LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP081820\
Data File : BP003025.D
Acq On : 18 Aug 2020 19:40
Operator : CG/JU
Sample : L3696-06
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :** BNA_P
**ClientSampleId :** W-1

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % 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:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP073020.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

| peak # | R.T. min | first scan | max scan | last scan | PK TY | peak height | corr. area | corr. % max. | % of total |
|---|---|---|---|---|---|---|---|---|---|
| 1 | 2.952 | 7 | 11 | 34 | rVB | 2990557 | 3757404 | 25.09% | 3.406% |
| 2 | 5.293 | 402 | 409 | 420 | rBV | 1870445 | 2751817 | 18.37% | 2.494% |
| 3 | 6.858 | 668 | 675 | 688 | rVB | 1089495 | 1756397 | 11.73% | 1.592% |
| 4 | 7.340 | 752 | 757 | 763 | rVB | 121569 | 177028 | 1.18% | 0.160% |
| 5 | 7.681 | 809 | 815 | 824 | rVB | 1061369 | 1668752 | 11.14% | 1.512% |
| 6 | 7.805 | 830 | 836 | 845 | rVB2 | 132935 | 239071 | 1.60% | 0.217% |
| 7 | 8.058 | 873 | 879 | 889 | rVB | 88431 | 150433 | 1.00% | 0.136% |
| 8 | 8.828 | 996 | 1010 | 1016 | rBV | 2416081 | 4357270 | 29.09% | 3.949% |
| 9 | 9.369 | 1096 | 1102 | 1107 | rBV | 100022 | 160032 | 1.07% | 0.145% |
| 10 | 9.875 | 1181 | 1188 | 1194 | rVB2 | 301195 | 587962 | 3.93% | 0.533% |
| 11 | 9.952 | 1194 | 1201 | 1209 | rVB3 | 121492 | 240292 | 1.60% | 0.218% |
| 12 | 10.463 | 1279 | 1288 | 1294 | rBV | 1522793 | 2630954 | 17.57% | 2.385% |
| 13 | 10.522 | 1294 | 1298 | 1305 | rVV4 | 107760 | 238897 | 1.60% | 0.217% |
| 14 | 10.587 | 1305 | 1309 | 1317 | rVB | 126391 | 247506 | 1.65% | 0.224% |
| 15 | 10.663 | 1317 | 1322 | 1325 | rBV | 107065 | 165709 | 1.11% | 0.150% |
| 16 | 10.934 | 1357 | 1368 | 1373 | rBV3 | 135152 | 281380 | 1.88% | 0.255% |
| 17 | 11.387 | 1440 | 1445 | 1449 | rBV | 146552 | 215602 | 1.44% | 0.195% |
| 18 | 11.528 | 1462 | 1469 | 1473 | rBV2 | 194134 | 379082 | 2.53% | 0.344% |
| 19 | 11.581 | 1474 | 1478 | 1485 | rVB | 110274 | 178452 | 1.19% | 0.162% |
| 20 | 11.663 | 1485 | 1492 | 1497 | rBV | 338495 | 569342 | 3.80% | 0.516% |
| 21 | 12.022 | 1548 | 1553 | 1558 | rVB2 | 200012 | 331624 | 2.21% | 0.301% |
| 22 | 12.128 | 1563 | 1571 | 1582 | rVB | 853822 | 1533662 | 10.24% | 1.390% |
| 23 | 12.352 | 1603 | 1609 | 1613 | rBV | 529517 | 831296 | 5.55% | 0.753% |
| 24 | 12.393 | 1613 | 1616 | 1622 | rVB3 | 272037 | 423160 | 2.83% | 0.384% |
| 25 | 12.675 | 1661 | 1664 | 1674 | rVB7 | 140083 | 373427 | 2.49% | 0.338% |
| 26 | 12.887 | 1695 | 1700 | 1703 | rBV2 | 251598 | 369111 | 2.46% | 0.335% |
| 27 | 12.940 | 1703 | 1709 | 1716 | rVB | 6690212 | 10010110 | 66.83% | 9.073% |
| 28 | 13.116 | 1731 | 1739 | 1742 | rBV5 | 87172 | 192197 | 1.28% | 0.174% |
| 29 | 13.269 | 1760 | 1765 | 1770 | rVB3 | 255287 | 401787 | 2.68% | 0.364% |
| 30 | 13.346 | 1770 | 1778 | 1788 | rBV3 | 230139 | 608916 | 4.07% | 0.552% |
| 31 | 13.481 | 1796 | 1801 | 1812 | rBV2 | 478236 | 1301328 | 8.69% | 1.179% |
| 32 | 13.640 | 1819 | 1828 | 1833 | rBV | 741862 | 1499452 | 10.01% | 1.359% |
| 33 | 13.698 | 1833 | 1838 | 1843 | rVV2 | 611280 | 955714 | 6.38% | 0.866% |
| 34 | 13.781 | 1844 | 1852 | 1857 | rVV | 493510 | 985008 | 6.58% | 0.893% |

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP081820\
Data File : BP003025.D
Acq On : 18 Aug 2020 19:40
Operator : CG/JU
Sample : L3696-06
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :** BNA_P
**ClientSampleId :** W-1

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % 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:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP073020.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 35 | 13.840 | 1857 | 1862 | 1865 | rVV3 | 455889 | 685851 | 4.58% | 0.622% |
| 36 | 13.881 | 1865 | 1869 | 1872 | rVV | 329847 | 569476 | 3.80% | 0.516% |
| 37 | 13.910 | 1872 | 1874 | 1879 | rVB2 | 238875 | 291014 | 1.94% | 0.264% |
| 38 | 13.987 | 1883 | 1887 | 1894 | rBV7 | 270163 | 457116 | 3.05% | 0.414% |
| 39 | 14.046 | 1894 | 1897 | 1901 | rVB | 216369 | 270923 | 1.81% | 0.246% |
| 40 | 14.322 | 1934 | 1944 | 1952 | rBV4 | 2550183 | 6335387 | 42.30% | 5.742% |
| 41 | 14.410 | 1956 | 1959 | 1964 | rVB2 | 205235 | 299889 | 2.00% | 0.272% |
| 42 | 14.540 | 1976 | 1981 | 1988 | rBV | 318523 | 712789 | 4.76% | 0.646% |
| 43 | 14.728 | 2007 | 2013 | 2017 | rBV2 | 376344 | 659327 | 4.40% | 0.598% |
| 44 | 14.804 | 2022 | 2026 | 2030 | rVB | 457660 | 616819 | 4.12% | 0.559% |
| 45 | 14.881 | 2035 | 2039 | 2045 | rVB6 | 191675 | 314390 | 2.10% | 0.285% |
| 46 | 14.951 | 2046 | 2051 | 2054 | rBV | 335534 | 591425 | 3.95% | 0.536% |
| 47 | 14.993 | 2054 | 2058 | 2065 | rVB3 | 1356028 | 2179648 | 14.55% | 1.976% |
| 48 | 15.122 | 2073 | 2080 | 2083 | rBV3 | 222686 | 487148 | 3.25% | 0.442% |
| 49 | 15.151 | 2083 | 2085 | 2094 | rVB5 | 227023 | 356674 | 2.38% | 0.323% |
| 50 | 15.369 | 2117 | 2122 | 2127 | rBV3 | 287764 | 447303 | 2.99% | 0.405% |
| 51 | 15.428 | 2127 | 2132 | 2136 | rBV4 | 218954 | 475894 | 3.18% | 0.431% |
| 52 | 15.498 | 2140 | 2144 | 2148 | rVB2 | 301579 | 394659 | 2.63% | 0.358% |
| 53 | 15.622 | 2157 | 2165 | 2169 | rBV3 | 530036 | 1053792 | 7.04% | 0.955% |
| 54 | 15.816 | 2191 | 2198 | 2207 | rVB | 4709062 | 7089596 | 47.33% | 6.426% |
| 55 | 15.898 | 2208 | 2212 | 2218 | rBV5 | 255474 | 427301 | 2.85% | 0.387% |
| 56 | 16.063 | 2236 | 2240 | 2242 | rBV3 | 204559 | 271592 | 1.81% | 0.246% |
| 57 | 16.110 | 2242 | 2248 | 2251 | rVV2 | 1424295 | 2693945 | 17.99% | 2.442% |
| 58 | 16.140 | 2251 | 2253 | 2259 | rVV | 1174276 | 1471820 | 9.83% | 1.334% |
| 59 | 16.192 | 2260 | 2262 | 2267 | rVB | 215194 | 258285 | 1.72% | 0.234% |
| 60 | 16.434 | 2298 | 2303 | 2308 | rBV4 | 419226 | 737131 | 4.92% | 0.668% |
| 61 | 16.487 | 2309 | 2312 | 2315 | rVB3 | 144914 | 178797 | 1.19% | 0.162% |
| 62 | 16.587 | 2326 | 2329 | 2333 | rBV | 132182 | 240920 | 1.61% | 0.218% |
| 63 | 16.922 | 2382 | 2386 | 2398 | rVB | 824630 | 1389431 | 9.28% | 1.259% |
| 64 | 17.039 | 2403 | 2406 | 2407 | rVV | 239155 | 248930 | 1.66% | 0.226% |
| 65 | 17.075 | 2407 | 2412 | 2416 | rVV | 2381303 | 3530731 | 23.57% | 3.200% |
| 66 | 17.116 | 2416 | 2419 | 2423 | rVB | 368789 | 491288 | 3.28% | 0.445% |
| 67 | 17.275 | 2442 | 2446 | 2451 | rBV3 | 170029 | 340895 | 2.28% | 0.309% |
| 68 | 17.528 | 2486 | 2489 | 2495 | rVB3 | 283189 | 428157 | 2.86% | 0.388% |
| 69 | 17.657 | 2507 | 2511 | 2514 | rBV | 260839 | 314044 | 2.10% | 0.285% |
| 70 | 17.945 | 2557 | 2560 | 2564 | rBV | 306289 | 380801 | 2.54% | 0.345% |
| 71 | 17.992 | 2564 | 2568 | 2576 | rVV2 | 1468990 | 2070974 | 13.83% | 1.877% |

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP081820\
Data File : BP003025.D
Acq On : 18 Aug 2020 19:40
Operator : CG/JU
Sample : L3696-06
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :** BNA_P
**ClientSampleId :** W-1

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % 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:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP073020.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 72 | 18.128 | 2588 | 2591 | 2596 | rBV2 | 220086 | 267961 | 1.79% | 0.243% |
| 73 | 18.334 | 2623 | 2626 | 2630 | rVB2 | 162208 | 217180 | 1.45% | 0.197% |
| 74 | 18.675 | 2682 | 2684 | 2687 | rVB3 | 199108 | 174766 | 1.17% | 0.158% |
| 75 | 18.722 | 2690 | 2692 | 2696 | rVV2 | 163170 | 202096 | 1.35% | 0.183% |
| 76 | 18.839 | 2708 | 2712 | 2716 | rVB | 467341 | 628666 | 4.20% | 0.570% |
| 77 | 18.892 | 2717 | 2721 | 2725 | rBV2 | 385982 | 526981 | 3.52% | 0.478% |
| 78 | 18.975 | 2731 | 2735 | 2740 | rBV3 | 156083 | 237580 | 1.59% | 0.215% |
| 79 | 19.116 | 2756 | 2759 | 2765 | rVB2 | 447770 | 533703 | 3.56% | 0.484% |
| 80 | 19.228 | 2775 | 2778 | 2784 | rBV2 | 240938 | 395905 | 2.64% | 0.359% |
| 81 | 19.469 | 2816 | 2819 | 2824 | rVB | 168240 | 236747 | 1.58% | 0.215% |
| 82 | 19.528 | 2826 | 2829 | 2834 | rVB | 213815 | 296434 | 1.98% | 0.269% |
| 83 | 19.651 | 2844 | 2850 | 2855 | rBV | 11657596 | 14977622 | 100.00% | 13.575% |
| 84 | 20.345 | 2965 | 2968 | 2977 | rVB | 201988 | 290597 | 1.94% | 0.263% |
| 85 | 20.904 | 3059 | 3063 | 3068 | rBV2 | 775411 | 1098505 | 7.33% | 0.996% |
| 86 | 20.951 | 3068 | 3071 | 3079 | rVB | 375405 | 452608 | 3.02% | 0.410% |
| 87 | 21.163 | 3102 | 3107 | 3112 | rBV | 3919891 | 4651287 | 31.05% | 4.216% |
| 88 | 23.404 | 3481 | 3488 | 3501 | rVB2 | 2473617 | 4809062 | 32.11% | 4.359% |

Sum of corrected areas: 110332036

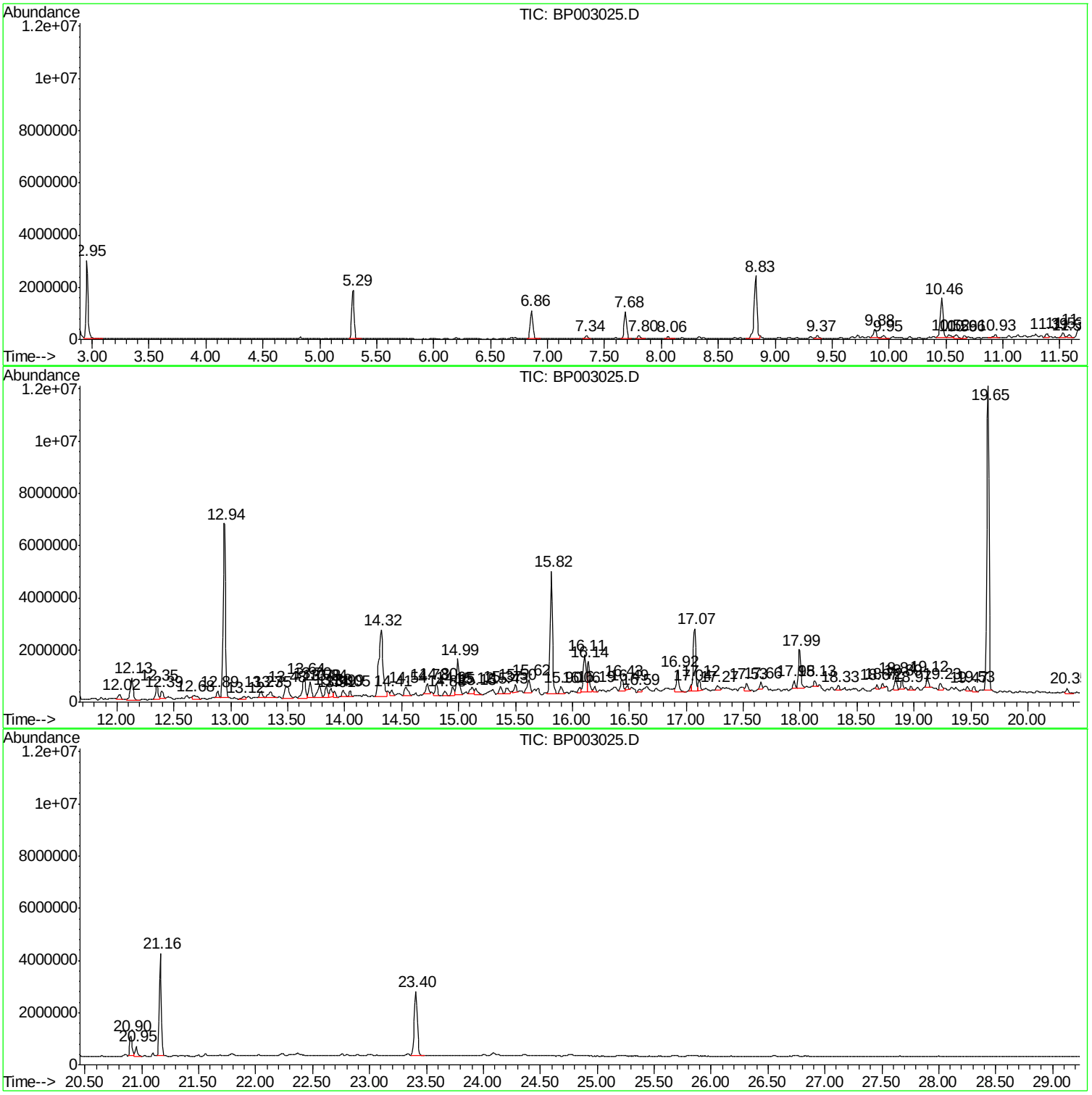
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP081820\
Data File : BP003025.D
Acq On : 18 Aug 2020 19:40
Operator : CG/JU
Sample : L3696-06
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :** BNA_P
**ClientSampleId :** W-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP073020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP081820\
Data File : BP003025.D
Acq On : 18 Aug 2020 19:40
Operator : CG/JU
Sample : L3696-06
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :** BNA_P
**ClientSampleId :** W-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP073020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

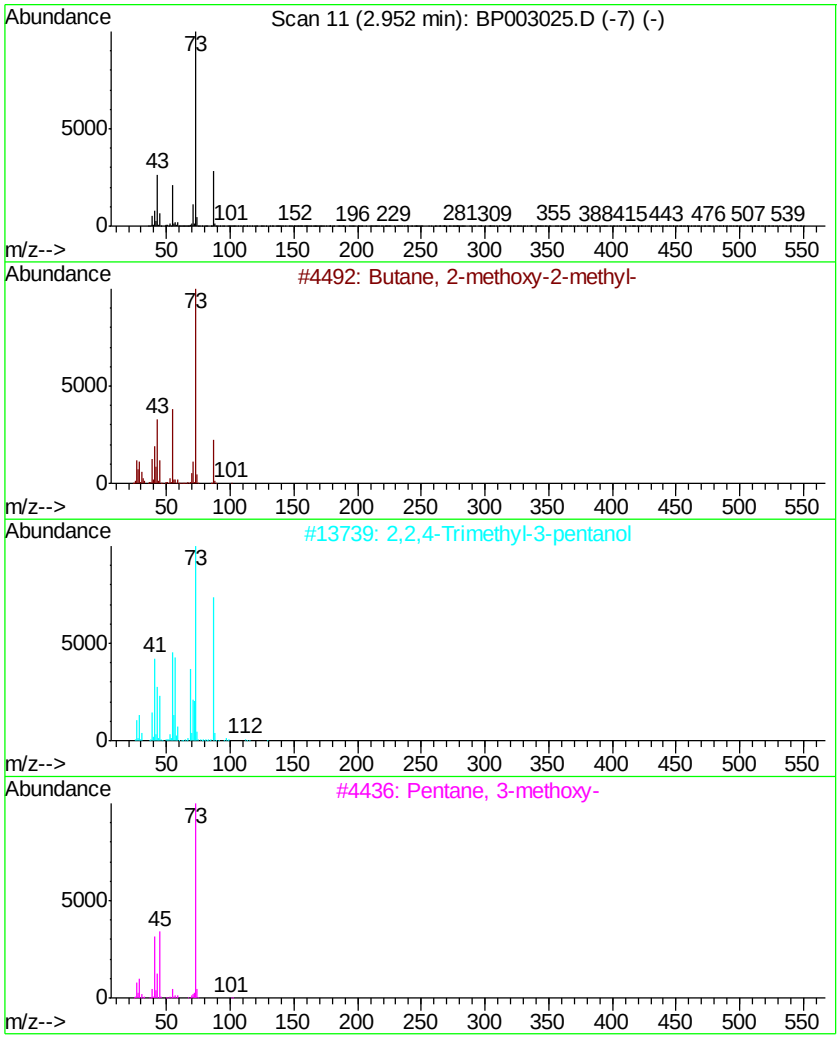
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

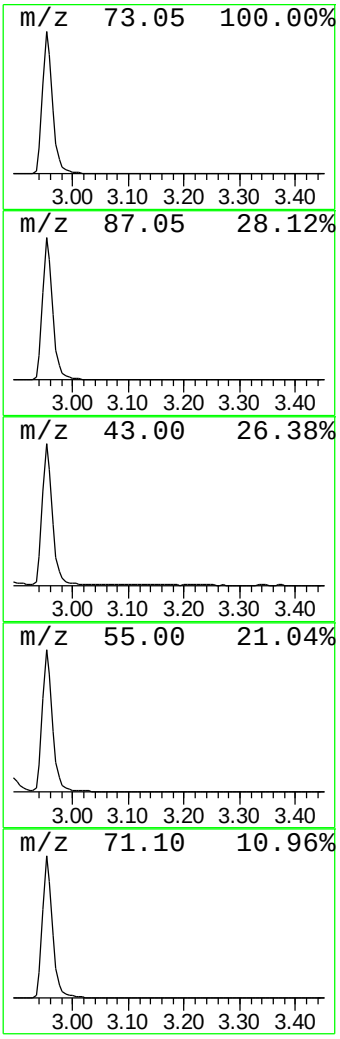
**********

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 2.95 | 45.03 ng | 3757400 | 1,4-Dichlorobenzene-d4 | 7.68 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Butane, 2-methoxy-2-methyl- | 102 | C6H14O | 000994-05-8 | 78 |
| 2 | 2,2,4-Trimethyl-3-pentanol | 130 | C8H18O | 005162-48-1 | 43 |
| 3 | Pentane, 3-methoxy- | 102 | C6H14O | 036839-67-5 | 17 |
| 4 | N-Ethylformamide | 73 | C3H7NO | 000627-45-2 | 9 |
| 5 | 3-Hexanol, 2-methyl- | 116 | C7H16O | 000617-29-8 | 9 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP081820\
Data File : BP003025.D
Acq On : 18 Aug 2020 19:40
Operator : CG/JU
Sample : L3696-06
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :** BNA_P
**ClientSampleId :** W-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP073020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

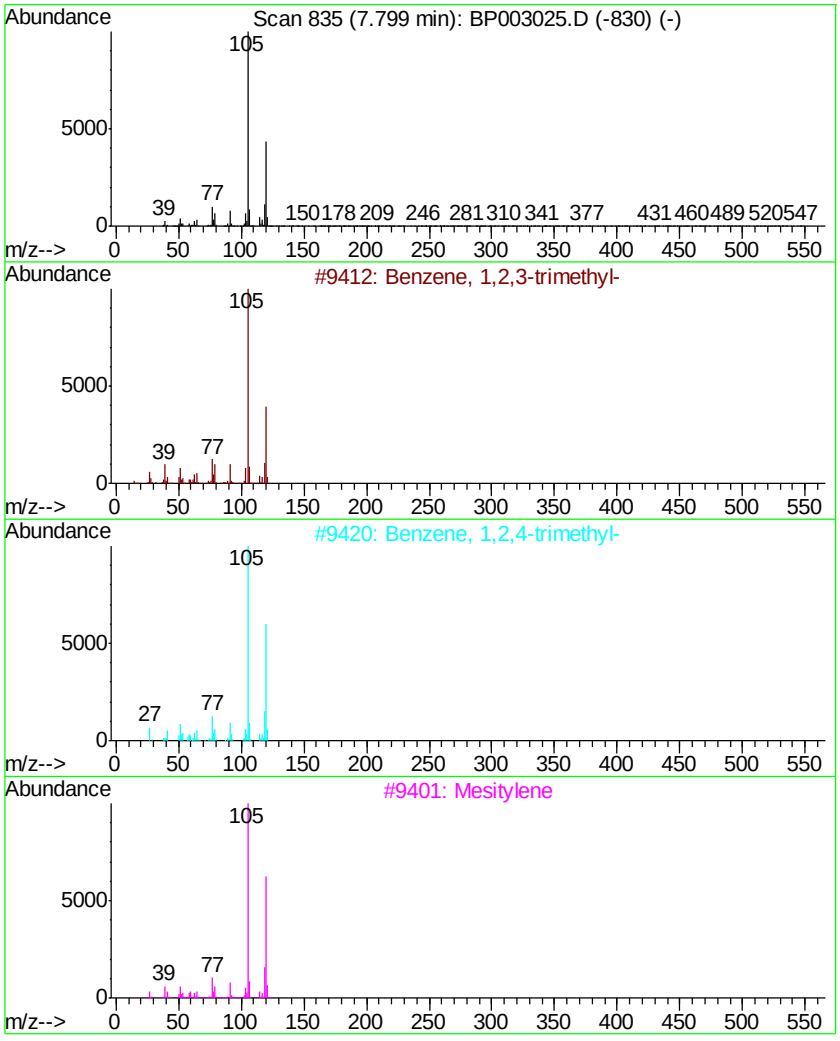
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

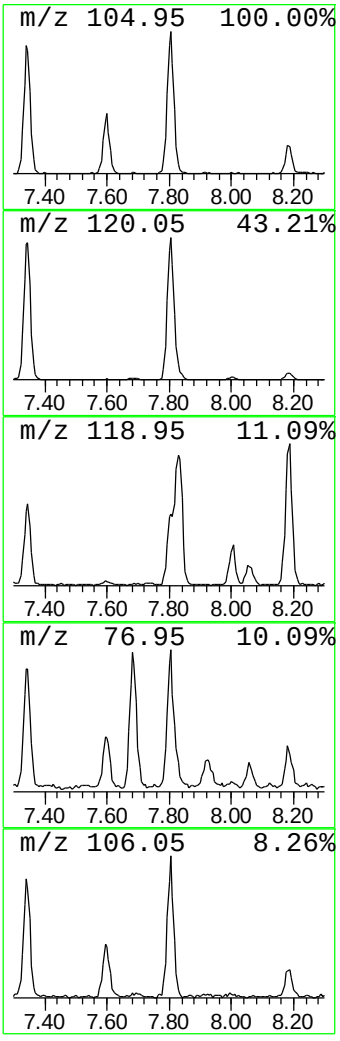
**********

Peak Number 2 Benzene, 1,2,3-trimethyl- Concentration Rank 19

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 7.80 | 2.87 ng | 239071 | 1,4-Dichlorobenzene-d4 | 7.68 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzene, 1,2,3-trimethyl- | 120 | C9H12 | 000526-73-8 | 95 |
| 2 | Benzene, 1,2,4-trimethyl- | 120 | C9H12 | 000095-63-6 | 95 |
| 3 | Mesitylene | 120 | C9H12 | 000108-67-8 | 94 |
| 4 | Benzene, 1-ethyl-3-methyl- | 120 | C9H12 | 000620-14-4 | 90 |
| 5 | Benzene, 1-ethyl-4-methyl- | 120 | C9H12 | 000622-96-8 | 87 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP081820\
Data File : BP003025.D
Acq On : 18 Aug 2020 19:40
Operator : CG/JU
Sample : L3696-06
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :** BNA_P
**ClientSampleId :** W-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP073020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

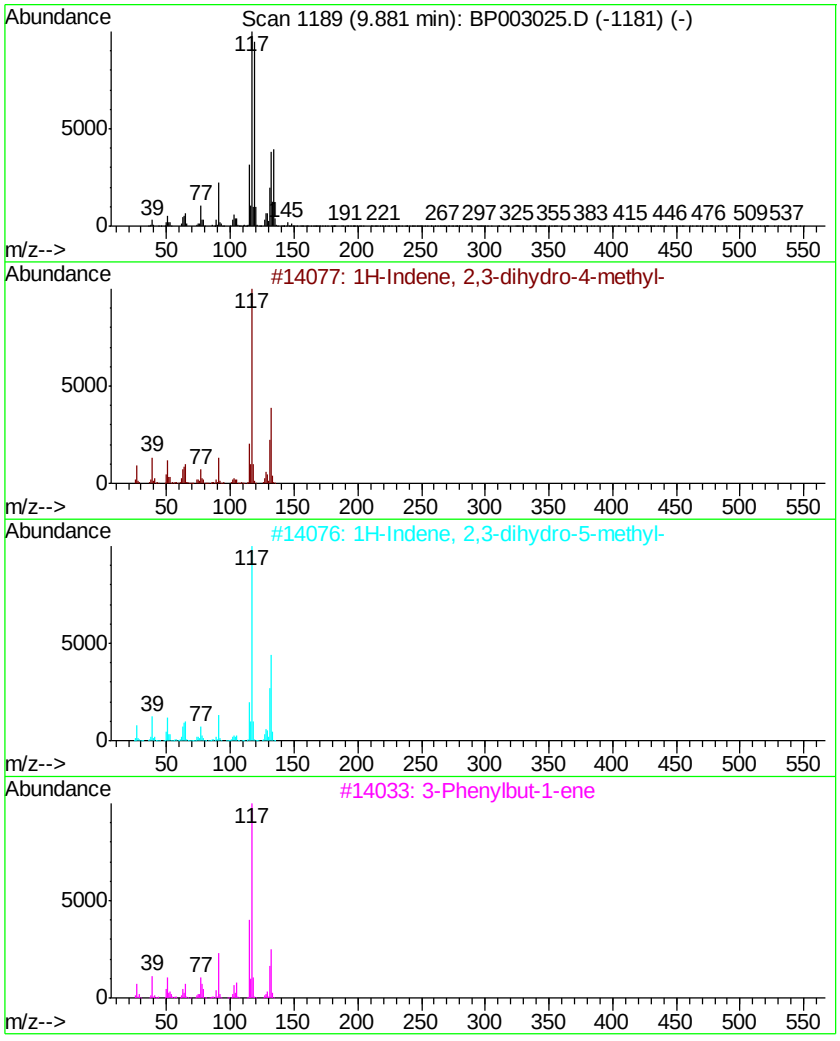
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

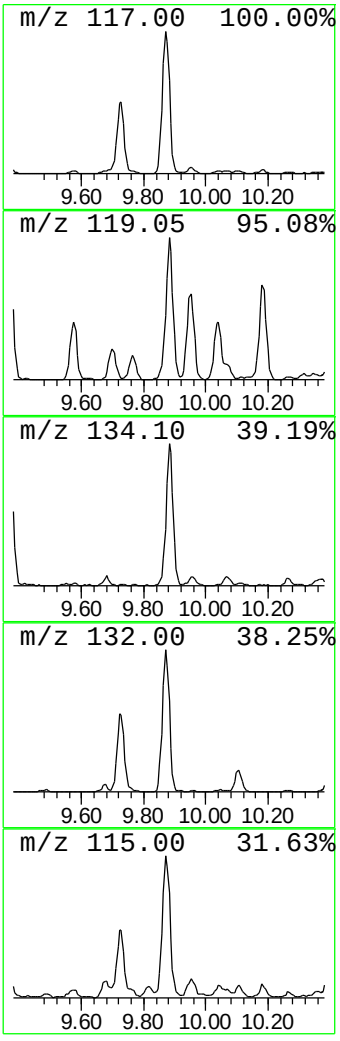
**********

Peak Number 3 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 9

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 9.88 | 4.47 ng | 587962 | Naphthalene-d8 | 10.46 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1H-Indene, 2,3-dihydro-4-methyl- | 132 | C10H12 | 000824-22-6 | 70 |
| 2 | 1H-Indene, 2,3-dihydro-5-methyl- | 132 | C10H12 | 000874-35-1 | 64 |
| 3 | 3-Phenylbut-1-ene | 132 | C10H12 | 000934-10-1 | 62 |
| 4 | Benzene, 1-ethenyl-4-ethyl- | 132 | C10H12 | 003454-07-7 | 60 |
| 5 | Indan, 1-methyl- | 132 | C10H12 | 000767-58-8 | 60 |



Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP081820\
Data File : BP003025.D
Acq On : 18 Aug 2020 19:40
Operator : CG/JU
Sample : L3696-06
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :** BNA_P
**ClientSampleId :** W-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP073020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

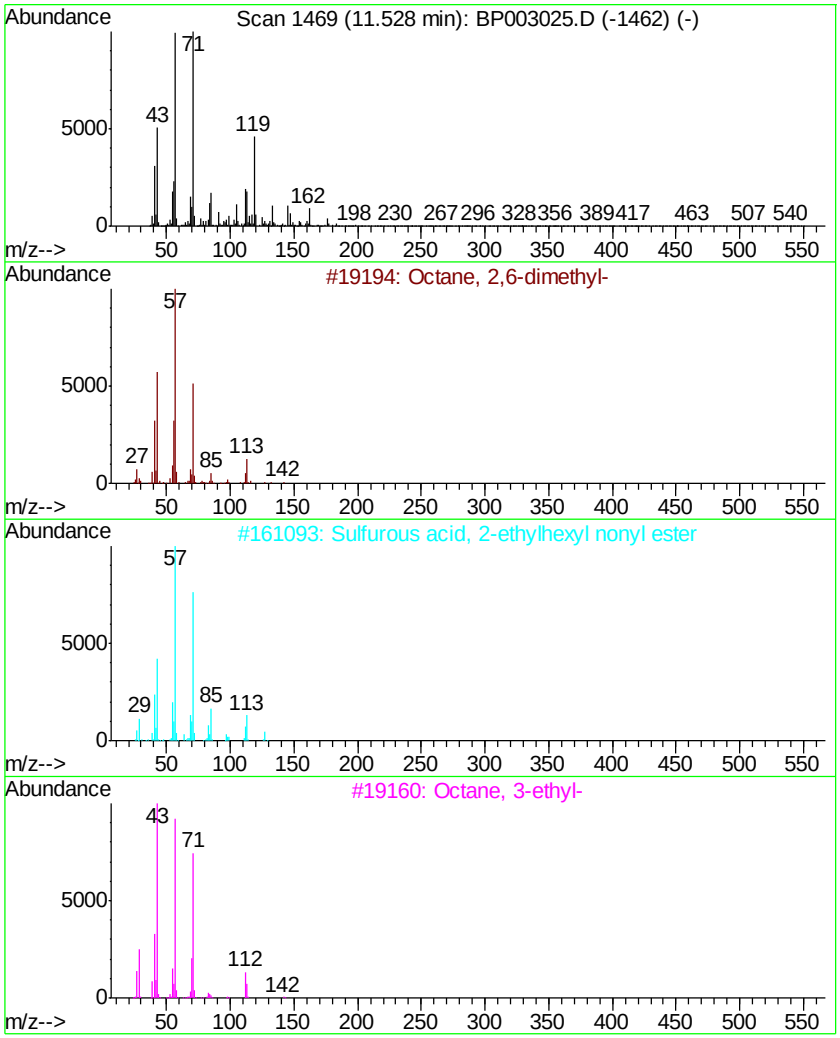
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

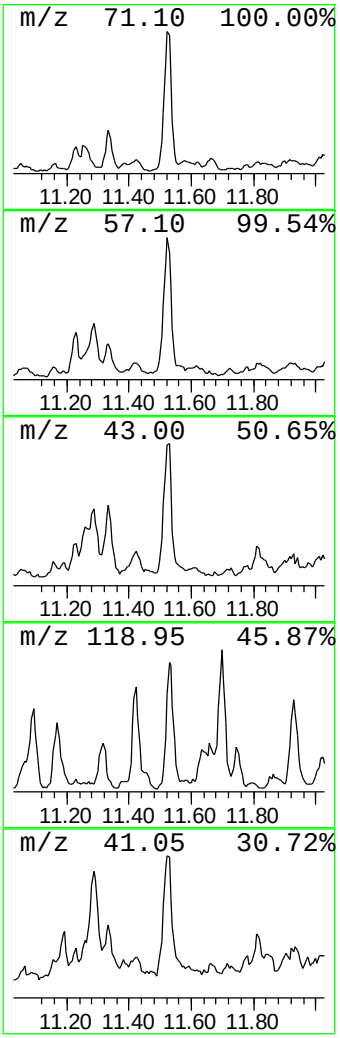
**********

Peak Number 4 Octane, 2,6-dimethyl- Concentration Rank 18

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.53 | 2.88 ng | 379082 | Naphthalene-d8 | 10.46 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Octane, 2,6-dimethyl- | 142 | C10H22 | 002051-30-1 | 50 |
| 2 | Sulfurous acid, 2-ethylhexyl non... | 320 | C17H36O3S | 1000309-19-2 | 49 |
| 3 | Octane, 3-ethyl- | 142 | C10H22 | 005881-17-4 | 47 |
| 4 | 3-Bromooctane | 192 | C8H17Br | 000999-64-4 | 47 |
| 5 | Dodecane, 4,6-dimethyl- | 198 | C14H30 | 061141-72-8 | 47 |



Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP081820\
Data File : BP003025.D
Acq On : 18 Aug 2020 19:40
Operator : CG/JU
Sample : L3696-06
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :** BNA_P
**ClientSampleId :** W-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP073020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

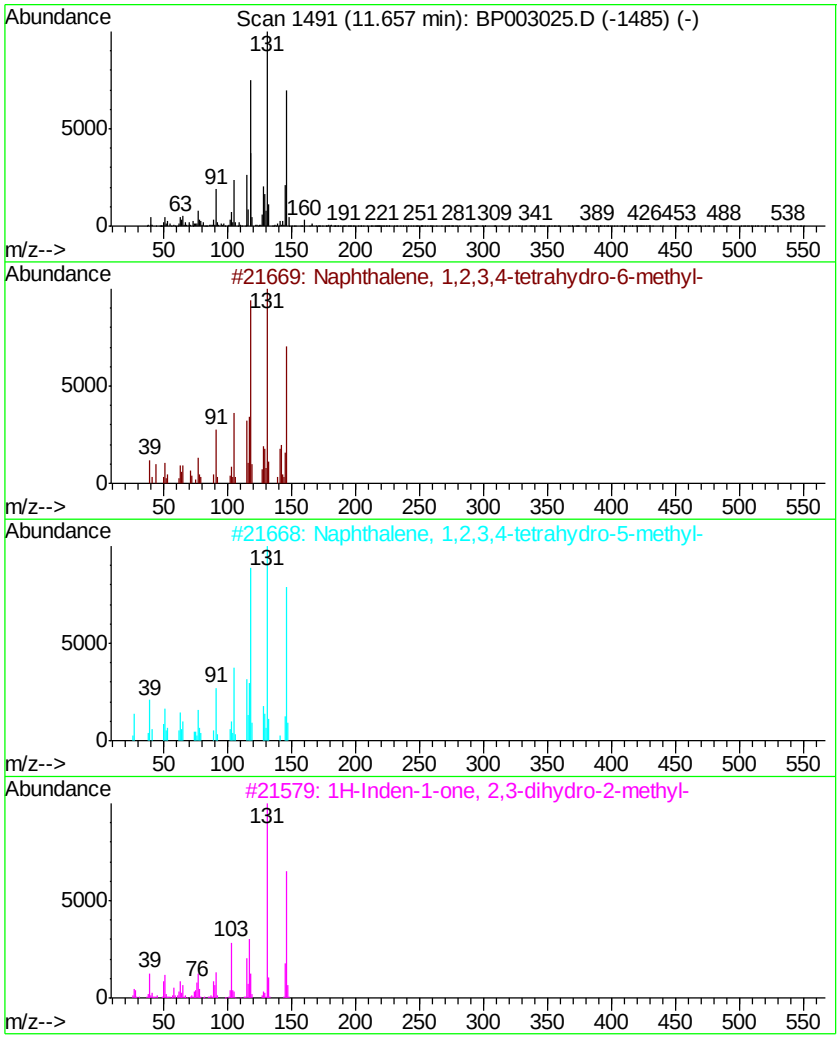
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

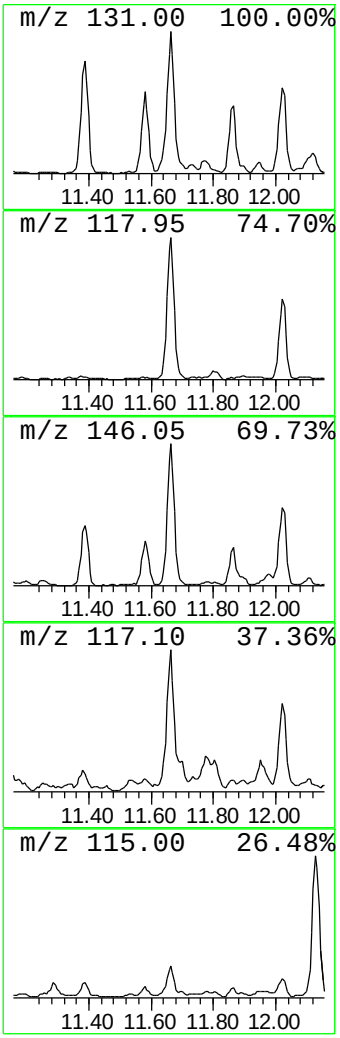
**********

Peak Number 5 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 10

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.66 | 4.33 ng | 569342 | Naphthalene-d8 | 10.46 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Naphthalene, 1,2,3,4-tetrahydro-... | 146 | C11H14 | 001680-51-9 | 94 |
| 2 | Naphthalene, 1,2,3,4-tetrahydro-... | 146 | C11H14 | 002809-64-5 | 94 |
| 3 | 1H-Inden-1-one, 2,3-dihydro-2-me... | 146 | C10H10O | 017496-14-9 | 74 |
| 4 | Benzene, (1-methyl-1-butenyl)- | 146 | C11H14 | 053172-84-2 | 70 |
| 5 | Benzene, 2-ethenyl-1,3,5-trimethyl- | 146 | C11H14 | 000769-25-5 | 58 |



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Data File : BP003025.D
Acq On : 18 Aug 2020 19:40
Operator : CG/JU
Sample : L3696-06
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :** BNA_P
**ClientSampleId :** W-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP073020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

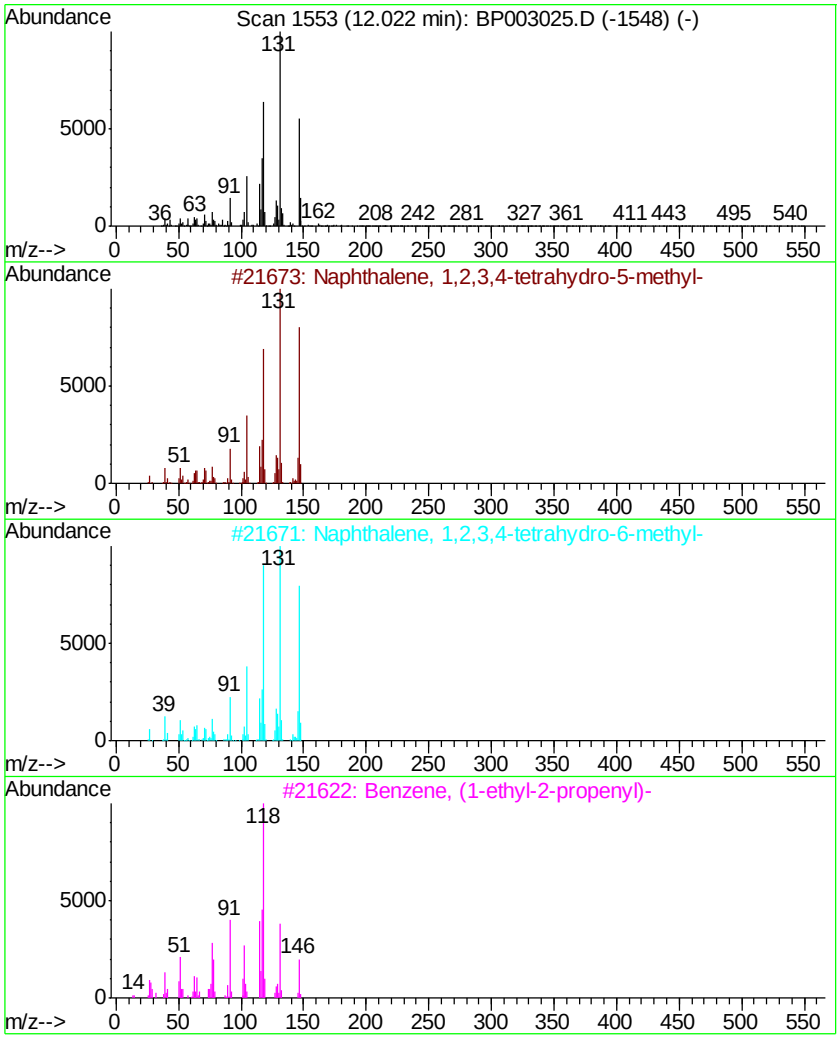
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

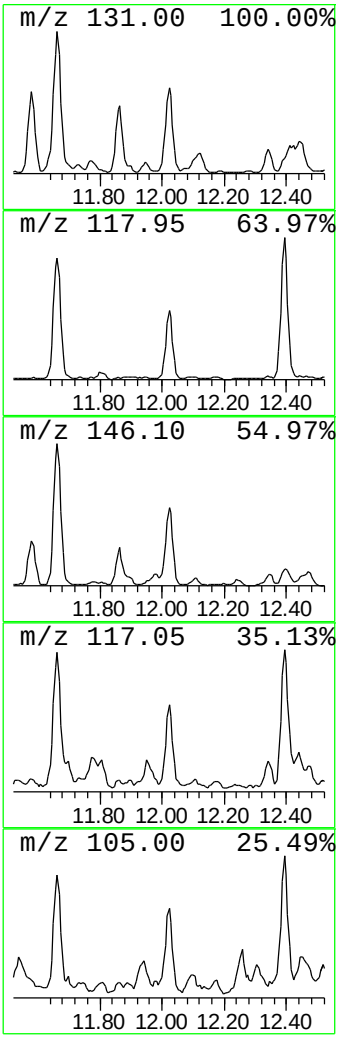
**********

Peak Number 6 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 20

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.02 | 2.52 ng | 331624 | Naphthalene-d8 | 10.46 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Naphthalene, 1,2,3,4-tetrahydro-... | 146 | C11H14 | 002809-64-5 | 94 |
| 2 | Naphthalene, 1,2,3,4-tetrahydro-... | 146 | C11H14 | 001680-51-9 | 94 |
| 3 | Benzene, (1-ethyl-2-propenyl)- | 146 | C11H14 | 019947-22-9 | 70 |
| 4 | 1H-Indene, 2,3-dihydro-5,6-dimet... | 146 | C11H14 | 001075-22-5 | 64 |
| 5 | 1H-Inden-1-one, 2,3-dihydro-2-me... | 146 | C10H10O | 017496-14-9 | 62 |



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Data File : BP003025.D
Acq On : 18 Aug 2020 19:40
Operator : CG/JU
Sample : L3696-06
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :** BNA_P
**ClientSampleId :** W-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP073020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

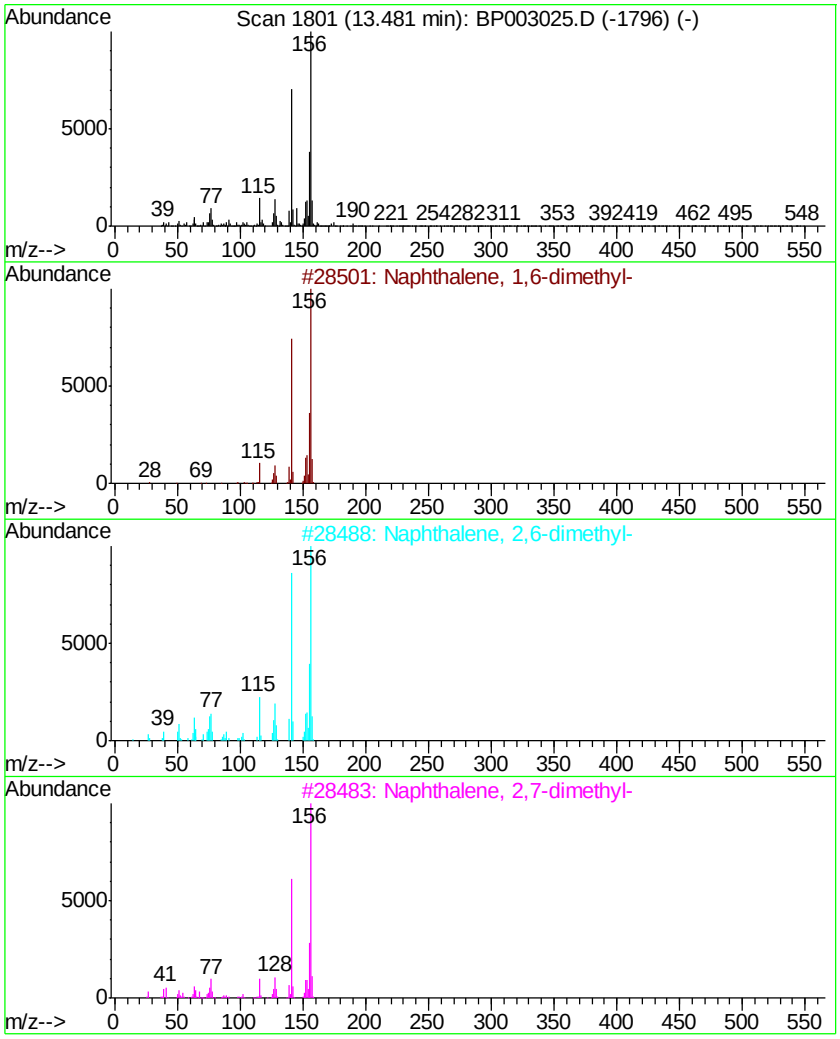
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

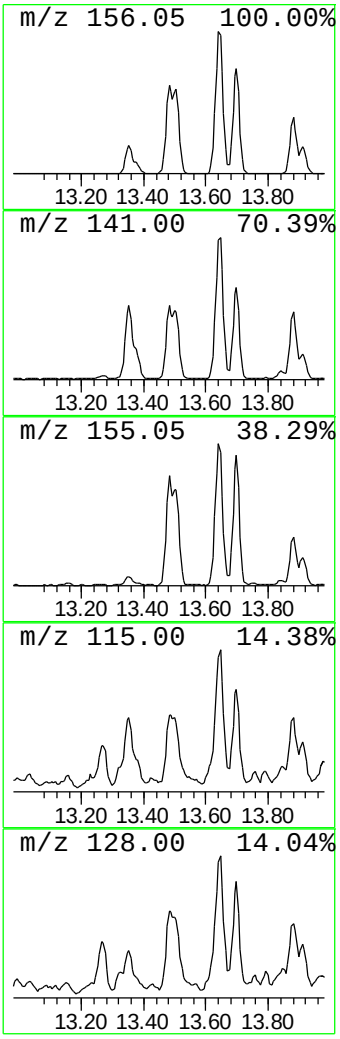
**********

Peak Number 7 Naphthalene, 1,6-dimethyl- Concentration Rank 12

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.48 | 4.11 ng | 1301330 | Acenaphthene-d10 | 14.32 |

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



Library Search Compound Report

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Data File : BP003025.D
Acq On : 18 Aug 2020 19:40
Operator : CG/JU
Sample : L3696-06
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :** BNA_P
**ClientSampleId :** W-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP073020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

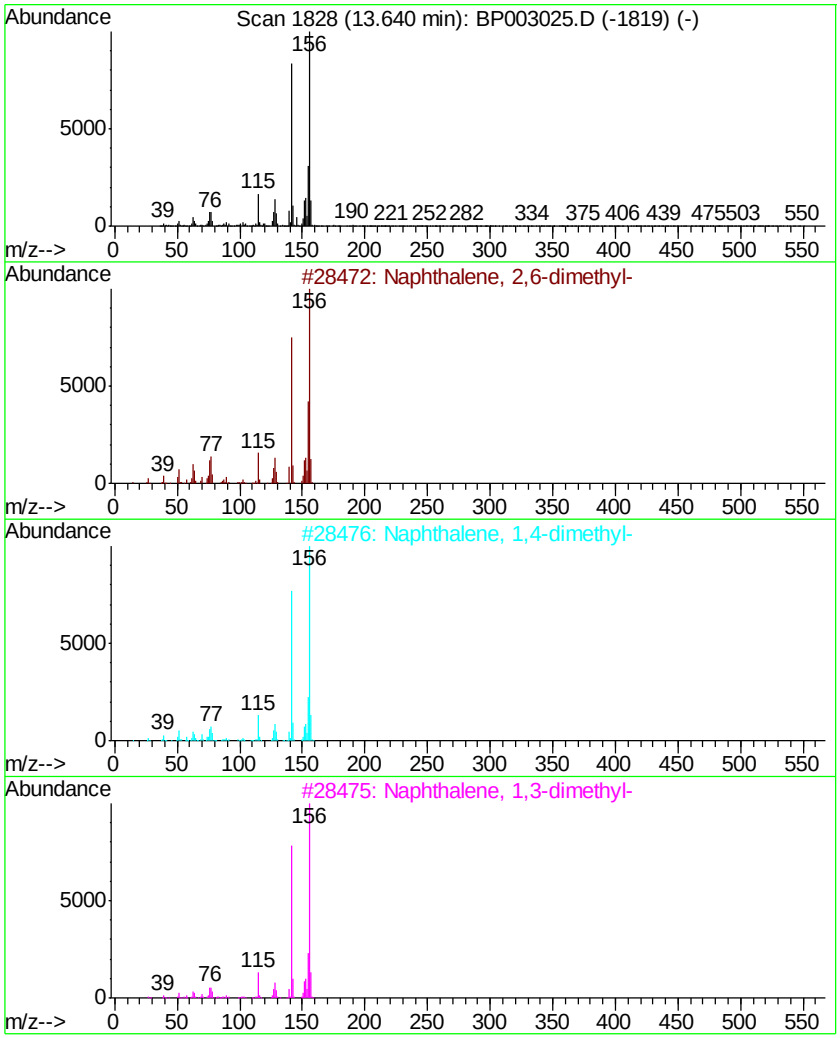
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

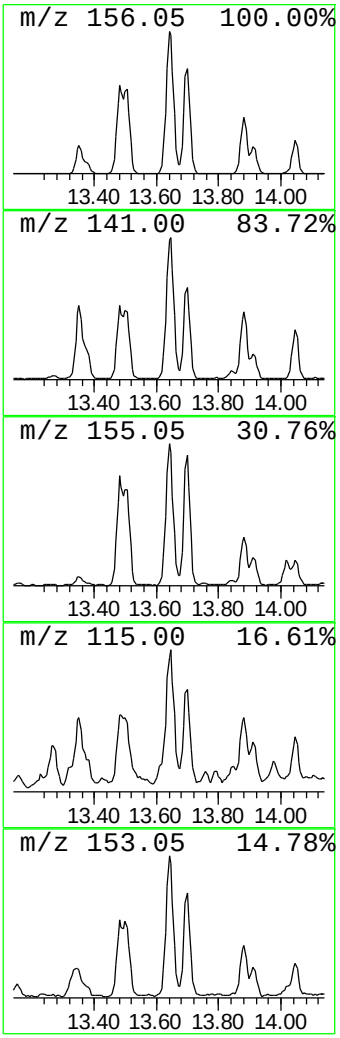
**********

Peak Number 8 Naphthalene, 2,6-dimethyl- Concentration Rank 7

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.64 | 4.73 ng | 1499450 | Acenaphthene-d10 | 14.32 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Naphthalene, 2,6-dimethyl- | 156 | C12H12 | 000581-42-0 | 97 |
| 2 | Naphthalene, 1,4-dimethyl- | 156 | C12H12 | 000571-58-4 | 97 |
| 3 | Naphthalene, 1,3-dimethyl- | 156 | C12H12 | 000575-41-7 | 97 |
| 4 | Naphthalene, 1,6-dimethyl- | 156 | C12H12 | 000575-43-9 | 96 |
| 5 | Naphthalene, 1,7-dimethyl- | 156 | C12H12 | 000575-37-1 | 96 |



Library Search Compound Report

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Data File : BP003025.D
Acq On : 18 Aug 2020 19:40
Operator : CG/JU
Sample : L3696-06
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :** BNA_P
**ClientSampleId :** W-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP073020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

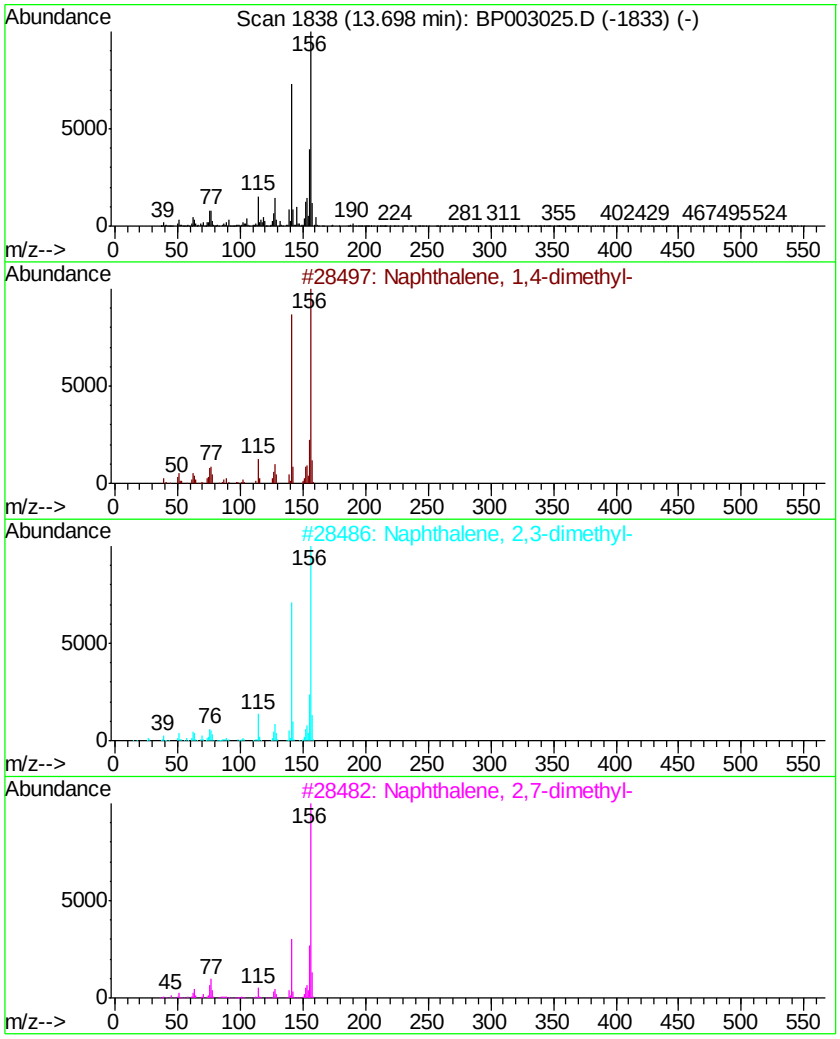
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

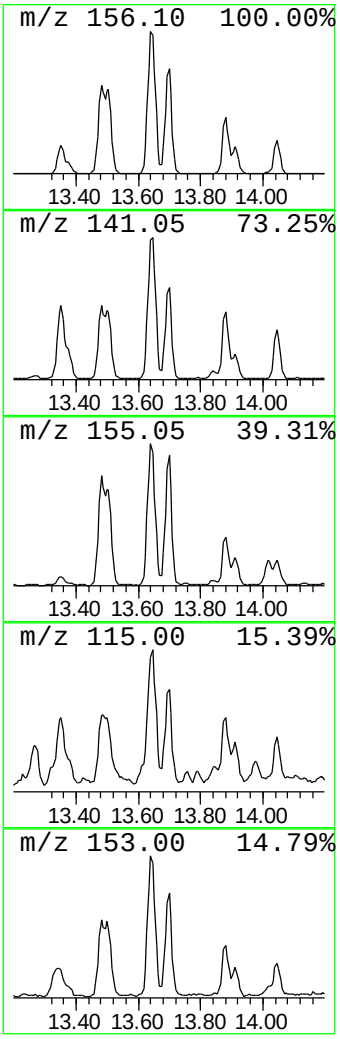
**********

Peak Number 9 Naphthalene, 1,4-dimethyl- Concentration Rank 16

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.70 | 3.02 ng | 955714 | Acenaphthene-d10 | 14.32 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Naphthalene, 1,4-dimethyl- | 156 | C12H12 | 000571-58-4 | 94 |
| 2 | Naphthalene, 2,3-dimethyl- | 156 | C12H12 | 000581-40-8 | 94 |
| 3 | Naphthalene, 2,7-dimethyl- | 156 | C12H12 | 000582-16-1 | 94 |
| 4 | Naphthalene, 2,6-dimethyl- | 156 | C12H12 | 000581-42-0 | 94 |
| 5 | Naphthalene, 1,5-dimethyl- | 156 | C12H12 | 000571-61-9 | 93 |



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Data File : BP003025.D
Acq On : 18 Aug 2020 19:40
Operator : CG/JU
Sample : L3696-06
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :** BNA_P
**ClientSampleId :** W-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP073020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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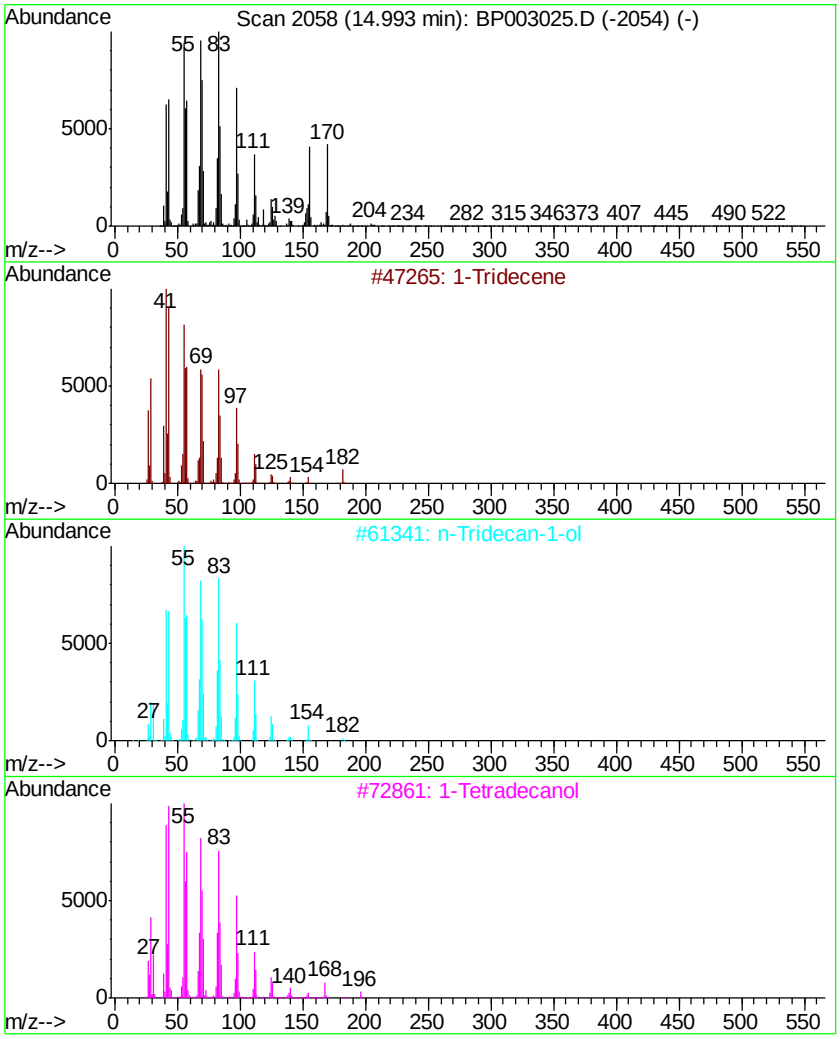
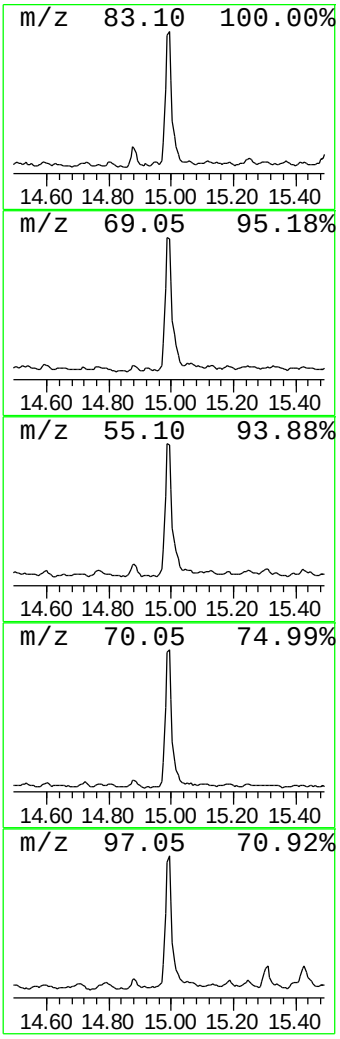
Peak Number 10 1-Tridecene Concentration Rank 6

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 14.99 | 6.88 ng | 2179650 | Acenaphthene-d10 | 14.32 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1-Tridecene | 182 | C13H26 | 002437-56-1 | 93 |
| 2 | n-Tridecan-1-ol | 200 | C13H28O | 000112-70-9 | 93 |
| 3 | 1-Tetradecanol | 214 | C14H30O | 000112-72-1 | 81 |
| 4 | 2-Butenedioic acid (Z)-, monodod... | 284 | C16H28O4 | 002424-61-5 | 76 |
| 5 | Acetic acid, trifluoro-, undecyl... | 268 | C13H23F3O2 | 053800-01-4 | 76 |



# Library Search Compound Report

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Data File : BP003025.D
Acq On : 18 Aug 2020 19:40
Operator : CG/JU
Sample : L3696-06
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :** BNA_P
**ClientSampleId :** W-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP073020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

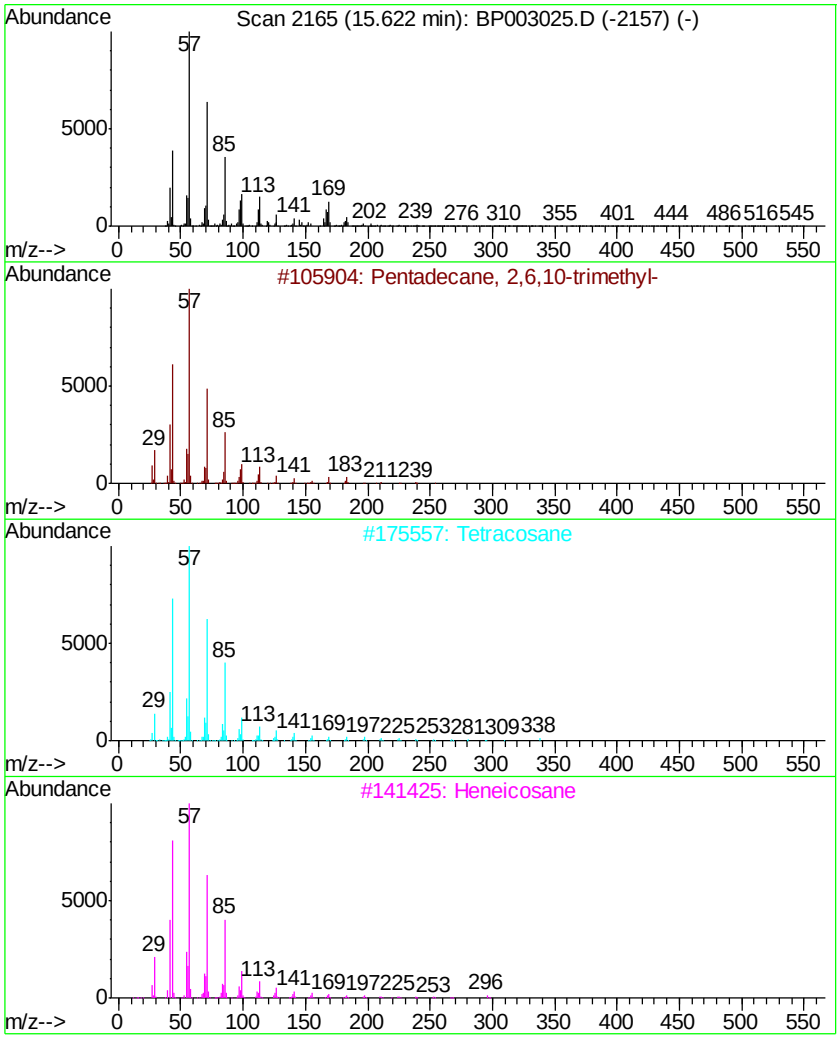
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

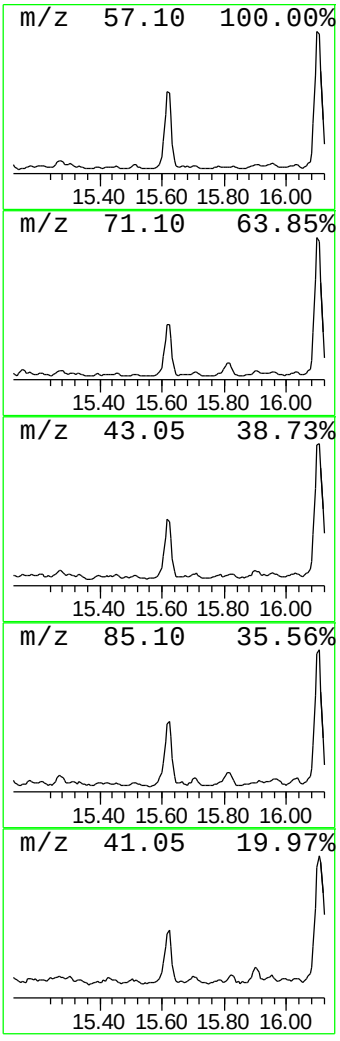
**********

Peak Number 11 Pentadecane, 2,6,10-trimethyl- Concentration Rank 14

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.62 | 3.33 ng | 1053790 | Acenaphthene-d10 | 14.32 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Pentadecane, 2,6,10-trimethyl- | 254 | C18H38 | 003892-00-0 | 95 |
| 2 | Tetracosane | 338 | C24H50 | 000646-31-1 | 95 |
| 3 | Heneicosane | 296 | C21H44 | 000629-94-7 | 93 |
| 4 | Tridecane, 5-propyl- | 226 | C16H34 | 055045-11-9 | 91 |
| 5 | 2-Bromo dodecane | 248 | C12H25Br | 013187-99-0 | 83 |



# Library Search Compound Report

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Data File : BP003025.D
Acq On : 18 Aug 2020 19:40
Operator : CG/JU
Sample : L3696-06
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :** BNA_P
**ClientSampleId :** W-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP073020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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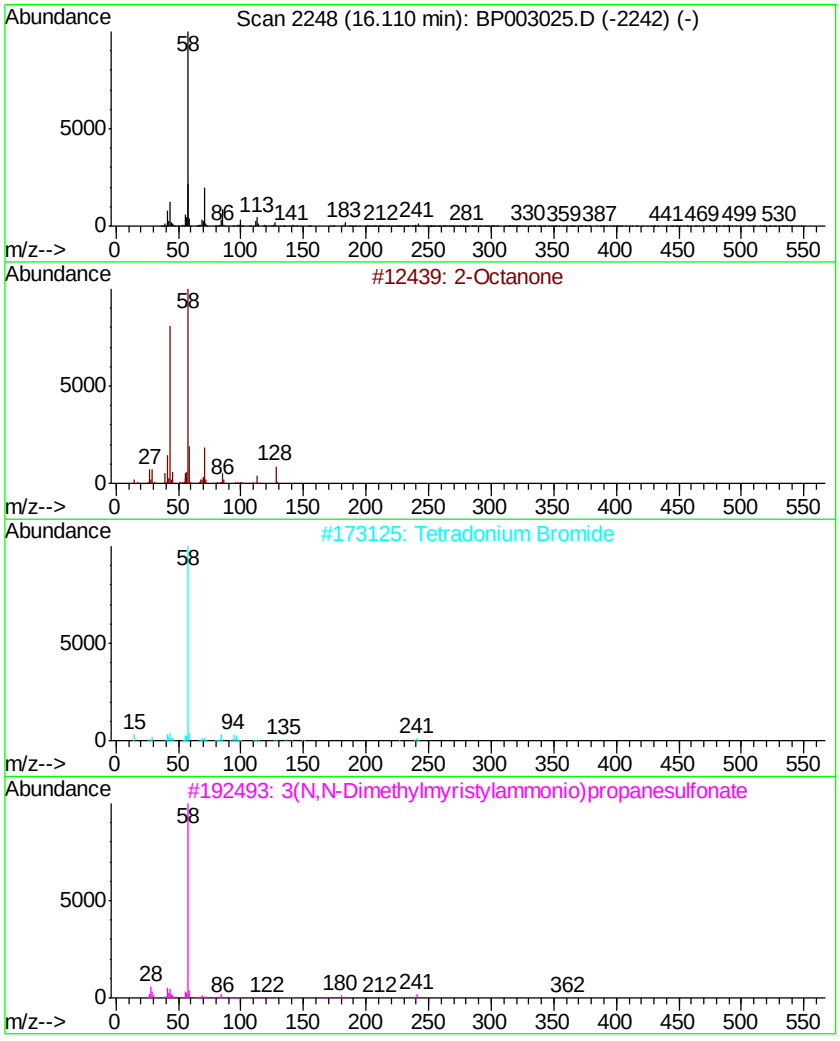
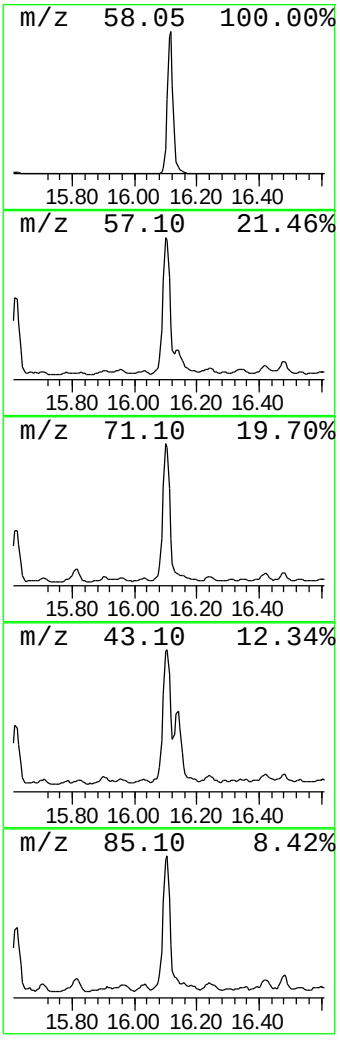
Peak Number 12 2-Octanone Concentration Rank 2

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.11 | 15.26 ng | 2693950 | Phenanthrene-d10 | 17.07 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-Octanone | 128 | C8H16O | 000111-13-7 | 56 |
| 2 | Tetradonium Bromide | 335 | C17H38BrN | 001119-97-7 | 53 |
| 3 | 3(N,N-Dimethylmyristylammonio)pr... | 363 | C19H41NO3S | 014933-09-6 | 53 |
| 4 | 2,6-Difluoro-3-methylbenzoic aci... | 243 | C12H15F2NO2 | 1000343-75-2 | 50 |
| 5 | Benzedrex | 155 | C10H21N | 000101-40-6 | 50 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP081820\
Data File : BP003025.D
Acq On : 18 Aug 2020 19:40
Operator : CG/JU
Sample : L3696-06
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :** BNA_P
**ClientSampleId :** W-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP073020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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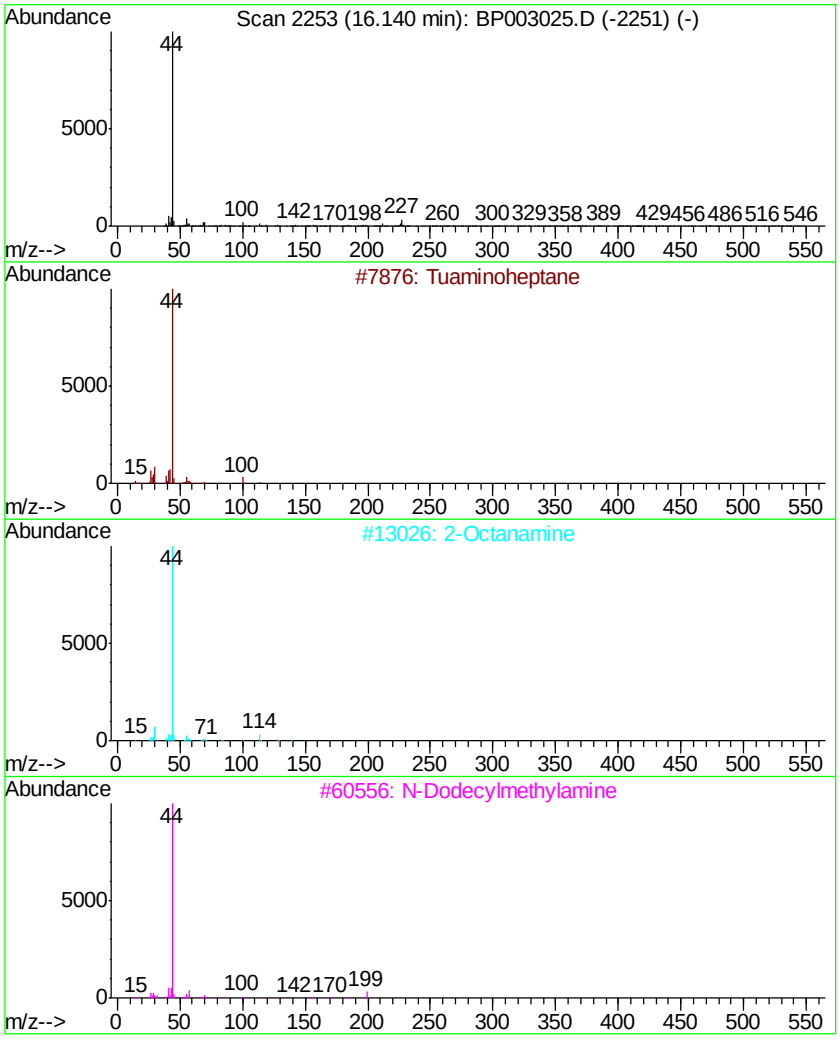
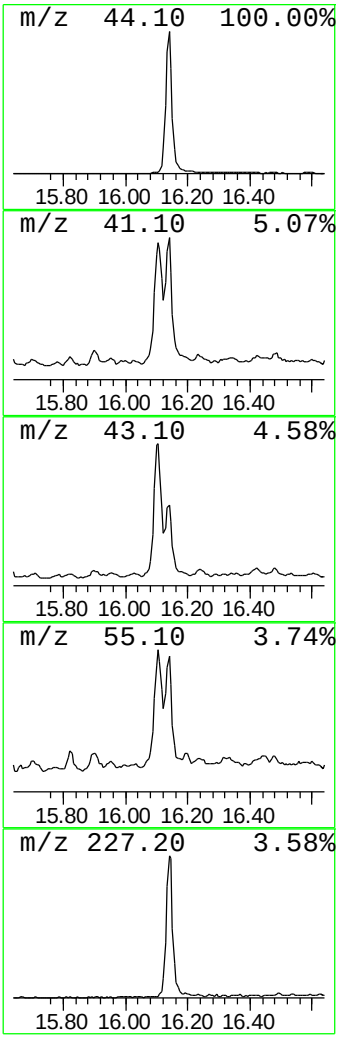
Peak Number 13 Tuaminoheptane Concentration Rank 4

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.14 | 8.34 ng | 1471820 | Phenanthrene-d10 | 17.07 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Tuaminoheptane | 115 | C7H17N | 000123-82-0 | 59 |
| 2 | 2-Octanamine | 129 | C8H19N | 000693-16-3 | 59 |
| 3 | N-Dodecylmethylamine | 199 | C13H29N | 1000342-26-1 | 59 |
| 4 | 12-Methylaminolauric acid | 229 | C13H27NO2 | 007408-81-3 | 38 |
| 5 | (+)-2-Aminoheptane | 115 | C7H17N | 006240-90-0 | 38 |



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Data File : BP003025.D
Acq On : 18 Aug 2020 19:40
Operator : CG/JU
Sample : L3696-06
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :** BNA_P
**ClientSampleId :** W-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP073020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

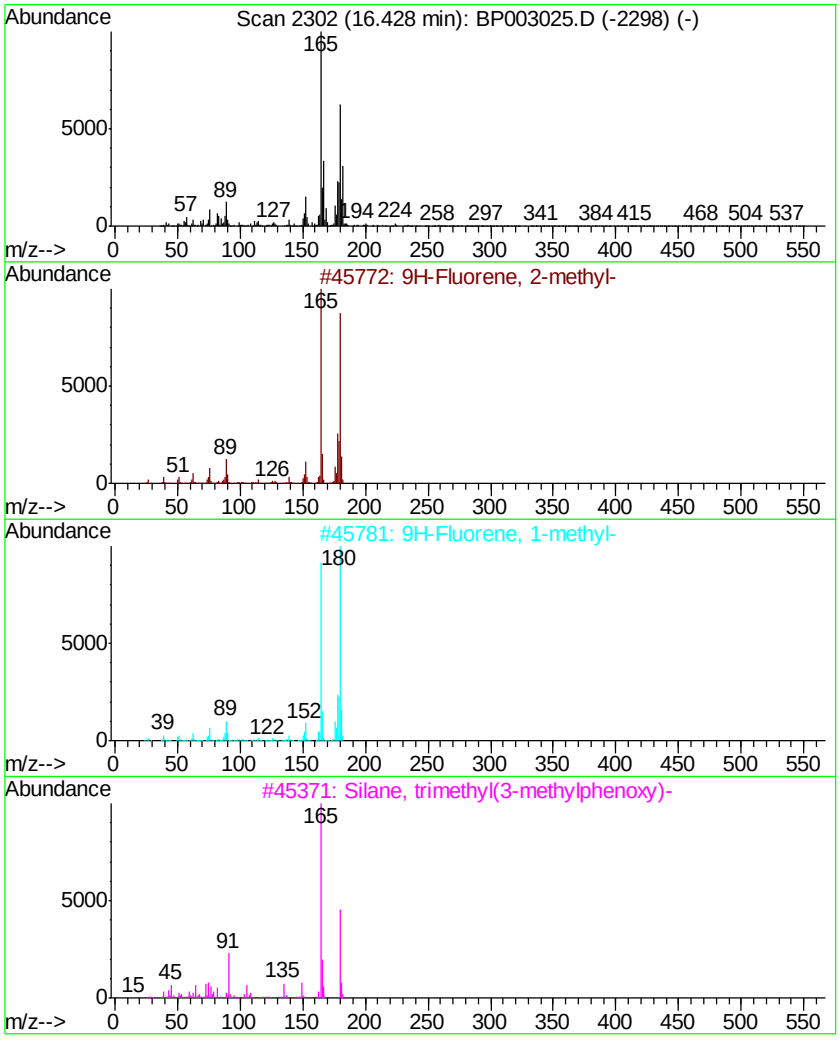
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

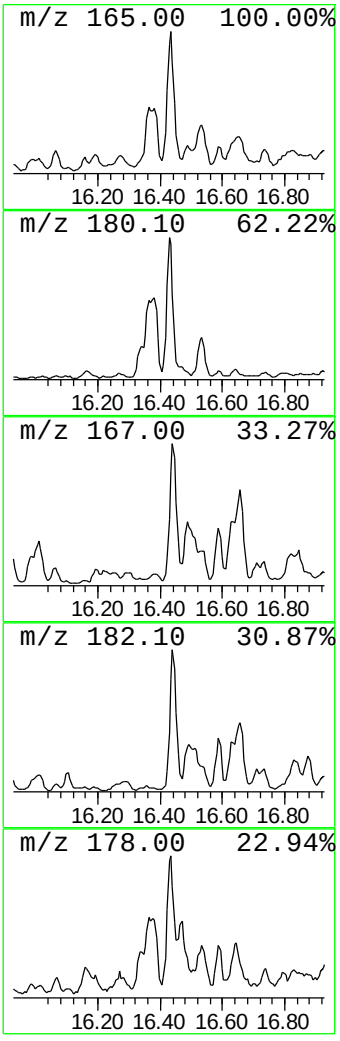
**********

Peak Number 14 9H-Fluorene, 2-methyl- Concentration Rank 11

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.43 | 4.18 ng | 737131 | Phenanthrene-d10 | 17.07 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 9H-Fluorene, 2-methyl- | 180 | C14H12 | 001430-97-3 | 55 |
| 2 | 9H-Fluorene, 1-methyl- | 180 | C14H12 | 001730-37-6 | 50 |
| 3 | Silane, trimethyl(3-methylphenoxy)- | 180 | C10H16OSi | 017902-31-7 | 38 |
| 4 | 1,1'-Biphenyl, 2,3'-dimethyl- | 182 | C14H14 | 000611-43-8 | 38 |
| 5 | 9H-Fluorene, 9-methyl- | 180 | C14H12 | 002523-37-7 | 38 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP081820\
Data File : BP003025.D
Acq On : 18 Aug 2020 19:40
Operator : CG/JU
Sample : L3696-06
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :** BNA_P
**ClientSampleId :** W-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP073020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

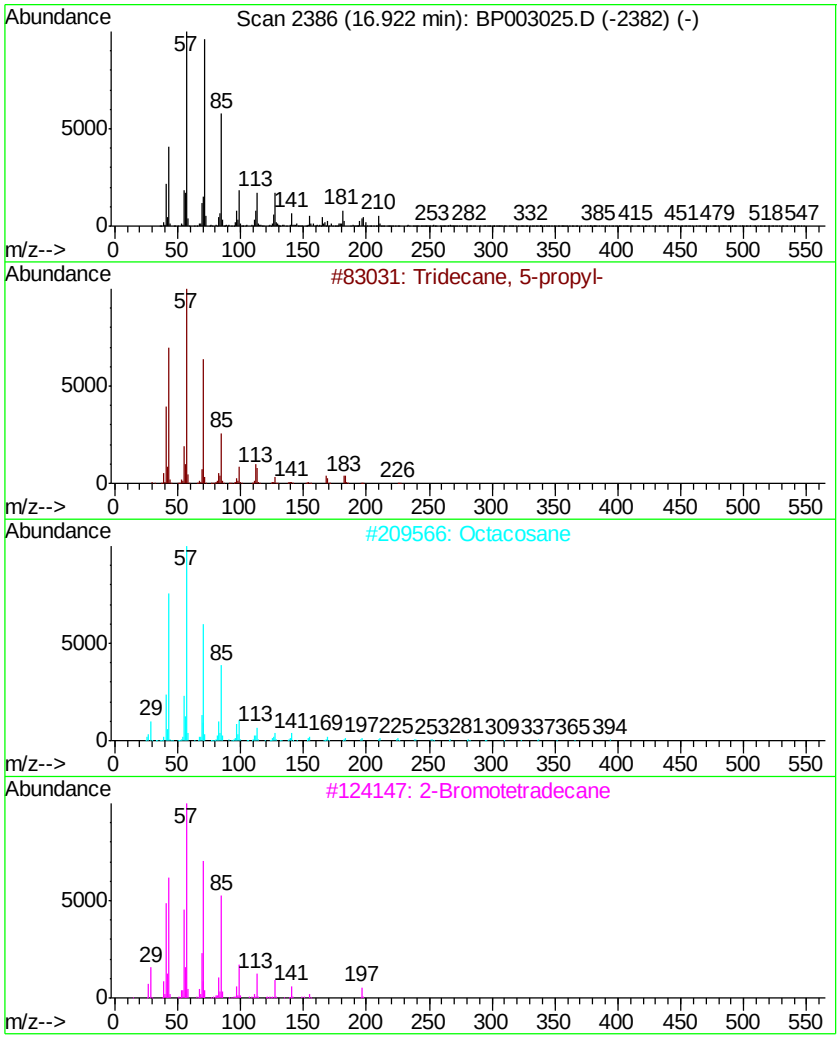
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

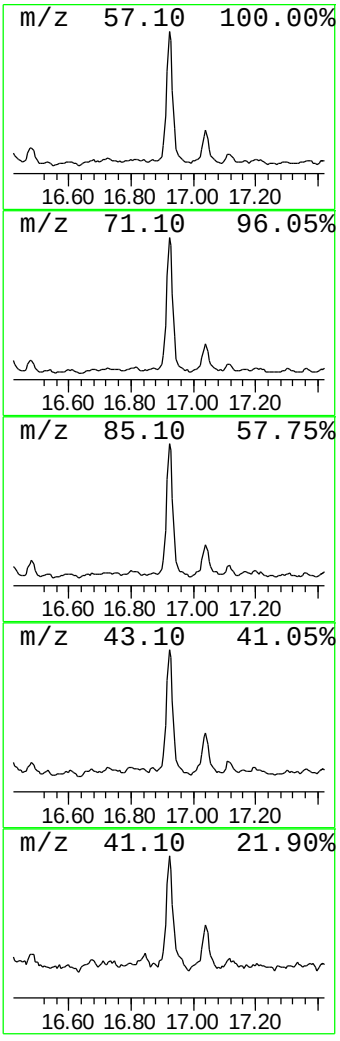
**********

Peak Number 15 Tridecane, 5-propyl- Concentration Rank 5

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.92 | 7.87 ng | 1389430 | Phenanthrene-d10 | 17.07 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Tridecane, 5-propyl- | 226 | C16H34 | 055045-11-9 | 91 |
| 2 | Octacosane | 394 | C28H58 | 000630-02-4 | 87 |
| 3 | 2-Bromotetradecane | 276 | C14H29Br | 074036-95-6 | 86 |
| 4 | Hexadecane | 226 | C16H34 | 000544-76-3 | 83 |
| 5 | Dodecane, 2-methyl-6-propyl- | 226 | C16H34 | 055045-08-4 | 81 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP081820\
Data File : BP003025.D
Acq On : 18 Aug 2020 19:40
Operator : CG/JU
Sample : L3696-06
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :** BNA_P
**ClientSampleId :** W-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP073020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

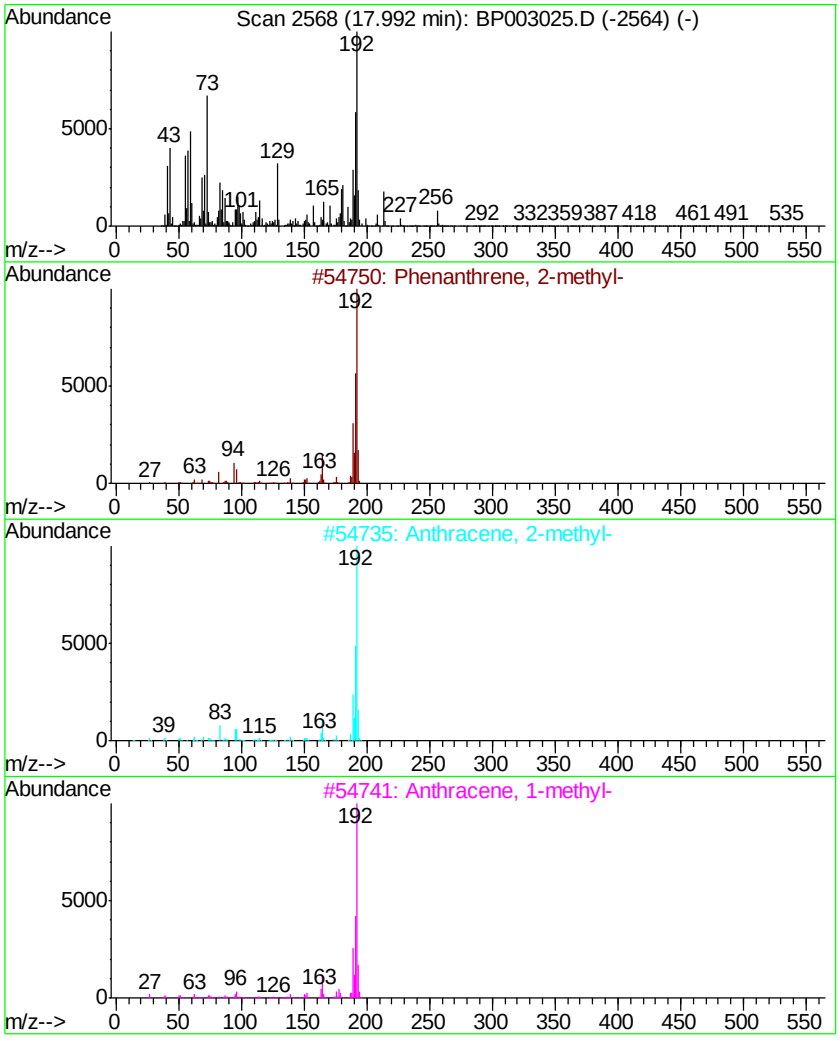
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

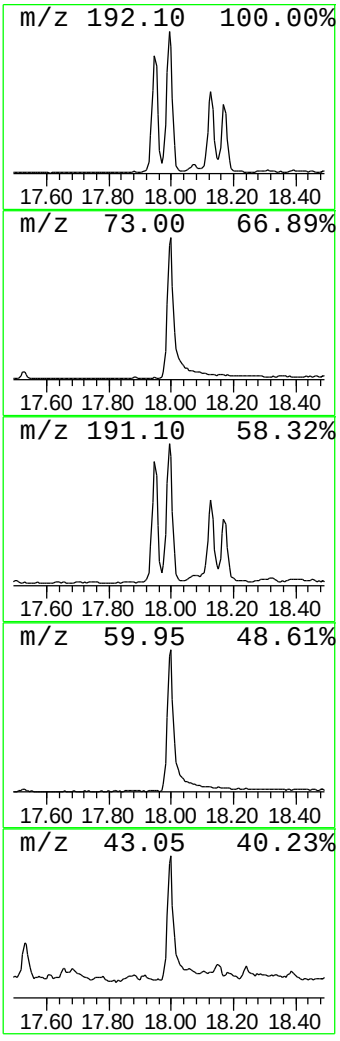
**********

Peak Number 16 Phenanthrene, 2-methyl- Concentration Rank 3

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 17.99 | 11.73 ng | 2070970 | Phenanthrene-d10 | 17.07 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenanthrene, 2-methyl- | 192 | C15H12 | 002531-84-2 | 95 |
| 2 | Anthracene, 2-methyl- | 192 | C15H12 | 000613-12-7 | 90 |
| 3 | Anthracene, 1-methyl- | 192 | C15H12 | 000610-48-0 | 90 |
| 4 | Phenanthrene, 1-methyl- | 192 | C15H12 | 000832-69-9 | 89 |
| 5 | 5H-Dibenzo[a,d]cycloheptene | 192 | C15H12 | 000256-81-5 | 86 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP081820\
Data File : BP003025.D
Acq On : 18 Aug 2020 19:40
Operator : CG/JU
Sample : L3696-06
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :** BNA_P
**ClientSampleId :** W-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP073020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

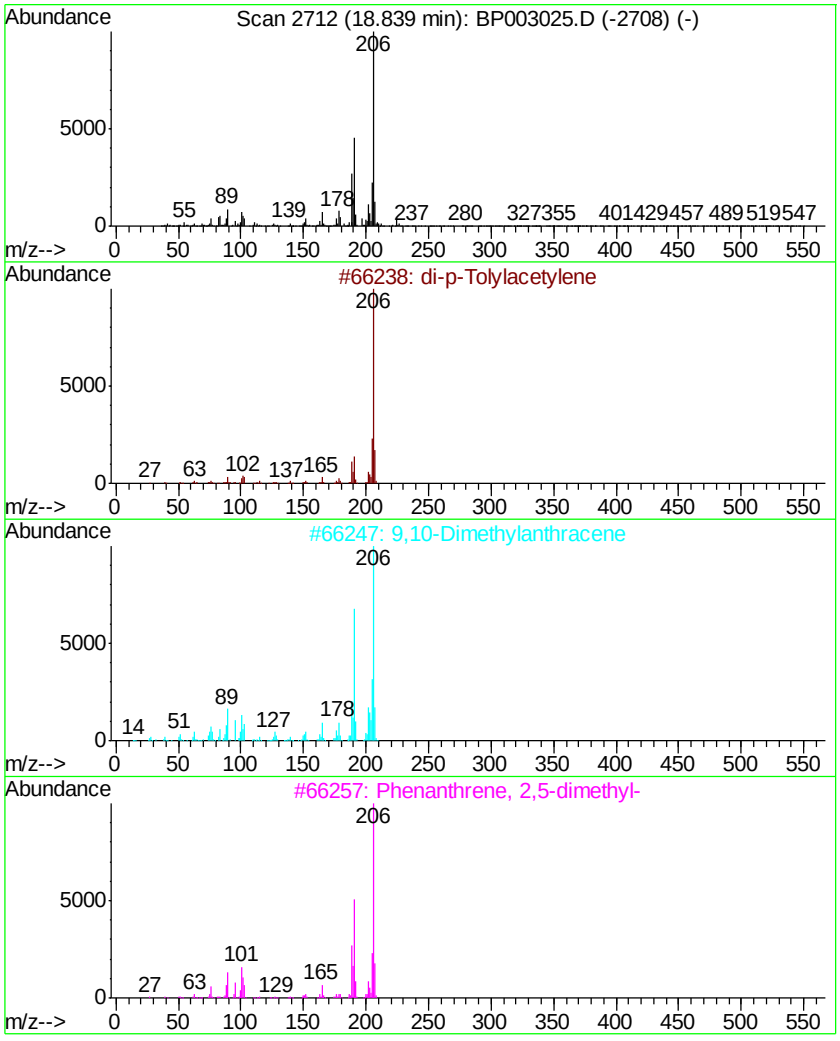
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

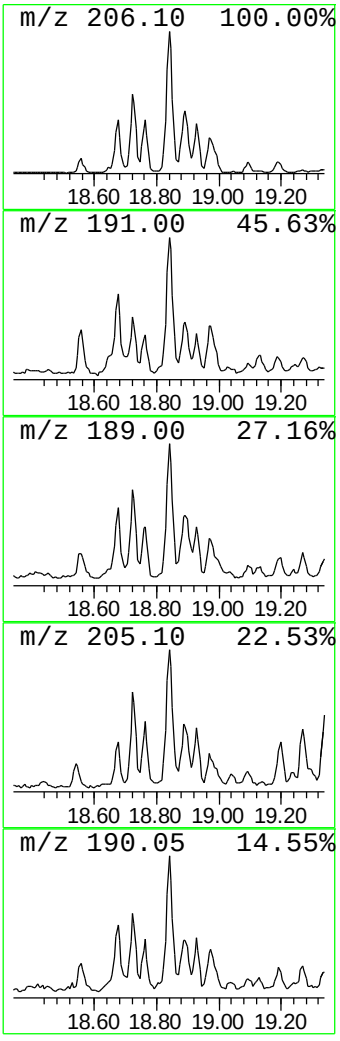
**********

Peak Number 17 di-p-Tolylacetylene Concentration Rank 13

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.84 | 3.56 ng | 628666 | Phenanthrene-d10 | 17.07 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | di-p-Tolylacetylene | 206 | C16H14 | 002789-88-0 | 91 |
| 2 | 9,10-Dimethylanthracene | 206 | C16H14 | 000781-43-1 | 91 |
| 3 | Phenanthrene, 2,5-dimethyl- | 206 | C16H14 | 003674-66-6 | 87 |
| 4 | Phenanthrene, 1,7-dimethyl- | 206 | C16H14 | 000483-87-4 | 86 |
| 5 | Anthracene, 1,4-dimethyl- | 206 | C16H14 | 000781-92-0 | 86 |



Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP081820\
Data File : BP003025.D
Acq On : 18 Aug 2020 19:40
Operator : CG/JU
Sample : L3696-06
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :** BNA_P
**ClientSampleId :** W-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP073020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

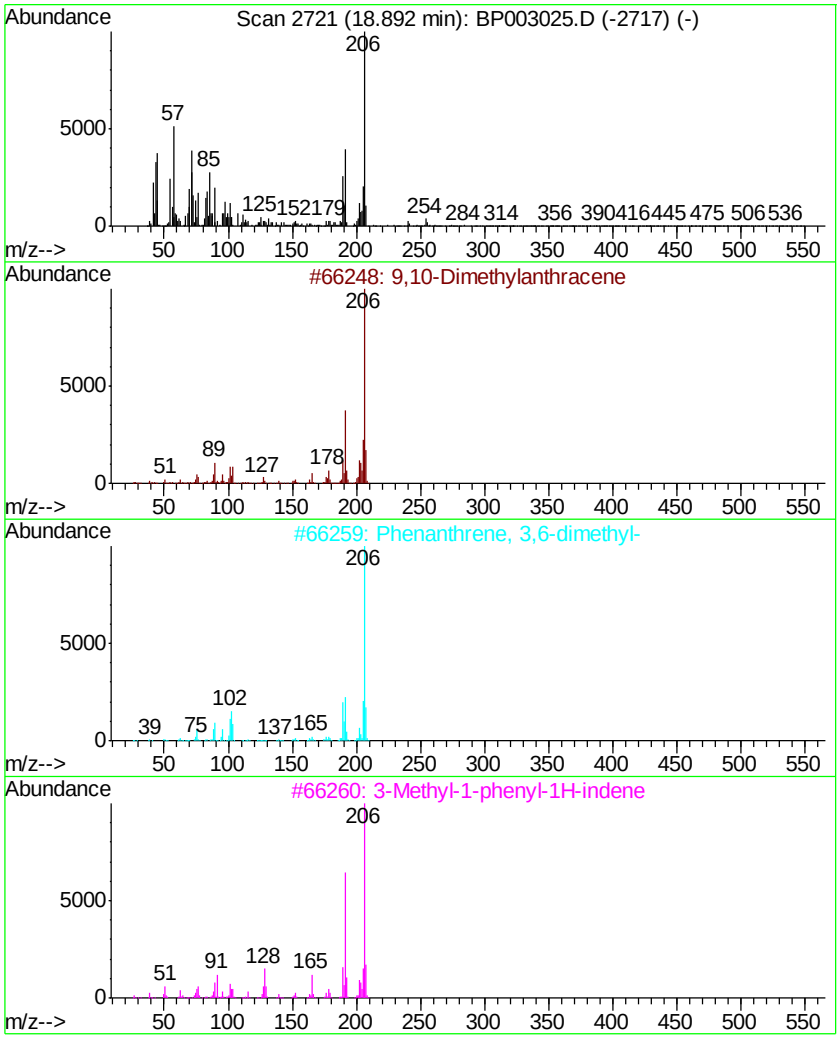
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

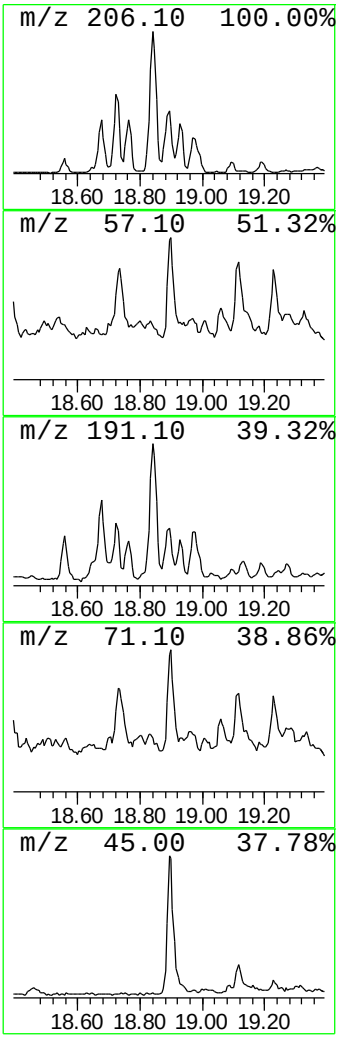
**********

Peak Number 18 9,10-Dimethylanthracene Concentration Rank 17

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.89 | 2.99 ng | 526981 | Phenanthrene-d10 | 17.07 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 9,10-Dimethylanthracene | 206 | C16H14 | 000781-43-1 | 60 |
| 2 | Phenanthrene, 3,6-dimethyl- | 206 | C16H14 | 001576-67-6 | 53 |
| 3 | 3-Methyl-1-phenyl-1H-indene | 206 | C16H14 | 022360-63-0 | 50 |
| 4 | Phenanthrene, 2,5-dimethyl- | 206 | C16H14 | 003674-66-6 | 49 |
| 5 | Anthracene, 1,4-dimethyl- | 206 | C16H14 | 000781-92-0 | 49 |



# Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP081820\
Data File : BP003025.D
Acq On : 18 Aug 2020 19:40
Operator : CG/JU
Sample : L3696-06
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :** BNA_P
**ClientSampleId :** W-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP073020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

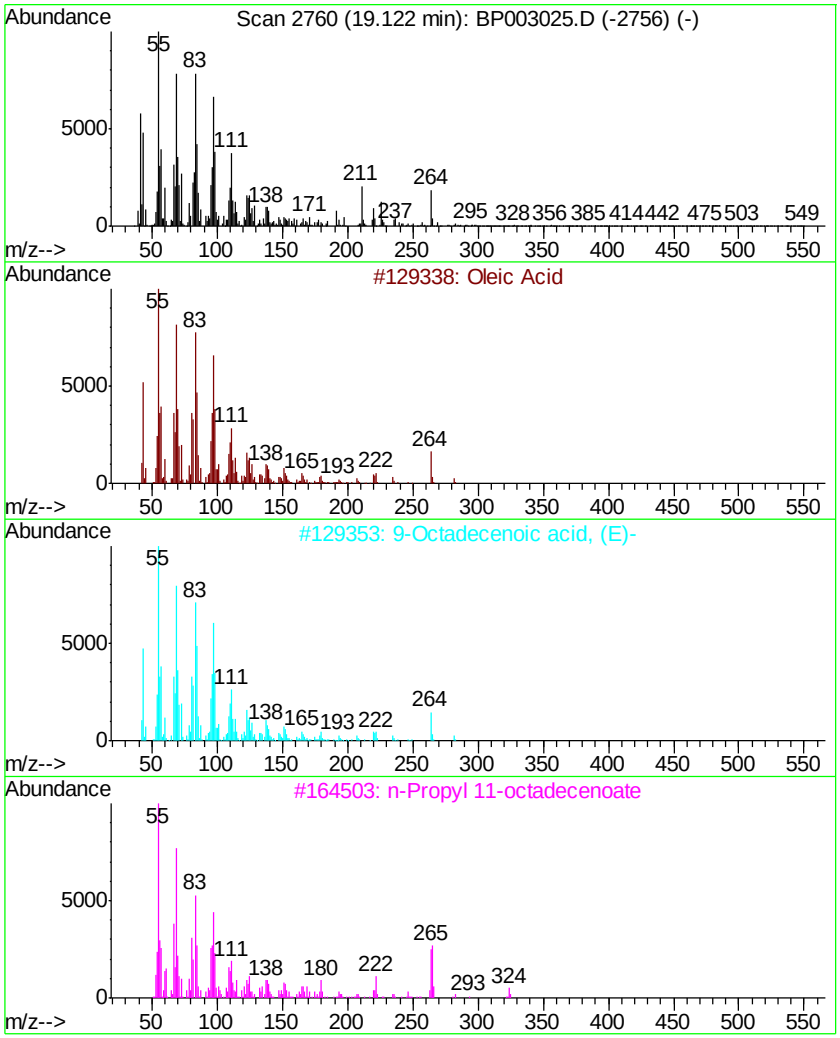
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

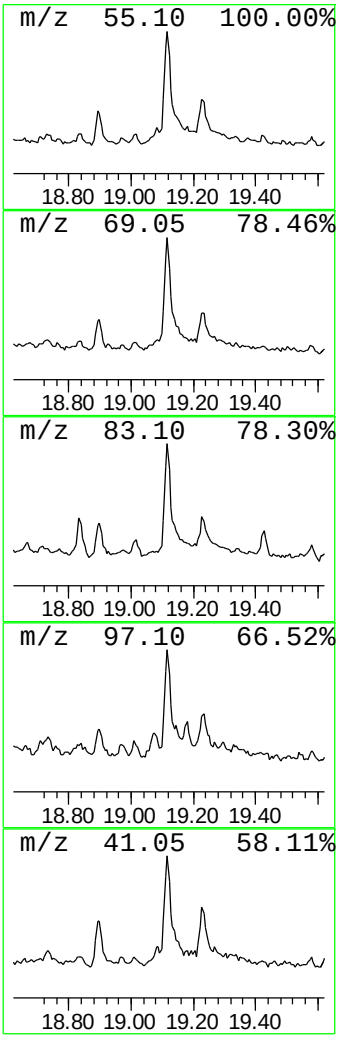
**********

Peak Number 19 Oleic Acid Concentration Rank 15

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 19.12 | 3.02 ng | 533703 | Phenanthrene-d10 | 17.07 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Oleic Acid | 282 | C18H34O2 | 000112-80-1 | 94 |
| 2 | 9-Octadecenoic acid, (E)- | 282 | C18H34O2 | 000112-79-8 | 94 |
| 3 | n-Propyl 11-octadecenoate | 324 | C21H40O2 | 1000336-71-7 | 94 |
| 4 | Octadec-9-enoic acid | 282 | C18H34O2 | 1000190-13-7 | 91 |
| 5 | 9-Nonadecene | 266 | C19H38 | 031035-07-1 | 90 |



# Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP081820\
Data File : BP003025.D
Acq On : 18 Aug 2020 19:40
Operator : CG/JU
Sample : L3696-06
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :** BNA_P
**ClientSampleId :** W-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP073020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

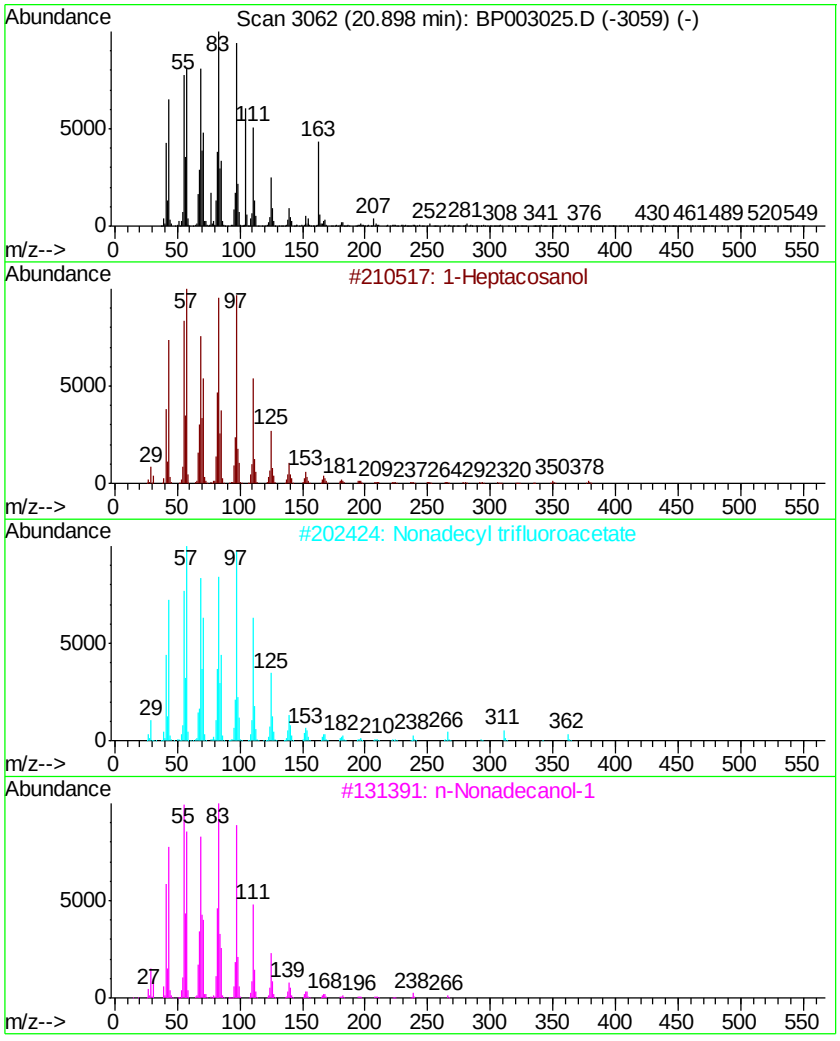
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

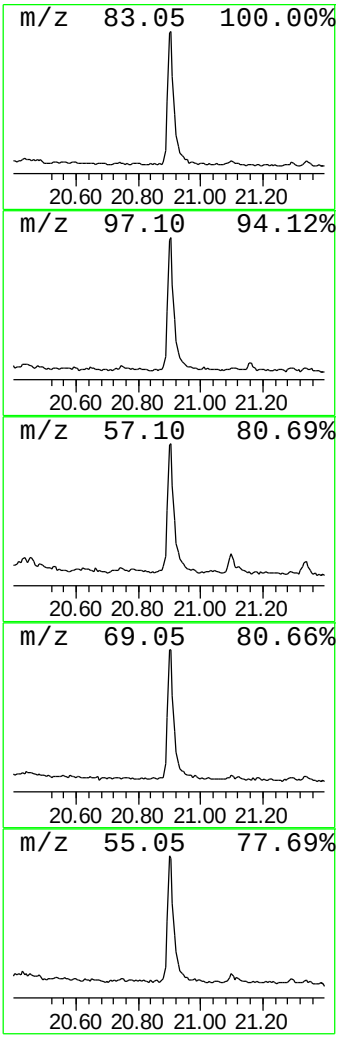
**********

Peak Number 20 1-Heptacosanol Concentration Rank 8

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 20.90 | 4.72 ng | 1098510 | Chrysene-d12 | 21.16 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1-Heptacosanol | 396 | C27H56O | 002004-39-9 | 93 |
| 2 | Nonadecyl trifluoroacetate | 380 | C21H39F3O2 | 1000351-76-3 | 93 |
| 3 | n-Nonadecanol-1 | 284 | C19H40O | 001454-84-8 | 93 |
| 4 | n-Tetracosanol-1 | 354 | C24H50O | 000506-51-4 | 93 |
| 5 | Trifluoroacetoxy hexadecane | 338 | C18H33F3O2 | 006222-03-3 | 90 |



Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP081820\
Data File : BP003025.D
Acq On : 18 Aug 2020 19:40
Operator : CG/JU
Sample : L3696-06
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :** BNA_P
**ClientSampleId :** W-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP073020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

| | | | | | Internal Standard | | | |
|---|---|---|---|---|---|---|---|---|
| TIC Top Hit name | RT | EstConc | Units | Response | # | RT | Resp | Conc |
| Butane, 2-methoxy... | 2.95 | 45.0 | ng | 3757400 | 1 | 7.68 | 1668750 | 20.0 |
| Benzene, 1,2,3-tr... | 7.80 | 2.9 | ng | 239071 | 1 | 7.68 | 1668750 | 20.0 |
| 1H-Indene, 2,3-di... | 9.88 | 4.5 | ng | 587962 | 2 | 10.46 | 2630950 | 20.0 |
| Octane, 2,6-dimet... | 11.53 | 2.9 | ng | 379082 | 2 | 10.46 | 2630950 | 20.0 |
| Naphthalene, 1,2,... | 11.66 | 4.3 | ng | 569342 | 2 | 10.46 | 2630950 | 20.0 |
| Naphthalene, 1,2,... | 12.02 | 2.5 | ng | 331624 | 2 | 10.46 | 2630950 | 20.0 |
| Naphthalene, 1,6-... | 13.48 | 4.1 | ng | 1301330 | 3 | 14.32 | 6335390 | 20.0 |
| Naphthalene, 2,6-... | 13.64 | 4.7 | ng | 1499450 | 3 | 14.32 | 6335390 | 20.0 |
| Naphthalene, 1,4-... | 13.70 | 3.0 | ng | 955714 | 3 | 14.32 | 6335390 | 20.0 |
| 1-Tridecene | 14.99 | 6.9 | ng | 2179650 | 3 | 14.32 | 6335390 | 20.0 |
| Pentadecane, 2,6,... | 15.62 | 3.3 | ng | 1053790 | 3 | 14.32 | 6335390 | 20.0 |
| 2-Octanone | 16.11 | 15.3 | ng | 2693950 | 4 | 17.07 | 3530730 | 20.0 |
| Tuaminoheptane | 16.14 | 8.3 | ng | 1471820 | 4 | 17.07 | 3530730 | 20.0 |
| 9H-Fluorene, 2-me... | 16.43 | 4.2 | ng | 737131 | 4 | 17.07 | 3530730 | 20.0 |
| Tridecane, 5-propyl- | 16.92 | 7.9 | ng | 1389430 | 4 | 17.07 | 3530730 | 20.0 |
| Phenanthrene, 2-m... | 17.99 | 11.7 | ng | 2070970 | 4 | 17.07 | 3530730 | 20.0 |
| di-p-Tolylacetylene | 18.84 | 3.6 | ng | 628666 | 4 | 17.07 | 3530730 | 20.0 |
| 9,10-Dimethylanth... | 18.89 | 3.0 | ng | 526981 | 4 | 17.07 | 3530730 | 20.0 |
| Oleic Acid | 19.12 | 3.0 | ng | 533703 | 4 | 17.07 | 3530730 | 20.0 |
| 1-Heptacosanol | 20.90 | 4.7 | ng | 1098510 | 5 | 21.16 | 4651290 | 20.0 |