LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080215\
Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
Operator : TP
Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** C01M4

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-BM072715.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 | 7.557 | 820 | 828 | 835 | rBV | 421682 | 658822 | 3.41% | 0.322% |
| 2 | 8.281 | 945 | 951 | 958 | rBV | 184341 | 293995 | 1.52% | 0.144% |
| 3 | 9.980 | 1234 | 1240 | 1248 | rBV | 198770 | 326364 | 1.69% | 0.160% |
| 4 | 10.345 | 1293 | 1302 | 1307 | rBV | 577093 | 968670 | 5.02% | 0.474% |
| 5 | 13.645 | 1858 | 1863 | 1867 | rVV | 329576 | 473841 | 2.45% | 0.232% |
| 6 | 13.921 | 1904 | 1910 | 1912 | rBV | 393576 | 597451 | 3.09% | 0.292% |
| 7 | 13.951 | 1912 | 1915 | 1928 | rVB | 484558 | 690145 | 3.57% | 0.337% |
| 8 | 14.233 | 1951 | 1963 | 1968 | rBV2 | 847731 | 1332977 | 6.90% | 0.652% |
| 9 | 14.292 | 1968 | 1973 | 1981 | rVB | 421449 | 623718 | 3.23% | 0.305% |
| 10 | 14.627 | 2020 | 2030 | 2039 | rVV | 220813 | 359203 | 1.86% | 0.176% |
| 11 | 15.227 | 2126 | 2132 | 2137 | rVV | 506886 | 761791 | 3.95% | 0.372% |
| 12 | 15.280 | 2137 | 2141 | 2152 | rVV | 451922 | 709066 | 3.67% | 0.347% |
| 13 | 15.574 | 2187 | 2191 | 2202 | rVB | 127746 | 262891 | 1.36% | 0.129% |
| 14 | 15.727 | 2213 | 2217 | 2224 | rVB2 | 129339 | 247026 | 1.28% | 0.121% |
| 15 | 16.286 | 2301 | 2312 | 2317 | rVV3 | 175023 | 460605 | 2.39% | 0.225% |
| 16 | 16.339 | 2317 | 2321 | 2326 | rVV2 | 149197 | 256106 | 1.33% | 0.125% |
| 17 | 16.556 | 2354 | 2358 | 2366 | rVV3 | 147520 | 295614 | 1.53% | 0.145% |
| 18 | 16.645 | 2368 | 2373 | 2379 | rVV | 272423 | 419443 | 2.17% | 0.205% |
| 19 | 16.733 | 2380 | 2388 | 2395 | rVV4 | 146143 | 391564 | 2.03% | 0.191% |
| 20 | 16.804 | 2395 | 2400 | 2409 | rVB | 389512 | 607919 | 3.15% | 0.297% |
| 21 | 16.986 | 2425 | 2431 | 2434 | rVV | 932127 | 1454460 | 7.53% | 0.711% |
| 22 | 17.039 | 2434 | 2440 | 2444 | rVV | 6047280 | 9019576 | 46.71% | 4.409% |
| 23 | 17.086 | 2444 | 2448 | 2451 | rVV | 593686 | 924946 | 4.79% | 0.452% |
| 24 | 17.121 | 2451 | 2454 | 2459 | rVV | 1600174 | 2006917 | 10.39% | 0.981% |
| 25 | 17.180 | 2459 | 2464 | 2468 | rVB3 | 134655 | 242829 | 1.26% | 0.119% |
| 26 | 17.398 | 2492 | 2501 | 2512 | rVV2 | 435308 | 872809 | 4.52% | 0.427% |
| 27 | 17.503 | 2515 | 2519 | 2523 | rVV | 225098 | 289270 | 1.50% | 0.141% |
| 28 | 17.562 | 2525 | 2529 | 2538 | rVB3 | 340471 | 589162 | 3.05% | 0.288% |
| 29 | 17.703 | 2549 | 2553 | 2561 | rVB | 194101 | 377118 | 1.95% | 0.184% |
| 30 | 17.862 | 2570 | 2580 | 2584 | rBV | 1184521 | 1838603 | 9.52% | 0.899% |
| 31 | 17.909 | 2584 | 2588 | 2593 | rVB | 1462217 | 2096425 | 10.86% | 1.025% |
| 32 | 17.992 | 2593 | 2602 | 2607 | rBV | 532446 | 823821 | 4.27% | 0.403% |
| 33 | 18.056 | 2607 | 2613 | 2617 | rBV2 | 3520285 | 5478973 | 28.38% | 2.678% |
| 34 | 18.098 | 2617 | 2620 | 2627 | rVB2 | 895730 | 1523596 | 7.89% | 0.745% |

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080215\
Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
Operator : TP
Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** C01M4

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-BM072715.M
Title : SVOA CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 35 | 18.168 | 2627 | 2632 | 2636 | rBV5 | 138500 | 249032 | 1.29% | 0.122% |
| 36 | 18.345 | 2653 | 2662 | 2663 | rBV3 | 750992 | 1124551 | 5.82% | 0.550% |
| 37 | 18.368 | 2663 | 2666 | 2670 | rVB | 801285 | 1030175 | 5.34% | 0.504% |
| 38 | 18.421 | 2671 | 2675 | 2682 | rVB | 534114 | 771413 | 4.00% | 0.377% |
| 39 | 18.521 | 2689 | 2692 | 2699 | rVB3 | 212483 | 309102 | 1.60% | 0.151% |
| 40 | 18.615 | 2704 | 2708 | 2713 | rVB | 336085 | 461515 | 2.39% | 0.226% |
| 41 | 18.668 | 2713 | 2717 | 2720 | rVB | 264335 | 307005 | 1.59% | 0.150% |
| 42 | 18.703 | 2720 | 2723 | 2728 | rVB2 | 263262 | 353942 | 1.83% | 0.173% |
| 43 | 18.792 | 2729 | 2738 | 2742 | rBV | 886665 | 1561758 | 8.09% | 0.763% |
| 44 | 18.845 | 2742 | 2747 | 2752 | rBV5 | 446015 | 869478 | 4.50% | 0.425% |
| 45 | 18.915 | 2752 | 2759 | 2770 | rVB6 | 2089815 | 4797074 | 24.84% | 2.345% |
| 46 | 18.997 | 2770 | 2773 | 2778 | rVB2 | 430008 | 516351 | 2.67% | 0.252% |
| 47 | 19.074 | 2778 | 2786 | 2790 | rBV | 9351513 | 14997605 | 77.67% | 7.331% |
| 48 | 19.127 | 2790 | 2795 | 2798 | rBV | 638360 | 935486 | 4.84% | 0.457% |
| 49 | 19.203 | 2803 | 2808 | 2814 | rVB2 | 3046388 | 4563944 | 23.64% | 2.231% |
| 50 | 19.268 | 2814 | 2819 | 2822 | rBV2 | 1227353 | 1560966 | 8.08% | 0.763% |
| 51 | 19.303 | 2822 | 2825 | 2829 | rVB | 407844 | 532666 | 2.76% | 0.260% |
| 52 | 19.345 | 2829 | 2832 | 2836 | rVB2 | 225379 | 306011 | 1.58% | 0.150% |
| 53 | 19.445 | 2836 | 2849 | 2855 | rBV | 9337009 | 19308693 | 100.00% | 9.439% |
| 54 | 19.497 | 2855 | 2858 | 2862 | rBV3 | 444261 | 612885 | 3.17% | 0.300% |
| 55 | 19.539 | 2862 | 2865 | 2870 | rVB | 1264700 | 1517392 | 7.86% | 0.742% |
| 56 | 19.592 | 2870 | 2874 | 2880 | rBV2 | 611031 | 1362375 | 7.06% | 0.666% |
| 57 | 19.650 | 2880 | 2884 | 2893 | rVB2 | 3103253 | 4289773 | 22.22% | 2.097% |
| 58 | 19.768 | 2896 | 2904 | 2910 | rBV2 | 903434 | 2535841 | 13.13% | 1.240% |
| 59 | 19.915 | 2919 | 2929 | 2934 | rBV2 | 1946323 | 4581153 | 23.73% | 2.239% |
| 60 | 19.968 | 2934 | 2938 | 2942 | rVB | 2518407 | 2746737 | 14.23% | 1.343% |
| 61 | 20.027 | 2944 | 2948 | 2952 | rBV | 1251627 | 1538080 | 7.97% | 0.752% |
| 62 | 20.086 | 2952 | 2958 | 2961 | rVV2 | 998503 | 1647368 | 8.53% | 0.805% |
| 63 | 20.115 | 2961 | 2963 | 2968 | rVB | 577267 | 752907 | 3.90% | 0.368% |
| 64 | 20.192 | 2968 | 2976 | 2979 | rBV3 | 1358784 | 1861962 | 9.64% | 0.910% |
| 65 | 20.239 | 2979 | 2984 | 2987 | rVV2 | 783227 | 1707480 | 8.84% | 0.835% |
| 66 | 20.274 | 2987 | 2990 | 2993 | rVB2 | 1752557 | 1983512 | 10.27% | 0.970% |
| 67 | 20.344 | 2998 | 3002 | 3006 | rBV5 | 299523 | 407274 | 2.11% | 0.199% |
| 68 | 20.439 | 3015 | 3018 | 3021 | rVB3 | 228173 | 265300 | 1.37% | 0.130% |
| 69 | 20.509 | 3021 | 3030 | 3035 | rBV2 | 1931011 | 3528071 | 18.27% | 1.725% |
| 70 | 20.680 | 3056 | 3059 | 3064 | rVB | 838136 | 1031154 | 5.34% | 0.504% |
| 71 | 20.815 | 3078 | 3082 | 3084 | rBV | 759274 | 1007747 | 5.22% | 0.493% |

LSC Area Percent Report

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Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
Operator : TP
Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** C01M4

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-BM072715.M
Title : SVOA CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 72 | 20.839 | 3084 | 3086 | 3090 | rVV2 | 1245839 | 1696953 | 8.79% | 0.830% |
| 73 | 20.880 | 3090 | 3093 | 3096 | rVV | 1055588 | 1403940 | 7.27% | 0.686% |
| 74 | 20.921 | 3096 | 3100 | 3104 | rVV2 | 990043 | 1644191 | 8.52% | 0.804% |
| 75 | 20.980 | 3107 | 3110 | 3115 | rVB | 861763 | 1113527 | 5.77% | 0.544% |
| 76 | 21.109 | 3121 | 3132 | 3136 | rBV | 9146744 | 10893150 | 56.42% | 5.325% |
| 77 | 21.191 | 3139 | 3146 | 3151 | rVV3 | 5524426 | 11140042 | 57.69% | 5.446% |
| 78 | 21.244 | 3151 | 3155 | 3163 | rVB | 5670163 | 7450266 | 38.59% | 3.642% |
| 79 | 21.315 | 3164 | 3167 | 3169 | rBV2 | 357940 | 406863 | 2.11% | 0.199% |
| 80 | 21.350 | 3169 | 3173 | 3179 | rVV2 | 935423 | 1472080 | 7.62% | 0.720% |
| 81 | 21.409 | 3180 | 3183 | 3185 | rVV3 | 409449 | 574212 | 2.97% | 0.281% |
| 82 | 21.439 | 3185 | 3188 | 3195 | rVB2 | 1427464 | 2082080 | 10.78% | 1.018% |
| 83 | 21.515 | 3196 | 3201 | 3205 | rBV3 | 581704 | 972678 | 5.04% | 0.475% |
| 84 | 21.556 | 3205 | 3208 | 3211 | rBV4 | 202229 | 264920 | 1.37% | 0.130% |
| 85 | 21.738 | 3233 | 3239 | 3243 | rBV | 981310 | 1535478 | 7.95% | 0.751% |
| 86 | 21.791 | 3245 | 3248 | 3252 | rVB | 538495 | 648001 | 3.36% | 0.317% |
| 87 | 21.850 | 3256 | 3258 | 3262 | rVB2 | 264886 | 281658 | 1.46% | 0.138% |
| 88 | 21.891 | 3262 | 3265 | 3268 | rVB3 | 315339 | 384449 | 1.99% | 0.188% |
| 89 | 21.933 | 3268 | 3272 | 3275 | rBV2 | 644489 | 920996 | 4.77% | 0.450% |
| 90 | 22.027 | 3284 | 3288 | 3297 | rVB2 | 500897 | 863623 | 4.47% | 0.422% |
| 91 | 22.221 | 3317 | 3321 | 3326 | rBV4 | 218331 | 466339 | 2.42% | 0.228% |
| 92 | 22.491 | 3363 | 3367 | 3373 | rBV3 | 326959 | 590779 | 3.06% | 0.289% |
| 93 | 22.756 | 3404 | 3412 | 3416 | rBV | 3559861 | 8490169 | 43.97% | 4.150% |
| 94 | 22.909 | 3433 | 3438 | 3443 | rVB | 745600 | 1140565 | 5.91% | 0.558% |
| 95 | 23.221 | 3484 | 3491 | 3496 | rBV | 2067338 | 4046403 | 20.96% | 1.978% |
| 96 | 23.321 | 3501 | 3508 | 3514 | rVB | 3347383 | 6190117 | 32.06% | 3.026% |
| 97 | 23.403 | 3517 | 3522 | 3526 | rVV | 622496 | 1184655 | 6.14% | 0.579% |
| 98 | 23.450 | 3526 | 3530 | 3537 | rVB2 | 1065599 | 1934966 | 10.02% | 0.946% |
| 99 | 25.638 | 3891 | 3902 | 3909 | rVB2 | 1541983 | 4712403 | 24.41% | 2.304% |
| 100 | 26.326 | 4008 | 4019 | 4031 | rVB | 1196971 | 3850763 | 19.94% | 1.882% |

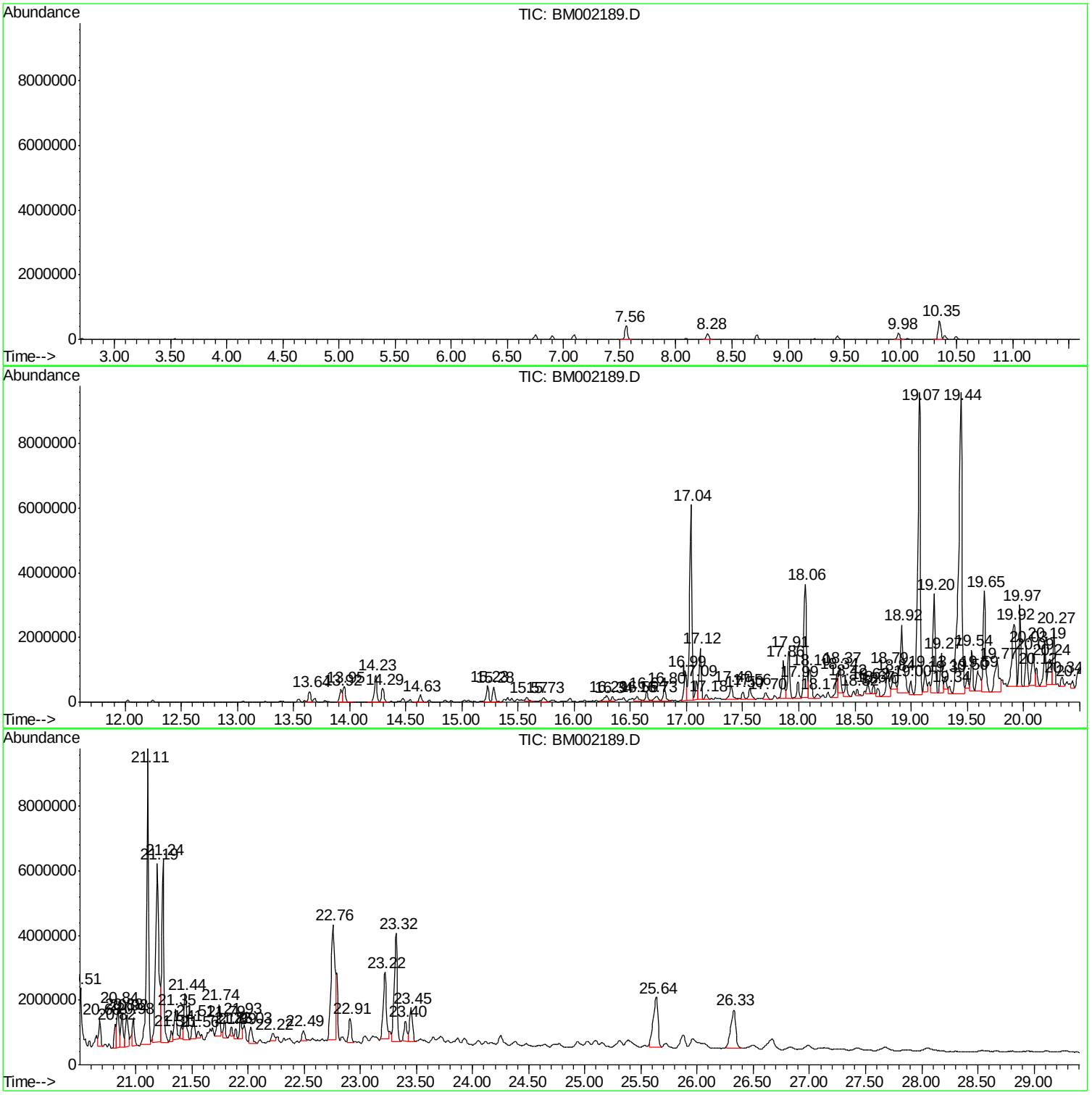
Sum of corrected areas: 204565760

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080215\
Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
Operator : TP
Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :**
BNA_M
**ClientSampleId :**
C01M4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080215\
Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
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Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** C01M4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
Quant Title : SVOA CALIBRATION

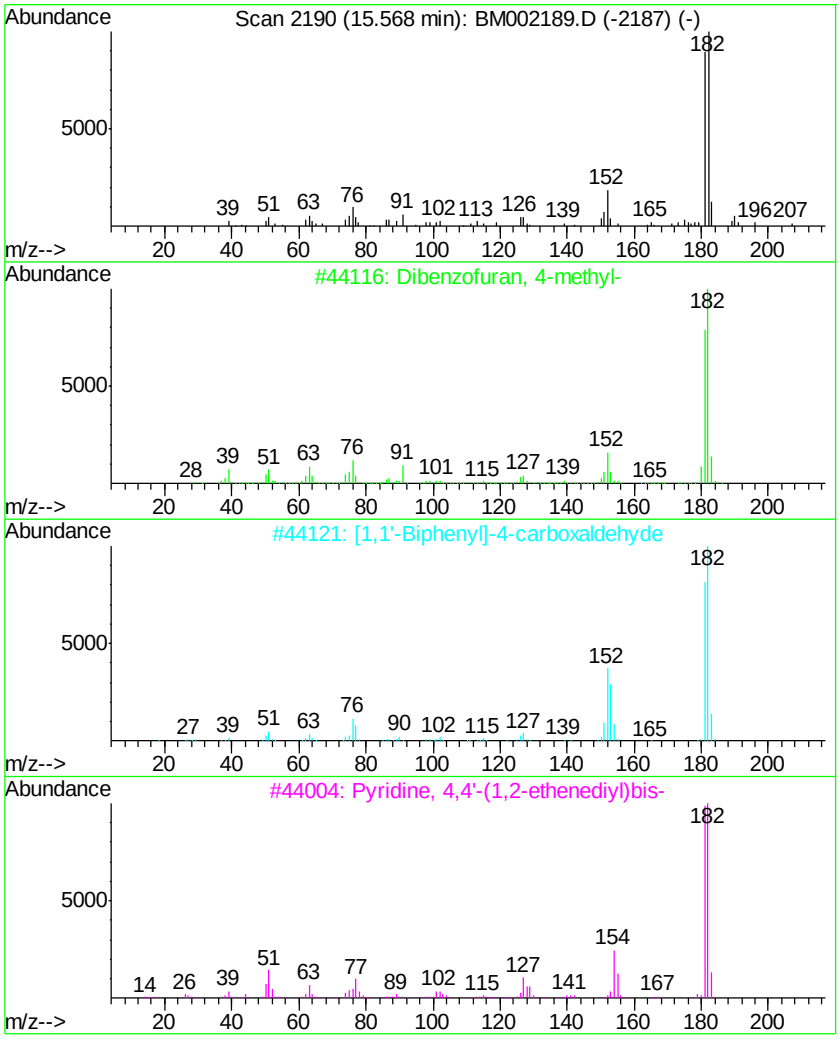
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

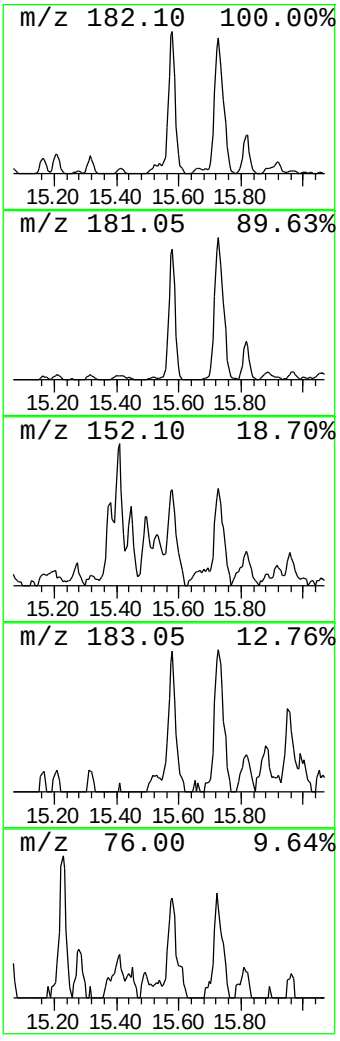
**********

Peak Number 1 Dibenzofuran, 4-methyl- Concentration Rank 33

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.57 | 3.94 ng/ul | 262891 | Acenaphthene-d10 | 14.23 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Dibenzofuran, 4-methyl- | 182 | C13H10O | 007320-53-8 | 90 |
| 2 | [1,1'-Biphenyl]-4-carboxaldehyde | 182 | C13H10O | 003218-36-8 | 83 |
| 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 | Pyridine, 4,4'-(1,2-ethenediyl)bis- | 182 | C12H10N2 | 001135-32-6 | 72 |



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080215\
Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
Operator : TP
Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** C01M4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
Quant Title : SVOA CALIBRATION

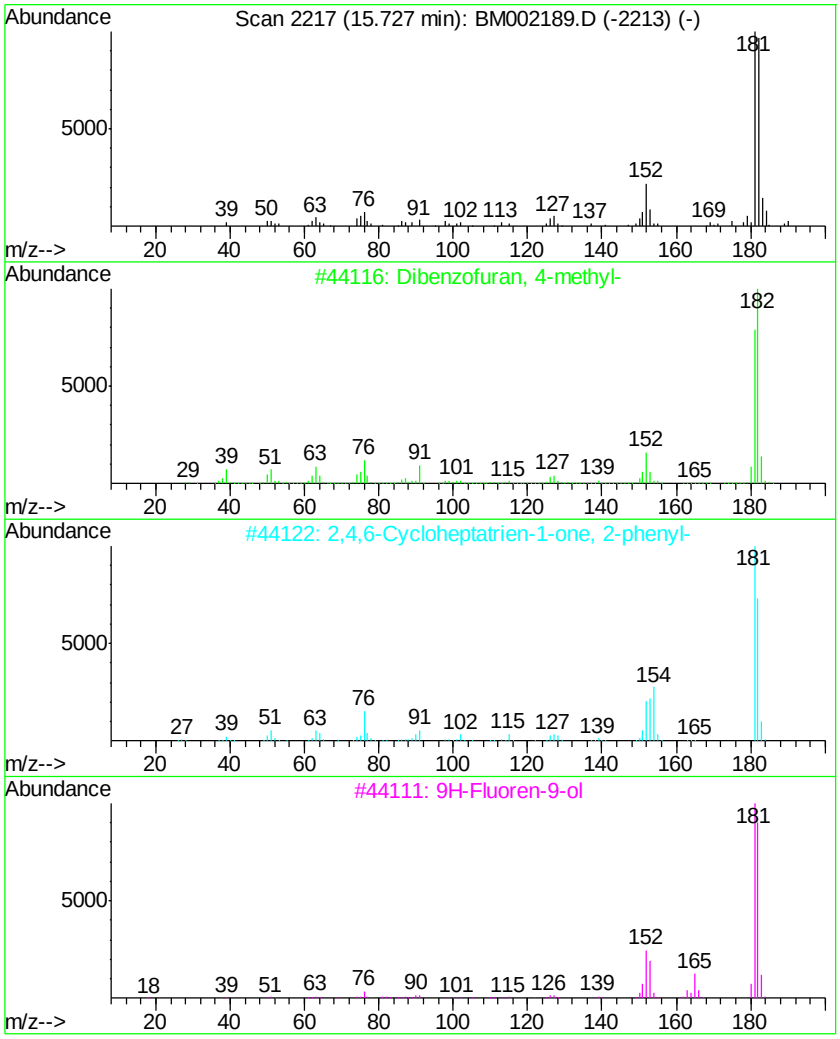
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

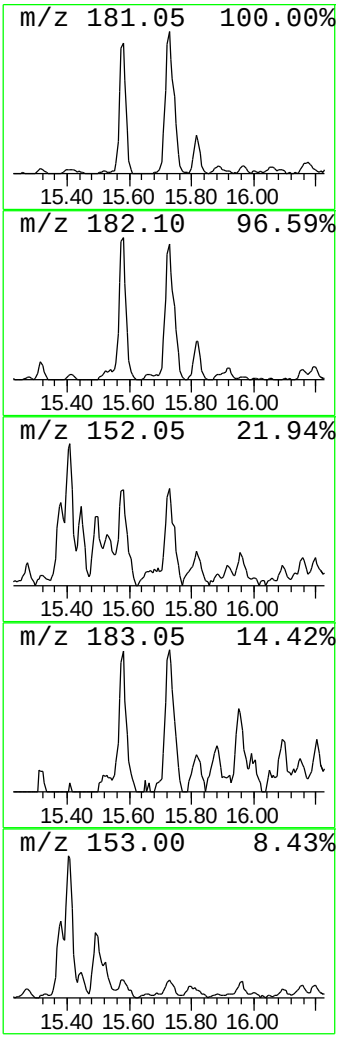
**********

Peak Number 2 2,4,6-Cycloheptatrien-1-one... Concentration Rank 38

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.73 | 3.40 ng/ul | 247026 | Phenanthrene-d10 | 16.99 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Dibenzofuran, 4-methyl- | 182 | C13H10O | 007320-53-8 | 91 |
| 2 | 2,4,6-Cycloheptatrien-1-one, 2-p... | 182 | C13H10O | 014562-09-5 | 86 |
| 3 | 9H-Fluoren-9-ol | 182 | C13H10O | 001689-64-1 | 78 |
| 4 | [1,1'-Biphenyl]-4-carboxaldehyde | 182 | C13H10O | 003218-36-8 | 72 |
| 5 | Pyrido[2,3-d]indole, 6-methyl- | 182 | C12H10N2 | 108349-67-3 | 72 |



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080215\
Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
Operator : TP
Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** C01M4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
Quant Title : SVOA CALIBRATION

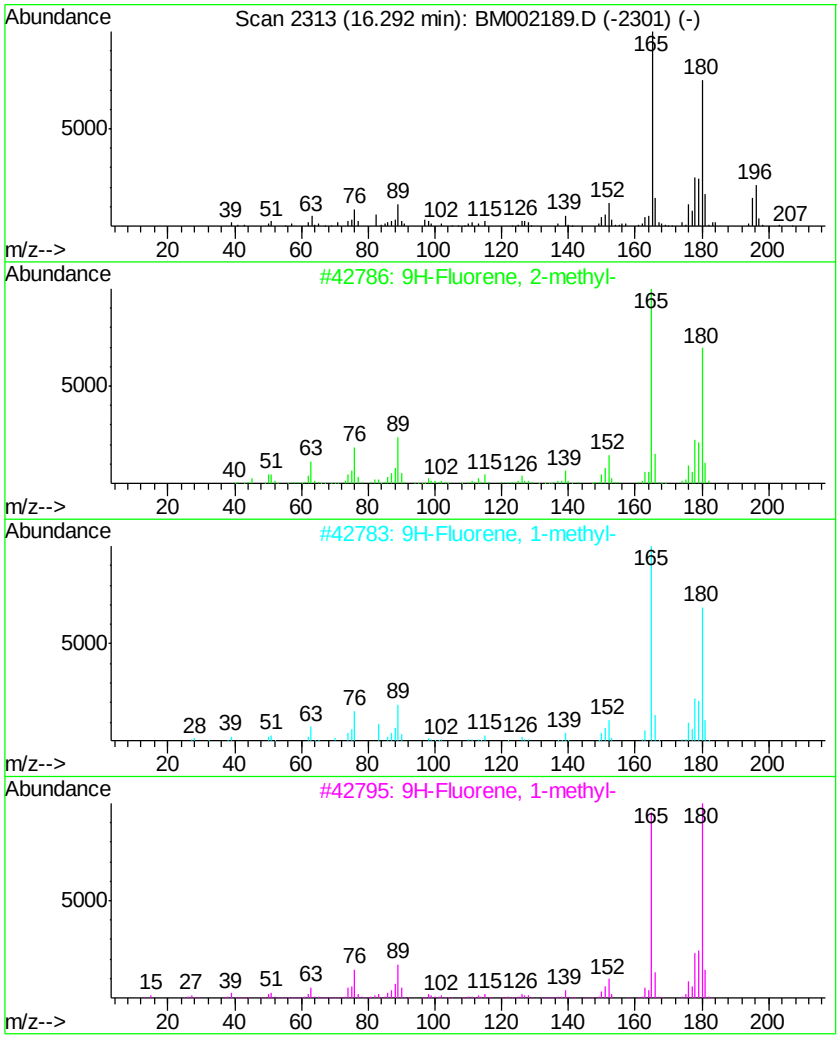
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

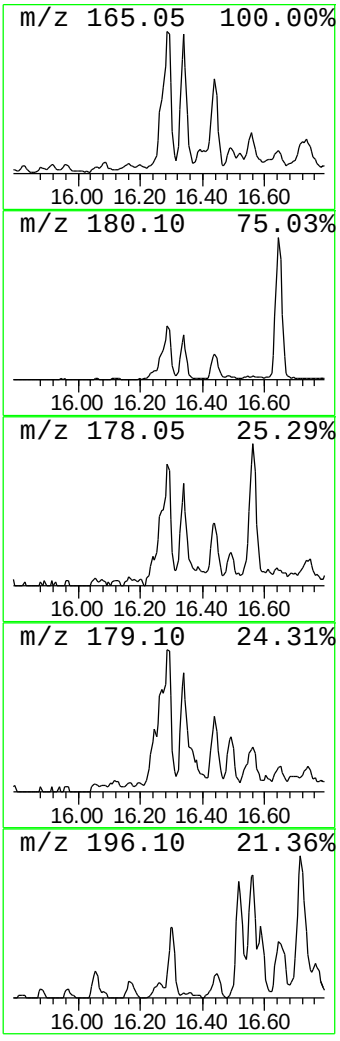
**********

Peak Number 3 9H-Fluorene, 2-methyl- Concentration Rank 22

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.29 | 6.33 ng/ul | 460605 | Phenanthrene-d10 | 16.99 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 9H-Fluorene, 2-methyl- | 180 | C14H12 | 001430-97-3 | 97 |
| 2 | 9H-Fluorene, 1-methyl- | 180 | C14H12 | 001730-37-6 | 96 |
| 3 | 9H-Fluorene, 1-methyl- | 180 | C14H12 | 001730-37-6 | 93 |
| 4 | 9H-Fluorene, 1-methyl- | 180 | C14H12 | 001730-37-6 | 93 |
| 5 | 9H-Fluorene, 2-methyl- | 180 | C14H12 | 001430-97-3 | 93 |



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080215\
Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
Operator : TP
Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** C01M4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
Quant Title : SVOA CALIBRATION

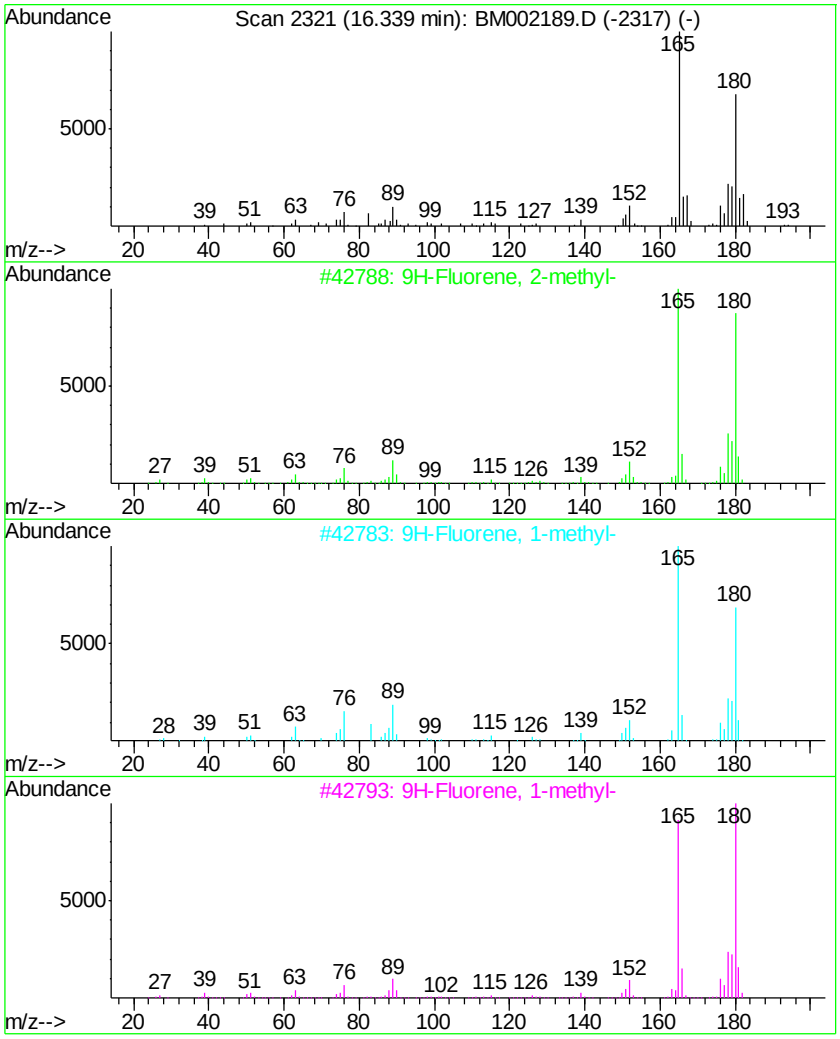
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

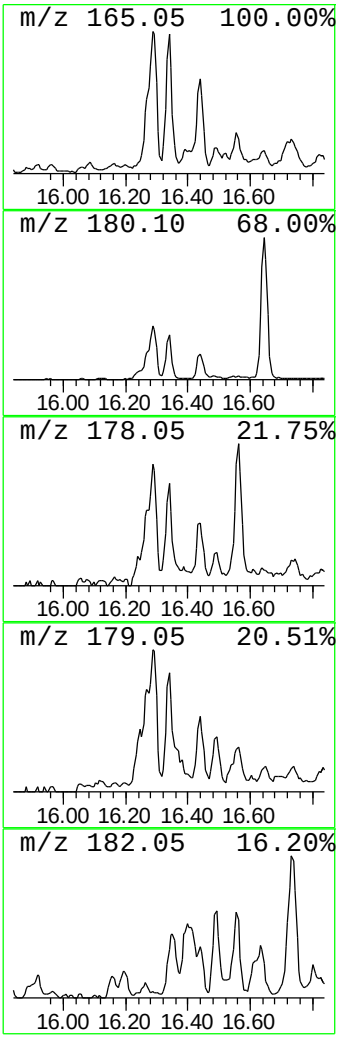
**********

Peak Number 4 9H-Fluorene, 1-methyl- Concentration Rank 36

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.34 | 3.52 ng/ul | 256106 | Phenanthrene-d10 | 16.99 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 9H-Fluorene, 2-methyl- | 180 | C14H12 | 001430-97-3 | 97 |
| 2 | 9H-Fluorene, 1-methyl- | 180 | C14H12 | 001730-37-6 | 94 |
| 3 | 9H-Fluorene, 1-methyl- | 180 | C14H12 | 001730-37-6 | 92 |
| 4 | 9H-Fluorene, 4-methyl- | 180 | C14H12 | 001556-99-6 | 76 |
| 5 | 9H-Fluorene, 1-methyl- | 180 | C14H12 | 001730-37-6 | 76 |



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080215\
Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
Operator : TP
Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** C01M4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
Quant Title : SVOA CALIBRATION

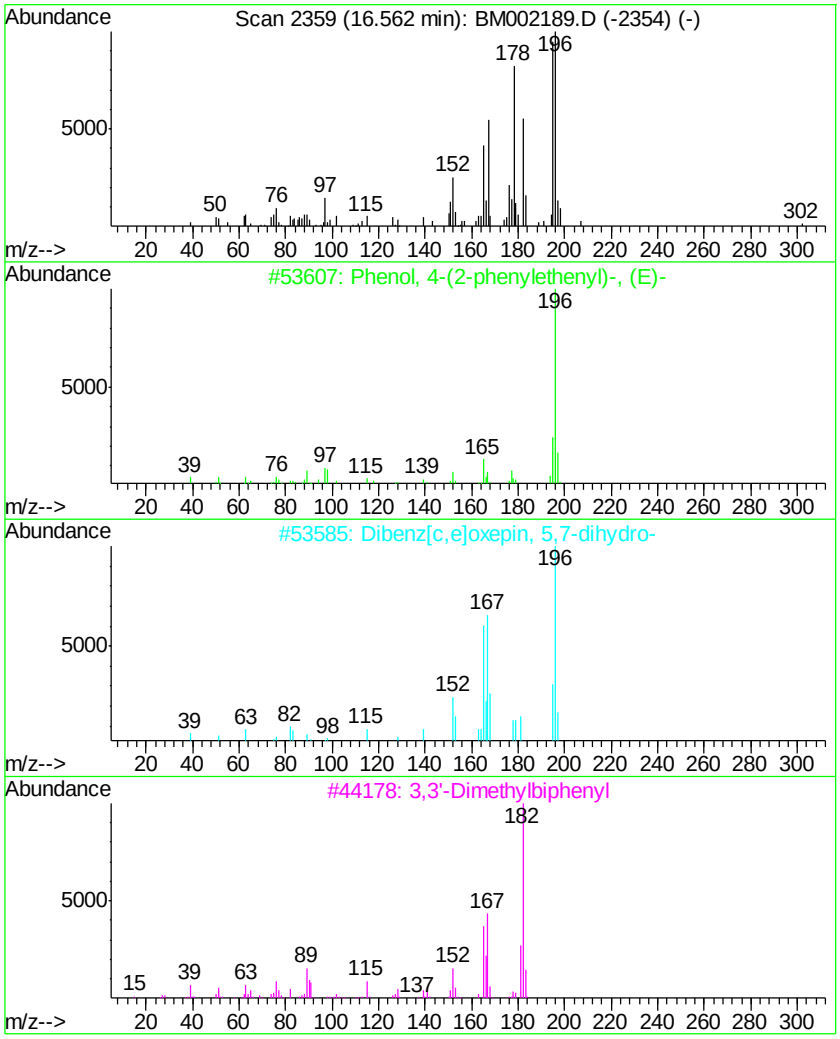
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

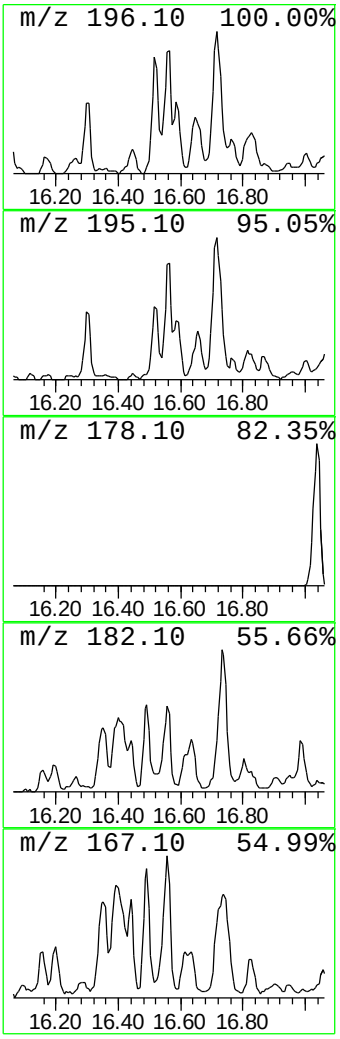
**********

Peak Number 5 Phenol, 4-(2-phenylethenyl)... Concentration Rank 31

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.56 | 4.06 ng/ul | 295614 | Phenanthrene-d10 | 16.99 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenol, 4-(2-phenylethenyl)-, (E)- | 196 | C14H12O | 006554-98-9 | 64 |
| 2 | Dibenz[c,e]oxepin, 5,7-dihydro- | 196 | C14H12O | 001136-22-7 | 45 |
| 3 | 3,3'-Dimethylbiphenyl | 182 | C14H14 | 000612-75-9 | 30 |
| 4 | Dibenz[b,f][1,4]oxazepine | 195 | C13H9NO | 000257-07-8 | 30 |
| 5 | 4,4'-Dimethylbiphenyl | 182 | C14H14 | 000613-33-2 | 30 |



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080215\
Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
Operator : TP
Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** C01M4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
Quant Title : SVOA CALIBRATION

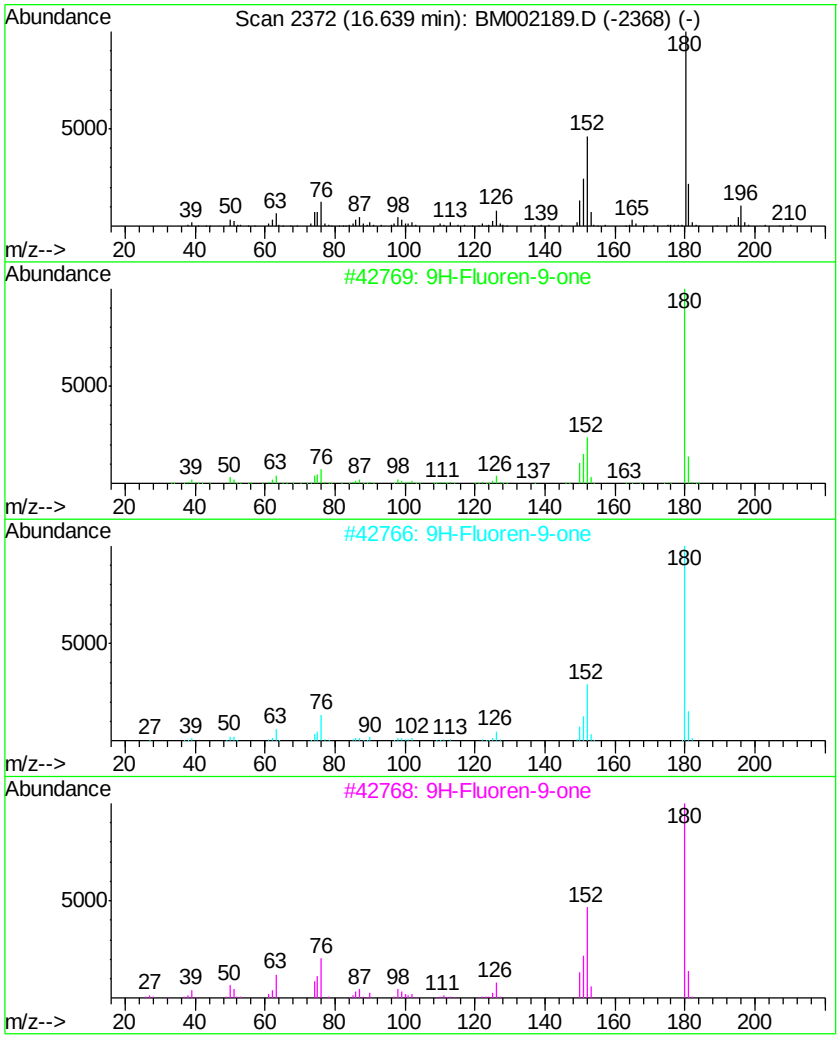
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

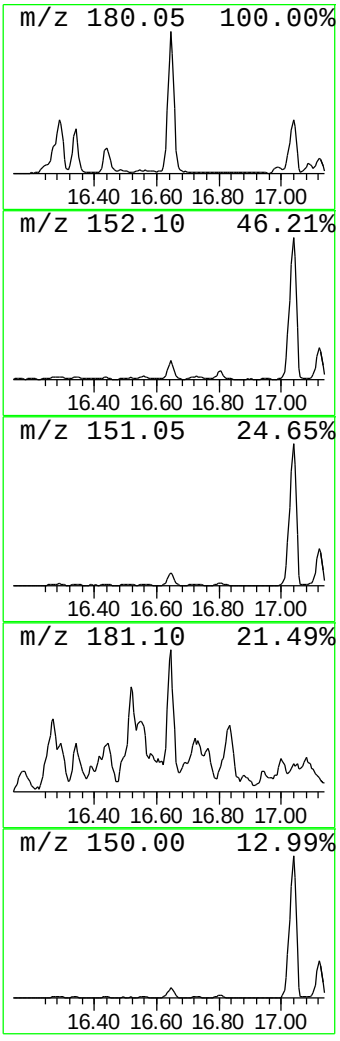
**********

Peak Number 6 9H-Fluoren-9-one Concentration Rank 23

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.64 | 5.77 ng/ul | 419443 | Phenanthrene-d10 | 16.99 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 9H-Fluoren-9-one | 180 | C13H8O | 000486-25-9 | 93 |
| 2 | 9H-Fluoren-9-one | 180 | C13H8O | 000486-25-9 | 93 |
| 3 | 9H-Fluoren-9-one | 180 | C13H8O | 000486-25-9 | 93 |
| 4 | Benzo[h]cinnoline | 180 | C12H8N2 | 000230-31-9 | 90 |
| 5 | 9H-Fluoren-9-one | 180 | C13H8O | 000486-25-9 | 90 |



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080215\
Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
Operator : TP
Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** C01M4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
Quant Title : SVOA CALIBRATION

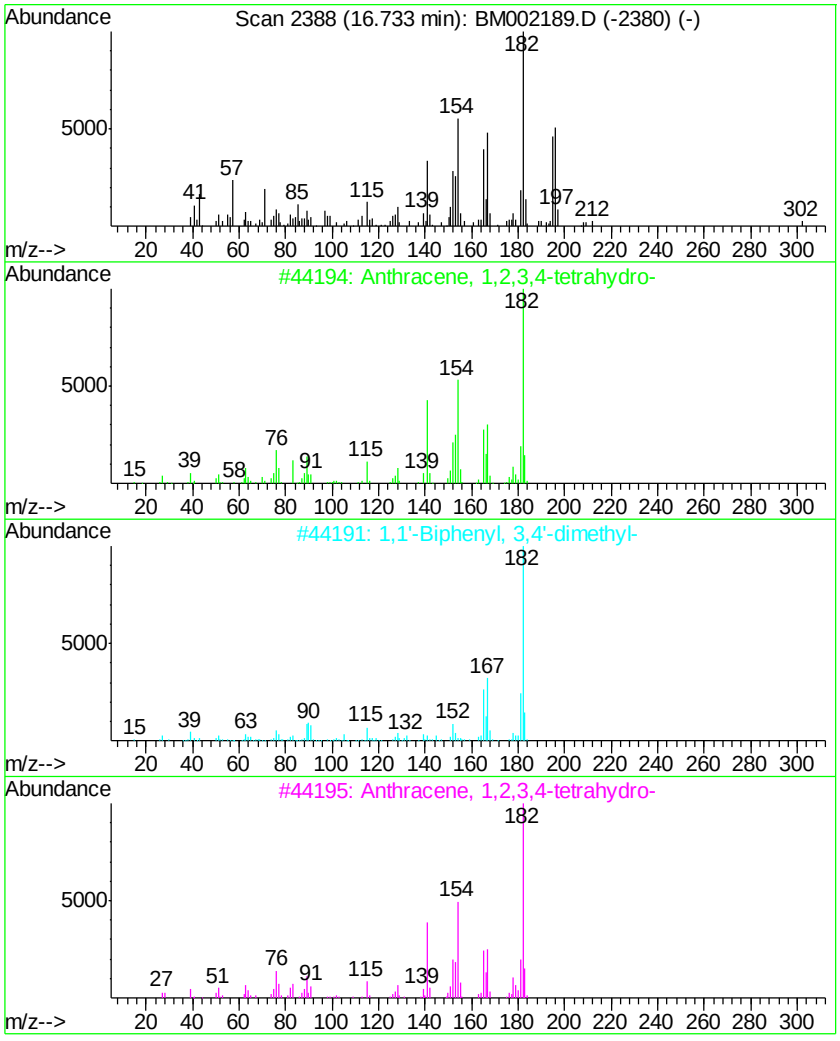
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

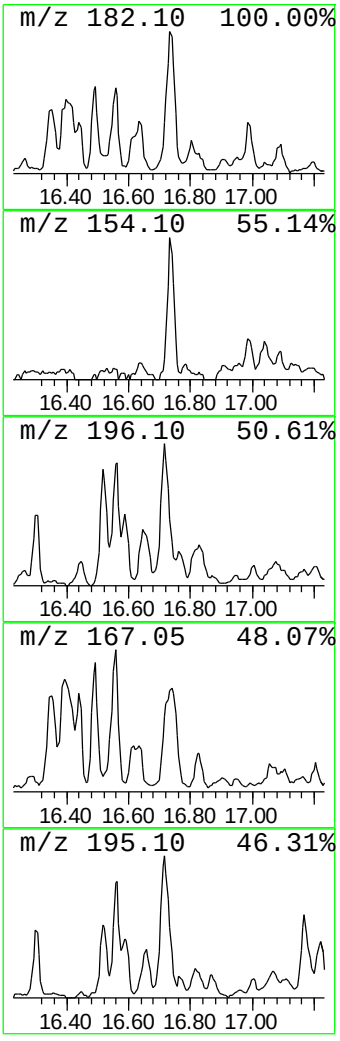
**********

Peak Number 7 Anthracene, 1,2,3,4-tetrahy... Concentration Rank 24

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.73 | 5.38 ng/ul | 391564 | Phenanthrene-d10 | 16.99 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Anthracene, 1,2,3,4-tetrahydro- | 182 | C14H14 | 002141-42-6 | 91 |
| 2 | 1,1'-Biphenyl, 3,4'-dimethyl- | 182 | C14H14 | 007383-90-6 | 64 |
| 3 | Anthracene, 1,2,3,4-tetrahydro- | 182 | C14H14 | 002141-42-6 | 60 |
| 4 | 7-Phenylbicyclo[3.2.1]octa-2,6-d... | 182 | C14H14 | 1000201-01-7 | 55 |
| 5 | 1,1'-Biphenyl, 3,4'-dimethyl- | 182 | C14H14 | 007383-90-6 | 50 |



Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080215\
Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
Operator : TP
Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** C01M4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
Quant Title : SVOA CALIBRATION

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

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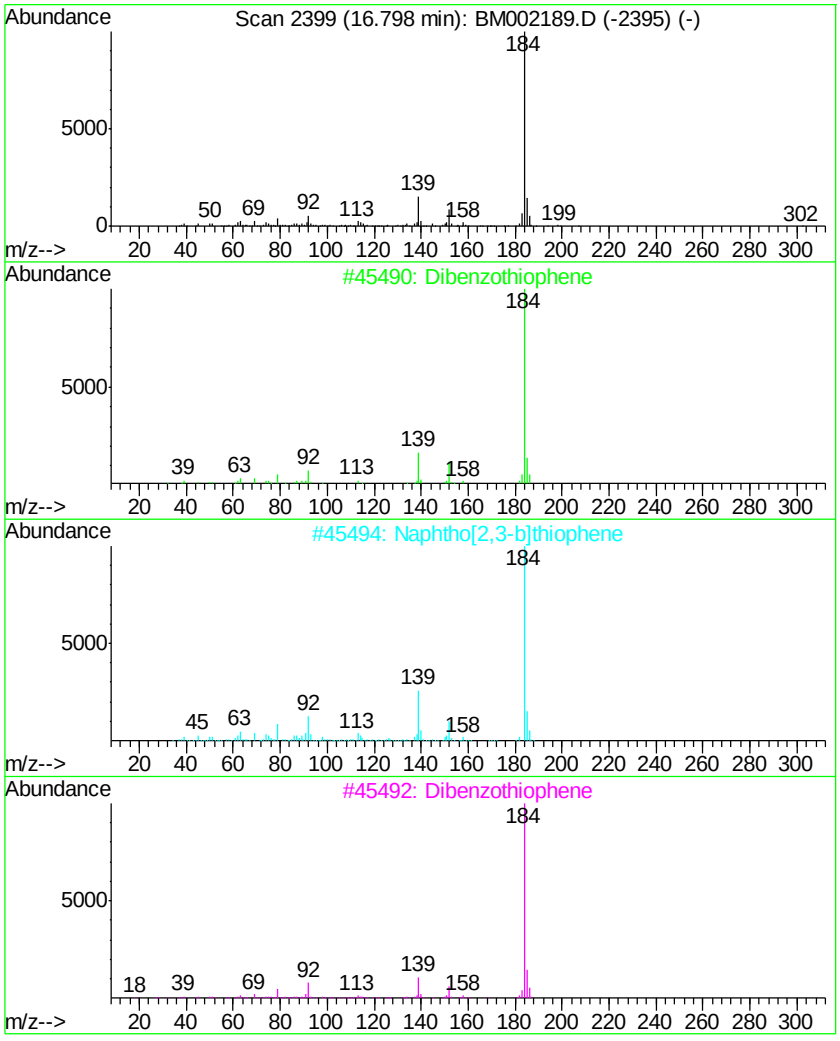
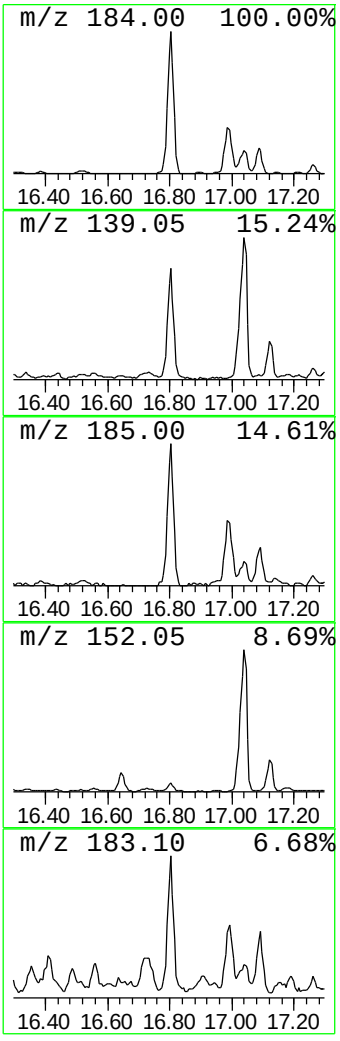
Peak Number 8 Dibenzothiophene Concentration Rank 14

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.80 | 8.36 ng/ul | 607919 | Phenanthrene-d10 | 16.99 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Dibenzothiophene | 184 | C12H8S | 000132-65-0 | 96 |
| 2 | Naphtho[2,3-b]thiophene | 184 | C12H8S | 000268-77-9 | 94 |
| 3 | Dibenzothiophene | 184 | C12H8S | 000132-65-0 | 94 |
| 4 | Dibenzothiophene | 184 | C12H8S | 000132-65-0 | 93 |
| 5 | Dibenzothiophene | 184 | C12H8S | 000132-65-0 | 93 |



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080215\
Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
Operator : TP
Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :**
BNA_M
**ClientSampleId :**
C01M4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
Quant Title : SVOA CALIBRATION

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

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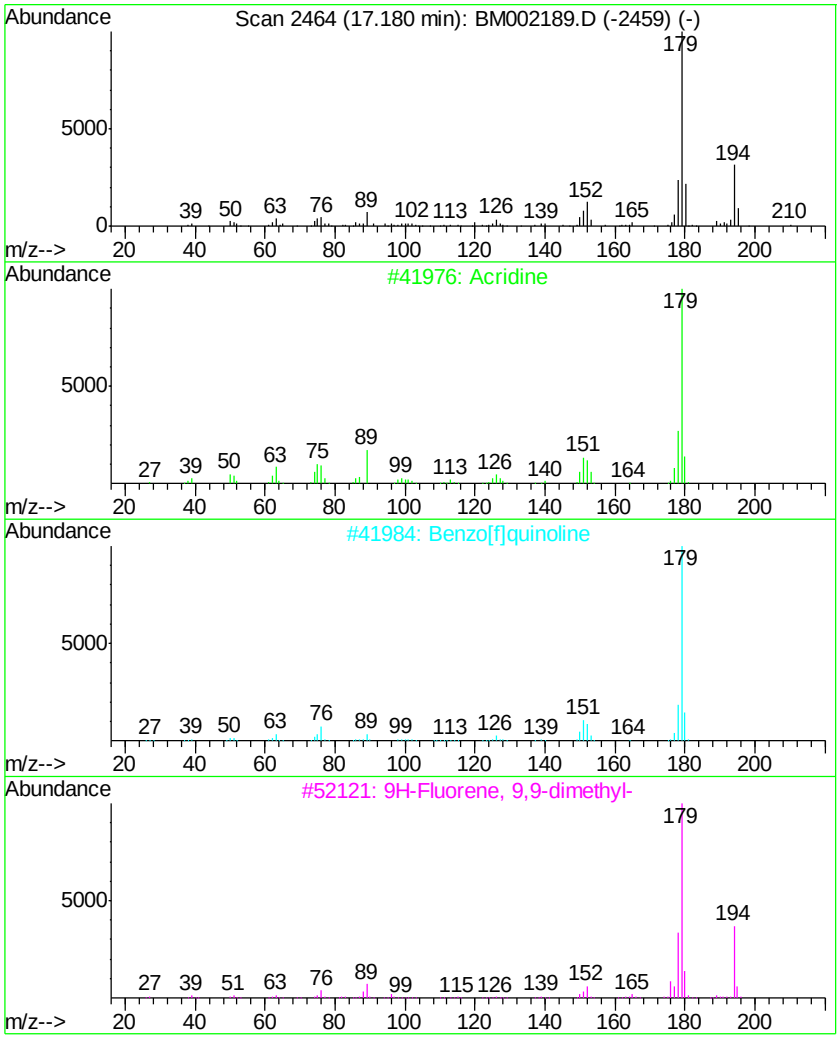
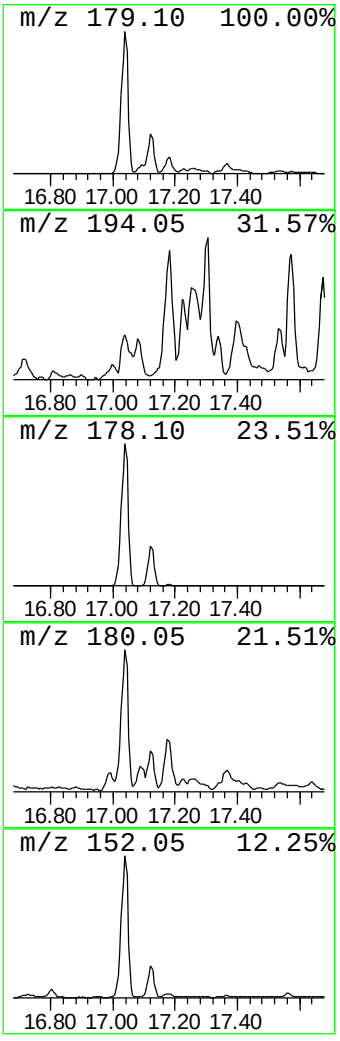
Peak Number 9 Acridine Concentration Rank 40

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 17.18 | 3.34 ng/ul | 242829 | Phenanthrene-d10 | 16.99 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Acridine | 179 | C13H9N | 000260-94-6 | 93 |
| 2 | Benzo[f]quinoline | 179 | C13H9N | 000085-02-9 | 90 |
| 3 | 9H-Fluorene, 9,9-dimethyl- | 194 | C15H14 | 004569-45-3 | 80 |
| 4 | Benzo[h]quinoline | 179 | C13H9N | 000230-27-3 | 76 |
| 5 | Phenanthridine | 179 | C13H9N | 000229-87-8 | 76 |



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080215\
Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
Operator : TP
Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** C01M4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
Quant Title : SVOA CALIBRATION

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

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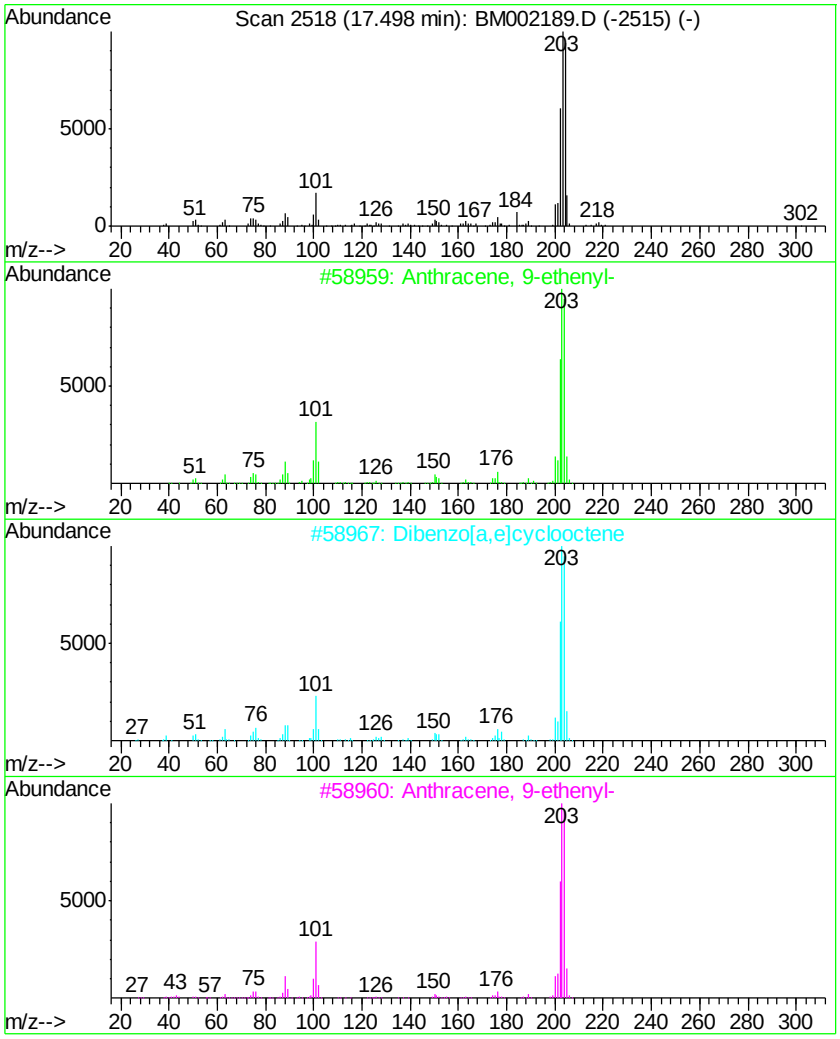
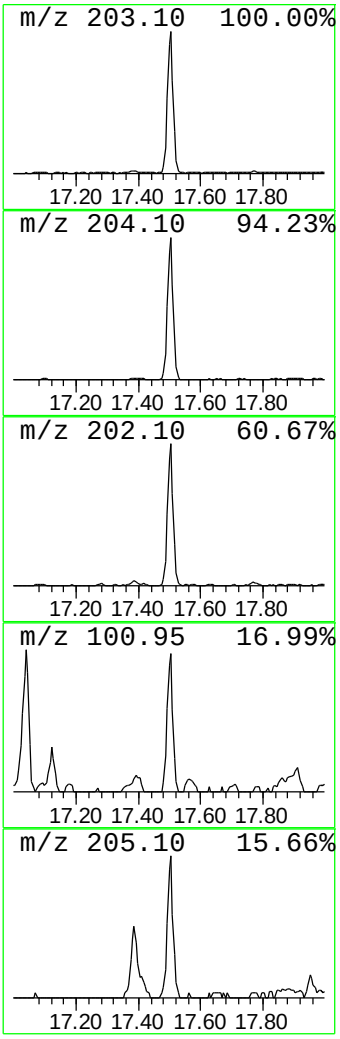
Peak Number 10 Anthracene, 9-ethenyl- Concentration Rank 32

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 17.50 | 3.98 ng/ul | 289270 | Phenanthrene-d10 | 16.99 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Anthracene, 9-ethenyl- | 204 | C16H12 | 002444-68-0 | 94 |
| 2 | Dibenzo[a,e]cyclooctene | 204 | C16H12 | 000262-89-5 | 93 |
| 3 | Anthracene, 9-ethenyl- | 204 | C16H12 | 002444-68-0 | 89 |
| 4 | 1,4-Ethenoanthracene, 1,4-dihydro- | 204 | C16H12 | 027765-96-4 | 81 |
| 5 | Naphthalene, 1-phenyl- | 204 | C16H12 | 000605-02-7 | 76 |



Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080215\
Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
Operator : TP
Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** C01M4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
Quant Title : SVOA CALIBRATION

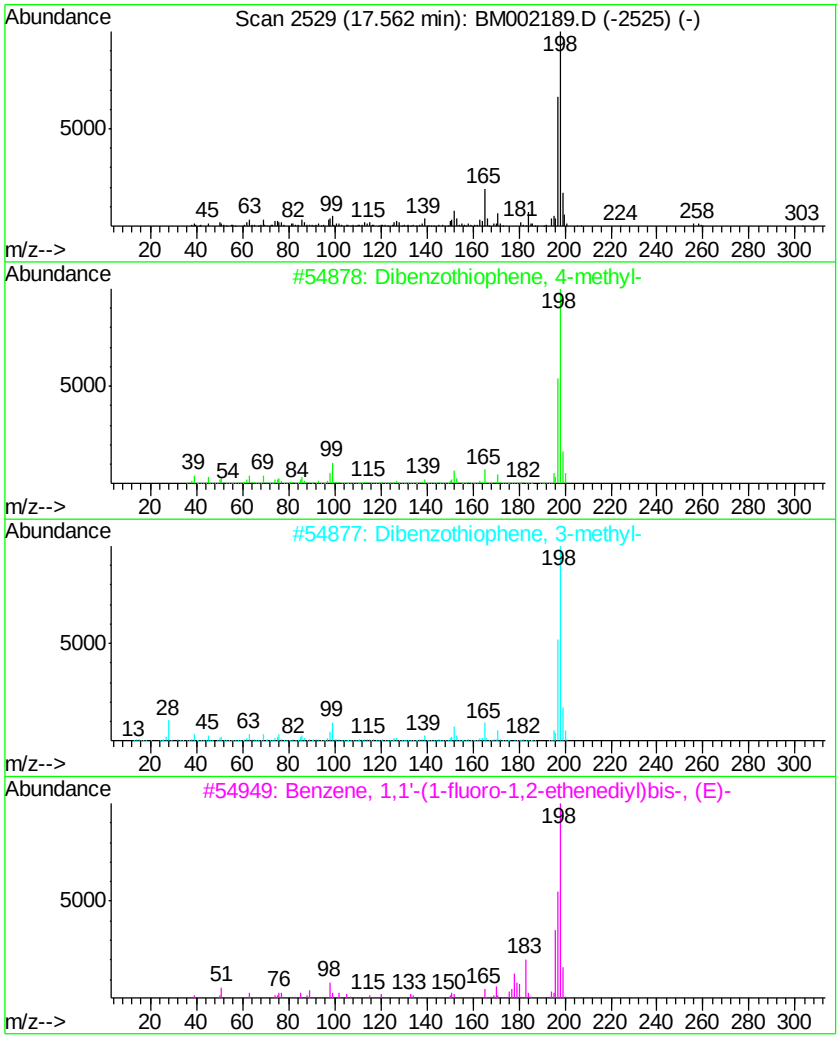
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

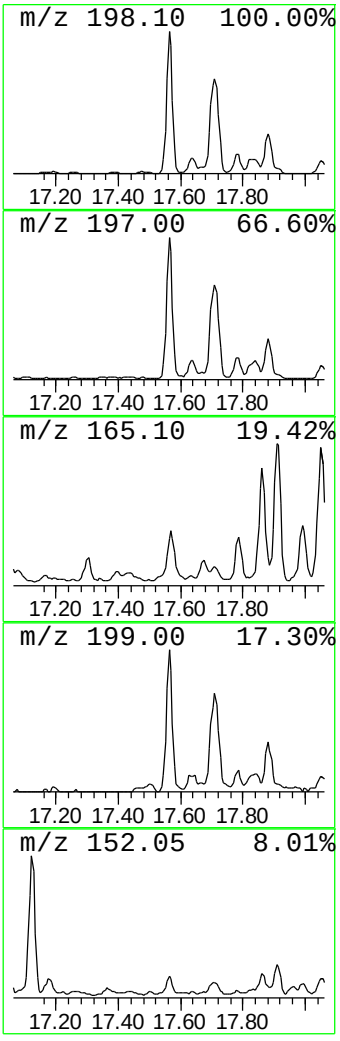
**********

Peak Number 11 Dibenzothiophene, 4-methyl- Concentration Rank 17

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 17.56 | 8.10 ng/ul | 589162 | Phenanthrene-d10 | 16.99 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Dibenzothiophene, 4-methyl- | 198 | C13H10S | 007372-88-5 | 91 |
| 2 | Dibenzothiophene, 3-methyl- | 198 | C13H10S | 016587-52-3 | 87 |
| 3 | Benzene, 1,1'-(1-fluoro-1,2-ethe... | 198 | C14H11F | 000671-19-2 | 72 |
| 4 | Benzene, 1-fluoro-4-(2-phenyleth... | 198 | C14H11F | 017404-69-2 | 72 |
| 5 | Benzenamine, 4,4'-methylenebis- | 198 | C13H14N2 | 000101-77-9 | 64 |



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080215\
Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
Operator : TP
Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** C01M4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
Quant Title : SVOA CALIBRATION

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

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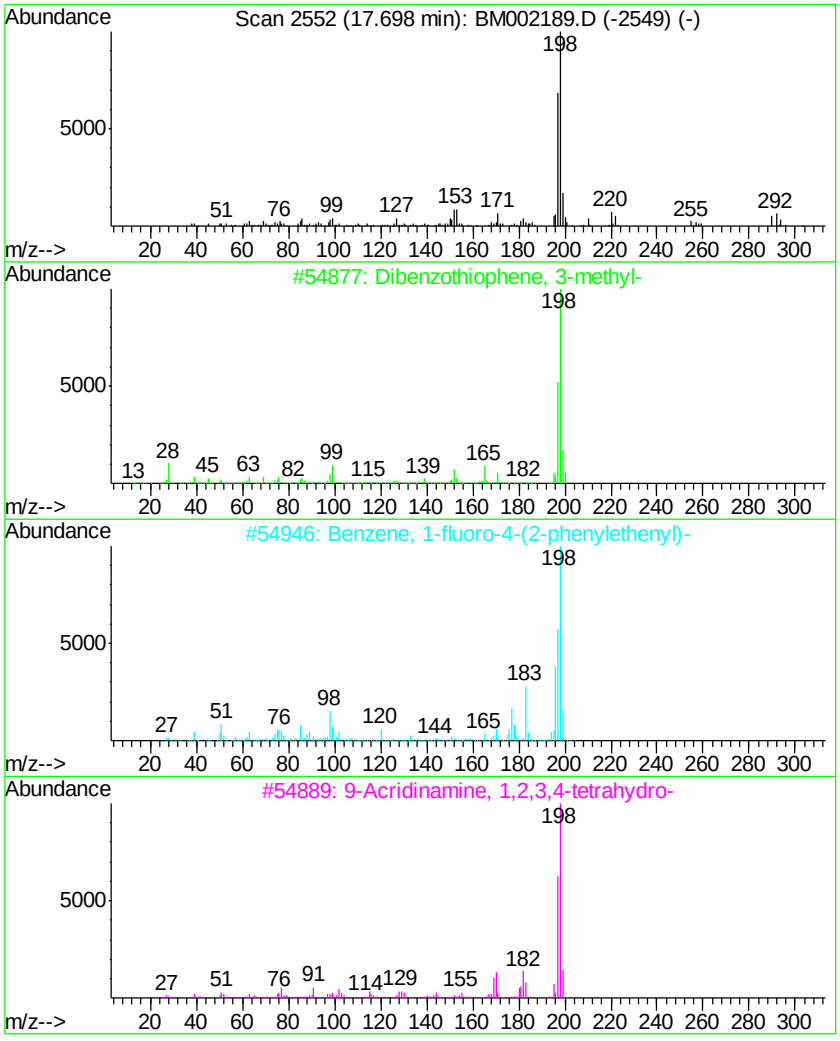
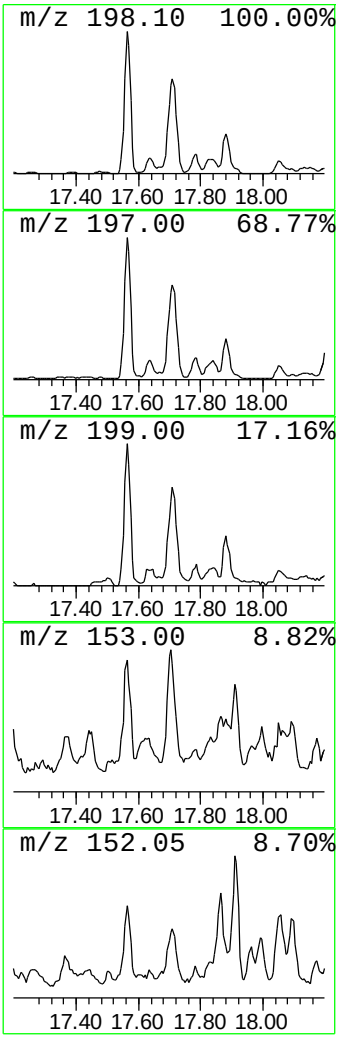
Peak Number 12 Dibenzothiophene, 3-methyl- Concentration Rank 25

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 17.70 | 5.19 ng/ul | 377118 | Phenanthrene-d10 | 16.99 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Dibenzothiophene, 3-methyl- | 198 | C13H10S | 016587-52-3 | 87 |
| 2 | Benzene, 1-fluoro-4-(2-phenyleth... | 198 | C14H11F | 017404-69-2 | 80 |
| 3 | 9-Acridinamine, 1,2,3,4-tetrahydro- | 198 | C13H14N2 | 000321-64-2 | 72 |
| 4 | .alpha.-[1-Indanylidene]cyanoace... | 198 | C12H10N2O | 058787-18-1 | 60 |
| 5 | Benzene, 1,1'-(1-fluoro-1,2-ethe... | 198 | C14H11F | 000671-19-2 | 56 |



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080215\
Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
Operator : TP
Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** C01M4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
Quant Title : SVOA CALIBRATION

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

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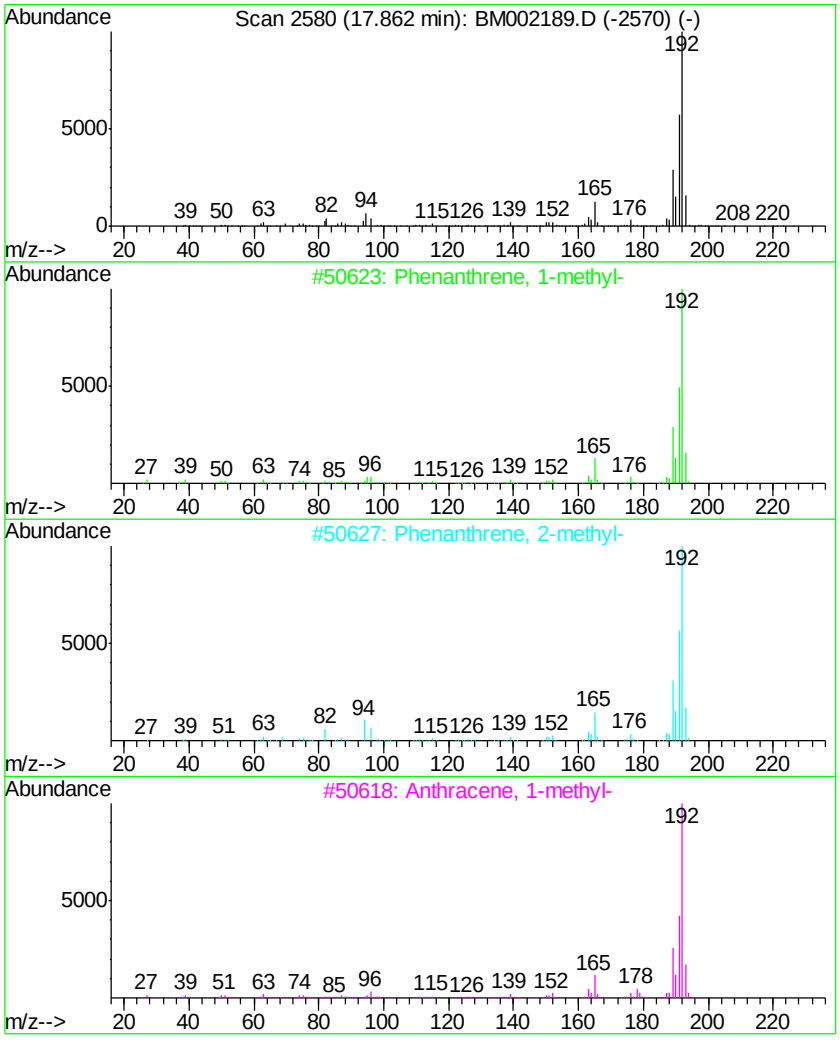
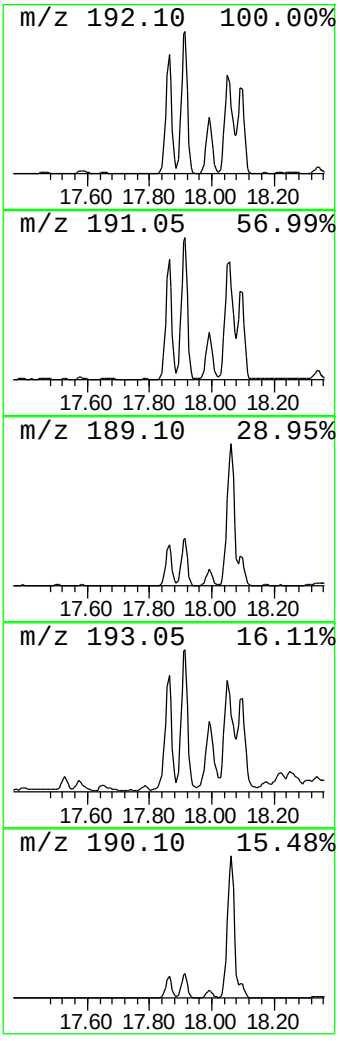
Peak Number 13 Phenanthrene, 1-methyl- Concentration Rank 4

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 17.86 | 25.28 ng/ul | 1838600 | Phenanthrene-d10 | 16.99 |

| 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 | 93 |
| 3 | Anthracene, 1-methyl- | 192 | C15H12 | 000610-48-0 | 93 |
| 4 | Anthracene, 2-methyl- | 192 | C15H12 | 000613-12-7 | 93 |
| 5 | Anthracene, 1-methyl- | 192 | C15H12 | 000610-48-0 | 91 |



# Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080215\
Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
Operator : TP
Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** C01M4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
Quant Title : SVOA CALIBRATION

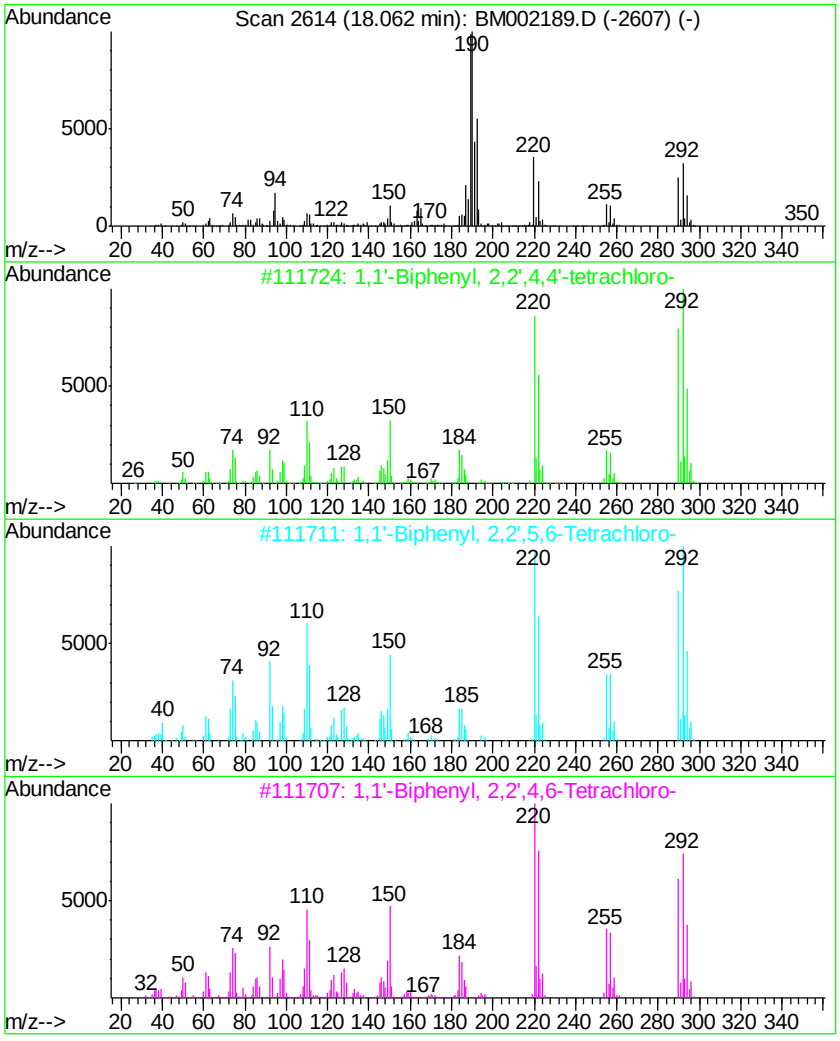
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

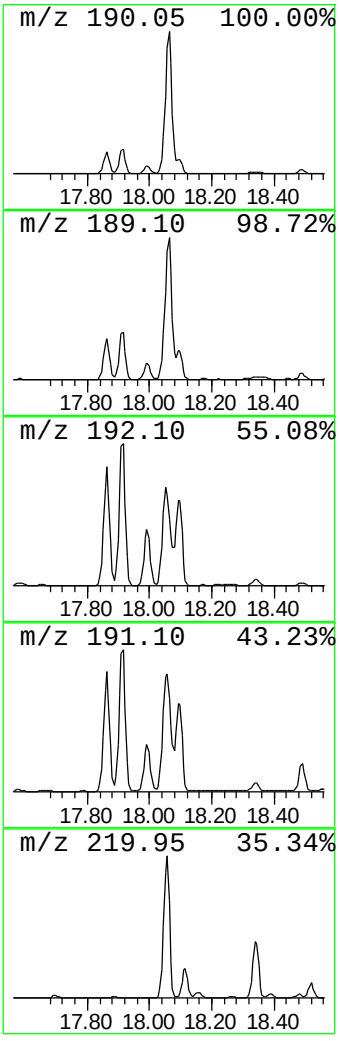
**********

Peak Number 16 1,1'-Biphenyl, 2,2',4,4'-te... Concentration Rank 1

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.06 | 75.34 ng/ul | 5478970 | Phenanthrene-d10 | 16.99 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1,1'-Biphenyl, 2,2',4,4'-tetrach... | 290 | C12H6Cl4 | 002437-79-8 | 94 |
| 2 | 1,1'-Biphenyl, 2,2',5,6-Tetrachl... | 290 | C12H6Cl4 | 041464-41-9 | 94 |
| 3 | 1,1'-Biphenyl, 2,2',4,6-Tetrachl... | 290 | C12H6Cl4 | 062796-65-0 | 94 |
| 4 | 1,1'-Biphenyl, 2,2',4,5'-tetrach... | 290 | C12H6Cl4 | 041464-40-8 | 94 |
| 5 | 1,1'-Biphenyl, 2,2',4,5'-tetrach... | 290 | C12H6Cl4 | 041464-40-8 | 93 |



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080215\
Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
Operator : TP
Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** C01M4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
Quant Title : SVOA CALIBRATION

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

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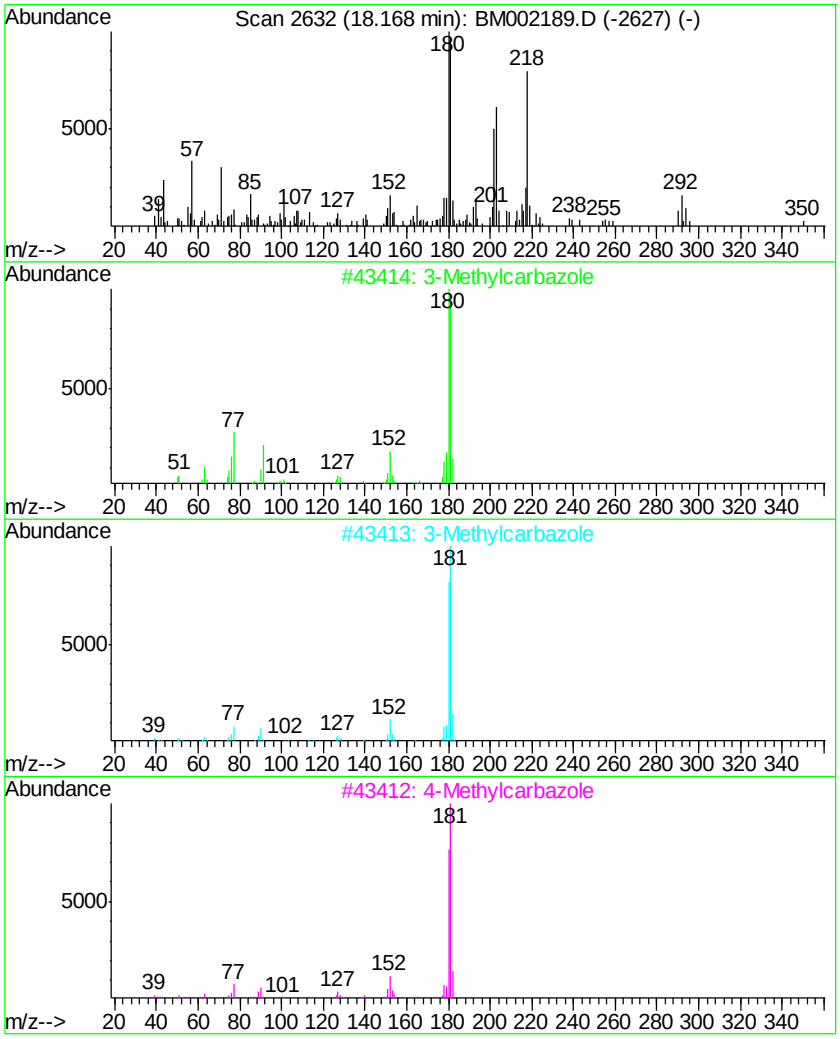
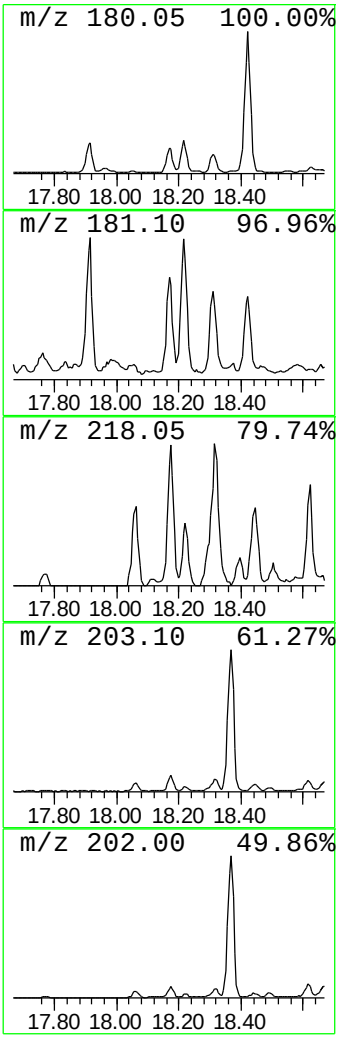
Peak Number 18 3-Methylcarbazole Concentration Rank 37

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.17 | 3.42 ng/ul | 249032 | Phenanthrene-d10 | 16.99 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 3-Methylcarbazole | 181 | C13H11N | 004630-20-0 | 83 |
| 2 | 3-Methylcarbazole | 181 | C13H11N | 004630-20-0 | 42 |
| 3 | 4-Methylcarbazole | 181 | C13H11N | 003770-48-7 | 42 |
| 4 | 9H-Carbazole, 2-methyl- | 181 | C13H11N | 003652-91-3 | 42 |
| 5 | 3-Methylcarbazole | 181 | C13H11N | 004630-20-0 | 42 |



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080215\
Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
Operator : TP
Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** C01M4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
Quant Title : SVOA CALIBRATION

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

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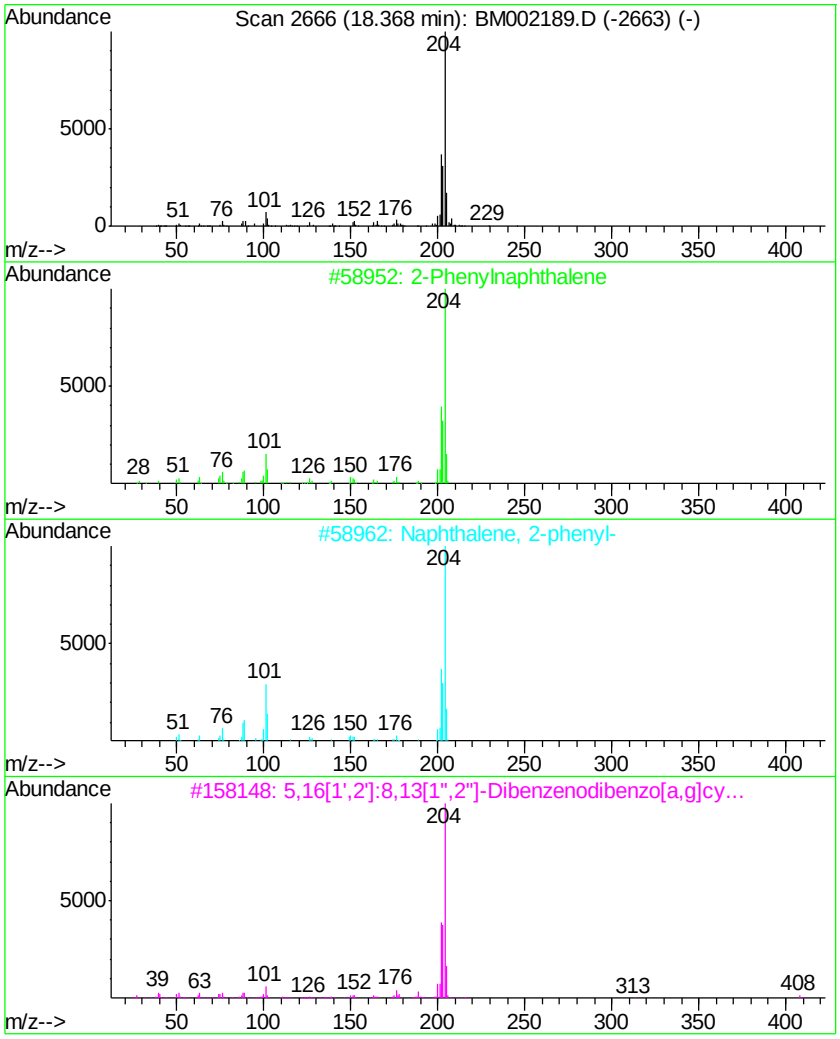
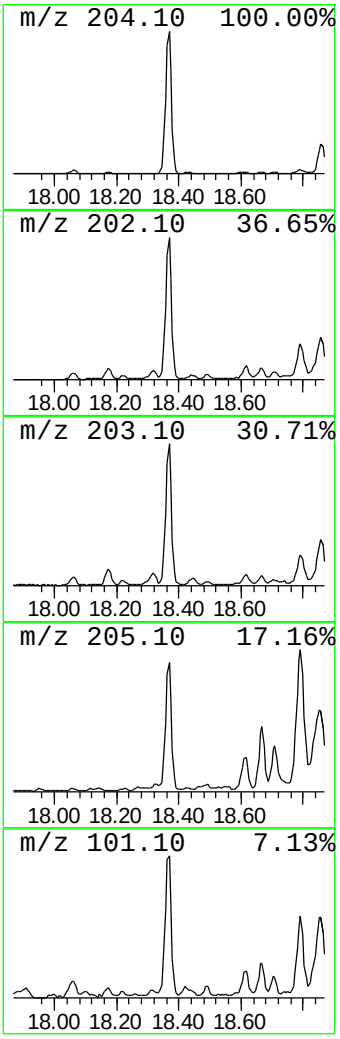
Peak Number 20 2-Phenylnaphthalene Concentration Rank 9

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.37 | 14.17 ng/ul | 1030180 | Phenanthrene-d10 | 16.99 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-Phenylnaphthalene | 204 | C16H12 | 035465-71-5 | 93 |
| 2 | Naphthalene, 2-phenyl- | 204 | C16H12 | 000612-94-2 | 90 |
| 3 | 5,16[1',2']:8,13[1'',2'']-Dibenz... | 408 | C32H24 | 005672-97-9 | 87 |
| 4 | Naphthalene, 2-phenyl- | 204 | C16H12 | 000612-94-2 | 81 |
| 5 | Naphthalene, 2-phenyl- | 204 | C16H12 | 000612-94-2 | 81 |



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Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
Operator : TP
Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** C01M4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
Quant Title : SVOA CALIBRATION

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

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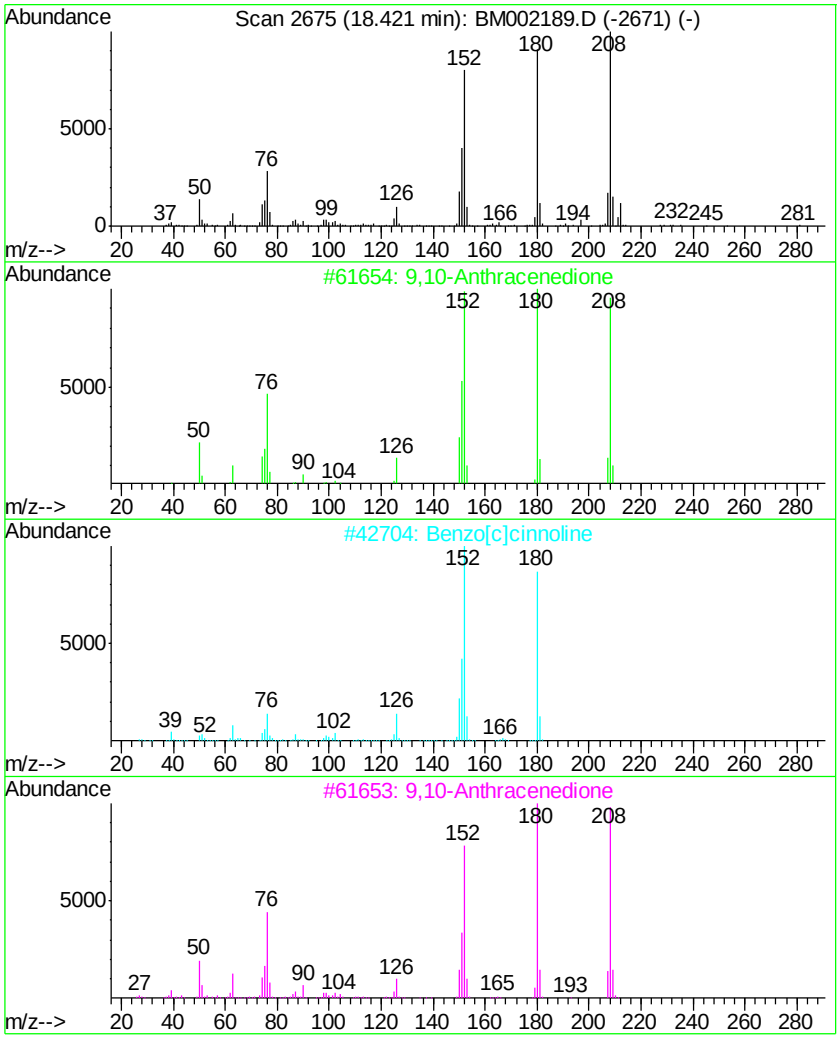
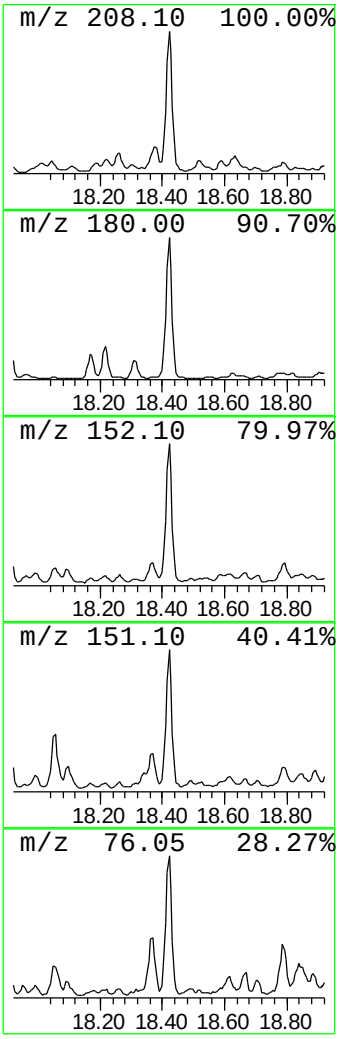
Peak Number 21 9,10-Anthracenedione Concentration Rank 12

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.42 | 10.61 ng/ul | 771413 | Phenanthrene-d10 | 16.99 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 9,10-Anthracenedione | 208 | C14H8O2 | 000084-65-1 | 97 |
| 2 | Benzo[c]cinnoline | 180 | C12H8N2 | 000230-17-1 | 96 |
| 3 | 9,10-Anthracenedione | 208 | C14H8O2 | 000084-65-1 | 94 |
| 4 | 9,10-Anthracenedione | 208 | C14H8O2 | 000084-65-1 | 90 |
| 5 | Benzo[c]cinnoline | 180 | C12H8N2 | 000230-17-1 | 64 |



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Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
Operator : TP
Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** C01M4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
Quant Title : SVOA CALIBRATION

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

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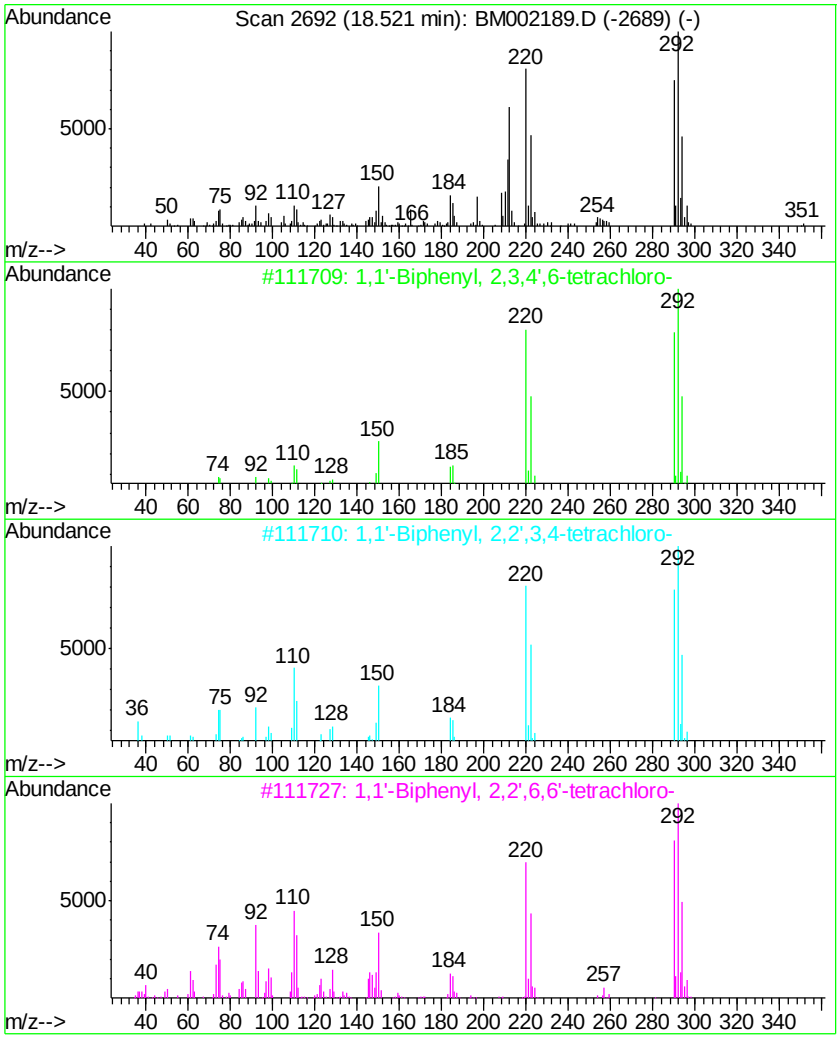
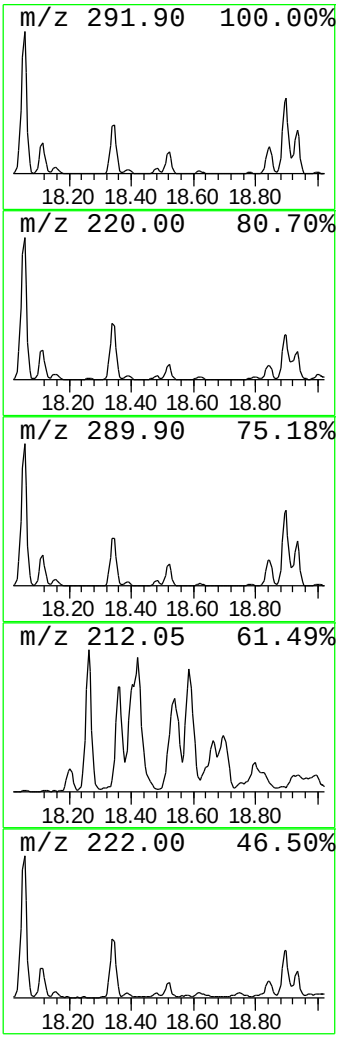
Peak Number 22 1,1'-Biphenyl, 2,3,4',6-tet... Concentration Rank 29

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.52 | 4.25 ng/ul | 309102 | Phenanthrene-d10 | 16.99 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1,1'-Biphenyl, 2,3,4',6-tetrachl... | 290 | C12H6Cl4 | 052663-58-8 | 99 |
| 2 | 1,1'-Biphenyl, 2,2',3,4-tetrachl... | 290 | C12H6Cl4 | 052663-59-9 | 98 |
| 3 | 1,1'-Biphenyl, 2,2',6,6'-tetrach... | 290 | C12H6Cl4 | 015968-05-5 | 98 |
| 4 | 1,1'-Biphenyl, 2,3',5,5'-tetrach... | 290 | C12H6Cl4 | 041464-42-0 | 98 |
| 5 | 1,1'-Biphenyl, 2,2',4,4'-tetrach... | 290 | C12H6Cl4 | 002437-79-8 | 95 |



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Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
Operator : TP
Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** C01M4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
Quant Title : SVOA CALIBRATION

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

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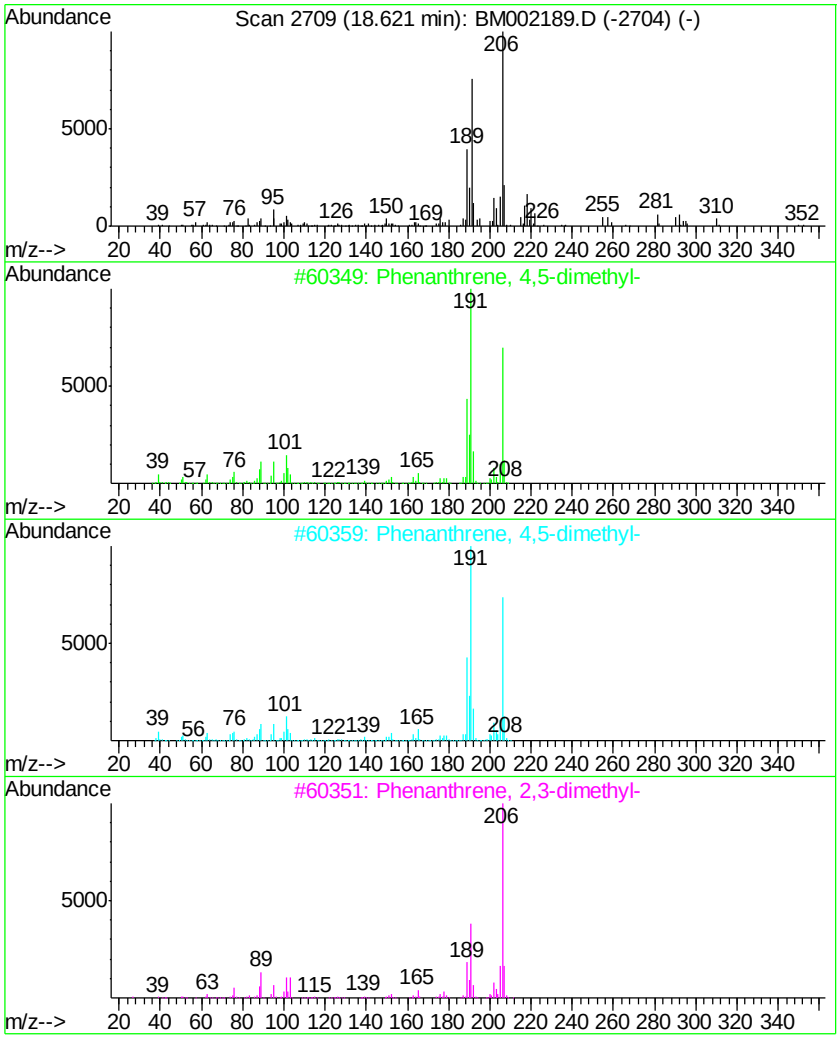
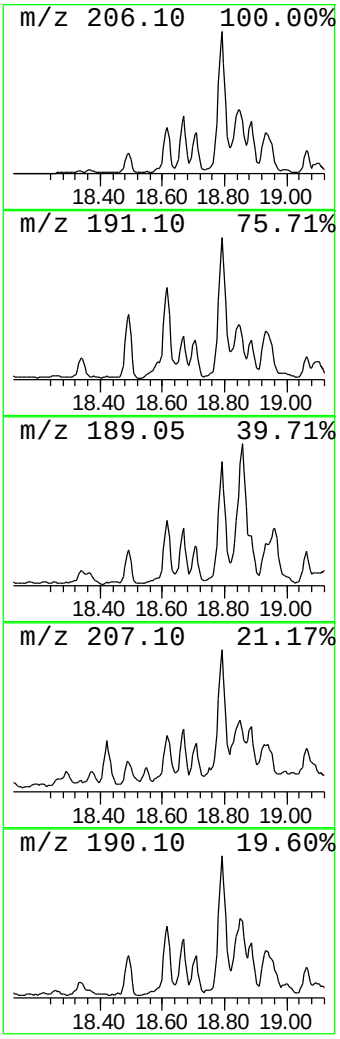
Peak Number 23 Phenanthrene, 4,5-dimethyl- Concentration Rank 20

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.62 | 6.35 ng/ul | 461515 | Phenanthrene-d10 | 16.99 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenanthrene, 4,5-dimethyl- | 206 | C16H14 | 003674-69-9 | 93 |
| 2 | Phenanthrene, 4,5-dimethyl- | 206 | C16H14 | 003674-69-9 | 93 |
| 3 | Phenanthrene, 2,3-dimethyl- | 206 | C16H14 | 003674-65-5 | 90 |
| 4 | Phenanthrene, 4,5-dimethyl- | 206 | C16H14 | 003674-69-9 | 90 |
| 5 | Phenanthrene, 2,5-dimethyl- | 206 | C16H14 | 003674-66-6 | 86 |



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Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
Operator : TP
Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** C01M4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
Quant Title : SVOA CALIBRATION

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

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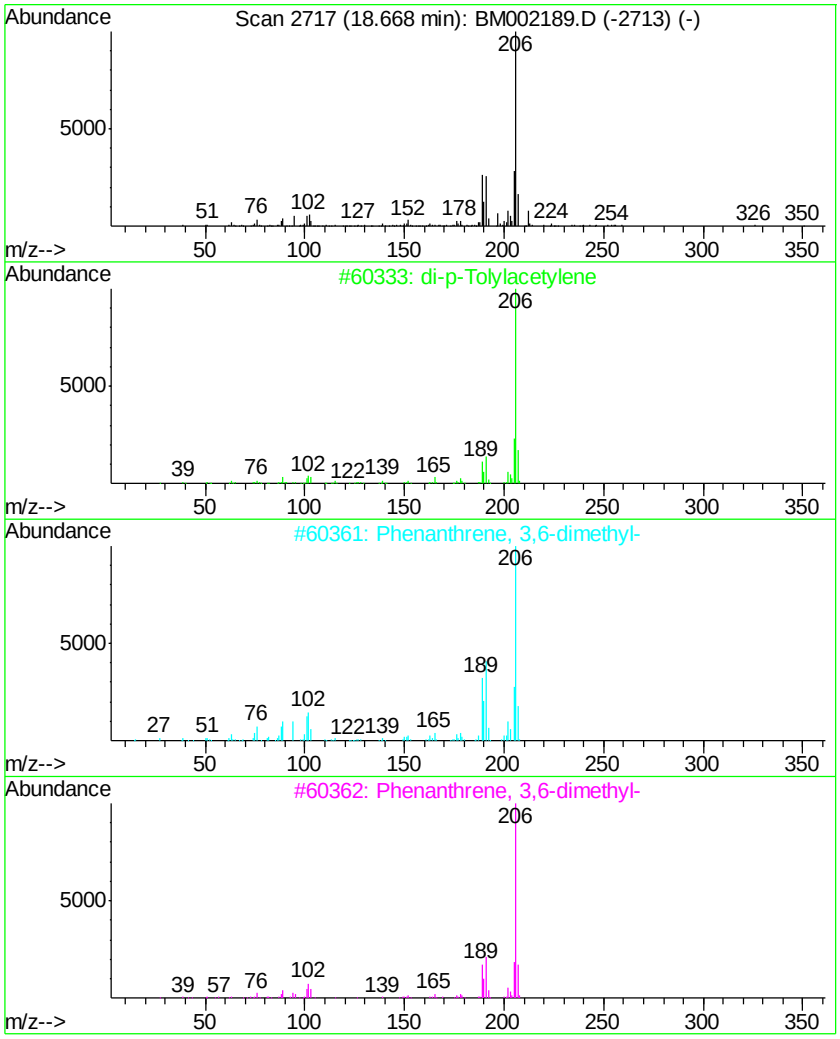
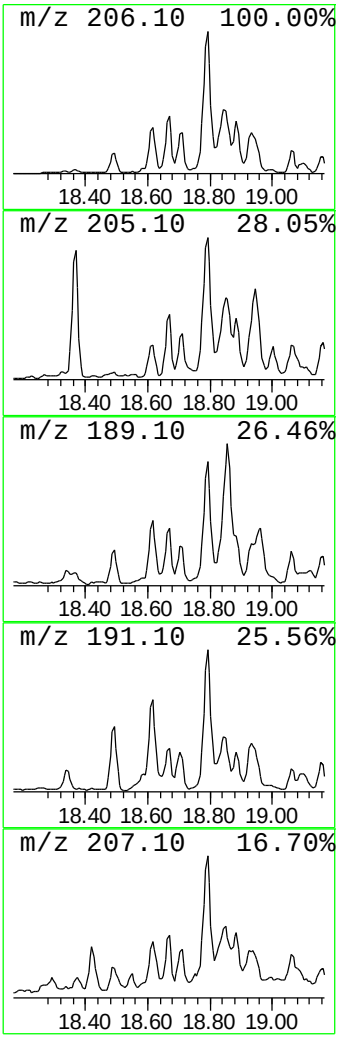
Peak Number 24 di-p-Tolylacetylene Concentration Rank 30

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.67 | 4.22 ng/ul | 307005 | Phenanthrene-d10 | 16.99 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | di-p-Tolylacetylene | 206 | C16H14 | 002789-88-0 | 94 |
| 2 | Phenanthrene, 3,6-dimethyl- | 206 | C16H14 | 001576-67-6 | 93 |
| 3 | Phenanthrene, 3,6-dimethyl- | 206 | C16H14 | 001576-67-6 | 91 |
| 4 | di-p-Tolylacetylene | 206 | C16H14 | 002789-88-0 | 90 |
| 5 | Phenanthrene, 2,7-dimethyl- | 206 | C16H14 | 001576-69-8 | 90 |



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080215\
Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
Operator : TP
Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** C01M4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
Quant Title : SVOA CALIBRATION

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

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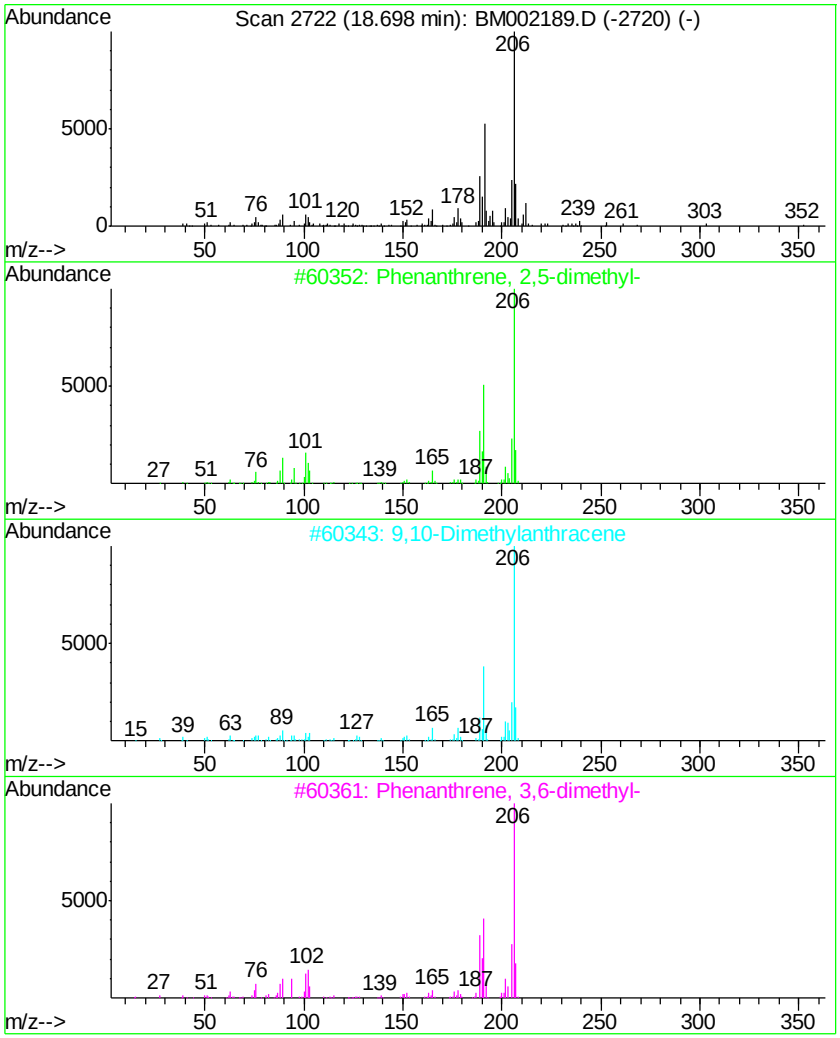
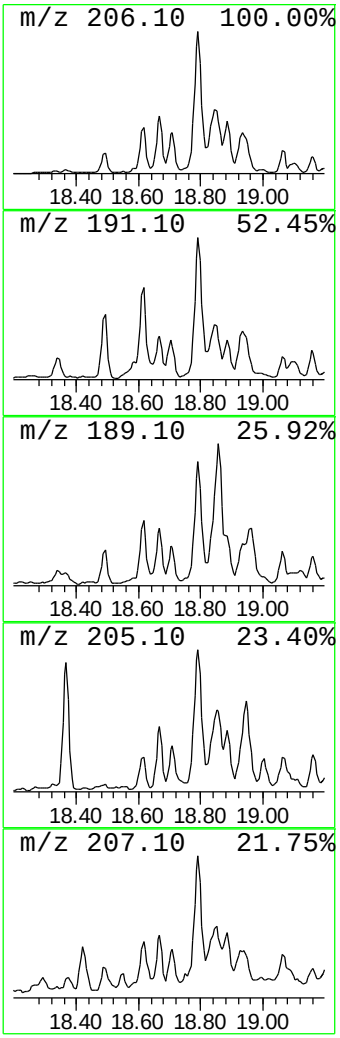
Peak Number 25 Phenanthrene, 2,5-dimethyl- Concentration Rank 27

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.70 | 4.87 ng/ul | 353942 | Phenanthrene-d10 | 16.99 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenanthrene, 2,5-dimethyl- | 206 | C16H14 | 003674-66-6 | 94 |
| 2 | 9,10-Dimethylanthracene | 206 | C16H14 | 000781-43-1 | 91 |
| 3 | Phenanthrene, 3,6-dimethyl- | 206 | C16H14 | 001576-67-6 | 91 |
| 4 | Phenanthrene, 3,6-dimethyl- | 206 | C16H14 | 001576-67-6 | 91 |
| 5 | Naphthalene, 1,2-dihydro-4-phenyl- | 206 | C16H14 | 007469-40-1 | 91 |



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080215\
Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
Operator : TP
Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** C01M4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
Quant Title : SVOA CALIBRATION

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

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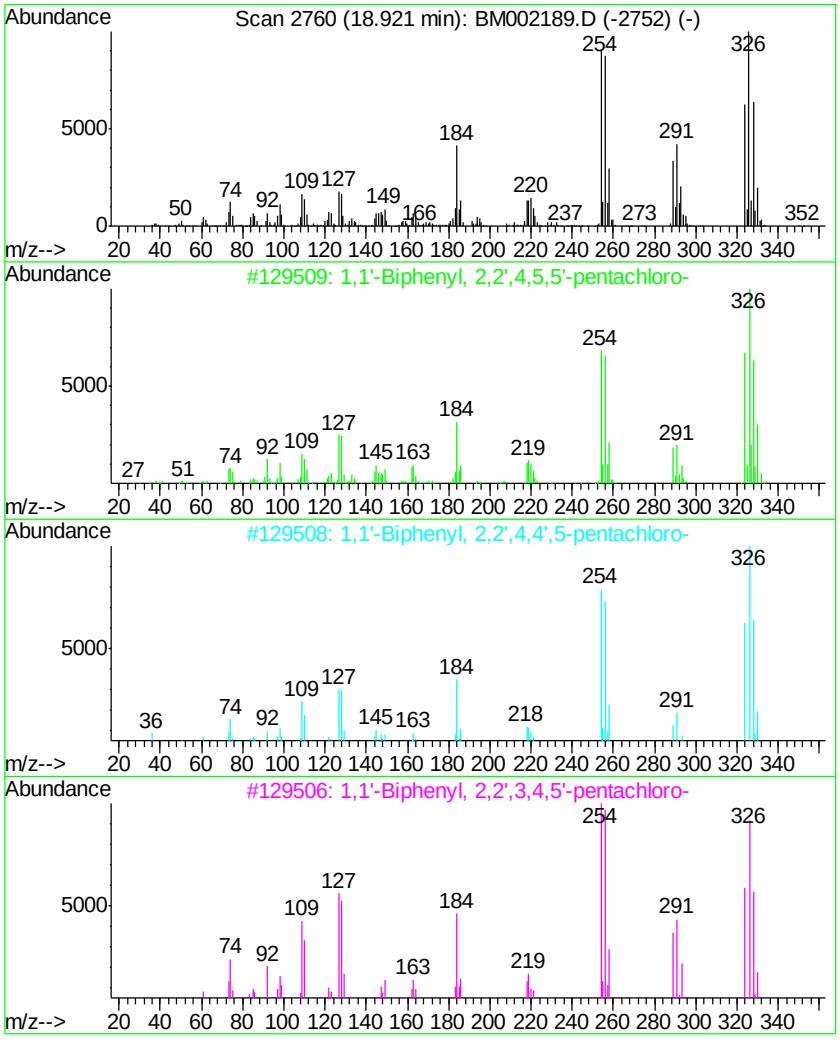
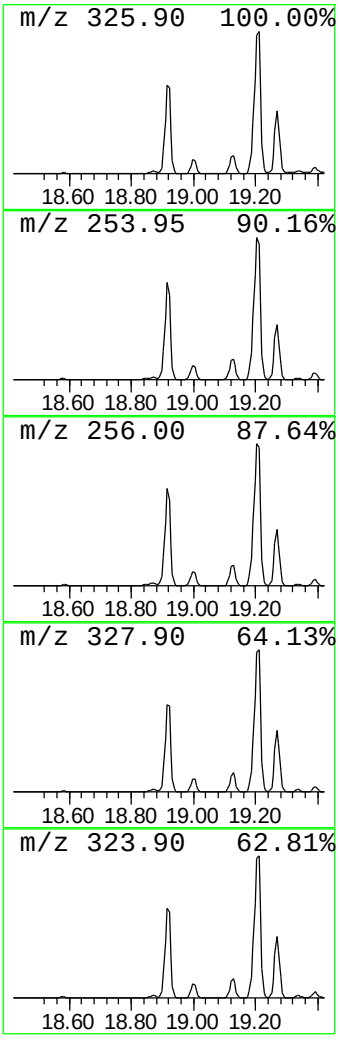
Peak Number 28 1,1'-Biphenyl, 2,2',4,5,5'-... Concentration Rank 2

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.92 | 65.96 ng/ul | 4797070 | Phenanthrene-d10 | 16.99 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1,1'-Biphenyl, 2,2',4,5,5'-penta... | 324 | C12H5Cl5 | 037680-73-2 | 99 |
| 2 | 1,1'-Biphenyl, 2,2',4,4',5-penta... | 324 | C12H5Cl5 | 038380-01-7 | 99 |
| 3 | 1,1'-Biphenyl, 2,2',3,4,5'-penta... | 324 | C12H5Cl5 | 038380-02-8 | 99 |
| 4 | 1,1'-Biphenyl, 2,2',3',4,5-Penta... | 324 | C12H5Cl5 | 041464-51-1 | 99 |
| 5 | 1,1'-Biphenyl, 2,2',3,5',6-penta... | 324 | C12H5Cl5 | 038379-99-6 | 99 |



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080215\
Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
Operator : TP
Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** C01M4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
Quant Title : SVOA CALIBRATION

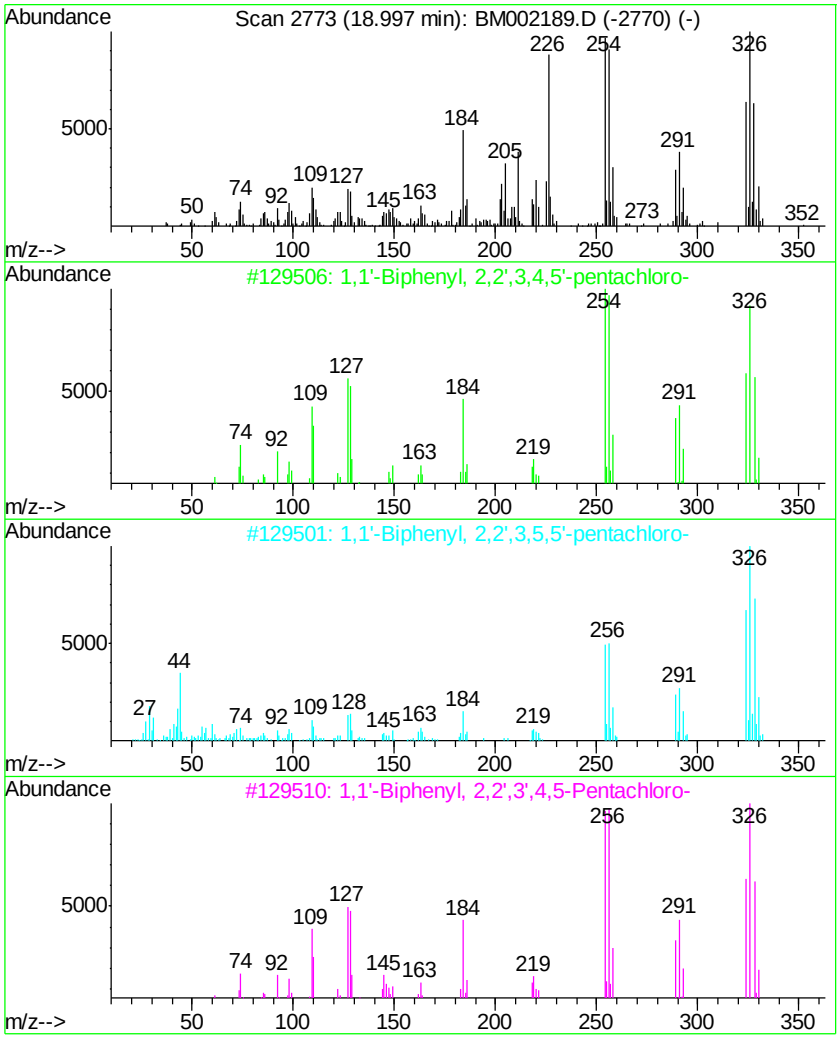
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

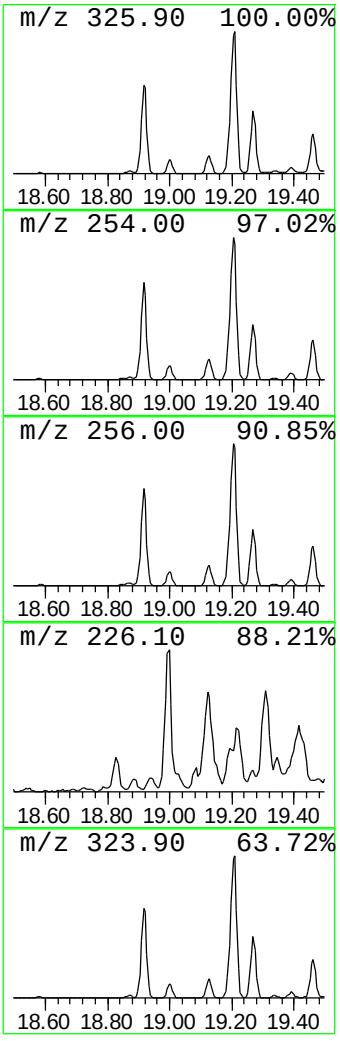
**********

Peak Number 29 1,1'-Biphenyl, 2,2',3,4,5'-... Concentration Rank 19

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 19.00 | 7.10 ng/ul | 516351 | Phenanthrene-d10 | 16.99 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1,1'-Biphenyl, 2,2',3,4,5'-penta... | 324 | C12H5Cl5 | 038380-02-8 | 99 |
| 2 | 1,1'-Biphenyl, 2,2',3,5,5'-penta... | 324 | C12H5Cl5 | 052663-61-3 | 99 |
| 3 | 1,1'-Biphenyl, 2,2',3',4,5-Penta... | 324 | C12H5Cl5 | 041464-51-1 | 99 |
| 4 | 1,1'-Biphenyl, 2,2',4,4',5-penta... | 324 | C12H5Cl5 | 038380-01-7 | 98 |
| 5 | 1,1'-Biphenyl, 2,2',3,5,5'-penta... | 324 | C12H5Cl5 | 052663-61-3 | 98 |



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Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
Operator : TP
Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** C01M4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
Quant Title : SVOA CALIBRATION

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

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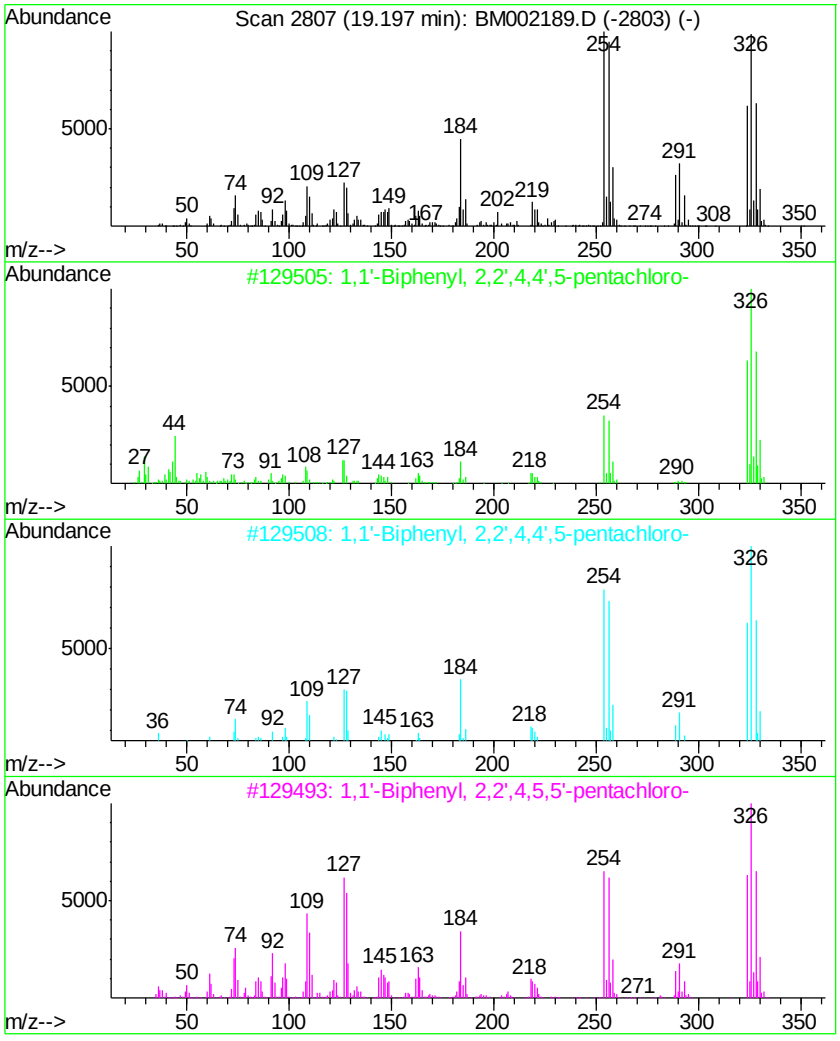
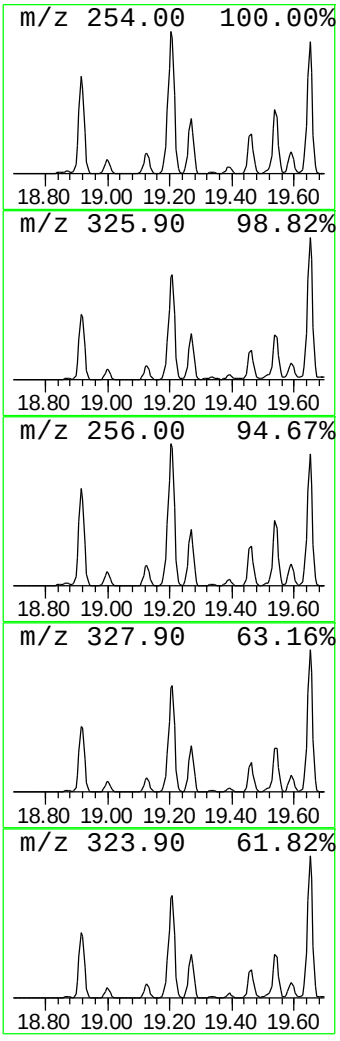
Peak Number 30 1,1'-Biphenyl, 2,2',4,4',5-... Concentration Rank 16

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 19.20 | 8.19 ng/ul | 4563940 | Chrysene-d12 | 21.20 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1,1'-Biphenyl, 2,2',4,4',5-penta... | 324 | C12H5Cl5 | 038380-01-7 | 98 |
| 2 | 1,1'-Biphenyl, 2,2',4,4',5-penta... | 324 | C12H5Cl5 | 038380-01-7 | 96 |
| 3 | 1,1'-Biphenyl, 2,2',4,5,5'-penta... | 324 | C12H5Cl5 | 037680-73-2 | 95 |
| 4 | 1,1'-Biphenyl, 2,3,3',4',6-penta... | 324 | C12H5Cl5 | 038380-03-9 | 95 |
| 5 | 1,1'-Biphenyl, 2,3,3',4,6-Pentac... | 324 | C12H5Cl5 | 074472-35-8 | 95 |



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080215\
Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
Operator : TP
Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** C01M4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
Quant Title : SVOA CALIBRATION

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

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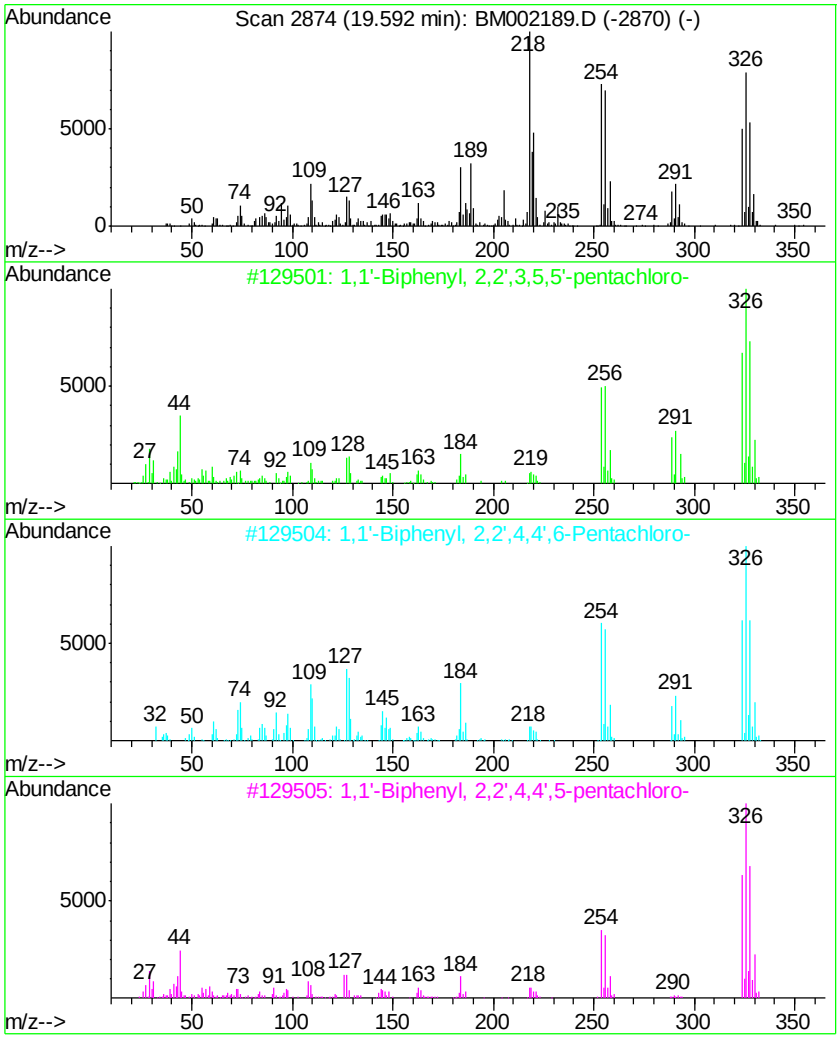
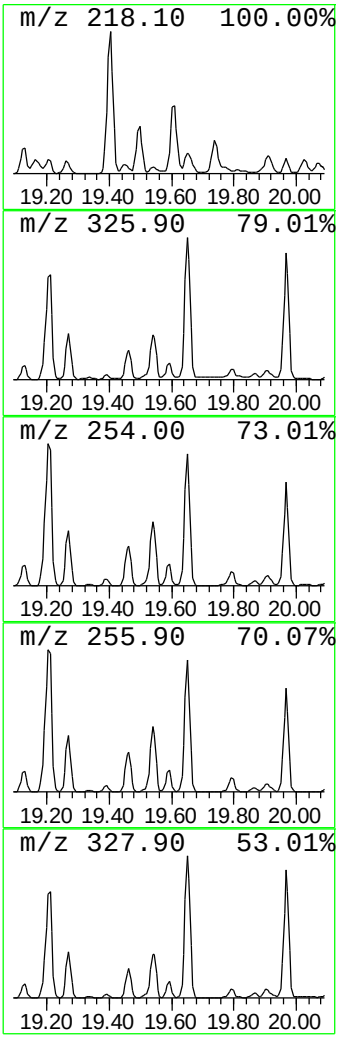
Peak Number 33 1,1'-Biphenyl, 2,2',3,5,5'-... Concentration Rank 51

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 19.59 | 2.45 ng/ul | 1362380 | Chrysene-d12 | 21.20 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1,1'-Biphenyl, 2,2',3,5,5'-penta... | 324 | C12H5Cl5 | 052663-61-3 | 97 |
| 2 | 1,1'-Biphenyl, 2,2',4,4',6-Penta... | 324 | C12H5Cl5 | 039485-83-1 | 91 |
| 3 | 1,1'-Biphenyl, 2,2',4,4',5-penta... | 324 | C12H5Cl5 | 038380-01-7 | 90 |
| 4 | 1,1'-Biphenyl, 2,2',3,3',6-penta... | 324 | C12H5Cl5 | 052663-60-2 | 90 |
| 5 | 1,1'-Biphenyl, 2,3,3',4,6-Pentac... | 324 | C12H5Cl5 | 074472-35-8 | 90 |



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080215\
Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
Operator : TP
Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** C01M4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
Quant Title : SVOA CALIBRATION

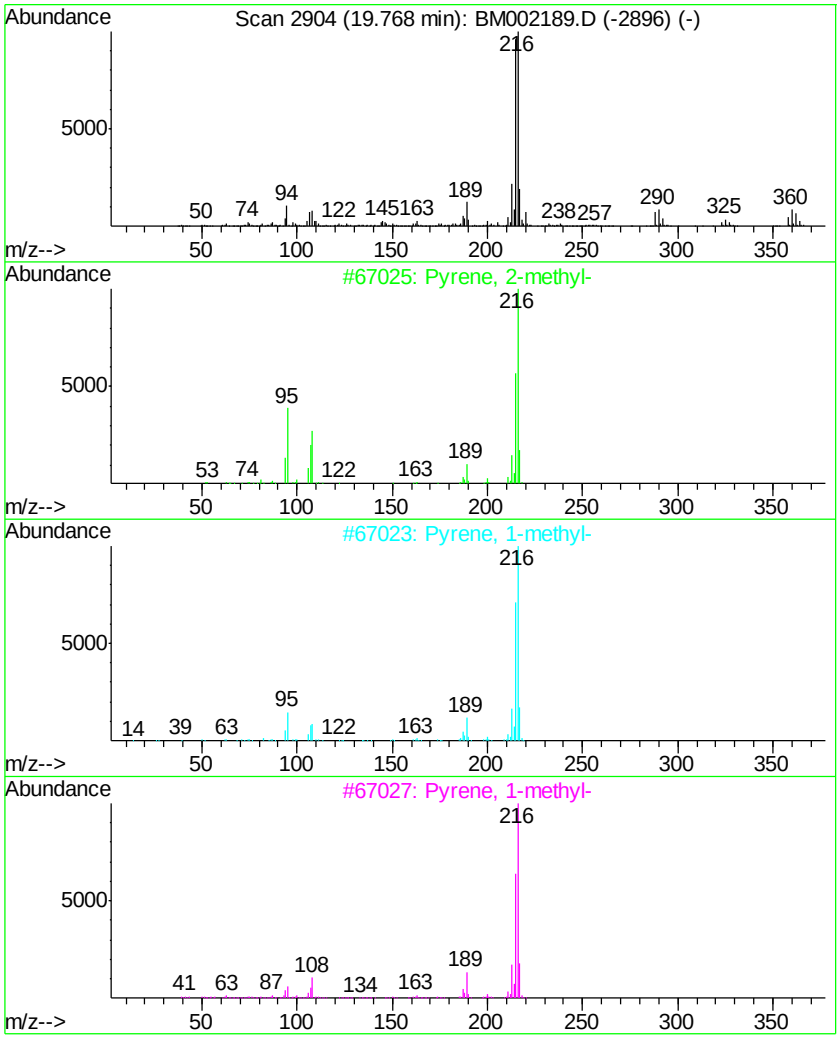
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

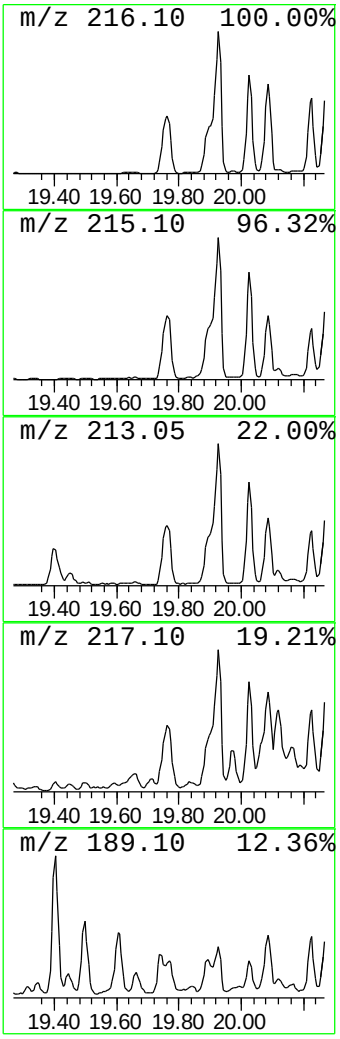
**********

Peak Number 35 Pyrene, 2-methyl- Concentration Rank 28

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 19.77 | 4.55 ng/ul | 2535840 | Chrysene-d12 | 21.20 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Pyrene, 2-methyl- | 216 | C17H12 | 003442-78-2 | 92 |
| 2 | Pyrene, 1-methyl- | 216 | C17H12 | 002381-21-7 | 92 |
| 3 | Pyrene, 1-methyl- | 216 | C17H12 | 002381-21-7 | 91 |
| 4 | Pyrene, 1-methyl- | 216 | C17H12 | 002381-21-7 | 90 |
| 5 | Fluoranthene, 2-methyl- | 216 | C17H12 | 033543-31-6 | 89 |



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080215\
Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
Operator : TP
Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** C01M4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
Quant Title : SVOA CALIBRATION

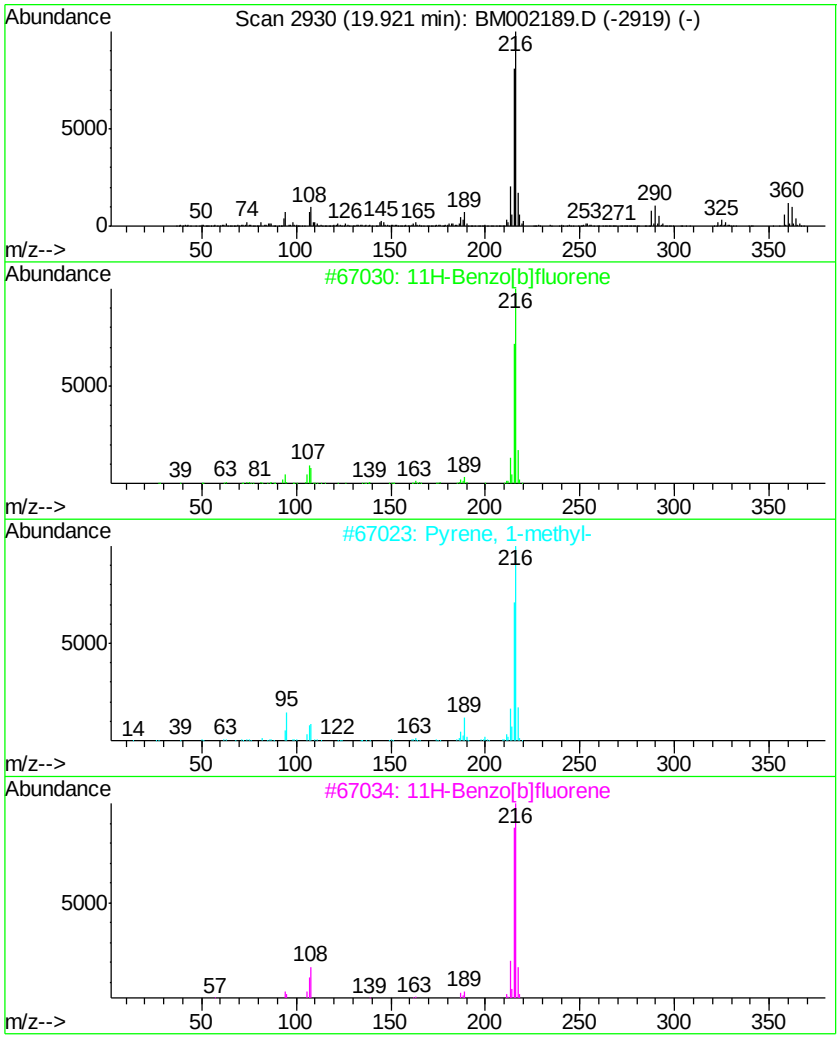
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

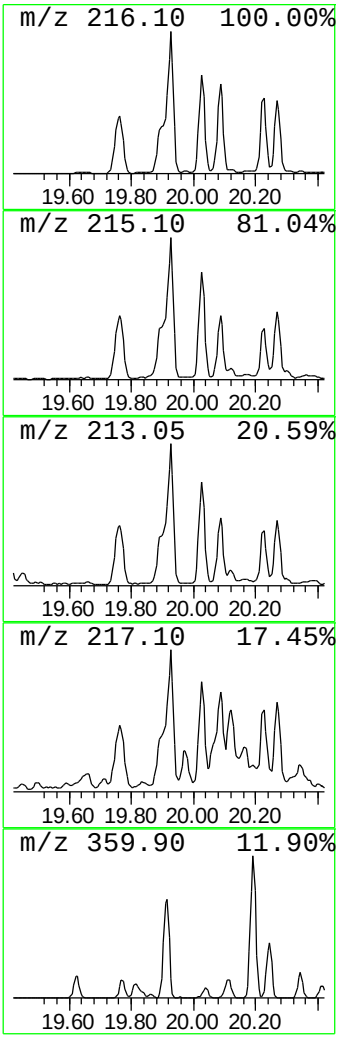
**********

Peak Number 36 11H-Benzo[b]fluorene Concentration Rank 15

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 19.92 | 8.22 ng/ul | 4581150 | Chrysene-d12 | 21.20 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 11H-Benzo[b]fluorene | 216 | C17H12 | 000243-17-4 | 78 |
| 2 | Pyrene, 1-methyl- | 216 | C17H12 | 002381-21-7 | 78 |
| 3 | 11H-Benzo[b]fluorene | 216 | C17H12 | 000243-17-4 | 70 |
| 4 | Pyrene, 1-methyl- | 216 | C17H12 | 002381-21-7 | 70 |
| 5 | Fluoranthene, 2-methyl- | 216 | C17H12 | 033543-31-6 | 70 |



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080215\
Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
Operator : TP
Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** C01M4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
Quant Title : SVOA CALIBRATION

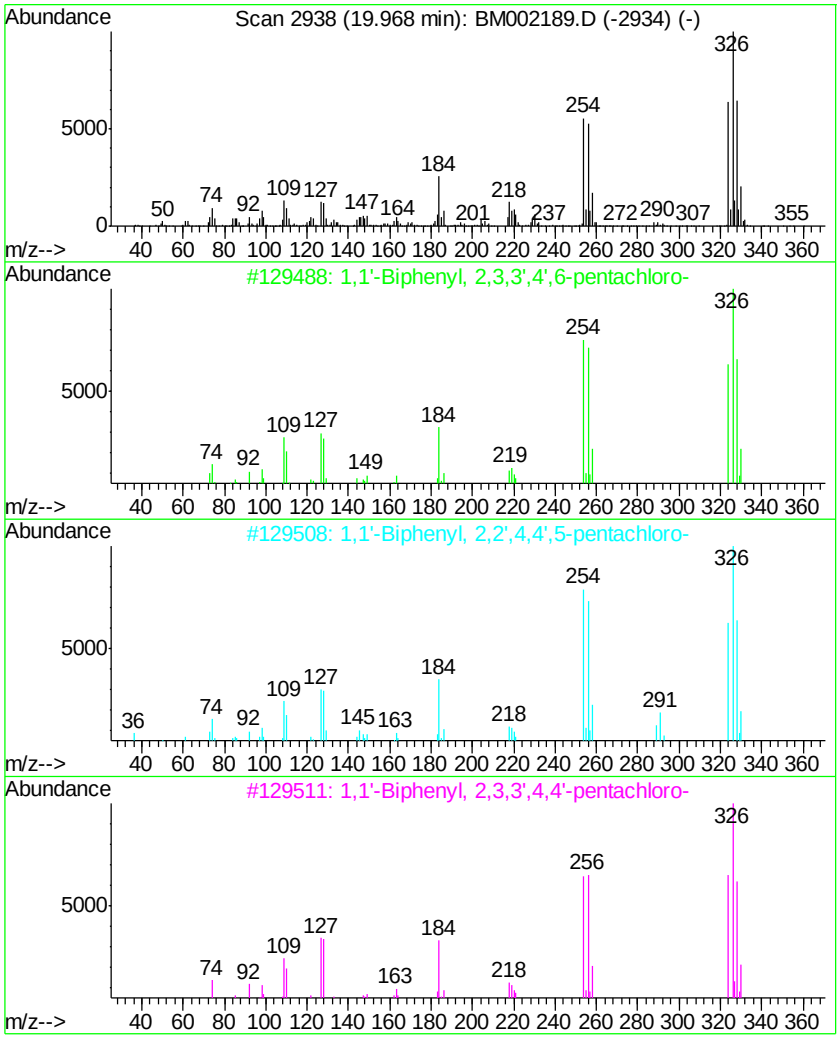
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

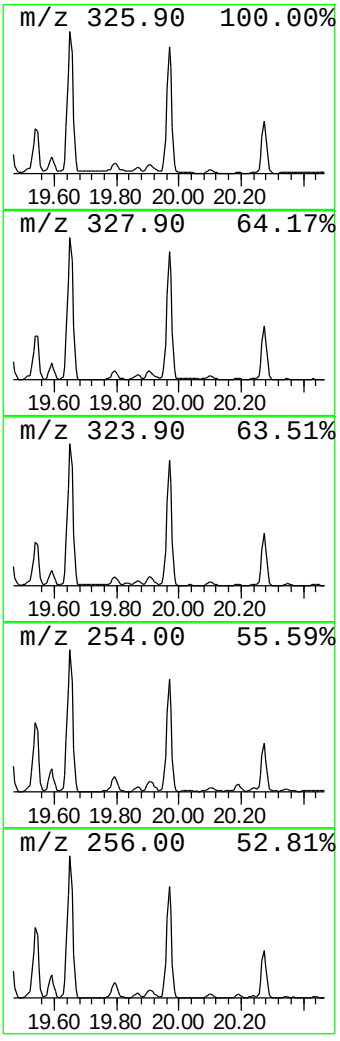
**********

Peak Number 37 1,1'-Biphenyl, 2,3,3',4',6-... Concentration Rank 26

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 19.97 | 4.93 ng/ul | 2746740 | Chrysene-d12 | 21.20 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1,1'-Biphenyl, 2,3,3',4',6-penta... | 324 | C12H5Cl5 | 038380-03-9 | 99 |
| 2 | 1,1'-Biphenyl, 2,2',4,4',5-penta... | 324 | C12H5Cl5 | 038380-01-7 | 99 |
| 3 | 1,1'-Biphenyl, 2,3,3',4,4'-penta... | 324 | C12H5Cl5 | 032598-14-4 | 99 |
| 4 | 1,1'-Biphenyl, 2,2',3,5,5'-penta... | 324 | C12H5Cl5 | 052663-61-3 | 96 |
| 5 | 1,1'-Biphenyl, 2,3,4,5,6-pentach... | 324 | C12H5Cl5 | 018259-05-7 | 96 |



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080215\
Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
Operator : TP
Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** C01M4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
Quant Title : SVOA CALIBRATION

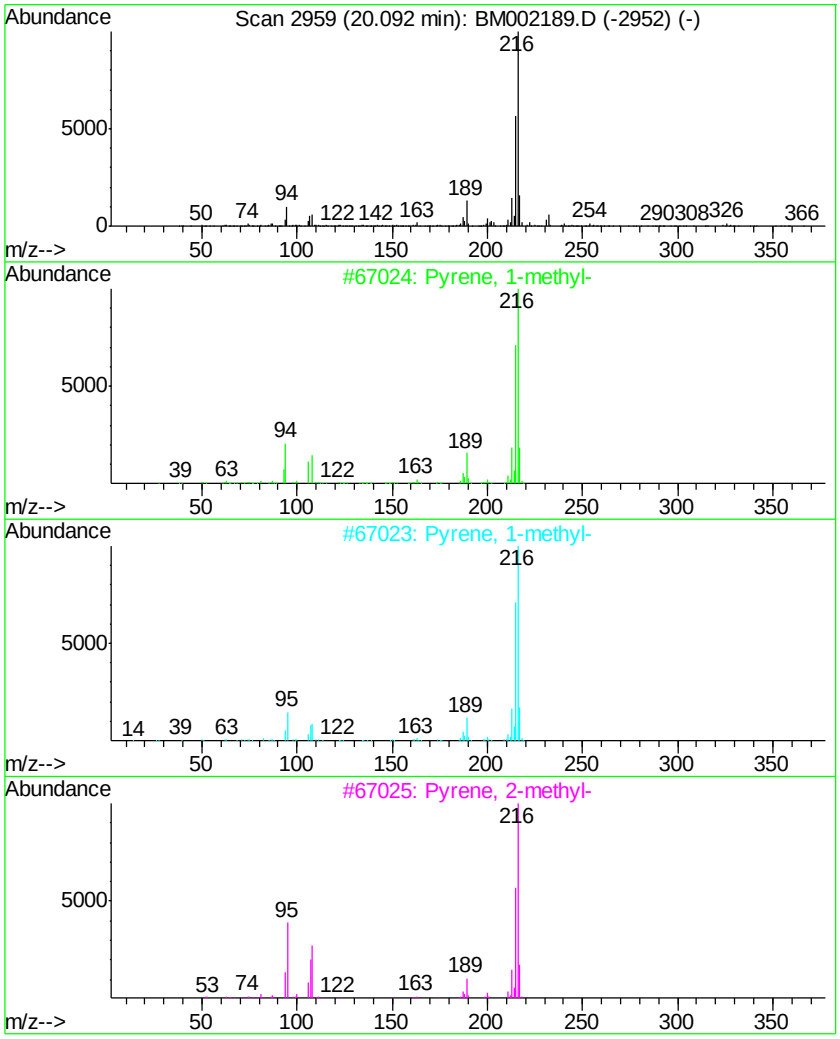
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

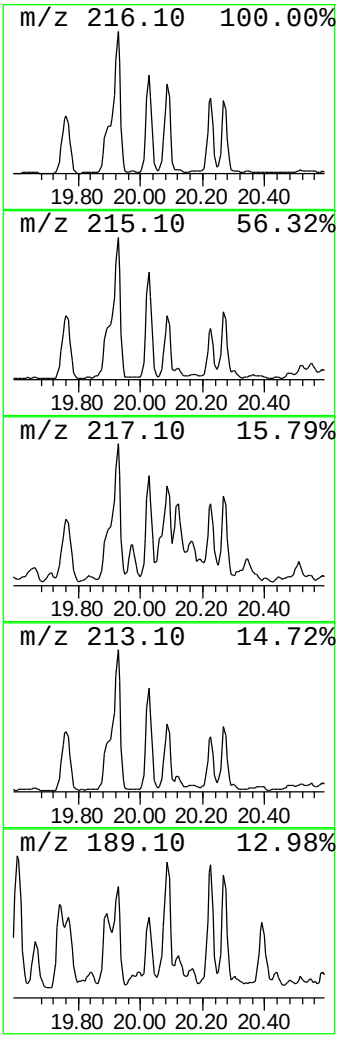
**********

Peak Number 39 Pyrene, 1-methyl- Concentration Rank 43

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 20.09 | 2.96 ng/ul | 1647370 | Chrysene-d12 | 21.20 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Pyrene, 1-methyl- | 216 | C17H12 | 002381-21-7 | 95 |
| 2 | Pyrene, 1-methyl- | 216 | C17H12 | 002381-21-7 | 95 |
| 3 | Pyrene, 2-methyl- | 216 | C17H12 | 003442-78-2 | 93 |
| 4 | Fluoranthene, 2-methyl- | 216 | C17H12 | 033543-31-6 | 90 |
| 5 | Pyrene, 2-methyl- | 216 | C17H12 | 003442-78-2 | 89 |



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080215\
Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
Operator : TP
Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** C01M4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
Quant Title : SVOA CALIBRATION

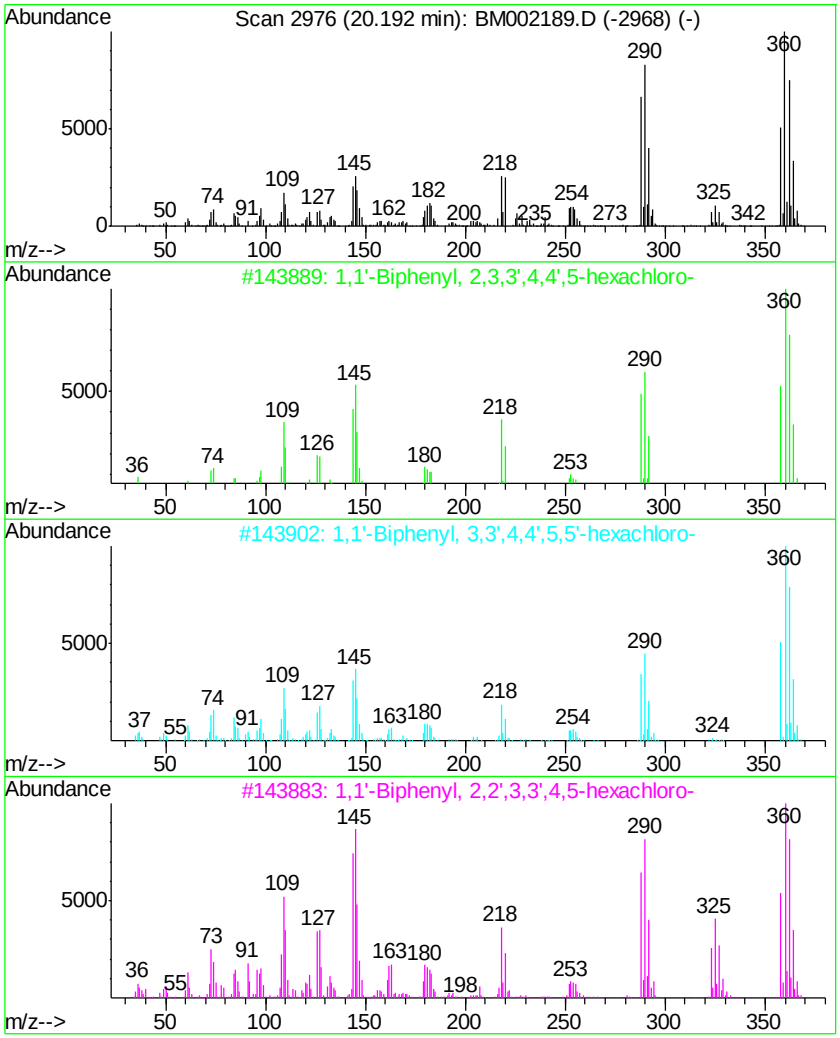
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

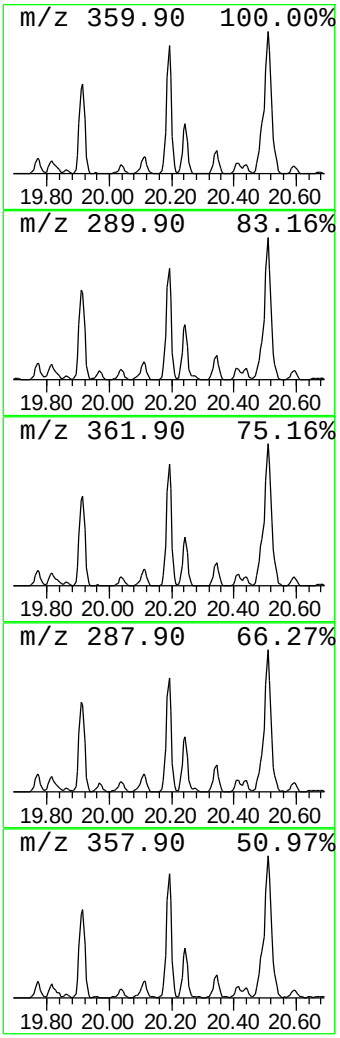
**********

Peak Number 40 1,1'-Biphenyl, 2,3,3',4,4',... Concentration Rank 39

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 20.19 | 3.34 ng/ul | 1861960 | Chrysene-d12 | 21.20 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1,1'-Biphenyl, 2,3,3',4,4',5-hex... | 358 | C12H4Cl6 | 038380-08-4 | 99 |
| 2 | 1,1'-Biphenyl, 3,3',4,4',5,5'-he... | 358 | C12H4Cl6 | 032774-16-6 | 99 |
| 3 | 1,1'-Biphenyl, 2,2',3,3',4,5-hex... | 358 | C12H4Cl6 | 055215-18-4 | 96 |
| 4 | 1,1'-Biphenyl, 2,2',3,4,5,6'-Hex... | 358 | C12H4Cl6 | 068194-15-0 | 95 |
| 5 | 1,1'-Biphenyl, 2,2',3,3',5,5'-He... | 358 | C12H4Cl6 | 035694-04-3 | 95 |



# Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080215\
Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
Operator : TP
Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** C01M4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
Quant Title : SVOA CALIBRATION

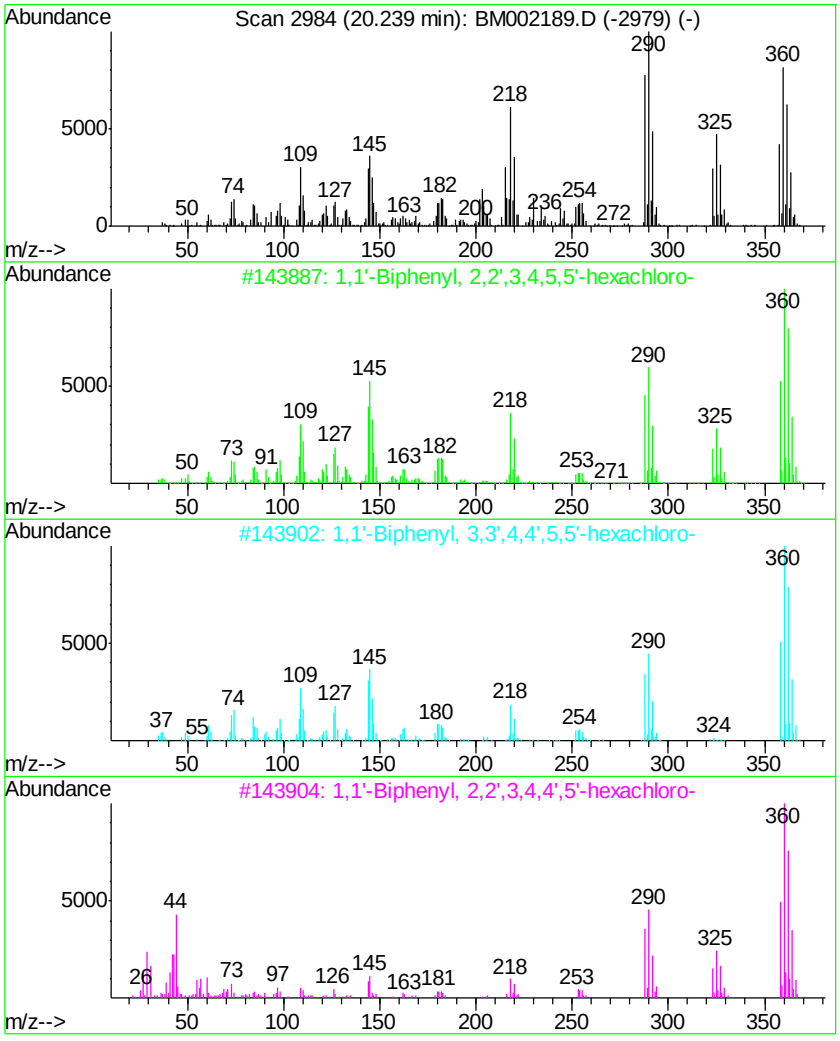
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

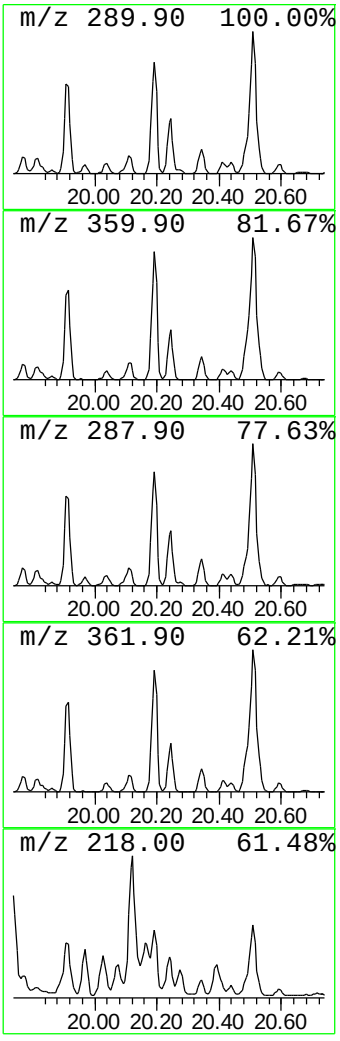
**********

Peak Number 41 1,1'-Biphenyl, 2,2',3,4,5,5... Concentration Rank 41

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 20.24 | 3.07 ng/ul | 1707480 | Chrysene-d12 | 21.20 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1,1'-Biphenyl, 2,2',3,4,5,5'-hex... | 358 | C12H4Cl6 | 052712-04-6 | 99 |
| 2 | 1,1'-Biphenyl, 3,3',4,4',5,5'-he... | 358 | C12H4Cl6 | 032774-16-6 | 99 |
| 3 | 1,1'-Biphenyl, 2,2',3,4,4',5'-he... | 358 | C12H4Cl6 | 035065-28-2 | 99 |
| 4 | 1,1'-Biphenyl, 2,2',3,4,5,6'-Hex... | 358 | C12H4Cl6 | 068194-15-0 | 99 |
| 5 | 1,1'-Biphenyl, 2,2',3,4,4',6-Hex... | 358 | C12H4Cl6 | 056030-56-9 | 99 |



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080215\
Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
Operator : TP
Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** C01M4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
Quant Title : SVOA CALIBRATION

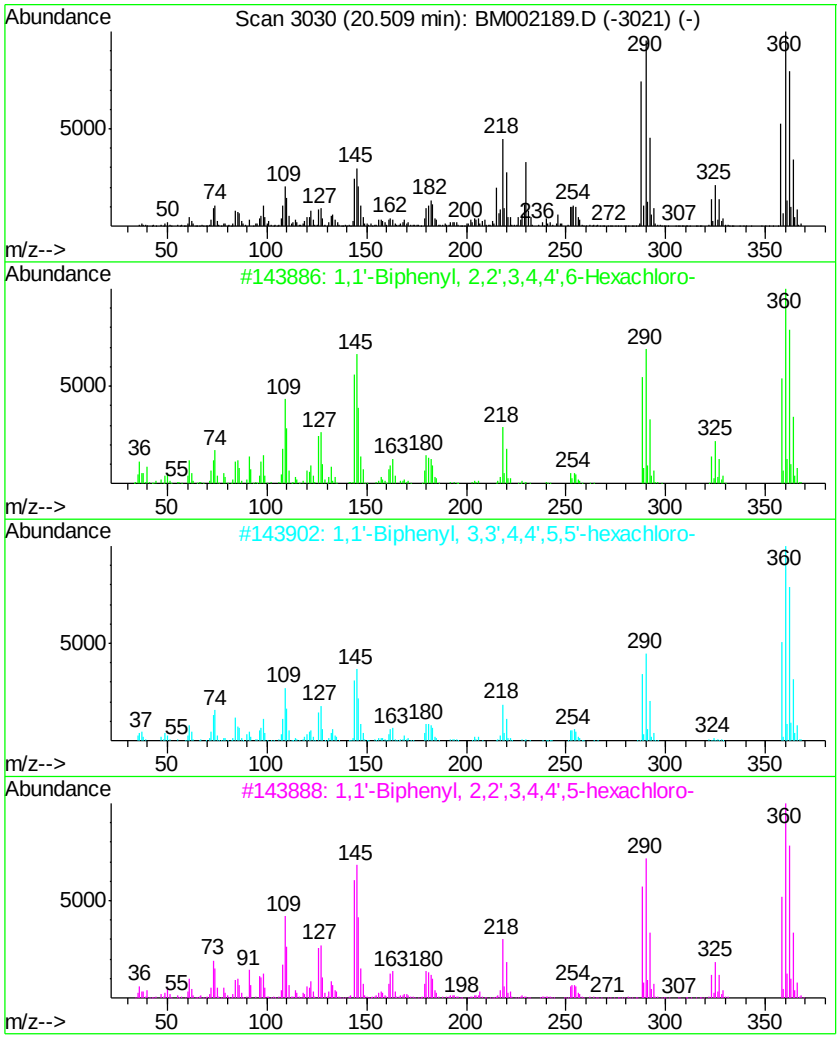
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

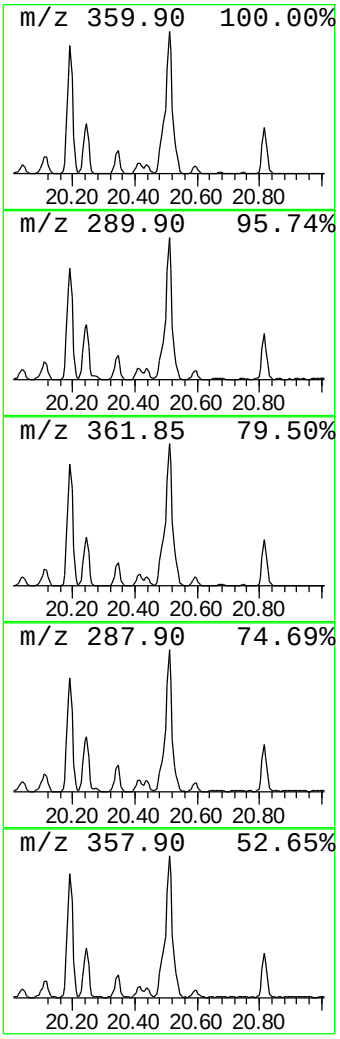
**********

Peak Number 43 1,1'-Biphenyl, 2,2',3,4,4',... Concentration Rank 21

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 20.51 | 6.33 ng/ul | 3528070 | Chrysene-d12 | 21.20 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1,1'-Biphenyl, 2,2',3,4,4',6-Hex... | 358 | C12H4Cl6 | 056030-56-9 | 99 |
| 2 | 1,1'-Biphenyl, 3,3',4,4',5,5'-he... | 358 | C12H4Cl6 | 032774-16-6 | 99 |
| 3 | 1,1'-Biphenyl, 2,2',3,4,4',5-hex... | 358 | C12H4Cl6 | 035694-06-5 | 99 |
| 4 | 1,1'-Biphenyl, 2,3,3',4,4',5-hex... | 358 | C12H4Cl6 | 038380-08-4 | 99 |
| 5 | 1,1'-Biphenyl, 2,2',4,4',6,6'-he... | 358 | C12H4Cl6 | 033979-03-2 | 98 |



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080215\
Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
Operator : TP
Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** C01M4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
Quant Title : SVOA CALIBRATION

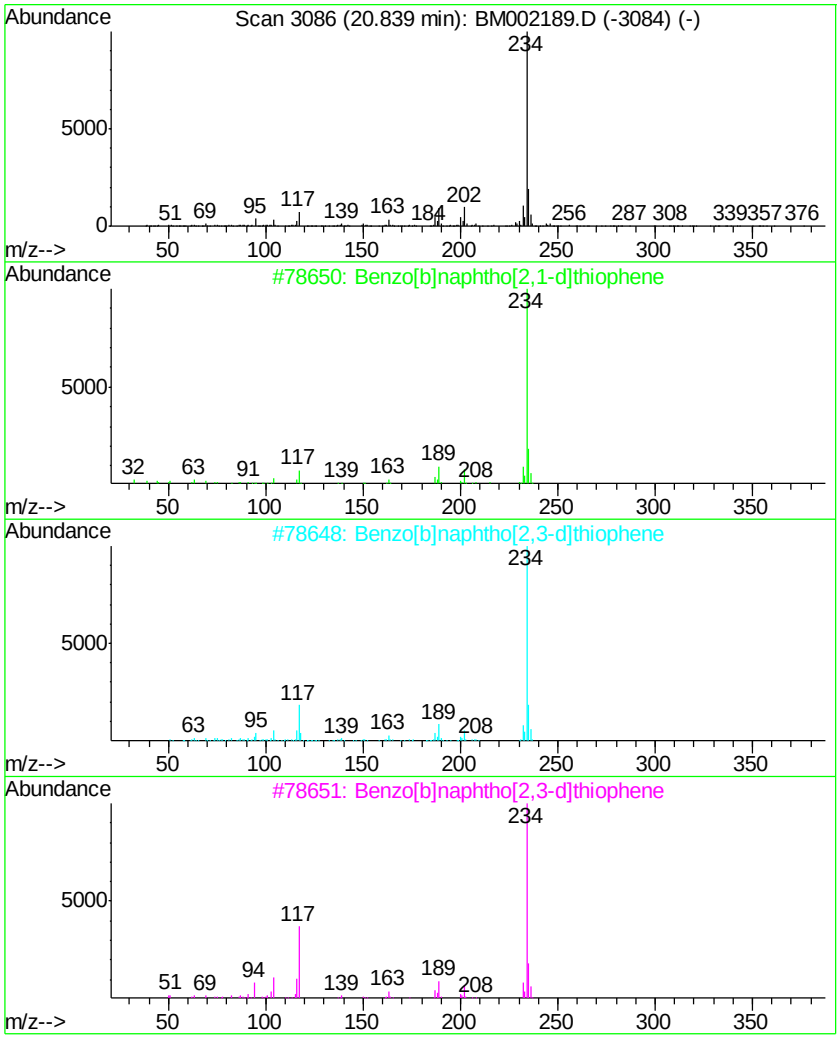
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

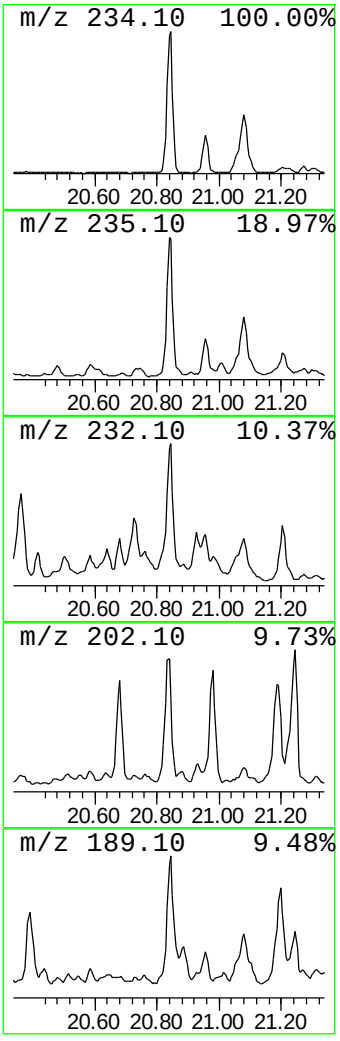
**********

Peak Number 44 Benzo[b]naphtho[2,1-d]thiop... Concentration Rank 42

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 20.84 | 3.05 ng/ul | 1696950 | Chrysene-d12 | 21.20 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzo[b]naphtho[2,1-d]thiophene | 234 | C16H10S | 000239-35-0 | 95 |
| 2 | Benzo[b]naphtho[2,3-d]thiophene | 234 | C16H10S | 000243-46-9 | 94 |
| 3 | Benzo[b]naphtho[2,3-d]thiophene | 234 | C16H10S | 000243-46-9 | 94 |
| 4 | Benzo[b]naphtho[2,1-d]thiophene | 234 | C16H10S | 000239-35-0 | 94 |
| 5 | Benzo[b]naphtho[2,1-d]thiophene | 234 | C16H10S | 000239-35-0 | 93 |



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080215\
Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
Operator : TP
Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** C01M4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
Quant Title : SVOA CALIBRATION

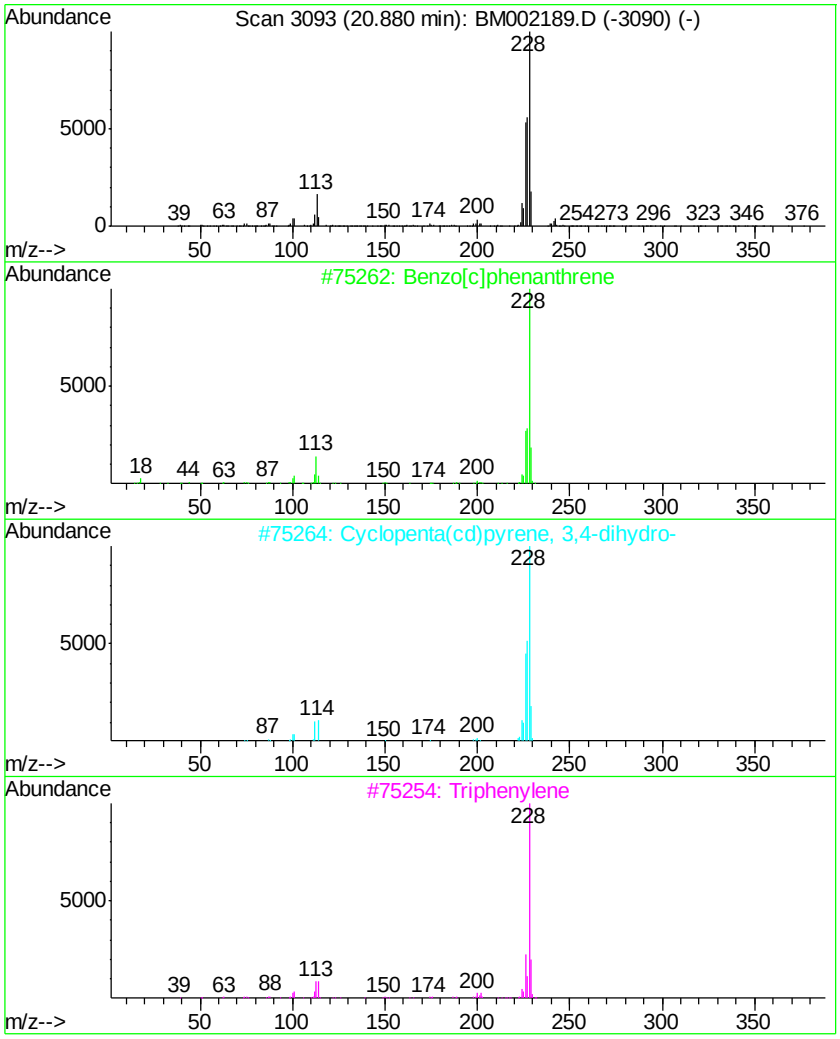
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

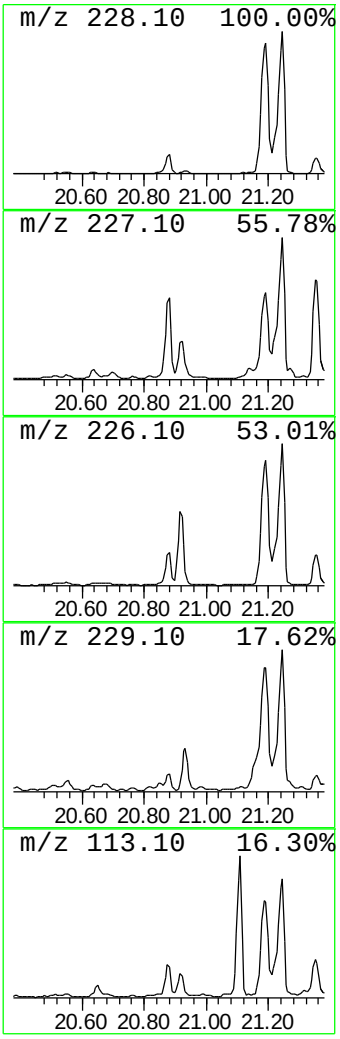
**********

Peak Number 45 Benzo[c]phenanthrene Concentration Rank 50

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 20.88 | 2.52 ng/ul | 1403940 | Chrysene-d12 | 21.20 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzo[c]phenanthrene | 228 | C18H12 | 000195-19-7 | 64 |
| 2 | Cyclopenta(cd)pyrene, 3,4-dihydro- | 228 | C18H12 | 025732-74-5 | 64 |
| 3 | Triphenylene | 228 | C18H12 | 000217-59-4 | 49 |
| 4 | Triphenylene | 228 | C18H12 | 000217-59-4 | 49 |
| 5 | Triphenylene | 228 | C18H12 | 000217-59-4 | 46 |



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080215\
Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
Operator : TP
Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** C01M4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
Quant Title : SVOA CALIBRATION

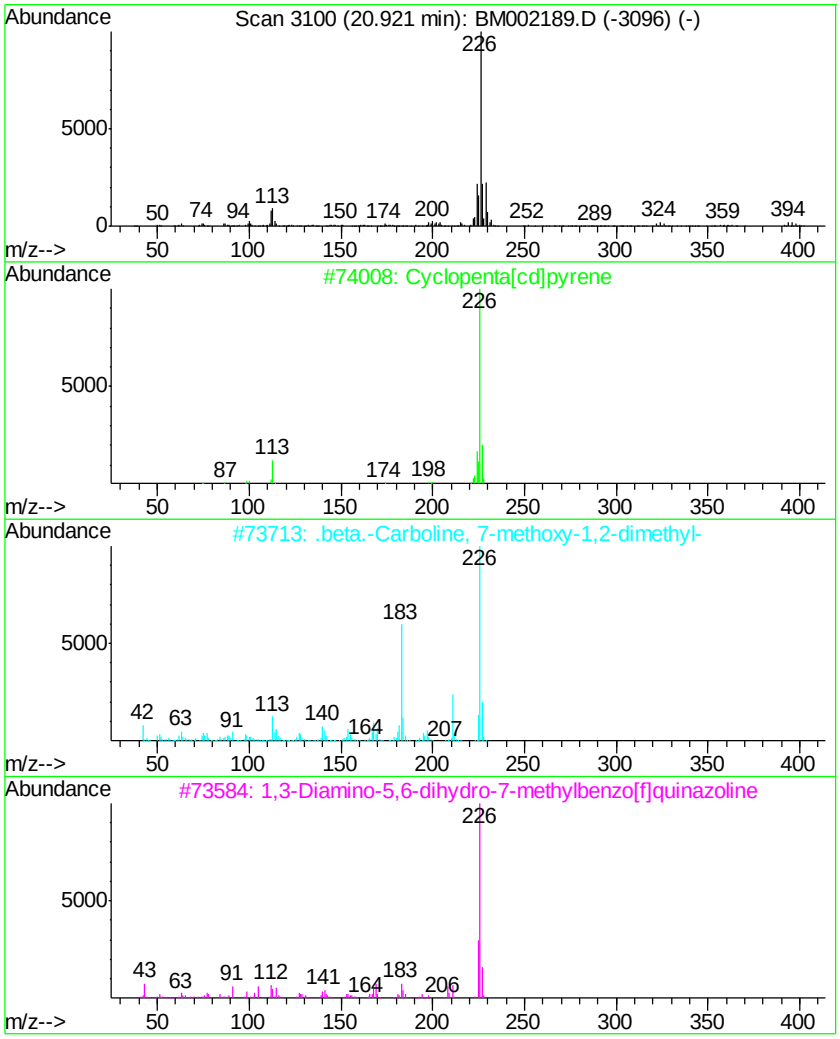
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

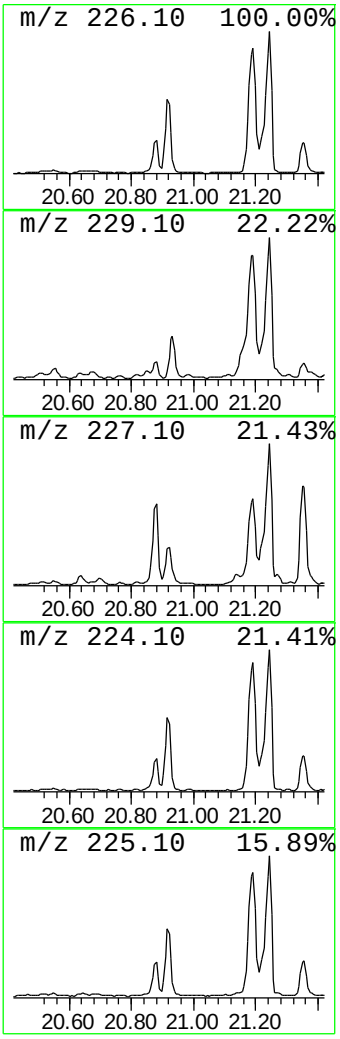
**********

Peak Number 46 Cyclopenta[cd]pyrene Concentration Rank 44

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 20.92 | 2.95 ng/ul | 1644190 | Chrysene-d12 | 21.20 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Cyclopenta[cd]pyrene | 226 | C18H10 | 027208-37-3 | 64 |
| 2 | .beta.-Carboline, 7-methoxy-1,2-... | 226 | C14H14N2O | 006519-18-2 | 52 |
| 3 | 1,3-Diamino-5,6-dihydro-7-methyl... | 226 | C13H14N4 | 037436-37-6 | 47 |
| 4 | 1-Bromo-2,4-dichlorobenzene | 224 | C6H3BrCl2 | 001193-72-2 | 47 |
| 5 | 1,3-Diamino-5,6-dihydro-7-methyl... | 226 | C13H14N4 | 037436-37-6 | 43 |



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080215\
Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
Operator : TP
Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** C01M4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
Quant Title : SVOA CALIBRATION

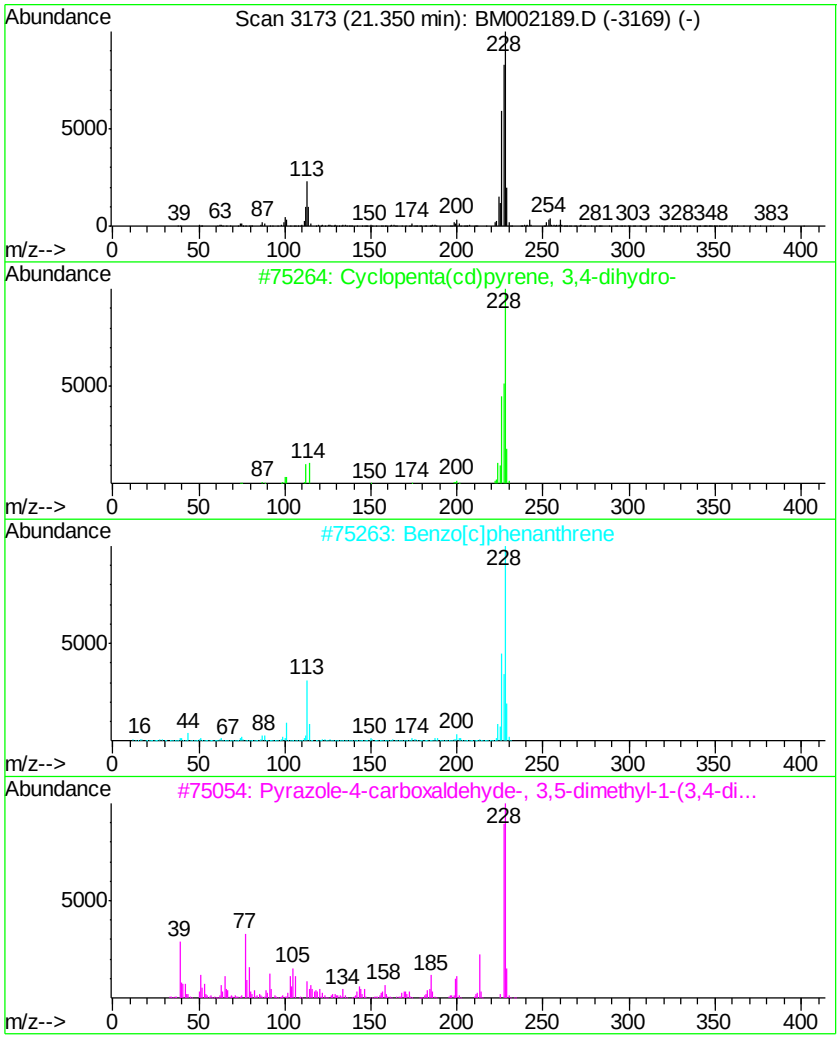
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

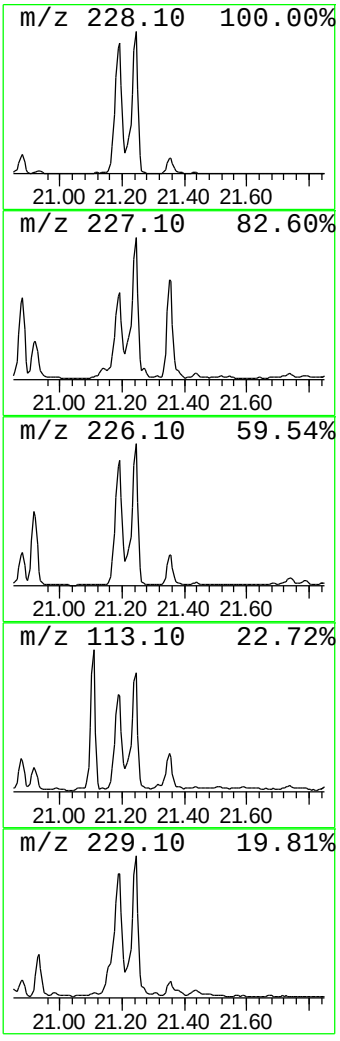
**********

Peak Number 47 Cyclopenta(cd)pyrene, 3,4-d... Concentration Rank 49

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 21.35 | 2.64 ng/ul | 1472080 | Chrysene-d12 | 21.20 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Cyclopenta(cd)pyrene, 3,4-dihydro- | 228 | C18H12 | 025732-74-5 | 96 |
| 2 | Benzo[c]phenanthrene | 228 | C18H12 | 000195-19-7 | 55 |
| 3 | Pyrazole-4-carboxaldehyde-, 3,5-... | 228 | C14H16N2O | 1000272-54-8 | 50 |
| 4 | Diphenylvinylphosphine oxide | 228 | C14H13OP | 002096-78-8 | 49 |
| 5 | Triphenylene | 228 | C18H12 | 000217-59-4 | 45 |



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080215\
Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
Operator : TP
Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** C01M4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
Quant Title : SVOA CALIBRATION

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

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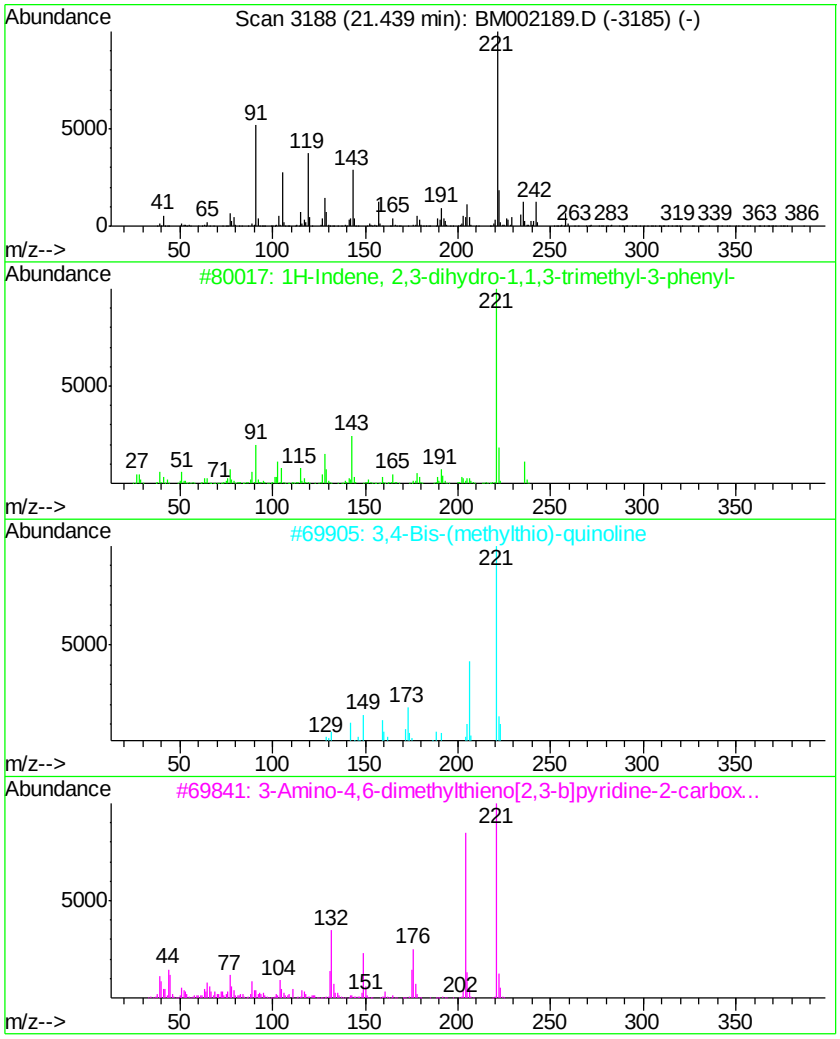
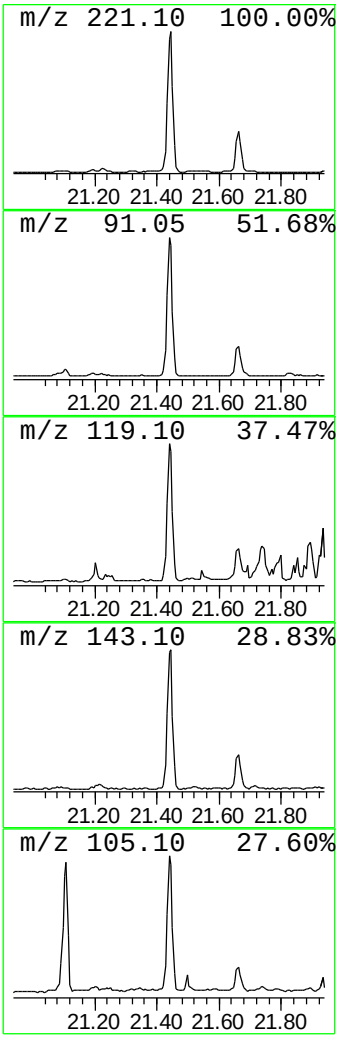
Peak Number 48 1H-Indene, 2,3-dihydro-1,1,... Concentration Rank 34

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 21.44 | 3.74 ng/ul | 2082080 | Chrysene-d12 | 21.20 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1H-Indene, 2,3-dihydro-1,1,3-tri... | 236 | C18H20 | 003910-35-8 | 52 |
| 2 | 3,4-Bis-(methylthio)-quinoline | 221 | C11H11NS2 | 074579-34-3 | 50 |
| 3 | 3-Amino-4,6-dimethylthieno[2,3-b... | 221 | C10H11N3OS | 067795-42-0 | 47 |
| 4 | 1,3-Cyclohexanedione, 2-[2-[4-(4... | 401 | C22H28FN3O3 | 339310-26-8 | 43 |
| 5 | Benz[j]isoquinoline, 3,6,8-trime... | 221 | C16H15N | 061171-16-2 | 43 |



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080215\
Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
Operator : TP
Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** C01M4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
Quant Title : SVOA CALIBRATION

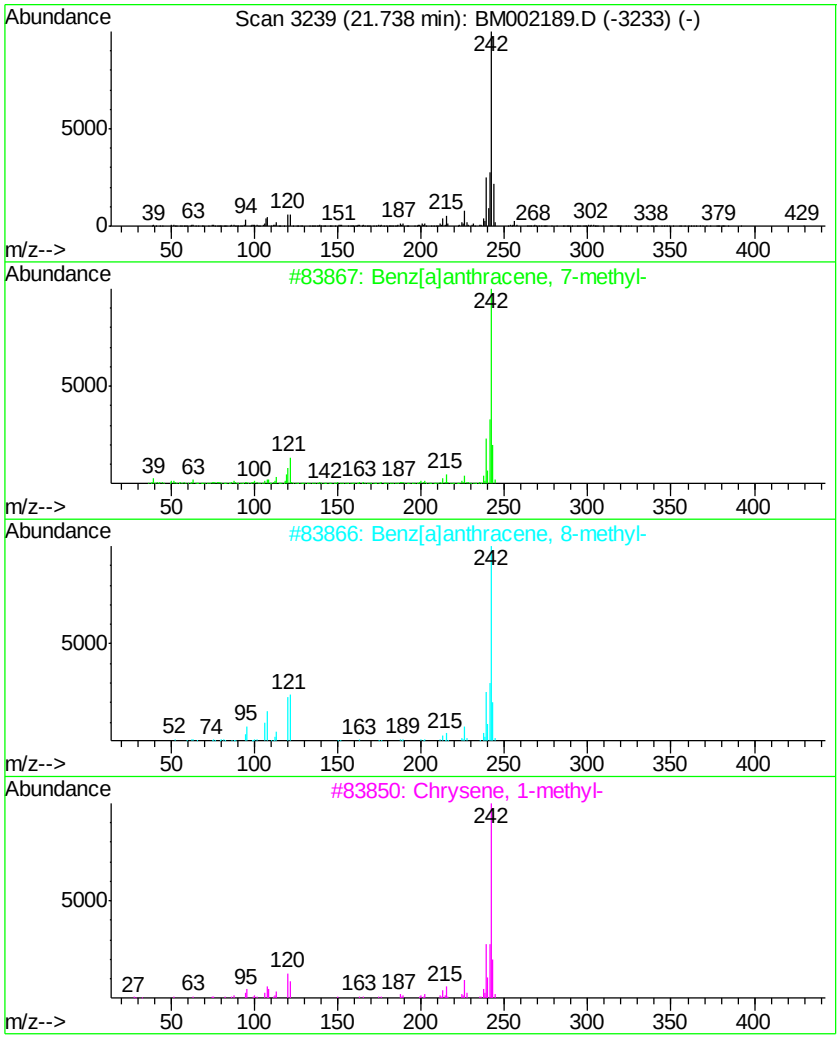
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

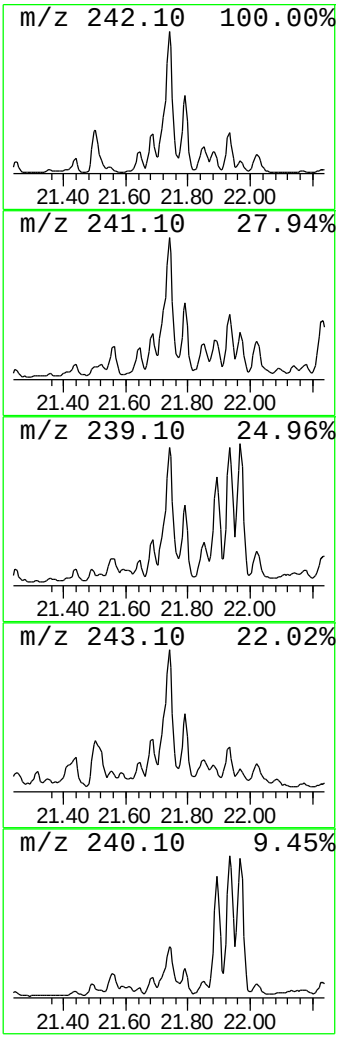
**********

Peak Number 49 Benz[a]anthracene, 7-methyl- Concentration Rank 47

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 21.74 | 2.76 ng/ul | 1535480 | Chrysene-d12 | 21.20 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benz[a]anthracene, 7-methyl- | 242 | C19H14 | 002541-69-7 | 95 |
| 2 | Benz[a]anthracene, 8-methyl- | 242 | C19H14 | 002381-31-9 | 95 |
| 3 | Chrysene, 1-methyl- | 242 | C19H14 | 003351-28-8 | 95 |
| 4 | Chrysene, 2-methyl- | 242 | C19H14 | 003351-32-4 | 94 |
| 5 | Triphenylene, 2-methyl- | 242 | C19H14 | 001705-84-6 | 93 |



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080215\
Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
Operator : TP
Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** C01M4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
Quant Title : SVOA CALIBRATION

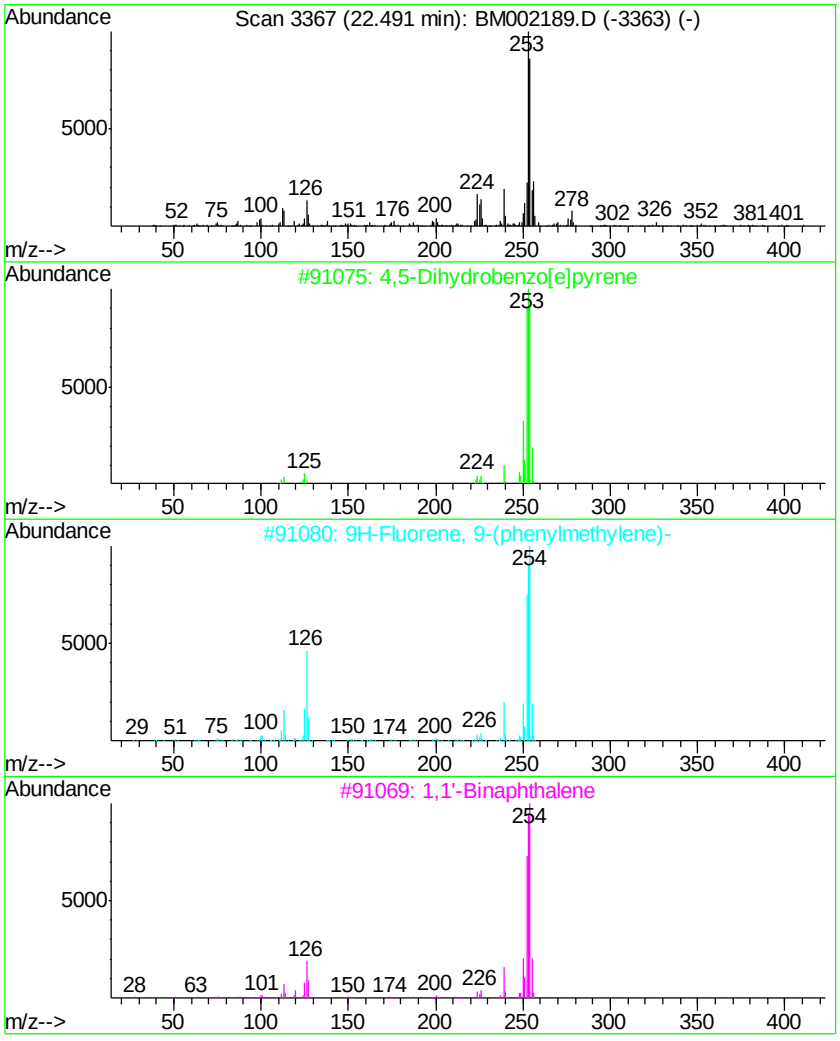
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

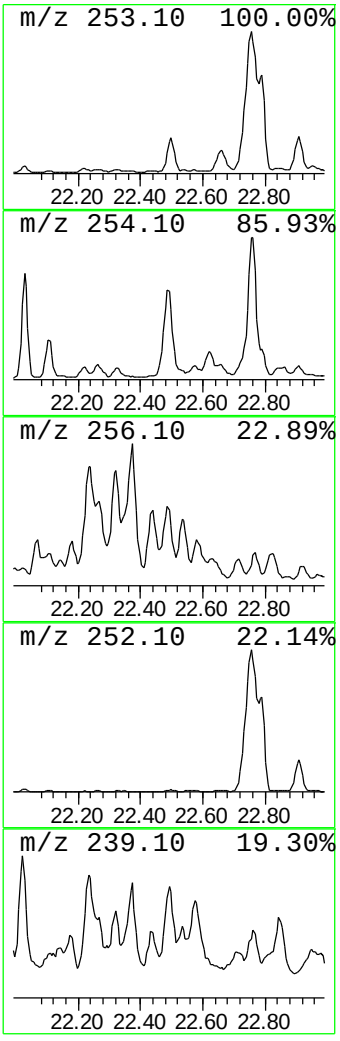
**********

Peak Number 50 4,5-Dihydrobenzo[e]pyrene Concentration Rank 13

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 22.49 | 9.97 ng/ul | 590779 | Perylene-d12 | 23.40 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 4,5-Dihydrobenzo[e]pyrene | 254 | C20H14 | 095676-42-9 | 81 |
| 2 | 9H-Fluorene, 9-(phenylmethylene)- | 254 | C20H14 | 001836-87-9 | 76 |
| 3 | 1,1'-Binaphthalene | 254 | C20H14 | 000604-53-5 | 64 |
| 4 | 9H-Fluorene, 9-(phenylmethylene)- | 254 | C20H14 | 001836-87-9 | 64 |
| 5 | Anthracene, 9-phenyl- | 254 | C20H14 | 000602-55-1 | 60 |



Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080215\
Data File : BM002189.D
Acq On : 03 Aug 2015 03:20
Operator : TP
Sample : G3095-20
Misc :
ALS Vial : 27 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** C01M4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.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 |
| Dibenzofuran, 4-m... | 15.57 | 3.9 | ng/ul | 262891 | 3 | 14.23 | 1332980 | 20.0 |
| 2,4,6-Cycloheptat... | 15.73 | 3.4 | ng/ul | 247026 | 4 | 16.99 | 1454460 | 20.0 |
| 9H-Fluorene, 2-me... | 16.29 | 6.3 | ng/ul | 460605 | 4 | 16.99 | 1454460 | 20.0 |
| 9H-Fluorene, 1-me... | 16.34 | 3.5 | ng/ul | 256106 | 4 | 16.99 | 1454460 | 20.0 |
| Phenol, 4-(2-phen... | 16.56 | 4.1 | ng/ul | 295614 | 4 | 16.99 | 1454460 | 20.0 |
| 9H-Fluoren-9-one | 16.64 | 5.8 | ng/ul | 419443 | 4 | 16.99 | 1454460 | 20.0 |
| Anthracene, 1,2,3... | 16.73 | 5.4 | ng/ul | 391564 | 4 | 16.99 | 1454460 | 20.0 |
| Dibenzothiophene | 16.80 | 8.4 | ng/ul | 607919 | 4 | 16.99 | 1454460 | 20.0 |
| Acridine | 17.18 | 3.3 | ng/ul | 242829 | 4 | 16.99 | 1454460 | 20.0 |
| Anthracene, 9-eth... | 17.50 | 4.0 | ng/ul | 289270 | 4 | 16.99 | 1454460 | 20.0 |
| Dibenzothiophene,... | 17.56 | 8.1 | ng/ul | 589162 | 4 | 16.99 | 1454460 | 20.0 |
| Dibenzothiophene,... | 17.70 | 5.2 | ng/ul | 377118 | 4 | 16.99 | 1454460 | 20.0 |
| Phenanthrene, 1-m... | 17.86 | 25.3 | ng/ul | 1838600 | 4 | 16.99 | 1454460 | 20.0 |
| 1,1'-Biphenyl, 2,... | 18.06 | 75.3 | ng/ul | 5478970 | 4 | 16.99 | 1454460 | 20.0 |
| 3-Methylcarbazole | 18.17 | 3.4 | ng/ul | 249032 | 4 | 16.99 | 1454460 | 20.0 |
| 2-Phenylnaphthalene | 18.37 | 14.2 | ng/ul | 1030180 | 4 | 16.99 | 1454460 | 20.0 |
| 9,10-Anthracenedione | 18.42 | 10.6 | ng/ul | 771413 | 4 | 16.99 | 1454460 | 20.0 |
| 1,1'-Biphenyl, 2,... | 18.52 | 4.3 | ng/ul | 309102 | 4 | 16.99 | 1454460 | 20.0 |
| Phenanthrene, 4,5... | 18.62 | 6.3 | ng/ul | 461515 | 4 | 16.99 | 1454460 | 20.0 |
| di-p-Tolylacetylene | 18.67 | 4.2 | ng/ul | 307005 | 4 | 16.99 | 1454460 | 20.0 |
| Phenanthrene, 2,5... | 18.70 | 4.9 | ng/ul | 353942 | 4 | 16.99 | 1454460 | 20.0 |
| 1,1'-Biphenyl, 2,... | 18.92 | 66.0 | ng/ul | 4797070 | 4 | 16.99 | 1454460 | 20.0 |
| 1,1'-Biphenyl, 2,... | 19.00 | 7.1 | ng/ul | 516351 | 4 | 16.99 | 1454460 | 20.0 |
| 1,1'-Biphenyl, 2,... | 19.20 | 8.2 | ng/ul | 4563940 | 5 | 21.20 | 11140000 | 20.0 |
| 1,1'-Biphenyl, 2,... | 19.59 | 2.5 | ng/ul | 1362380 | 5 | 21.20 | 11140000 | 20.0 |
| Pyrene, 2-methyl- | 19.77 | 4.5 | ng/ul | 2535840 | 5 | 21.20 | 11140000 | 20.0 |
| 11H-Benzo[b]fluorene | 19.92 | 8.2 | ng/ul | 4581150 | 5 | 21.20 | 11140000 | 20.0 |
| 1,1'-Biphenyl, 2,... | 19.97 | 4.9 | ng/ul | 2746740 | 5 | 21.20 | 11140000 | 20.0 |
| Pyrene, 1-methyl- | 20.09 | 3.0 | ng/ul | 1647370 | 5 | 21.20 | 11140000 | 20.0 |
| 1,1'-Biphenyl, 2,... | 20.19 | 3.3 | ng/ul | 1861960 | 5 | 21.20 | 11140000 | 20.0 |
| 1,1'-Biphenyl, 2,... | 20.24 | 3.1 | ng/ul | 1707480 | 5 | 21.20 | 11140000 | 20.0 |
| 1,1'-Biphenyl, 2,... | 20.51 | 6.3 | ng/ul | 3528070 | 5 | 21.20 | 11140000 | 20.0 |
| Benzo[b]naphtho[2... | 20.84 | 3.0 | ng/ul | 1696950 | 5 | 21.20 | 11140000 | 20.0 |
| Benzo[c]phenanthrene | 20.88 | 2.5 | ng/ul | 1403940 | 5 | 21.20 | 11140000 | 20.0 |
| Cyclopenta[cd]pyrene | 20.92 | 3.0 | ng/ul | 1644190 | 5 | 21.20 | 11140000 | 20.0 |
| Cyclopenta(cd)pyr... | 21.35 | 2.6 | ng/ul | 1472080 | 5 | 21.20 | 11140000 | 20.0 |
| 1H-Indene, 2,3-di... | 21.44 | 3.7 | ng/ul | 2082080 | 5 | 21.20 | 11140000 | 20.0 |
| Benz[a]anthracene... | 21.74 | 2.8 | ng/ul | 1535480 | 5 | 21.20 | 11140000 | 20.0 |
| 4,5-Dihydrobenzo[... | 22.49 | 10.0 | ng/ul | 590779 | 6 | 23.40 | 1184660 | 20.0 |