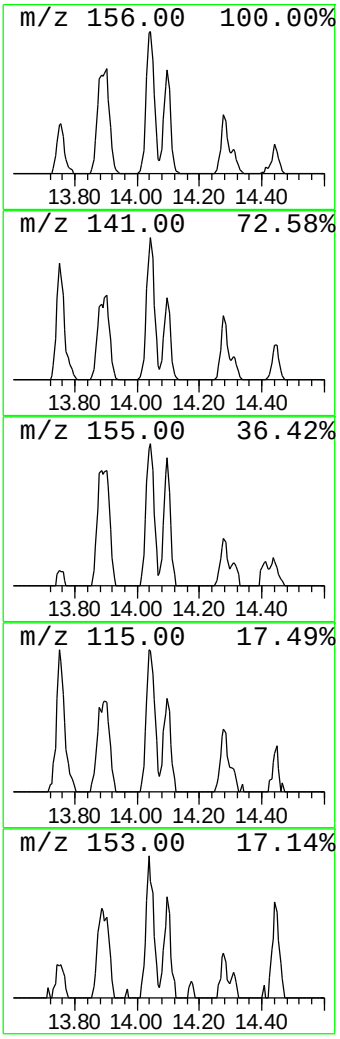
**********

Peak Number 8 Naphthalene, 2,6-dimethyl- Concentration Rank 13

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 14.10 | 6.56 ng/ul | 72854 | Acenaphthene-d10 | 14.72 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Naphthalene, 2,6-dimethyl- | 156 | C12H12 | 000581-42-0 | 97 |
| 2 | Naphthalene, 1,5-dimethyl- | 156 | C12H12 | 000571-61-9 | 97 |
| 3 | Naphthalene, 2,7-dimethyl- | 156 | C12H12 | 000582-16-1 | 96 |
| 4 | Naphthalene, 1,5-dimethyl- | 156 | C12H12 | 000571-61-9 | 96 |
| 5 | Naphthalene, 2,3-dimethyl- | 156 | C12H12 | 000581-40-8 | 96 |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG122115\
Data File : BG020400.D
Acq On : 22 Dec 2015 6:10
Operator : UM/SJ
Sample : G4848-04DL 30X
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** D9N50DL

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG121415.M
Quant Title : SVOA CALIBRATION

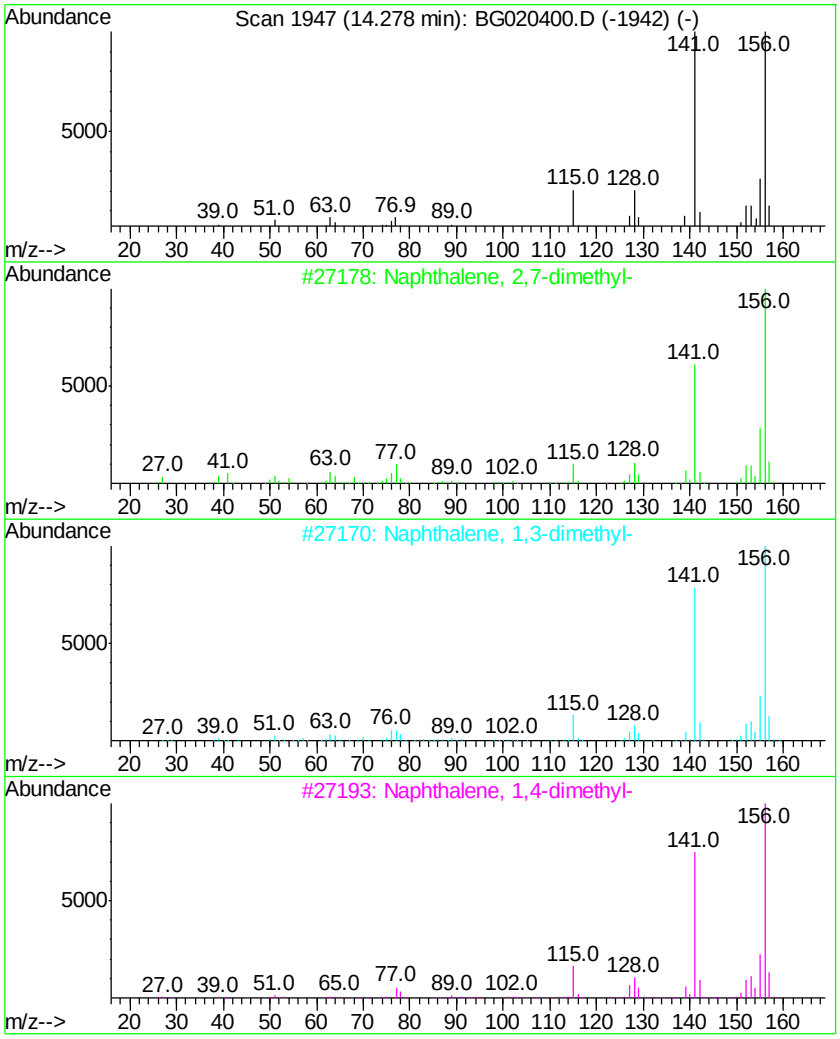
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

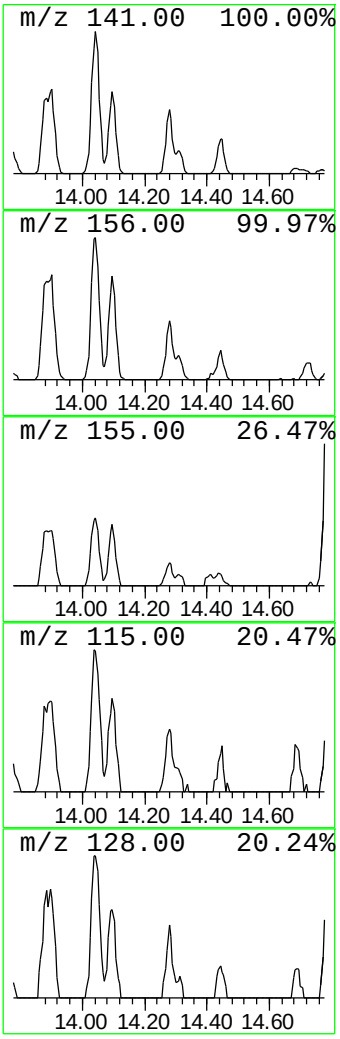
**********

Peak Number 9 Naphthalene, 2,7-dimethyl- Concentration Rank 27

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 14.28 | 3.24 ng/ul | 35925 | Acenaphthene-d10 | 14.72 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Naphthalene, 2,7-dimethyl- | 156 | C12H12 | 000582-16-1 | 96 |
| 2 | Naphthalene, 1,3-dimethyl- | 156 | C12H12 | 000575-41-7 | 95 |
| 3 | Naphthalene, 1,4-dimethyl- | 156 | C12H12 | 000571-58-4 | 95 |
| 4 | Naphthalene, 1,6-dimethyl- | 156 | C12H12 | 000575-43-9 | 95 |
| 5 | Naphthalene, 2,6-dimethyl- | 156 | C12H12 | 000581-42-0 | 95 |



# Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG122115\
Data File : BG020400.D
Acq On : 22 Dec 2015 6:10
Operator : UM/SJ
Sample : G4848-04DL 30X
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** D9N50DL

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG121415.M
Quant Title : SVOA CALIBRATION

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

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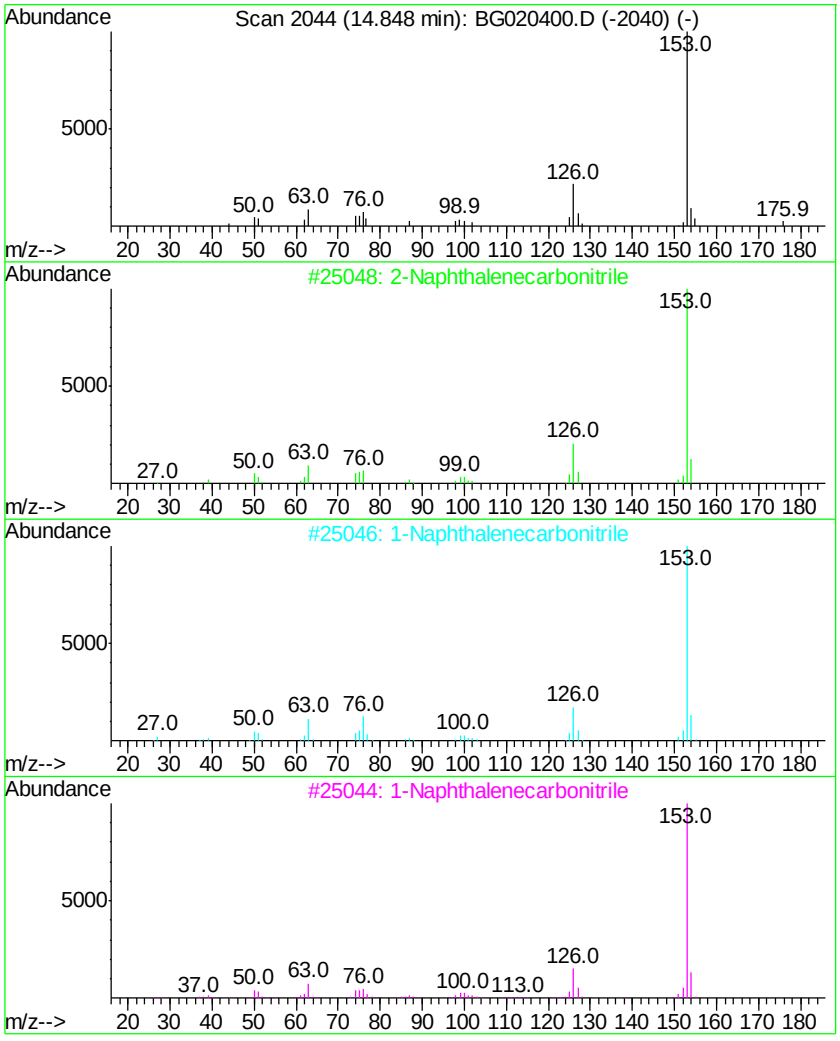
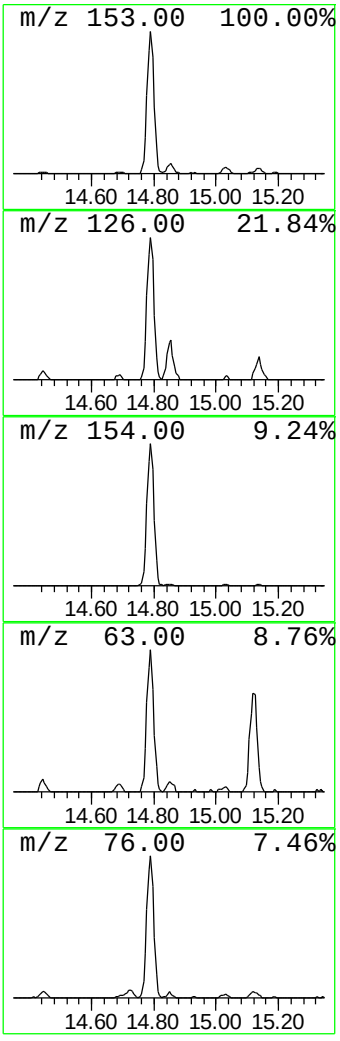
Peak Number 10 2-Naphthalenecarbonitrile Concentration Rank 26

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 14.85 | 3.34 ng/ul | 37029 | Acenaphthene-d10 | 14.72 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-Naphthalenecarbonitrile | 153 | C11H7N | 000613-46-7 | 91 |
| 2 | 1-Naphthalenecarbonitrile | 153 | C11H7N | 000086-53-3 | 91 |
| 3 | 1-Naphthalenecarbonitrile | 153 | C11H7N | 000086-53-3 | 91 |
| 4 | 2-Naphthalenecarbonitrile | 153 | C11H7N | 000613-46-7 | 80 |
| 5 | 1-Naphthalenecarbonitrile | 153 | C11H7N | 000086-53-3 | 80 |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG122115\
Data File : BG020400.D
Acq On : 22 Dec 2015 6:10
Operator : UM/SJ
Sample : G4848-04DL 30X
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** D9N50DL

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG121415.M
Quant Title : SVOA CALIBRATION

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

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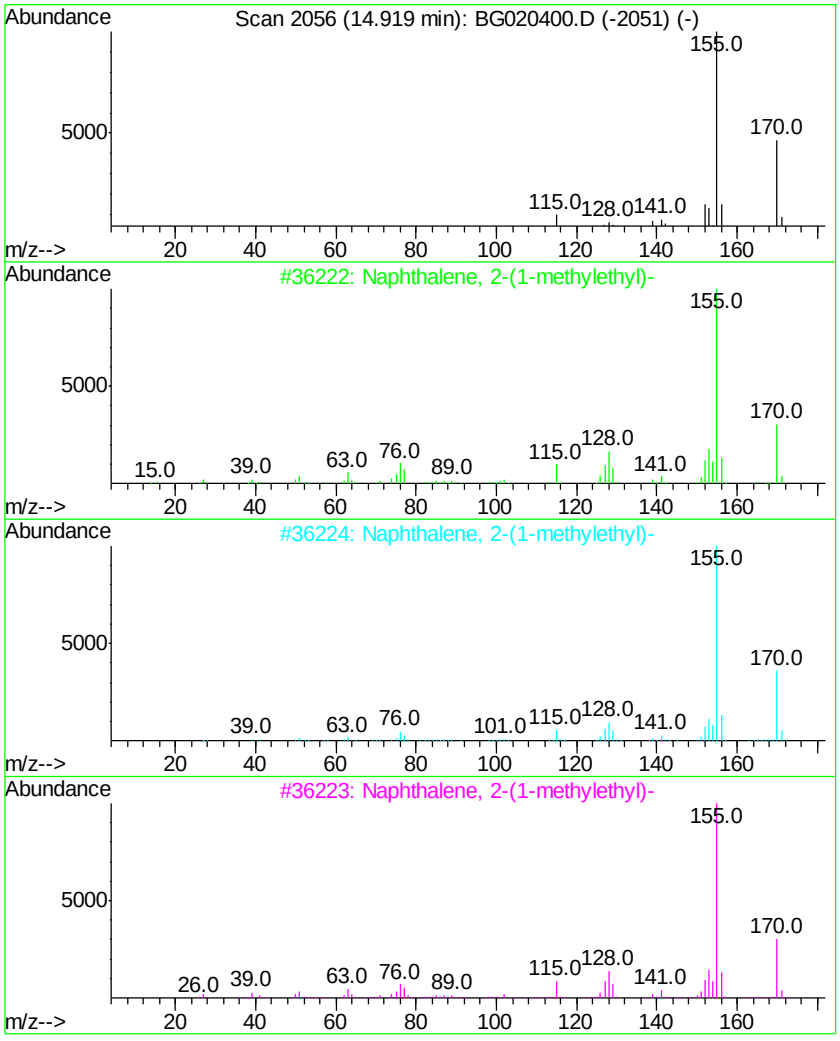
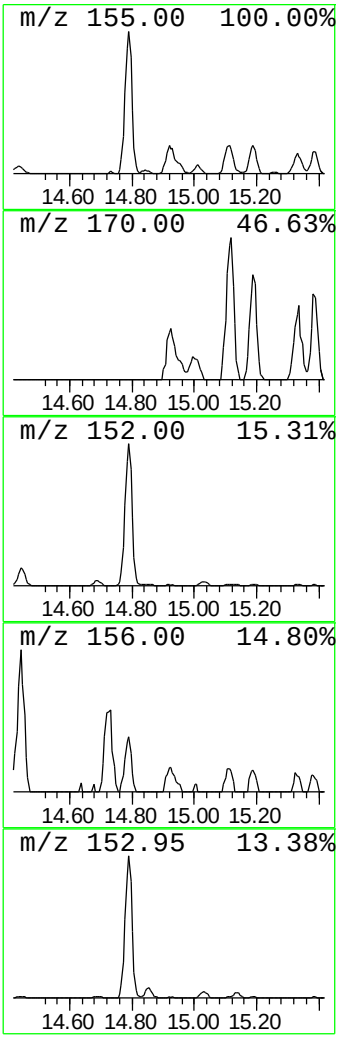
Peak Number 11 Naphthalene, 2-(1-methyleth... Concentration Rank 36

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 14.92 | 2.01 ng/ul | 22338 | Acenaphthene-d10 | 14.72 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Naphthalene, 2-(1-methylethyl)- | 170 | C13H14 | 002027-17-0 | 86 |
| 2 | Naphthalene, 2-(1-methylethyl)- | 170 | C13H14 | 002027-17-0 | 86 |
| 3 | Naphthalene, 2-(1-methylethyl)- | 170 | C13H14 | 002027-17-0 | 86 |
| 4 | 3-(2-Methyl-propenyl)-1H-indene | 170 | C13H14 | 1000187-78-5 | 83 |
| 5 | 1'-Acetonaphthone | 170 | C12H10O | 000941-98-0 | 64 |



# Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG122115\
Data File : BG020400.D
Acq On : 22 Dec 2015 6:10
Operator : UM/SJ
Sample : G4848-04DL 30X
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** D9N50DL

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG121415.M
Quant Title : SVOA CALIBRATION

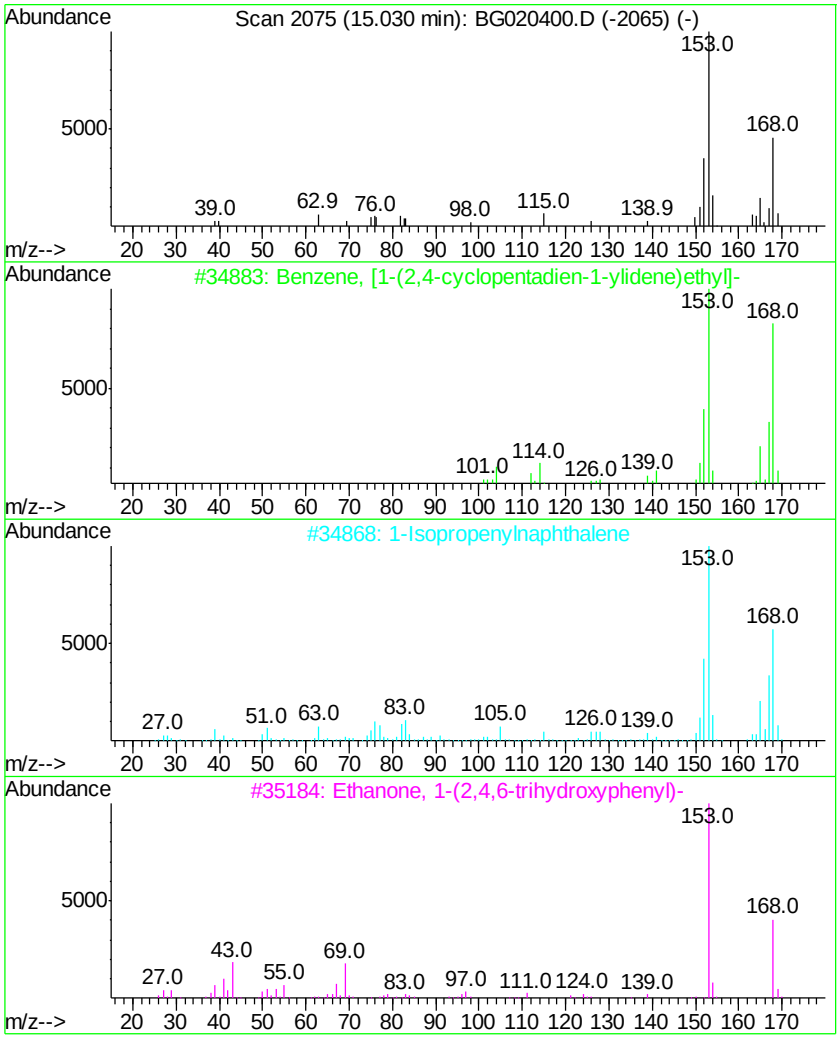
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

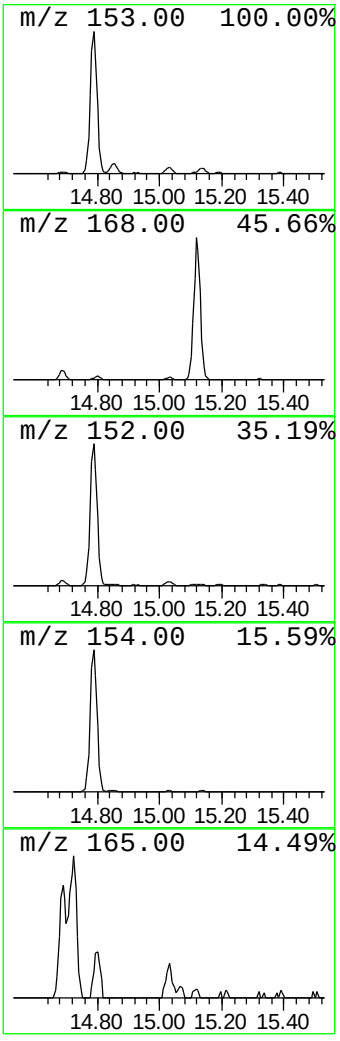
**********

Peak Number 12 Benzene, [1-(2,4-cyclopenta... Concentration Rank 24

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.03 | 3.55 ng/ul | 39462 | Acenaphthene-d10 | 14.72 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzene, [1-(2,4-cyclopentadien-... | 168 | C13H12 | 002320-32-3 | 72 |
| 2 | 1-Isopropenylnaphthalene | 168 | C13H12 | 001855-47-6 | 58 |
| 3 | Ethanone, 1-(2,4,6-trihydroxyphe... | 168 | C8H8O4 | 000480-66-0 | 50 |
| 4 | Ethanone, 1-(2,4,6-trihydroxyphe... | 168 | C8H8O4 | 000480-66-0 | 50 |
| 5 | Ethanone, 1-(2,3,4-trihydroxyphe... | 168 | C8H8O4 | 000528-21-2 | 50 |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG122115\
Data File : BG020400.D
Acq On : 22 Dec 2015 6:10
Operator : UM/SJ
Sample : G4848-04DL 30X
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** D9N50DL

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG121415.M
Quant Title : SVOA CALIBRATION

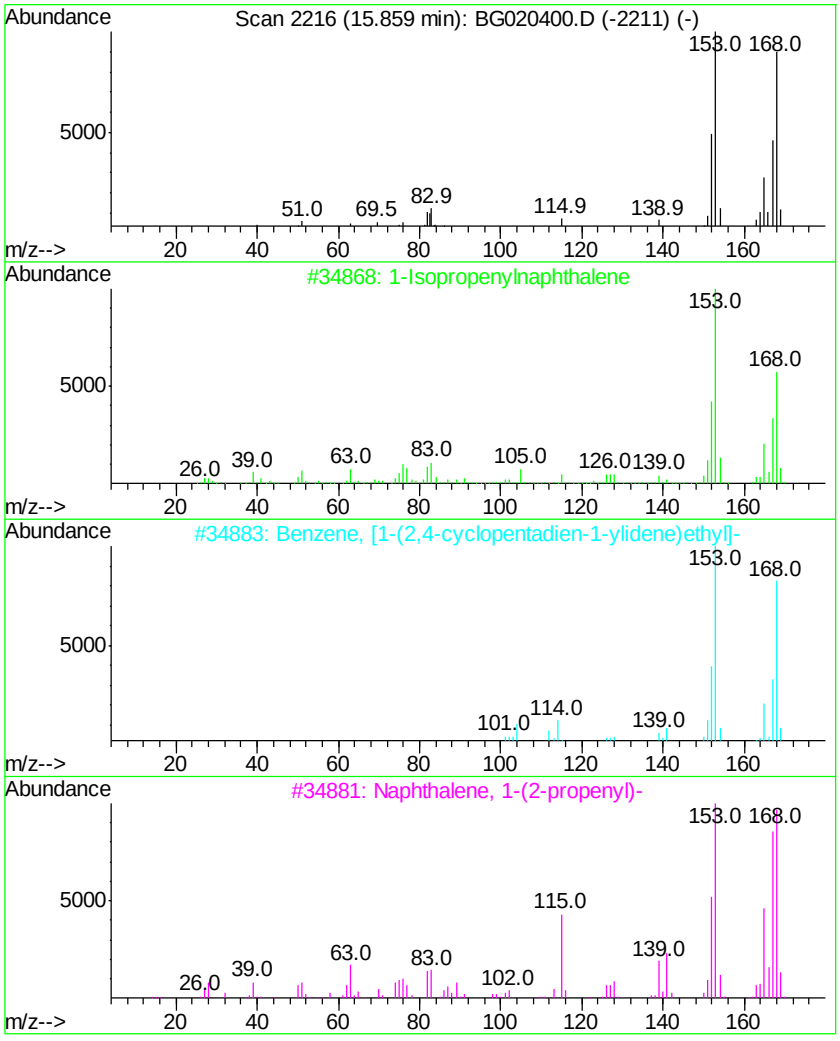
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

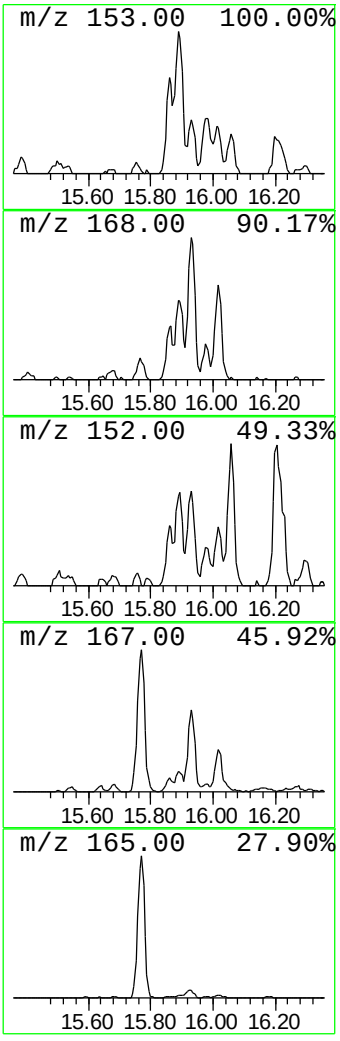
**********

Peak Number 13 1-Isopropenylnaphthalene Concentration Rank 28

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.86 | 2.85 ng/ul | 31649 | Acenaphthene-d10 | 14.72 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1-Isopropenylnaphthalene | 168 | C13H12 | 001855-47-6 | 81 |
| 2 | Benzene, [1-(2,4-cyclopentadien-... | 168 | C13H12 | 002320-32-3 | 72 |
| 3 | Naphthalene, 1-(2-propenyl)- | 168 | C13H12 | 002489-86-3 | 58 |
| 4 | Naphthalene, 1-(2-propenyl)- | 168 | C13H12 | 002489-86-3 | 53 |
| 5 | Naphthalene, 2-(1-methylethenyl)- | 168 | C13H12 | 003710-23-4 | 50 |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG122115\
Data File : BG020400.D
Acq On : 22 Dec 2015 6:10
Operator : UM/SJ
Sample : G4848-04DL 30X
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** D9N50DL

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG121415.M
Quant Title : SVOA CALIBRATION

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

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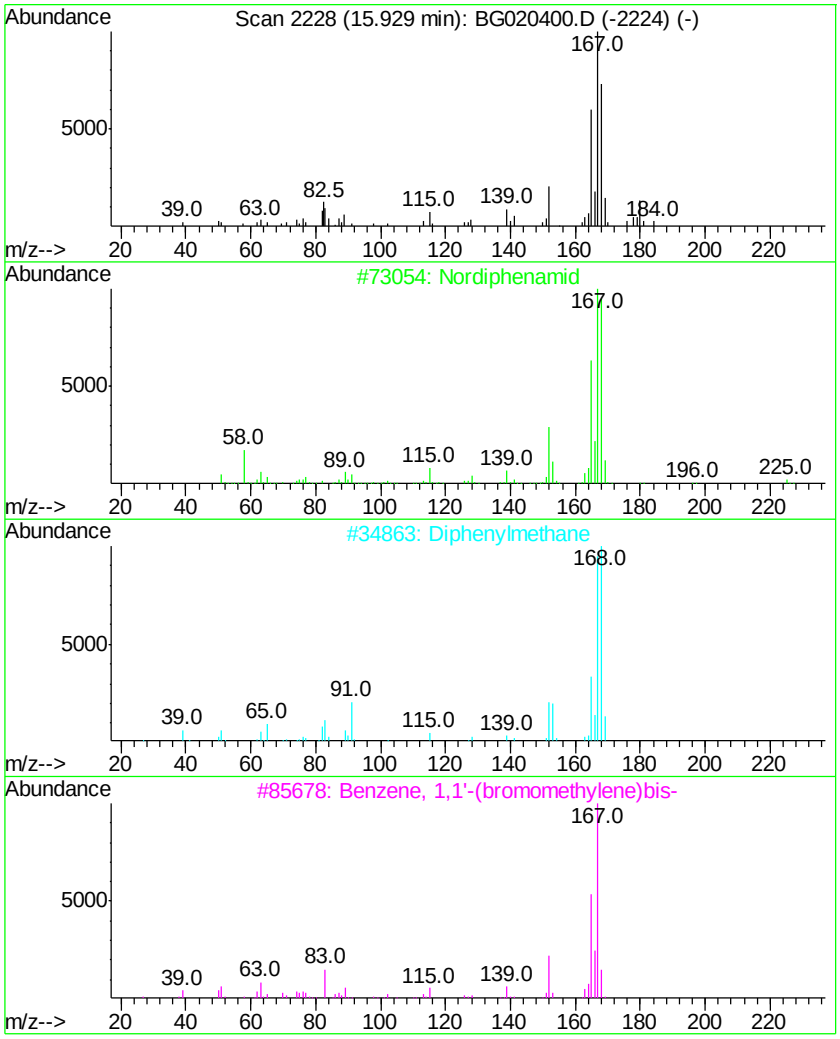
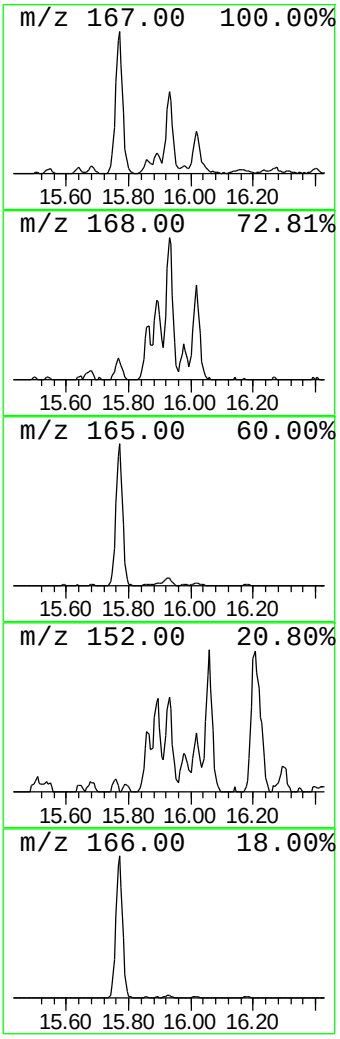
Peak Number 15 Nordiphenamid Concentration Rank 9

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.93 | 9.69 ng/ul | 107532 | Acenaphthene-d10 | 14.72 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Nordiphenamid | 225 | C15H15NO | 000954-21-2 | 80 |
| 2 | Diphenylmethane | 168 | C13H12 | 000101-81-5 | 52 |
| 3 | Benzene, 1,1'-(bromomethylene)bis- | 246 | C13H11Br | 000776-74-9 | 50 |
| 4 | Benzhydryl isothiocyanate | 225 | C14H11NS | 003550-21-8 | 50 |
| 5 | Benzene, 1,1'-(bromomethylene)bis- | 246 | C13H11Br | 000776-74-9 | 47 |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG122115\
Data File : BG020400.D
Acq On : 22 Dec 2015 6:10
Operator : UM/SJ
Sample : G4848-04DL 30X
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** D9N50DL

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG121415.M
Quant Title : SVOA CALIBRATION

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

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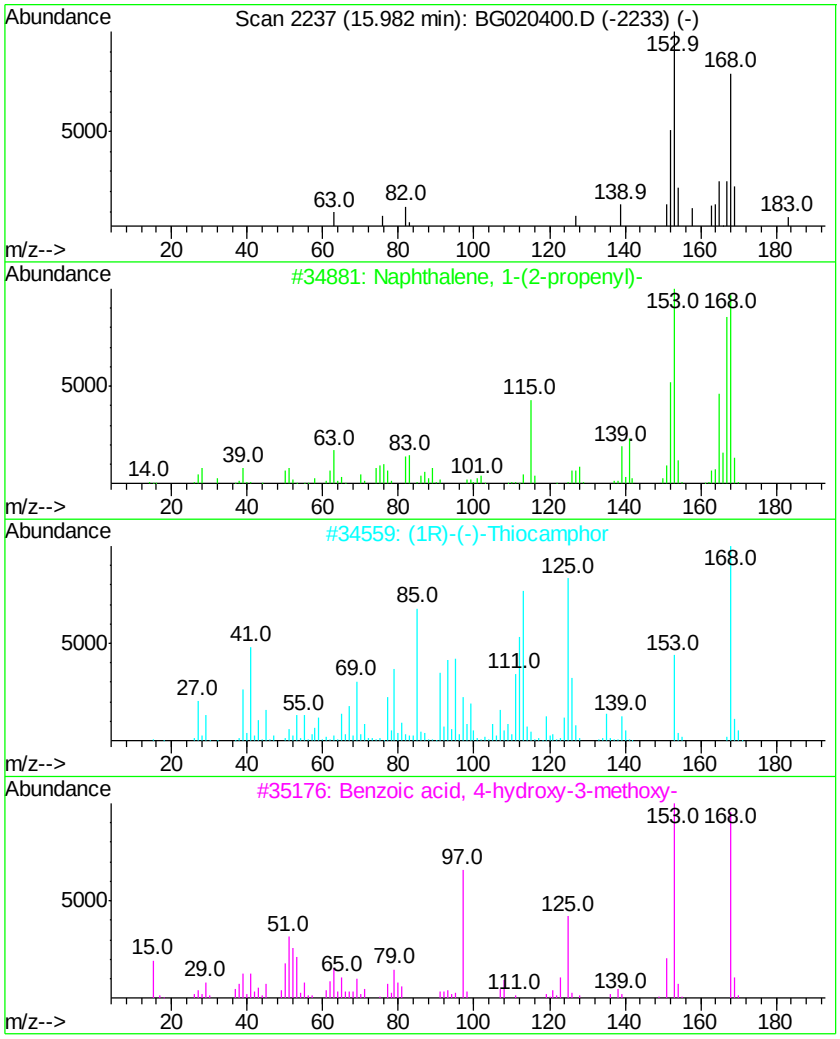
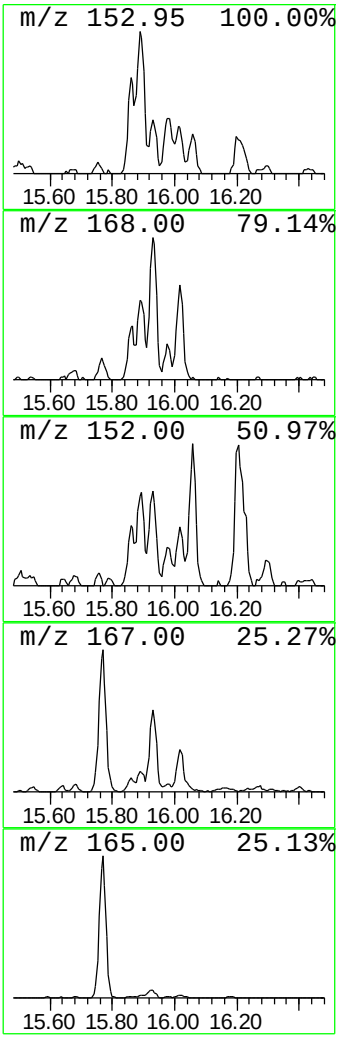
Peak Number 16 Naphthalene, 1-(2-propenyl)- Concentration Rank 35

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.98 | 2.09 ng/ul | 23190 | Acenaphthene-d10 | 14.72 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Naphthalene, 1-(2-propenyl)- | 168 | C13H12 | 002489-86-3 | 72 |
| 2 | (1R)-(-)-Thiocamphor | 168 | C10H16S | 053402-10-1 | 53 |
| 3 | Benzoic acid, 4-hydroxy-3-methoxy- | 168 | C8H8O4 | 000121-34-6 | 47 |
| 4 | 3-Hydroxy-4-methoxybenzoic acid | 168 | C8H8O4 | 000645-08-9 | 47 |
| 5 | 1-Isopropenylnaphthalene | 168 | C13H12 | 001855-47-6 | 46 |



# Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG122115\
Data File : BG020400.D
Acq On : 22 Dec 2015 6:10
Operator : UM/SJ
Sample : G4848-04DL 30X
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** D9N50DL

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG121415.M
Quant Title : SVOA CALIBRATION

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

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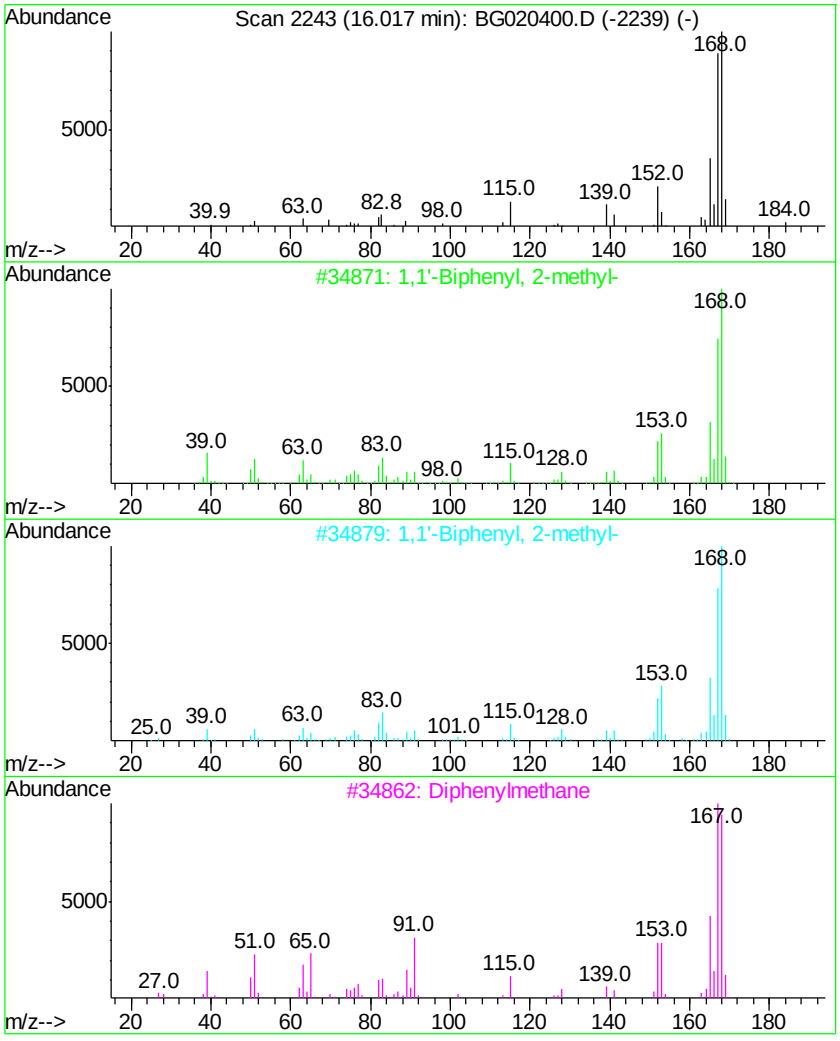
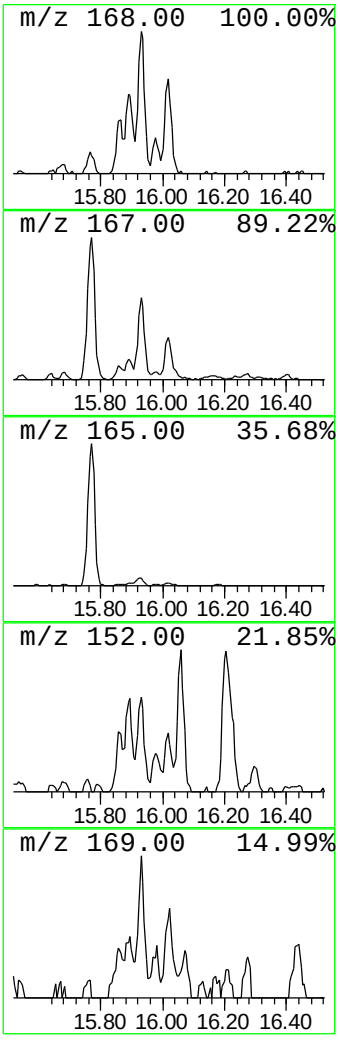
Peak Number 17 1,1'-Biphenyl, 2-methyl- Concentration Rank 20

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.02 | 4.65 ng/ul | 51639 | Acenaphthene-d10 | 14.72 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1,1'-Biphenyl, 2-methyl- | 168 | C13H12 | 000643-58-3 | 87 |
| 2 | 1,1'-Biphenyl, 2-methyl- | 168 | C13H12 | 000643-58-3 | 87 |
| 3 | Diphenylmethane | 168 | C13H12 | 000101-81-5 | 80 |
| 4 | Diphenylmethane | 168 | C13H12 | 000101-81-5 | 80 |
| 5 | 1,1'-Biphenyl, 4-methyl- | 168 | C13H12 | 000644-08-6 | 80 |



# Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG122115\
Data File : BG020400.D
Acq On : 22 Dec 2015 6:10
Operator : UM/SJ
Sample : G4848-04DL 30X
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** D9N50DL

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG121415.M
Quant Title : SVOA CALIBRATION

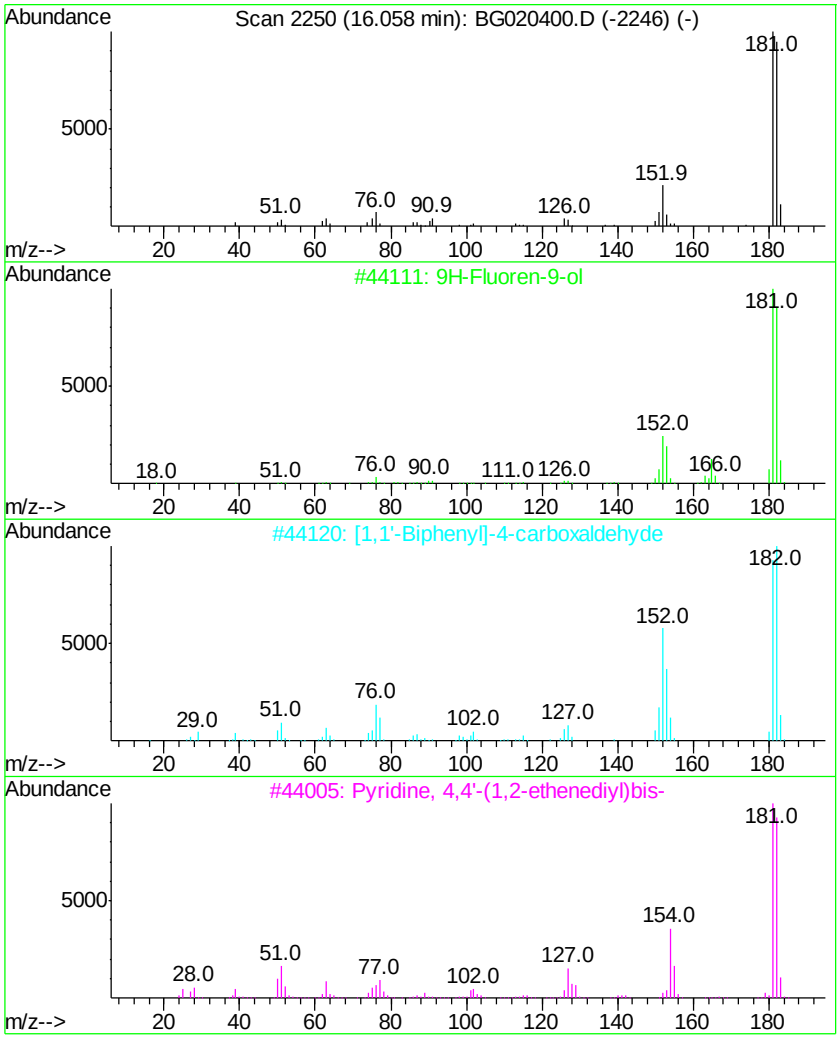
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

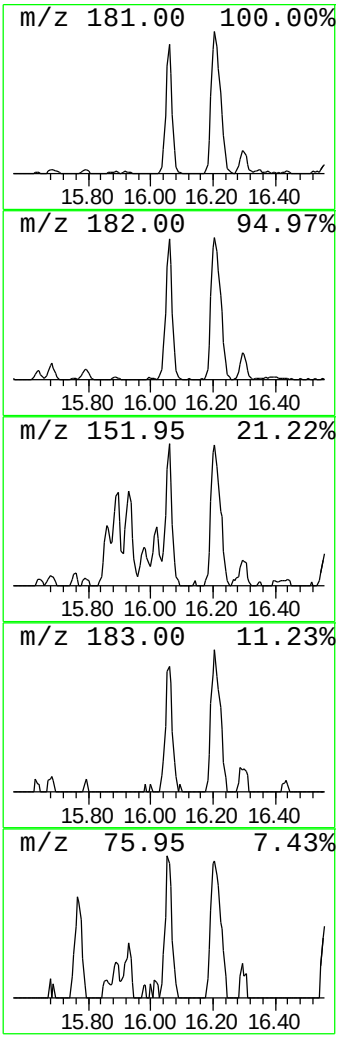
**********

Peak Number 18 9H-Fluoren-9-ol Concentration Rank 8

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.06 | 9.93 ng/ul | 110275 | Acenaphthene-d10 | 14.72 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 9H-Fluoren-9-ol | 182 | C13H10O | 001689-64-1 | 86 |
| 2 | [1,1'-Biphenyl]-4-carboxaldehyde | 182 | C13H10O | 003218-36-8 | 78 |
| 3 | Pyridine, 4,4'-(1,2-ethenediyl)bis- | 182 | C12H10N2 | 001135-32-6 | 72 |
| 4 | Pyrido[2,3-d]indole, 6-methyl- | 182 | C12H10N2 | 108349-67-3 | 72 |
| 5 | 9H-Fluoren-9-ol | 182 | C13H10O | 001689-64-1 | 72 |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG122115\
Data File : BG020400.D
Acq On : 22 Dec 2015 6:10
Operator : UM/SJ
Sample : G4848-04DL 30X
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** D9N50DL

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG121415.M
Quant Title : SVOA CALIBRATION

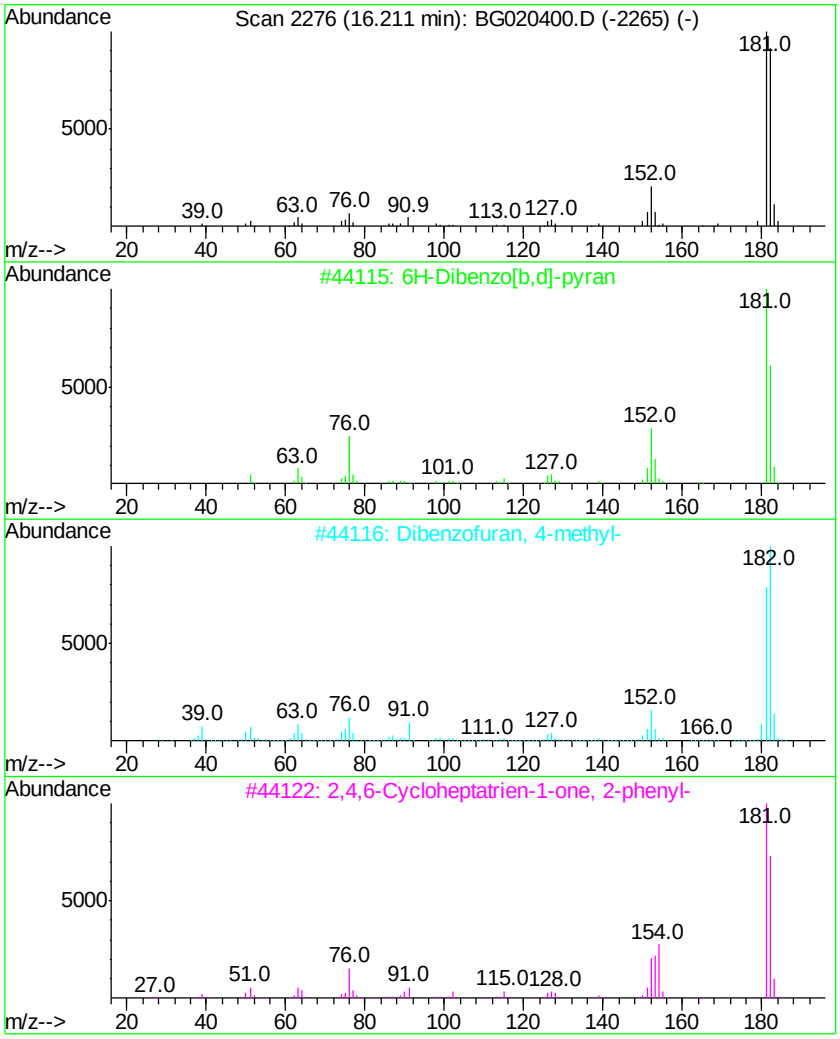
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

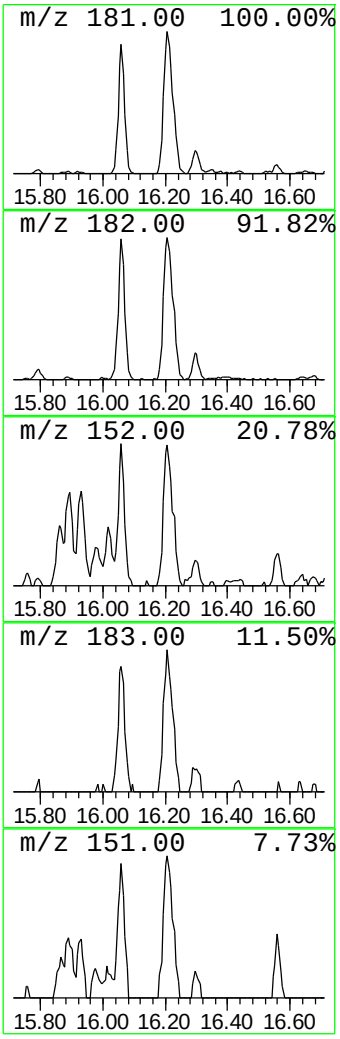
**********

Peak Number 19 6H-Dibenzo[b,d]-pyran Concentration Rank 5

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.21 | 10.94 ng/ul | 166043 | Phenanthrene-d10 | 17.47 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 6H-Dibenzo[b,d]-pyran | 182 | C13H10O | 000229-95-8 | 91 |
| 2 | Dibenzofuran, 4-methyl- | 182 | C13H10O | 007320-53-8 | 90 |
| 3 | 2,4,6-Cycloheptatrien-1-one, 2-p... | 182 | C13H10O | 014562-09-5 | 90 |
| 4 | 9H-Fluoren-9-ol | 182 | C13H10O | 001689-64-1 | 86 |
| 5 | [1,1'-Biphenyl]-4-carboxaldehyde | 182 | C13H10O | 003218-36-8 | 83 |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG122115\
Data File : BG020400.D
Acq On : 22 Dec 2015 6:10
Operator : UM/SJ
Sample : G4848-04DL 30X
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** D9N50DL

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG121415.M
Quant Title : SVOA CALIBRATION

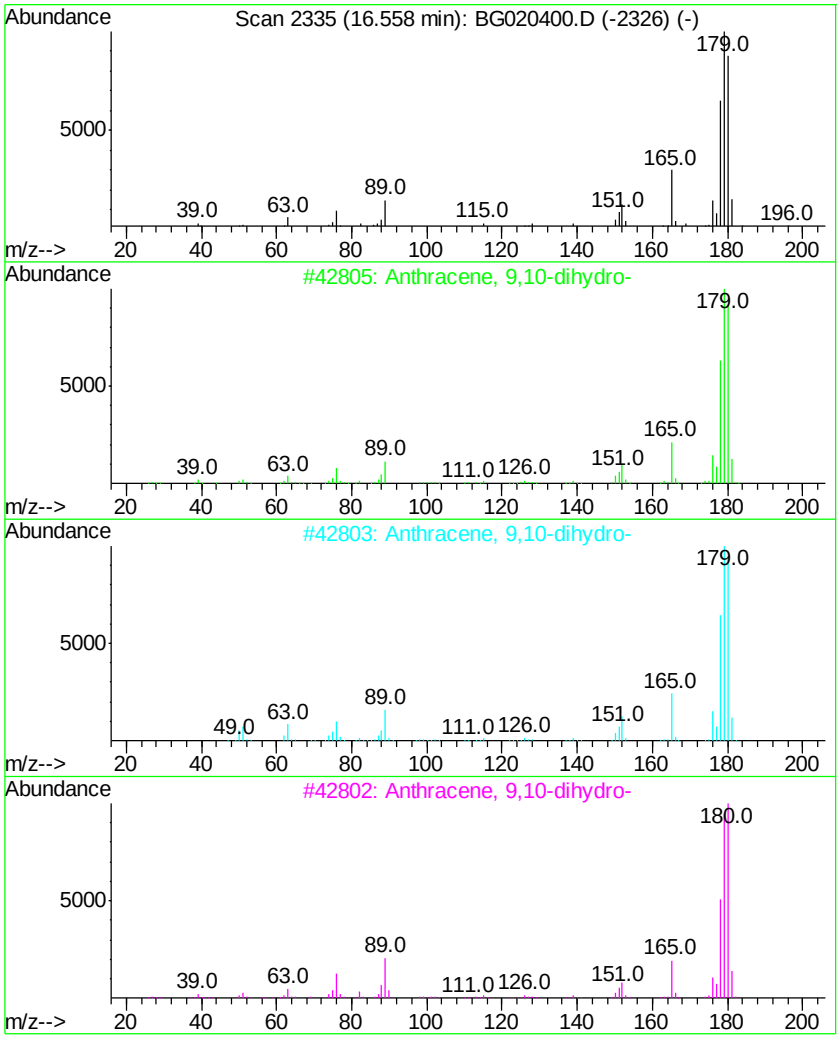
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

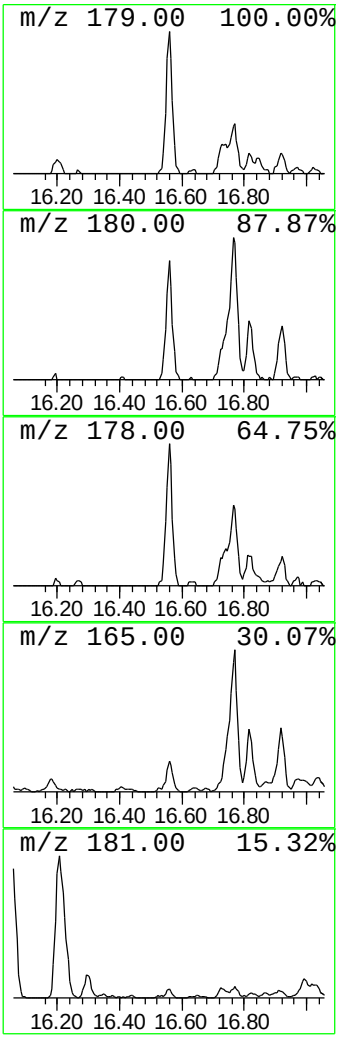
**********

Peak Number 20 Anthracene, 9,10-dihydro- Concentration Rank 22

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.56 | 4.11 ng/ul | 62337 | Phenanthrene-d10 | 17.47 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Anthracene, 9,10-dihydro- | 180 | C14H12 | 000613-31-0 | 96 |
| 2 | Anthracene, 9,10-dihydro- | 180 | C14H12 | 000613-31-0 | 96 |
| 3 | Anthracene, 9,10-dihydro- | 180 | C14H12 | 000613-31-0 | 95 |
| 4 | Anthracene, 9,10-dihydro- | 180 | C14H12 | 000613-31-0 | 94 |
| 5 | 1,2-Diphenylethylene | 180 | C14H12 | 000588-59-0 | 92 |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG122115\
Data File : BG020400.D
Acq On : 22 Dec 2015 6:10
Operator : UM/SJ
Sample : G4848-04DL 30X
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** D9N50DL

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG121415.M
Quant Title : SVOA CALIBRATION

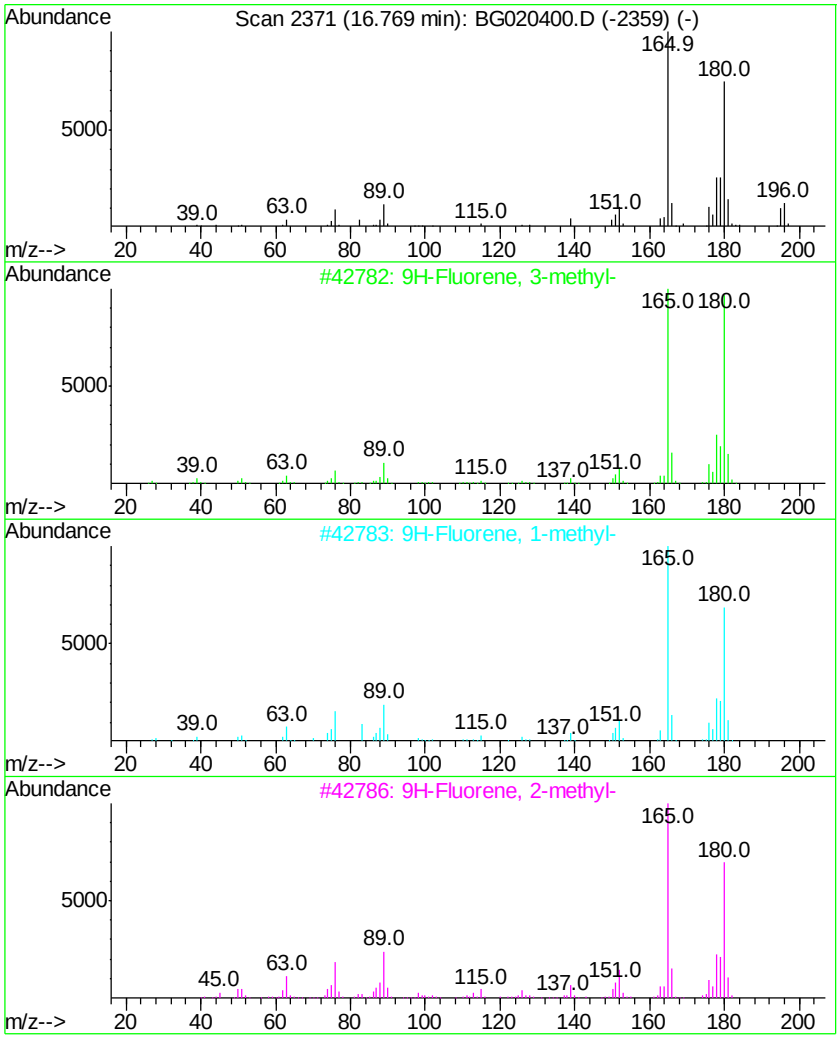
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

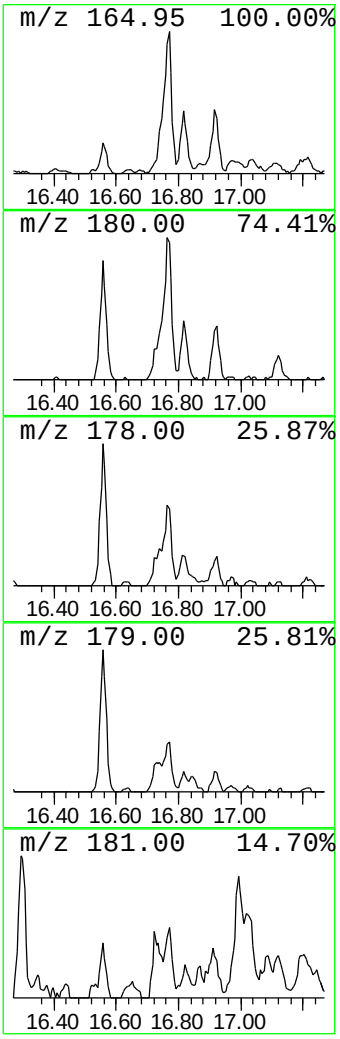
**********

Peak Number 21 9H-Fluorene, 3-methyl- Concentration Rank 11

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.77 | 8.14 ng/ul | 123601 | Phenanthrene-d10 | 17.47 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 9H-Fluorene, 3-methyl- | 180 | C14H12 | 002523-39-9 | 96 |
| 2 | 9H-Fluorene, 1-methyl- | 180 | C14H12 | 001730-37-6 | 96 |
| 3 | 9H-Fluorene, 2-methyl- | 180 | C14H12 | 001430-97-3 | 96 |
| 4 | 4a,9a-Methano-9H-fluorene | 180 | C14H12 | 019540-84-2 | 92 |
| 5 | 9H-Fluorene, 1-methyl- | 180 | C14H12 | 001730-37-6 | 91 |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG122115\
Data File : BG020400.D
Acq On : 22 Dec 2015 6:10
Operator : UM/SJ
Sample : G4848-04DL 30X
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** D9N50DL

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG121415.M
Quant Title : SVOA CALIBRATION

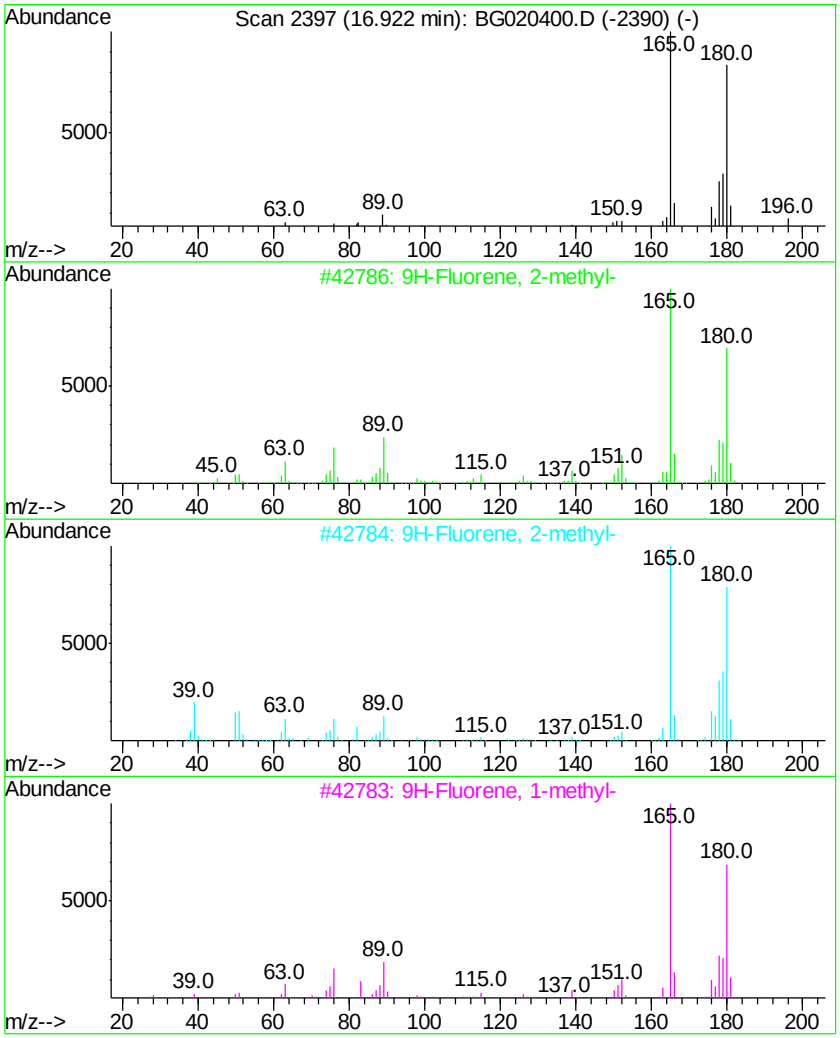
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

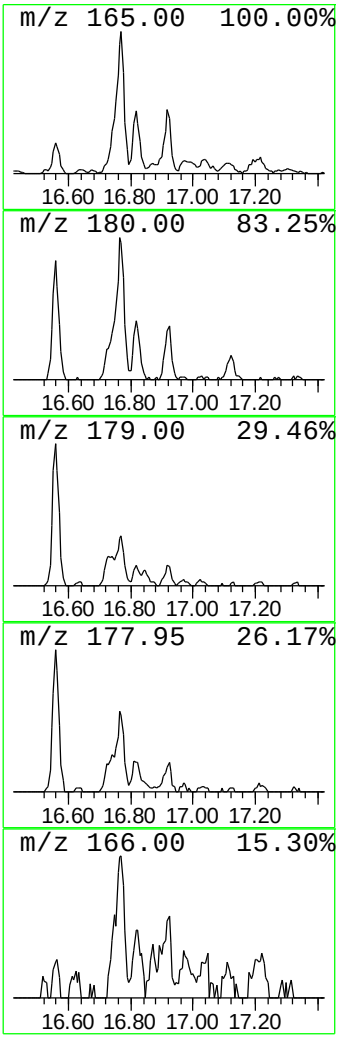
**********

Peak Number 22 9H-Fluorene, 2-methyl- Concentration Rank 31

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.92 | 2.42 ng/ul | 36712 | Phenanthrene-d10 | 17.47 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 9H-Fluorene, 2-methyl- | 180 | C14H12 | 001430-97-3 | 81 |
| 2 | 9H-Fluorene, 2-methyl- | 180 | C14H12 | 001430-97-3 | 81 |
| 3 | 9H-Fluorene, 1-methyl- | 180 | C14H12 | 001730-37-6 | 81 |
| 4 | 9H-Fluorene, 9-methyl- | 180 | C14H12 | 002523-37-7 | 81 |
| 5 | 9H-Fluorene, 3-methyl- | 180 | C14H12 | 002523-39-9 | 81 |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG122115\
Data File : BG020400.D
Acq On : 22 Dec 2015 6:10
Operator : UM/SJ
Sample : G4848-04DL 30X
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** D9N50DL

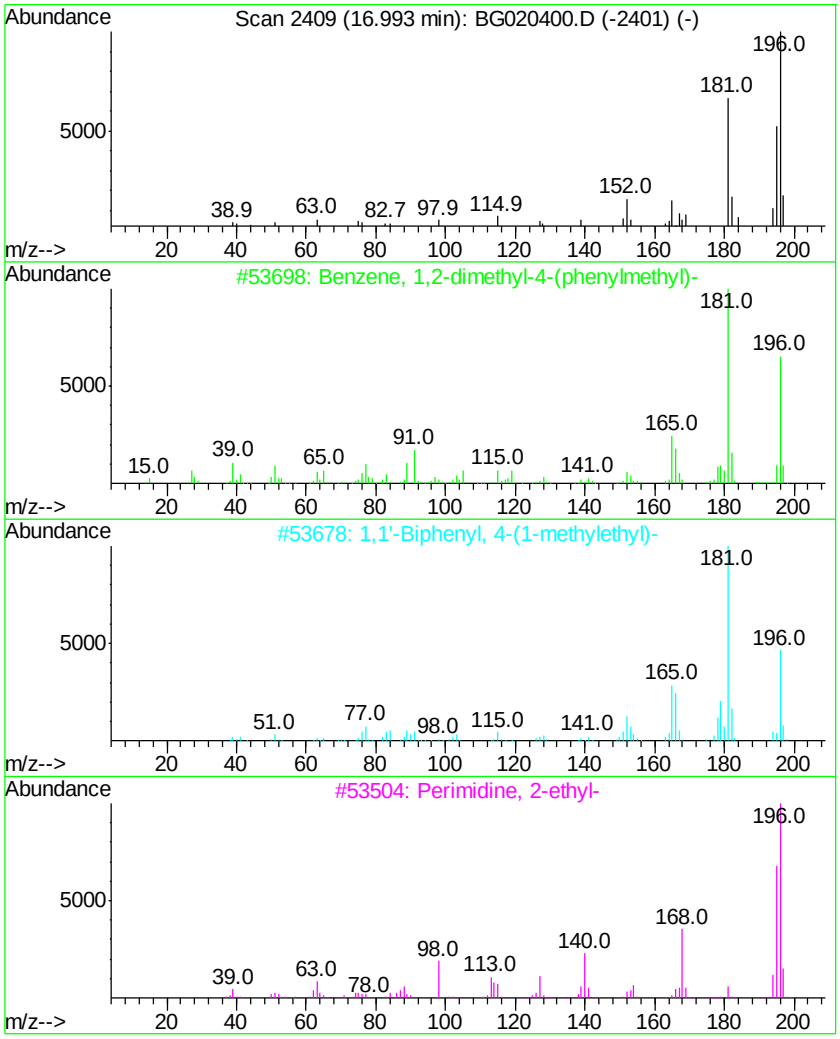
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG121415.M
Quant Title : SVOA CALIBRATION

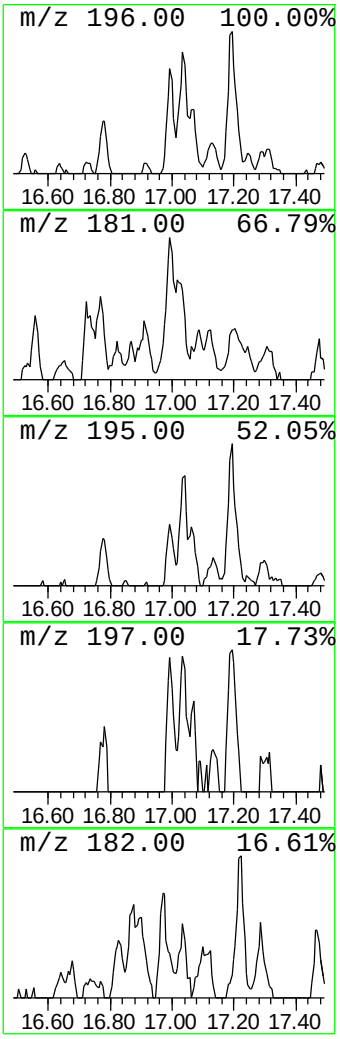
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 23 Benzene, 1,2-dimethyl-4-(ph... Concentration Rank 30

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.99 | 2.50 ng/ul | 37947 | Phenanthrene-d10 | 17.47 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzene, 1,2-dimethyl-4-(phenylm... | 196 | C15H16 | 013540-56-2 | 58 |
| 2 | 1,1'-Biphenyl, 4-(1-methylethyl)- | 196 | C15H16 | 007116-95-2 | 53 |
| 3 | Perimidine, 2-ethyl- | 196 | C13H12N2 | 028478-13-9 | 52 |
| 4 | 9H-Fluorene, 2-methoxy- | 196 | C14H12O | 002523-46-8 | 50 |
| 5 | 9H-Fluorene-2,7-diamine | 196 | C13H12N2 | 000525-64-4 | 49 |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG122115\
Data File : BG020400.D
Acq On : 22 Dec 2015 6:10
Operator : UM/SJ
Sample : G4848-04DL 30X
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** D9N50DL

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG121415.M
Quant Title : SVOA CALIBRATION

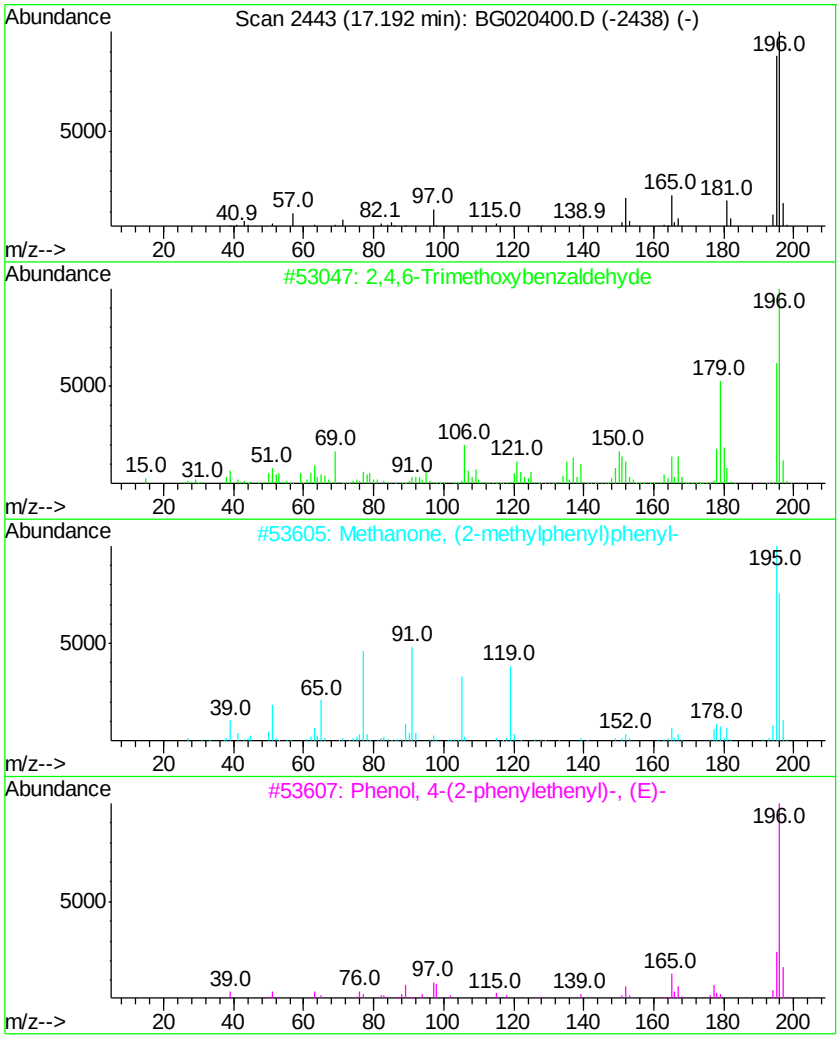
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

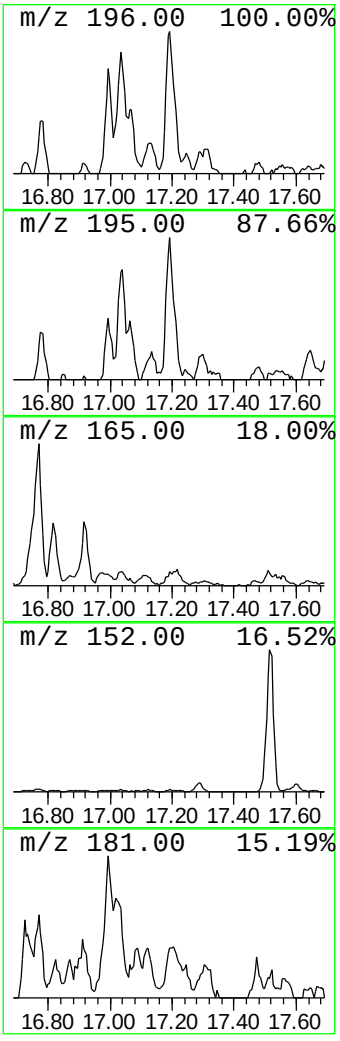
**********

Peak Number 24 2,4,6-Trimethoxybenzaldehyde Concentration Rank 21

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 17.19 | 4.22 ng/ul | 64046 | Phenanthrene-d10 | 17.47 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2,4,6-Trimethoxybenzaldehyde | 196 | C10H12O4 | 000830-79-5 | 72 |
| 2 | Methanone, (2-methylphenyl)phenyl- | 196 | C14H12O | 000131-58-8 | 72 |
| 3 | Phenol, 4-(2-phenylethenyl)-, (E)- | 196 | C14H12O | 006554-98-9 | 52 |
| 4 | Phenol, 4-(2-phenylethenyl)- | 196 | C14H12O | 003839-46-1 | 50 |
| 5 | Phenol, 4-(2-phenylethenyl)-, (E)- | 196 | C14H12O | 006554-98-9 | 50 |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG122115\
Data File : BG020400.D
Acq On : 22 Dec 2015 6:10
Operator : UM/SJ
Sample : G4848-04DL 30X
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** D9N50DL

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG121415.M
Quant Title : SVOA CALIBRATION

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

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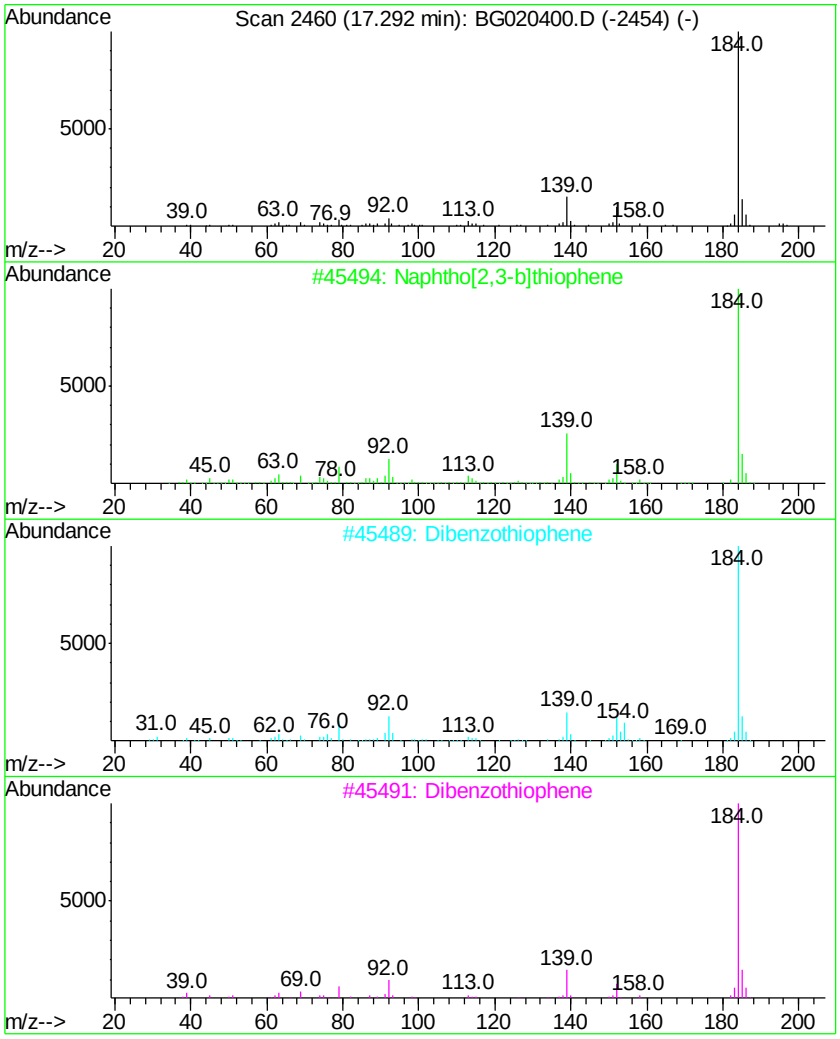
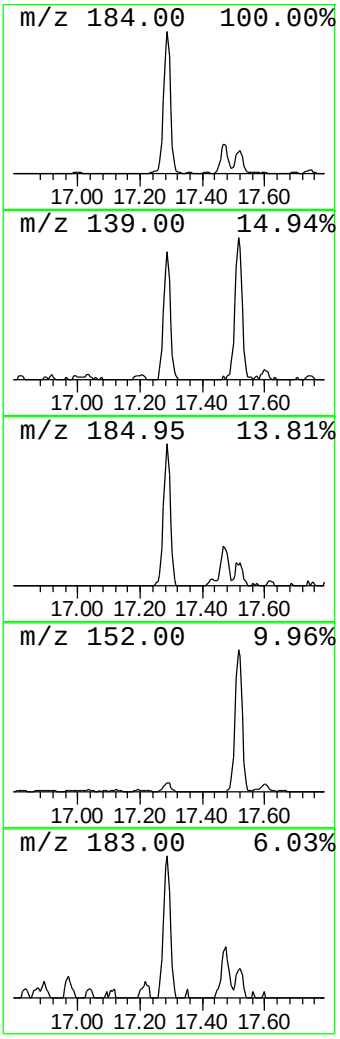
Peak Number 25 Naphtho[2,3-b]thiophene Concentration Rank 4

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 17.29 | 11.59 ng/ul | 175935 | Phenanthrene-d10 | 17.47 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Naphtho[2,3-b]thiophene | 184 | C12H8S | 000268-77-9 | 94 |
| 2 | Dibenzothiophene | 184 | C12H8S | 000132-65-0 | 94 |
| 3 | Dibenzothiophene | 184 | C12H8S | 000132-65-0 | 93 |
| 4 | Dibenzothiophene | 184 | C12H8S | 000132-65-0 | 93 |
| 5 | Dibenzothiophene | 184 | C12H8S | 000132-65-0 | 93 |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG122115\
Data File : BG020400.D
Acq On : 22 Dec 2015 6:10
Operator : UM/SJ
Sample : G4848-04DL 30X
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** D9N50DL

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG121415.M
Quant Title : SVOA CALIBRATION

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

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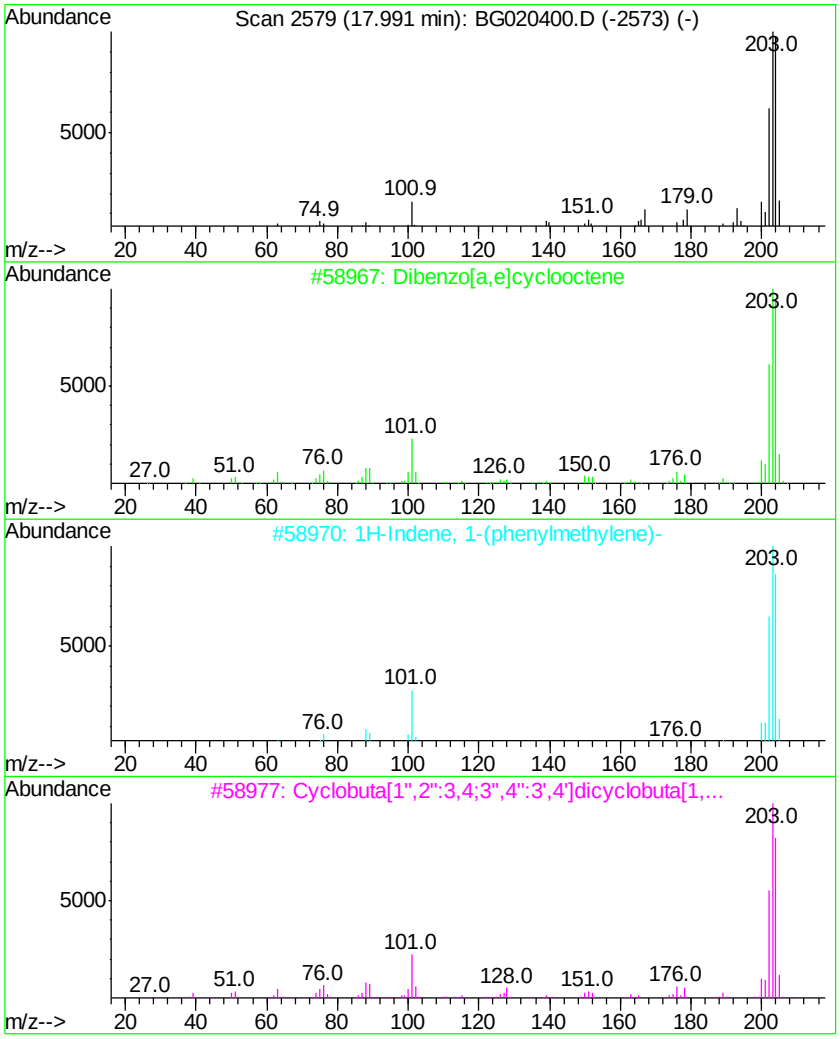
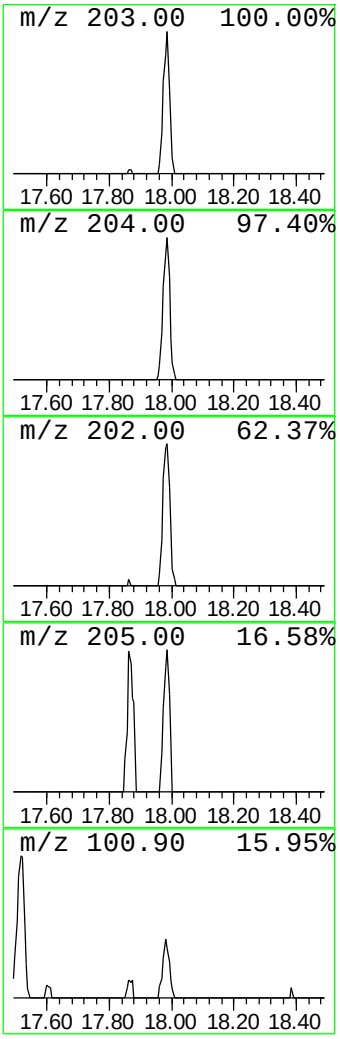
Peak Number 26 Dibenzo[a,e]cyclooctene Concentration Rank 29

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 17.99 | 2.61 ng/ul | 39608 | Phenanthrene-d10 | 17.47 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Dibenzo[a,e]cyclooctene | 204 | C16H12 | 000262-89-5 | 86 |
| 2 | 1H-Indene, 1-(phenylmethylene)- | 204 | C16H12 | 005394-86-5 | 70 |
| 3 | Cyclobuta[1'',2'':3,4;3'',4'':3'... | 204 | C16H12 | 006574-36-3 | 70 |
| 4 | 5H-Dibenzo[a,d]cycloheptene, 5-m... | 204 | C16H12 | 002975-79-3 | 64 |
| 5 | Anthracene, 9-ethenyl- | 204 | C16H12 | 002444-68-0 | 62 |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG122115\
Data File : BG020400.D
Acq On : 22 Dec 2015 6:10
Operator : UM/SJ
Sample : G4848-04DL 30X
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** D9N50DL

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG121415.M
Quant Title : SVOA CALIBRATION

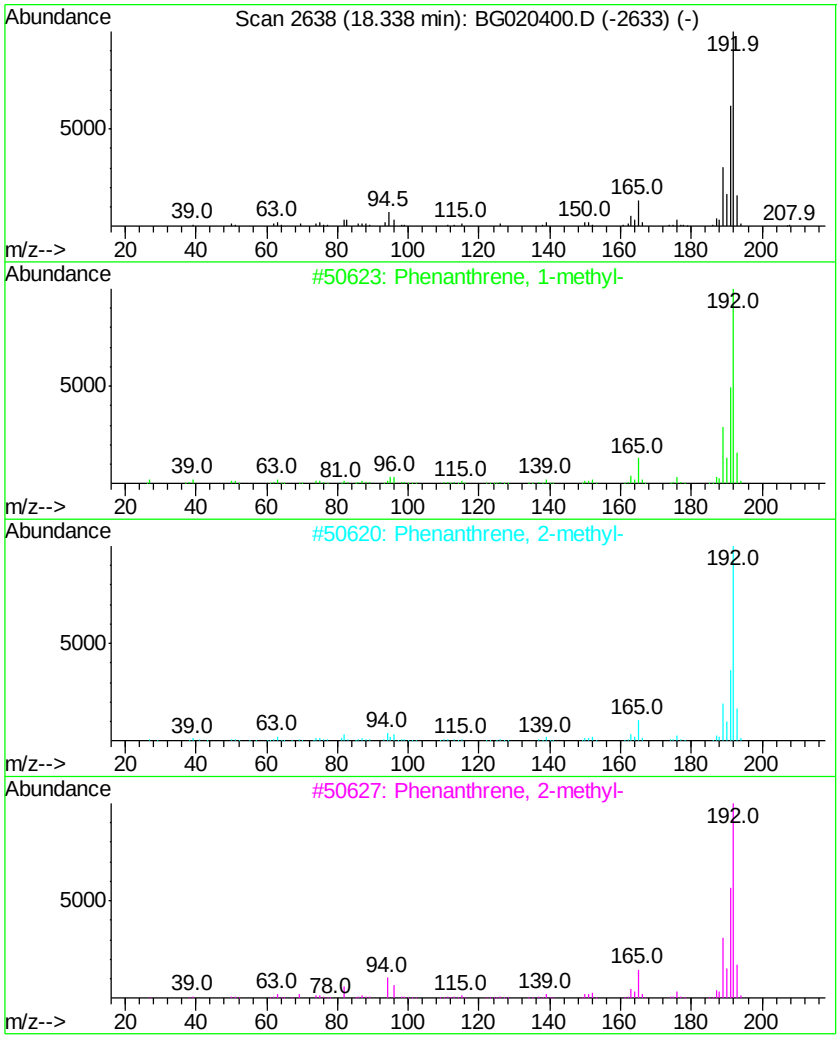
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

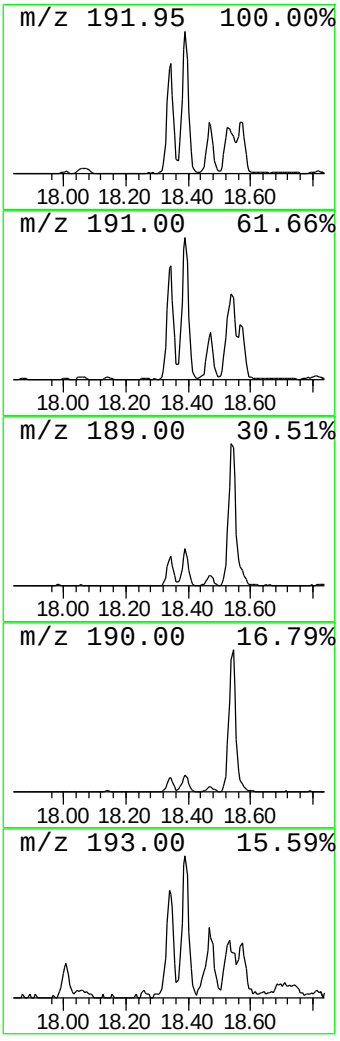
**********

Peak Number 27 Phenanthrene, 1-methyl- Concentration Rank 7

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.34 | 9.95 ng/ul | 151031 | Phenanthrene-d10 | 17.47 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenanthrene, 1-methyl- | 192 | C15H12 | 000832-69-9 | 96 |
| 2 | Phenanthrene, 2-methyl- | 192 | C15H12 | 002531-84-2 | 95 |
| 3 | Phenanthrene, 2-methyl- | 192 | C15H12 | 002531-84-2 | 95 |
| 4 | Anthracene, 1-methyl- | 192 | C15H12 | 000610-48-0 | 94 |
| 5 | Anthracene, 1-methyl- | 192 | C15H12 | 000610-48-0 | 94 |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG122115\
Data File : BG020400.D
Acq On : 22 Dec 2015 6:10
Operator : UM/SJ
Sample : G4848-04DL 30X
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** D9N50DL

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG121415.M
Quant Title : SVOA CALIBRATION

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

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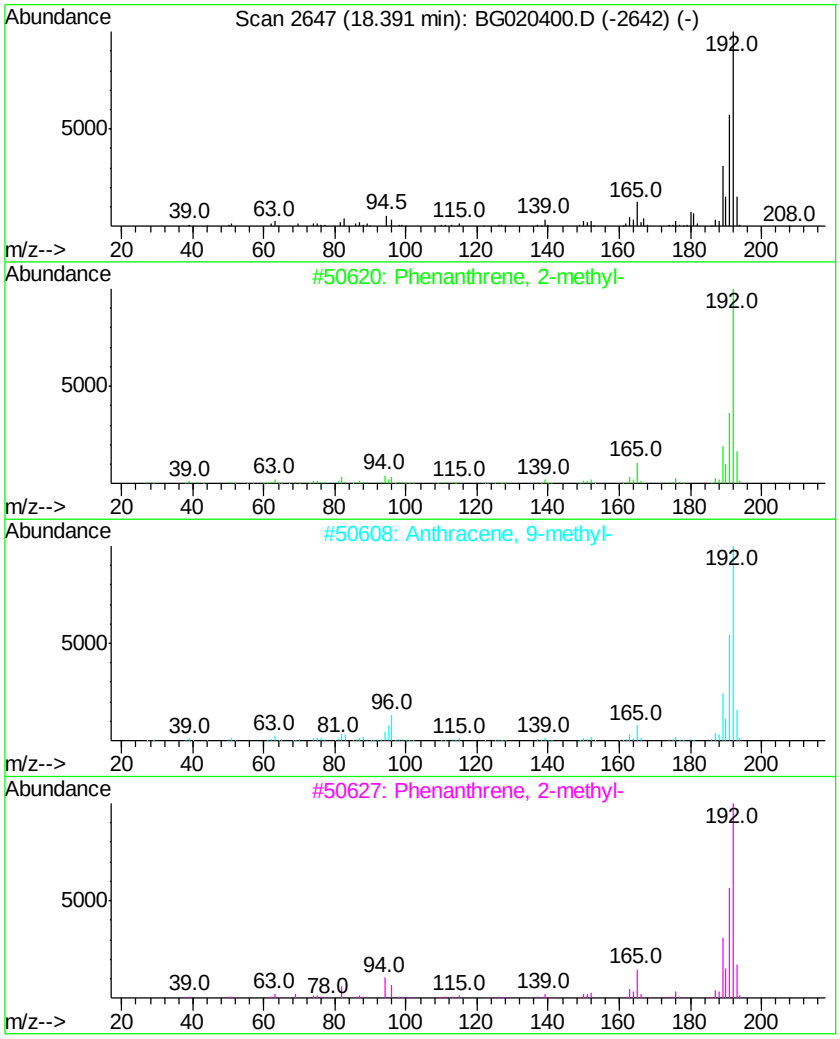
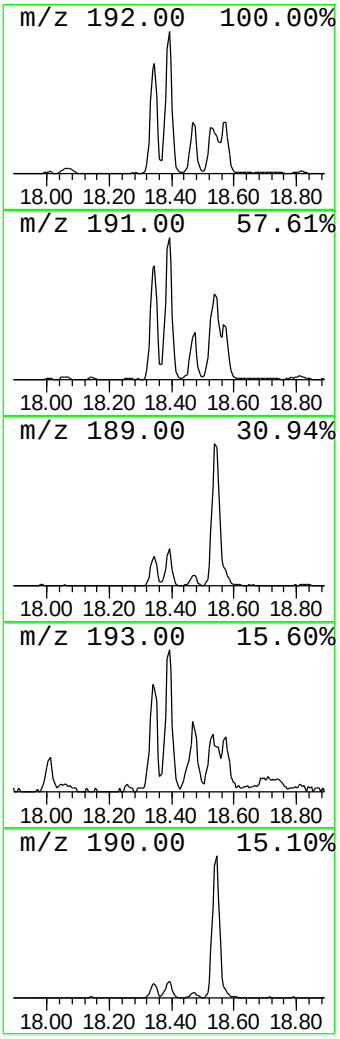
Peak Number 28 Phenanthrene, 2-methyl- Concentration Rank 3

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.39 | 14.64 ng/ul | 222216 | Phenanthrene-d10 | 17.47 |

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



Data Path : Z:\HPCHEM1\BNA G\DATA\BG122115\
Data File : BG020400.D
Acq On : 22 Dec 2015 6:10
Operator : UM/SJ
Sample : G4848-04DL 30X
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** D9N50DL

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG121415.M
Quant Title : SVOA CALIBRATION

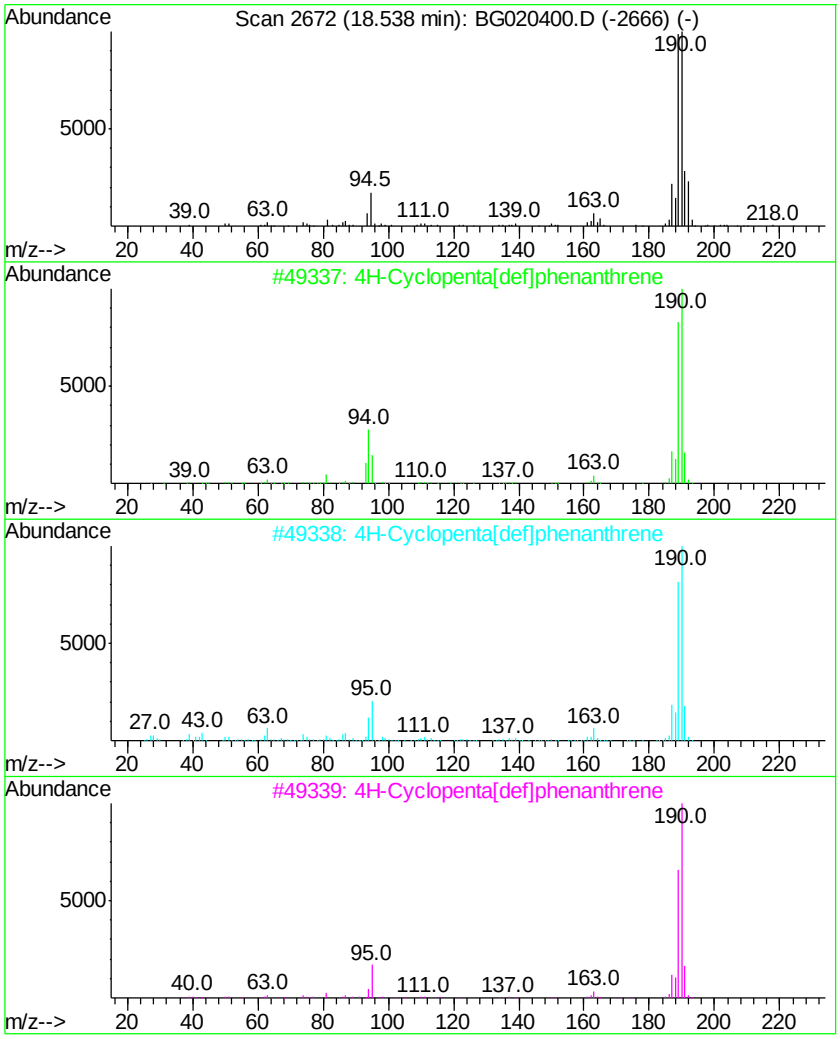
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

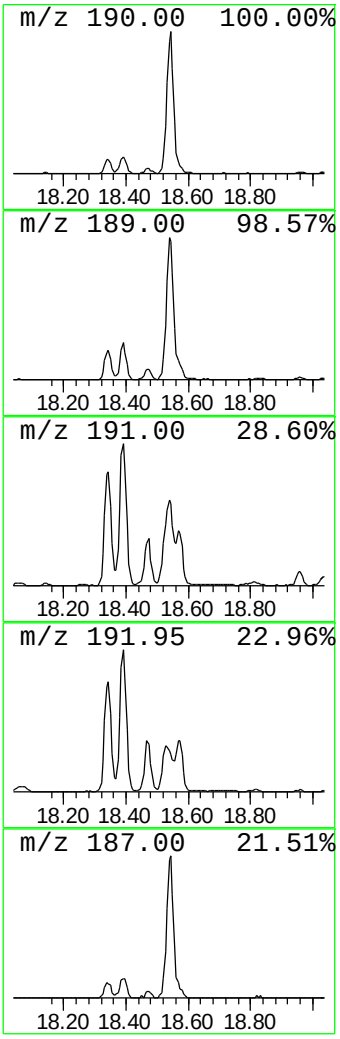
**********

Peak Number 30 4H-Cyclopenta[def]phenanthrene Concentration Rank 2

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.54 | 23.15 ng/ul | 351373 | Phenanthrene-d10 | 17.47 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 4H-Cyclopenta[def]phenanthrene | 190 | C15H10 | 000203-64-5 | 83 |
| 2 | 4H-Cyclopenta[def]phenanthrene | 190 | C15H10 | 000203-64-5 | 76 |
| 3 | 4H-Cyclopenta[def]phenanthrene | 190 | C15H10 | 000203-64-5 | 74 |
| 4 | 6H-Cyclobuta[jk]phenanthrene | 190 | C15H10 | 083469-43-6 | 64 |
| 5 | 2,3,5,6-Tetramethylterephthalald... | 190 | C12H14O2 | 007072-01-7 | 50 |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG122115\
Data File : BG020400.D
Acq On : 22 Dec 2015 6:10
Operator : UM/SJ
Sample : G4848-04DL 30X
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** D9N50DL

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG121415.M
Quant Title : SVOA CALIBRATION

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

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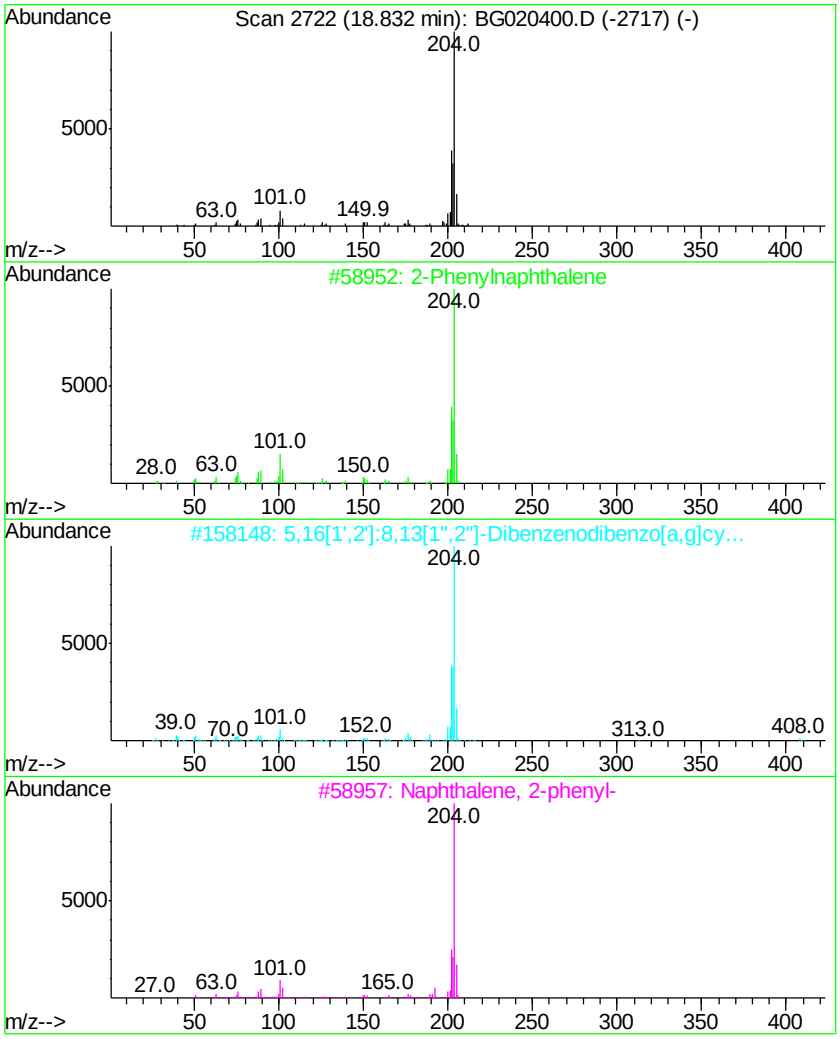
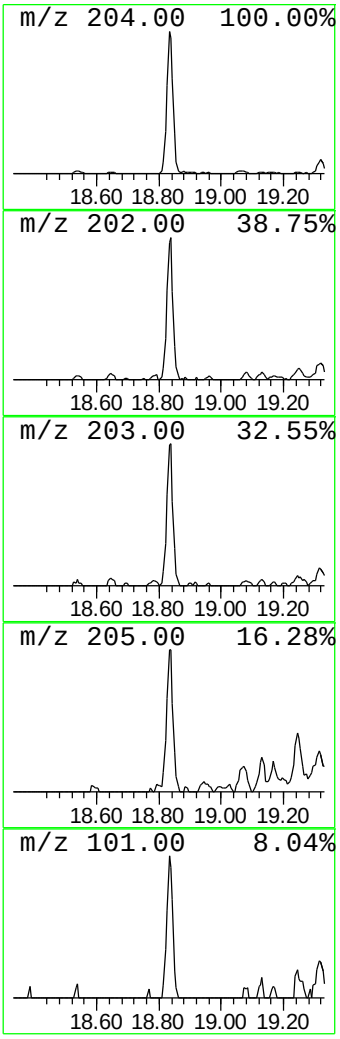
Peak Number 31 2-Phenylnaphthalene Concentration Rank 10

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.83 | 8.39 ng/ul | 127316 | Phenanthrene-d10 | 17.47 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-Phenylnaphthalene | 204 | C16H12 | 035465-71-5 | 94 |
| 2 | 5,16[1',2']:8,13[1'',2'']-Dibenz... | 408 | C32H24 | 005672-97-9 | 91 |
| 3 | Naphthalene, 2-phenyl- | 204 | C16H12 | 000612-94-2 | 64 |
| 4 | Naphthalene, 2-phenyl- | 204 | C16H12 | 000612-94-2 | 64 |
| 5 | Naphthalene, 1-phenyl- | 204 | C16H12 | 000605-02-7 | 60 |



Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG122115\
Data File : BG020400.D
Acq On : 22 Dec 2015 6:10
Operator : UM/SJ
Sample : G4848-04DL 30X
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** D9N50DL

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG121415.M
Quant Title : SVOA CALIBRATION

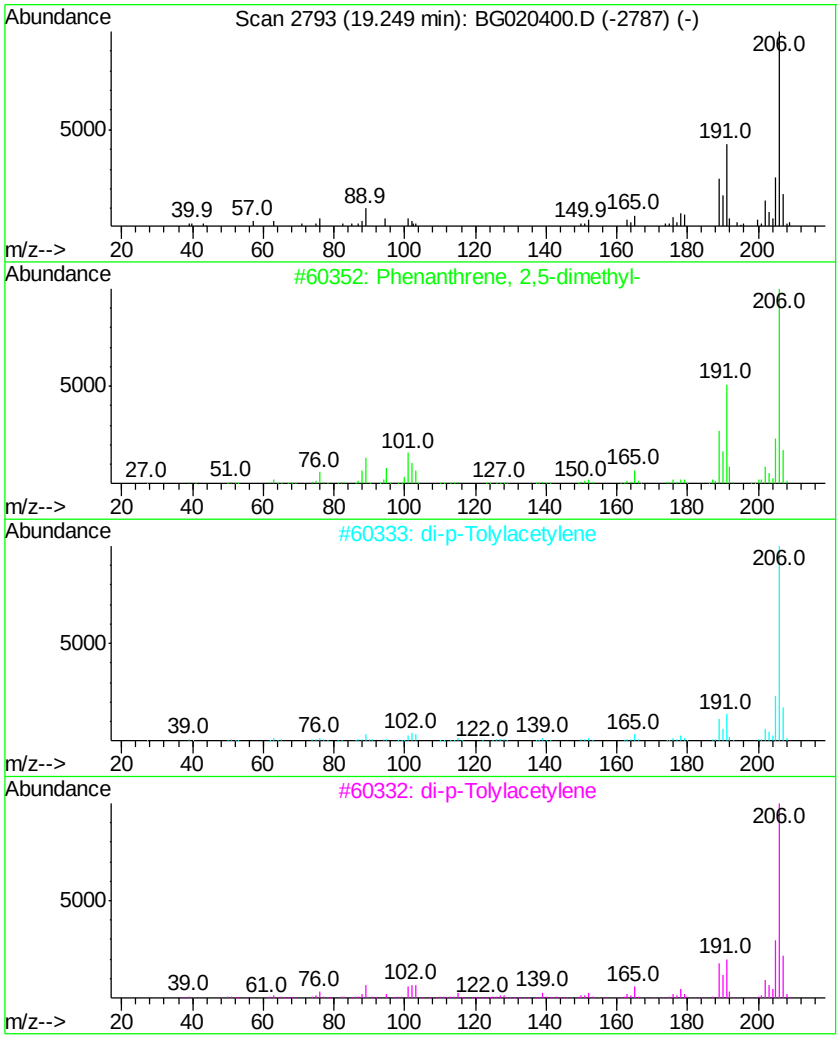
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

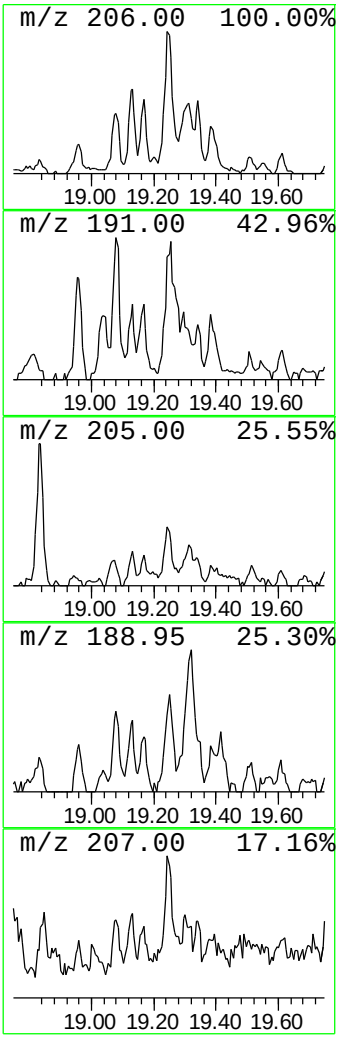
**********

Peak Number 32 Phenanthrene, 2,5-dimethyl- Concentration Rank 25

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 19.25 | 3.46 ng/ul | 52572 | Phenanthrene-d10 | 17.47 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenanthrene, 2,5-dimethyl- | 206 | C16H14 | 003674-66-6 | 93 |
| 2 | di-p-Tolylacetylene | 206 | C16H14 | 002789-88-0 | 90 |
| 3 | di-p-Tolylacetylene | 206 | C16H14 | 002789-88-0 | 90 |
| 4 | Naphthalene, 1,2-dihydro-4-phenyl- | 206 | C16H14 | 007469-40-1 | 87 |
| 5 | Phenanthrene, 2,7-dimethyl- | 206 | C16H14 | 001576-69-8 | 87 |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG122115\
Data File : BG020400.D
Acq On : 22 Dec 2015 6:10
Operator : UM/SJ
Sample : G4848-04DL 30X
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** D9N50DL

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG121415.M
Quant Title : SVOA CALIBRATION

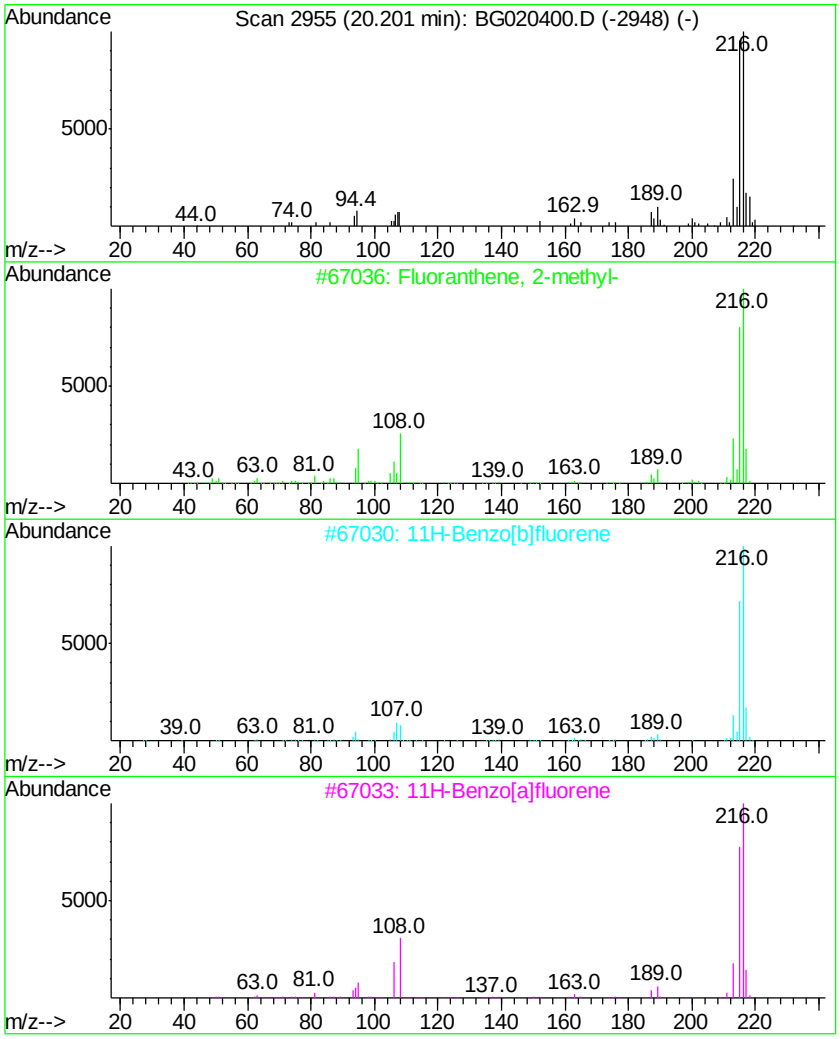
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

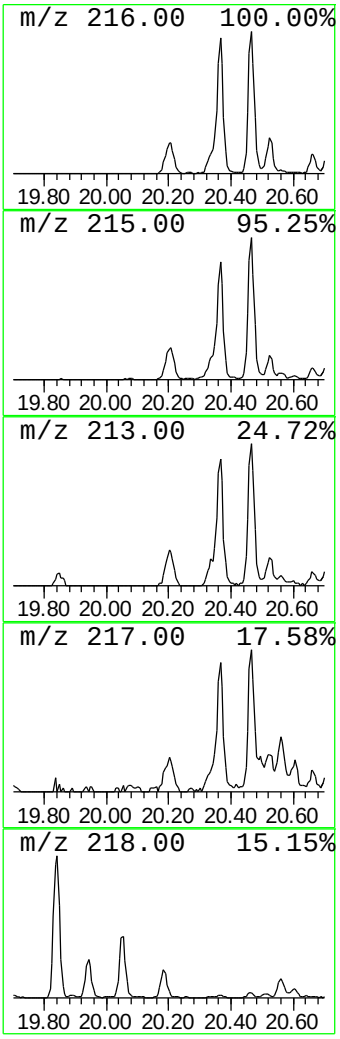
**********

Peak Number 33 Fluoranthene, 2-methyl- Concentration Rank 32

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 20.20 | 2.41 ng/ul | 85737 | Chrysene-d12 | 21.75 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Fluoranthene, 2-methyl- | 216 | C17H12 | 033543-31-6 | 91 |
| 2 | 11H-Benzo[b]fluorene | 216 | C17H12 | 000243-17-4 | 90 |
| 3 | 11H-Benzo[a]fluorene | 216 | C17H12 | 000238-84-6 | 87 |
| 4 | 11H-Benzo[a]fluorene | 216 | C17H12 | 000238-84-6 | 87 |
| 5 | 11H-Benzo[b]fluorene | 216 | C17H12 | 000243-17-4 | 83 |



Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG122115\
Data File : BG020400.D
Acq On : 22 Dec 2015 6:10
Operator : UM/SJ
Sample : G4848-04DL 30X
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** D9N50DL

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG121415.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 |
| Indene | 8.63 | 8.1 | ng/ul | 38714 | 1 | 8.09 | 96126 | 20.0 |
| 2-Benzothiophene # | 11.08 | 3.8 | ng/ul | 29212 | 2 | 10.91 | 154628 | 20.0 |
| Quinoline | 11.73 | 2.3 | ng/ul | 17545 | 2 | 10.91 | 154628 | 20.0 |
| Naphthalene, 1-et... | 13.75 | 5.1 | ng/ul | 56858 | 3 | 14.72 | 222035 | 20.0 |
| Naphthalene, 1,6-... | 13.88 | 5.4 | ng/ul | 60011 | 3 | 14.72 | 222035 | 20.0 |
| Naphthalene, 1,7-... | 14.04 | 10.2 | ng/ul | 112874 | 3 | 14.72 | 222035 | 20.0 |
| Naphthalene, 2,6-... | 14.10 | 6.6 | ng/ul | 72854 | 3 | 14.72 | 222035 | 20.0 |
| Naphthalene, 2,7-... | 14.28 | 3.2 | ng/ul | 35925 | 3 | 14.72 | 222035 | 20.0 |
| 2-Naphthalenecarb... | 14.85 | 3.3 | ng/ul | 37029 | 3 | 14.72 | 222035 | 20.0 |
| Naphthalene, 2-(1... | 14.92 | 2.0 | ng/ul | 22338 | 3 | 14.72 | 222035 | 20.0 |
| Benzene, [1-(2,4-... | 15.03 | 3.5 | ng/ul | 39462 | 3 | 14.72 | 222035 | 20.0 |
| 1-Isopropenylnaph... | 15.86 | 2.9 | ng/ul | 31649 | 3 | 14.72 | 222035 | 20.0 |
| Nordiphenamid | 15.93 | 9.7 | ng/ul | 107532 | 3 | 14.72 | 222035 | 20.0 |
| Naphthalene, 1-(2... | 15.98 | 2.1 | ng/ul | 23190 | 3 | 14.72 | 222035 | 20.0 |
| 1,1'-Biphenyl, 2-... | 16.02 | 4.7 | ng/ul | 51639 | 3 | 14.72 | 222035 | 20.0 |
| 9H-Fluoren-9-ol | 16.06 | 9.9 | ng/ul | 110275 | 3 | 14.72 | 222035 | 20.0 |
| 6H-Dibenzo[b,d]-p... | 16.21 | 10.9 | ng/ul | 166043 | 4 | 17.47 | 303589 | 20.0 |
| Anthracene, 9,10-... | 16.56 | 4.1 | ng/ul | 62337 | 4 | 17.47 | 303589 | 20.0 |
| 9H-Fluorene, 3-me... | 16.77 | 8.1 | ng/ul | 123601 | 4 | 17.47 | 303589 | 20.0 |
| 9H-Fluorene, 2-me... | 16.92 | 2.4 | ng/ul | 36712 | 4 | 17.47 | 303589 | 20.0 |
| Benzene, 1,2-dime... | 16.99 | 2.5 | ng/ul | 37947 | 4 | 17.47 | 303589 | 20.0 |
| 2,4,6-Trimethoxyb... | 17.19 | 4.2 | ng/ul | 64046 | 4 | 17.47 | 303589 | 20.0 |
| Naphtho[2,3-b]thi... | 17.29 | 11.6 | ng/ul | 175935 | 4 | 17.47 | 303589 | 20.0 |
| Dibenzo[a,e]cyclo... | 17.99 | 2.6 | ng/ul | 39608 | 4 | 17.47 | 303589 | 20.0 |
| Phenanthrene, 1-m... | 18.34 | 9.9 | ng/ul | 151031 | 4 | 17.47 | 303589 | 20.0 |
| Phenanthrene, 2-m... | 18.39 | 14.6 | ng/ul | 222216 | 4 | 17.47 | 303589 | 20.0 |
| 4H-Cyclopenta[def... | 18.54 | 23.1 | ng/ul | 351373 | 4 | 17.47 | 303589 | 20.0 |
| 2-Phenylnaphthalene | 18.83 | 8.4 | ng/ul | 127316 | 4 | 17.47 | 303589 | 20.0 |
| Phenanthrene, 2,5... | 19.25 | 3.5 | ng/ul | 52572 | 4 | 17.47 | 303589 | 20.0 |
| Fluoranthene, 2-m... | 20.20 | 2.4 | ng/ul | 85737 | 5 | 21.75 | 711969 | 20.0 |