LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
Data File : BP017348.D
Acq On : 26 Sep 2023 07:11
Operator : MA/JU
Sample : 04472-08
Misc :
ALS Vial : 40 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EZYE8

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:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BP017348.D\data.ms

| peak # | R.T. min | first scan | max scan | last scan | PK TY | peak height | corr. area | corr. % max. | % of total |
|---|---|---|---|---|---|---|---|---|---|
| 1 | 3.299 | 32 | 36 | 43 | rVB | 28817 | 39686 | 1.09% | 0.073% |
| 2 | 3.740 | 107 | 111 | 124 | rVV | 105469 | 158557 | 4.35% | 0.290% |
| 3 | 3.840 | 124 | 128 | 134 | rVV | 232306 | 303158 | 8.31% | 0.555% |
| 4 | 3.922 | 137 | 142 | 151 | rVB2 | 121352 | 182013 | 4.99% | 0.333% |
| 5 | 4.040 | 158 | 162 | 173 | rVB3 | 27911 | 46039 | 1.26% | 0.084% |
| 6 | 4.181 | 183 | 186 | 191 | rVB4 | 30804 | 42386 | 1.16% | 0.078% |
| 7 | 4.234 | 192 | 195 | 202 | rVB3 | 22294 | 28546 | 0.78% | 0.052% |
| 8 | 4.422 | 223 | 227 | 233 | rVB2 | 29471 | 44580 | 1.22% | 0.082% |
| 9 | 4.622 | 257 | 261 | 267 | rVB | 37682 | 50655 | 1.39% | 0.093% |
| 10 | 4.969 | 316 | 320 | 325 | rBV | 58702 | 81851 | 2.24% | 0.150% |
| 11 | 5.099 | 338 | 342 | 350 | rVV | 34968 | 51408 | 1.41% | 0.094% |
| 12 | 5.534 | 408 | 416 | 423 | rBV2 | 52077 | 96536 | 2.65% | 0.177% |
| 13 | 5.875 | 465 | 474 | 481 | rBV3 | 38351 | 65782 | 1.80% | 0.120% |
| 14 | 6.334 | 545 | 552 | 556 | rBV | 29303 | 45010 | 1.23% | 0.082% |
| 15 | 6.893 | 642 | 647 | 653 | rVB6 | 15003 | 31160 | 0.85% | 0.057% |
| 16 | 7.081 | 672 | 679 | 693 | rBV3 | 532464 | 1030892 | 28.27% | 1.888% |
| 17 | 7.269 | 704 | 711 | 719 | rBV | 374293 | 580132 | 15.91% | 1.063% |
| 18 | 7.446 | 734 | 741 | 748 | rBV | 421584 | 657320 | 18.02% | 1.204% |
| 19 | 7.522 | 750 | 754 | 762 | rVV2 | 22941 | 47661 | 1.31% | 0.087% |
| 20 | 7.922 | 814 | 822 | 831 | rBV | 942552 | 1455933 | 39.92% | 2.667% |
| 21 | 8.640 | 937 | 944 | 951 | rBV | 137554 | 242655 | 6.65% | 0.444% |
| 22 | 8.775 | 960 | 967 | 975 | rVB | 494872 | 763898 | 20.95% | 1.399% |
| 23 | 8.910 | 984 | 990 | 1001 | rBV | 37718 | 81259 | 2.23% | 0.149% |
| 24 | 9.122 | 1020 | 1026 | 1032 | rVV | 426755 | 663971 | 18.21% | 1.216% |
| 25 | 9.204 | 1036 | 1040 | 1045 | rVV | 26255 | 37747 | 1.04% | 0.069% |
| 26 | 9.834 | 1138 | 1147 | 1163 | rBV | 269931 | 469046 | 12.86% | 0.859% |
| 27 | 10.010 | 1166 | 1177 | 1188 | rBV | 207983 | 510828 | 14.01% | 0.936% |
| 28 | 10.204 | 1203 | 1210 | 1217 | rBV | 97943 | 163919 | 4.49% | 0.300% |
| 29 | 10.363 | 1231 | 1237 | 1245 | rVV | 404041 | 671906 | 18.42% | 1.231% |
| 30 | 10.634 | 1279 | 1283 | 1289 | rVB | 24322 | 38952 | 1.07% | 0.071% |
| 31 | 10.746 | 1295 | 1302 | 1309 | rBV | 1304555 | 2144069 | 58.79% | 3.927% |
| 32 | 10.799 | 1309 | 1311 | 1318 | rVB3 | 24400 | 40879 | 1.12% | 0.075% |
| 33 | 10.916 | 1323 | 1331 | 1342 | rBV | 144862 | 288277 | 7.90% | 0.528% |
| 34 | 11.304 | 1390 | 1397 | 1406 | rBV6 | 37189 | 107017 | 2.93% | 0.196% |
| 35 | 11.569 | 1436 | 1442 | 1451 | rVV | 40259 | 79316 | 2.17% | 0.145% |
| 36 | 11.681 | 1455 | 1461 | 1467 | rBV4 | 14276 | 28080 | 0.77% | 0.051% |

LSC Area Percent Report

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Operator : MA/JU
Sample : 04472-08
Misc :
ALS Vial : 40 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EZYE8

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:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Title : SVOA CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 37 | 12.087 | 1524 | 1530 | 1535 | rBV | 37218 | 56252 | 1.54% | 0.103% |
| 38 | 12.146 | 1535 | 1540 | 1547 | rVB | 25743 | 47147 | 1.29% | 0.086% |
| 39 | 12.404 | 1578 | 1584 | 1589 | rBV2 | 23438 | 37350 | 1.02% | 0.068% |
| 40 | 12.622 | 1617 | 1621 | 1628 | rVB | 16432 | 26908 | 0.74% | 0.049% |
| 41 | 12.904 | 1662 | 1669 | 1676 | rBV | 9051 | 21636 | 0.59% | 0.040% |
| 42 | 13.034 | 1687 | 1691 | 1699 | rVB4 | 20913 | 31726 | 0.87% | 0.058% |
| 43 | 13.334 | 1732 | 1742 | 1747 | rBV | 53106 | 81523 | 2.24% | 0.149% |
| 44 | 13.428 | 1751 | 1758 | 1767 | rVV2 | 68337 | 125660 | 3.45% | 0.230% |
| 45 | 13.757 | 1812 | 1814 | 1820 | rVV6 | 14657 | 22745 | 0.62% | 0.042% |
| 46 | 13.898 | 1832 | 1838 | 1842 | rBV5 | 16116 | 29136 | 0.80% | 0.053% |
| 47 | 13.998 | 1849 | 1855 | 1866 | rVB | 968837 | 1311575 | 35.96% | 2.402% |
| 48 | 14.287 | 1896 | 1904 | 1914 | rBV | 990944 | 1509863 | 41.40% | 2.765% |
| 49 | 14.410 | 1921 | 1925 | 1933 | rVB | 66135 | 92007 | 2.52% | 0.169% |
| 50 | 14.587 | 1946 | 1955 | 1963 | rBV | 1879032 | 2689126 | 73.74% | 4.925% |
| 51 | 14.839 | 1993 | 1998 | 2002 | rBV7 | 11926 | 24695 | 0.68% | 0.045% |
| 52 | 14.992 | 2015 | 2024 | 2027 | rVV4 | 37983 | 89174 | 2.45% | 0.163% |
| 53 | 15.028 | 2028 | 2030 | 2039 | rVV7 | 24945 | 55041 | 1.51% | 0.101% |
| 54 | 15.369 | 2084 | 2088 | 2100 | rVB | 76655 | 118328 | 3.24% | 0.217% |
| 55 | 15.581 | 2118 | 2124 | 2133 | rBV | 1403501 | 2008696 | 55.08% | 3.679% |
| 56 | 15.763 | 2152 | 2155 | 2161 | rVB | 29330 | 46577 | 1.28% | 0.085% |
| 57 | 15.928 | 2179 | 2183 | 2190 | rVB6 | 15376 | 27860 | 0.76% | 0.051% |
| 58 | 16.239 | 2231 | 2236 | 2243 | rBV2 | 118200 | 203536 | 5.58% | 0.373% |
| 59 | 16.875 | 2340 | 2344 | 2352 | rVV9 | 17736 | 38615 | 1.06% | 0.071% |
| 60 | 17.039 | 2364 | 2372 | 2375 | rBV2 | 84558 | 111897 | 3.07% | 0.205% |
| 61 | 17.081 | 2375 | 2379 | 2384 | rVB | 55016 | 77000 | 2.11% | 0.141% |
| 62 | 17.357 | 2420 | 2426 | 2431 | rBV | 2225748 | 3155321 | 86.52% | 5.779% |
| 63 | 17.398 | 2431 | 2433 | 2437 | rVV | 90342 | 107141 | 2.94% | 0.196% |
| 64 | 17.457 | 2437 | 2443 | 2454 | rVV | 1414424 | 2014756 | 55.25% | 3.690% |
| 65 | 17.786 | 2495 | 2499 | 2506 | rVB | 85701 | 104911 | 2.88% | 0.192% |
| 66 | 17.851 | 2506 | 2510 | 2512 | rBV5 | 19686 | 30731 | 0.84% | 0.056% |
| 67 | 17.975 | 2529 | 2531 | 2537 | rVB4 | 22011 | 26940 | 0.74% | 0.049% |
| 68 | 18.180 | 2562 | 2566 | 2567 | rBV | 114142 | 133429 | 3.66% | 0.244% |
| 69 | 18.345 | 2591 | 2594 | 2599 | rVB2 | 53081 | 68796 | 1.89% | 0.126% |
| 70 | 18.463 | 2610 | 2614 | 2619 | rVB3 | 78710 | 95019 | 2.61% | 0.174% |
| 71 | 18.516 | 2620 | 2623 | 2624 | rBV2 | 51820 | 53074 | 1.46% | 0.097% |
| 72 | 18.545 | 2624 | 2628 | 2632 | rVB | 148171 | 197461 | 5.41% | 0.362% |
| 73 | 18.733 | 2657 | 2660 | 2664 | rBV3 | 71909 | 94024 | 2.58% | 0.172% |
| 74 | 18.792 | 2665 | 2670 | 2676 | rVB2 | 446355 | 608442 | 16.68% | 1.114% |
| 75 | 18.910 | 2685 | 2690 | 2697 | rVB | 1009475 | 1273241 | 34.91% | 2.332% |
| 76 | 19.022 | 2705 | 2709 | 2712 | rBV | 168600 | 215908 | 5.92% | 0.395% |

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Operator : MA/JU
Sample : 04472-08
Misc :
ALS Vial : 40 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EZYE8

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:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Title : SVOA CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 77 | 19.075 | 2716 | 2718 | 2722 | rVV | 83978 | 101784 | 2.79% | 0.186% |
| 78 | 19.151 | 2726 | 2731 | 2736 | rVV | 744923 | 945948 | 25.94% | 1.733% |
| 79 | 19.257 | 2742 | 2749 | 2750 | rBV2 | 211992 | 368425 | 10.10% | 0.675% |
| 80 | 19.369 | 2764 | 2768 | 2776 | rVB | 383404 | 489586 | 13.42% | 0.897% |
| 81 | 19.433 | 2777 | 2779 | 2785 | rVB4 | 53618 | 62767 | 1.72% | 0.115% |
| 82 | 19.575 | 2799 | 2803 | 2807 | rBV | 514411 | 606187 | 16.62% | 1.110% |
| 83 | 19.639 | 2810 | 2814 | 2817 | rVV | 205462 | 241993 | 6.64% | 0.443% |
| 84 | 19.698 | 2819 | 2824 | 2835 | rVB | 1989561 | 2832316 | 77.66% | 5.188% |
| 85 | 19.886 | 2851 | 2856 | 2861 | rBV | 630743 | 794290 | 21.78% | 1.455% |
| 86 | 19.945 | 2861 | 2866 | 2869 | rVV | 744736 | 1002135 | 27.48% | 1.835% |
| 87 | 19.980 | 2869 | 2872 | 2877 | rVB | 697889 | 842502 | 23.10% | 1.543% |
| 88 | 20.098 | 2887 | 2892 | 2897 | rBV | 1677844 | 2025566 | 55.54% | 3.710% |
| 89 | 20.163 | 2899 | 2903 | 2908 | rBV4 | 281714 | 418048 | 11.46% | 0.766% |
| 90 | 20.439 | 2948 | 2950 | 2953 | rBV3 | 72754 | 72135 | 1.98% | 0.132% |
| 91 | 20.786 | 3006 | 3009 | 3014 | rVB | 400670 | 491876 | 13.49% | 0.901% |
| 92 | 20.845 | 3015 | 3019 | 3024 | rVB | 859272 | 997241 | 27.34% | 1.827% |
| 93 | 20.910 | 3025 | 3030 | 3034 | rBV | 1631728 | 1981490 | 54.33% | 3.629% |
| 94 | 21.104 | 3061 | 3063 | 3065 | rBV2 | 130581 | 125425 | 3.44% | 0.230% |
| 95 | 21.133 | 3065 | 3068 | 3079 | rVB2 | 526935 | 788698 | 21.63% | 1.445% |
| 96 | 21.251 | 3085 | 3088 | 3094 | rBV | 666423 | 834479 | 22.88% | 1.528% |
| 97 | 21.457 | 3118 | 3123 | 3128 | rBV | 2657936 | 3646946 | 100.00% | 6.680% |
| 98 | 22.010 | 3213 | 3217 | 3227 | rVB2 | 700405 | 1008144 | 27.64% | 1.846% |
| 99 | 23.804 | 3517 | 3522 | 3529 | rVB2 | 1157959 | 2120020 | 58.13% | 3.883% |
| 100 | 23.974 | 3545 | 3551 | 3559 | rVB | 1633447 | 3365232 | 92.28% | 6.164% |

Sum of corrected areas: 54597624

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
Data File : BP017348.D
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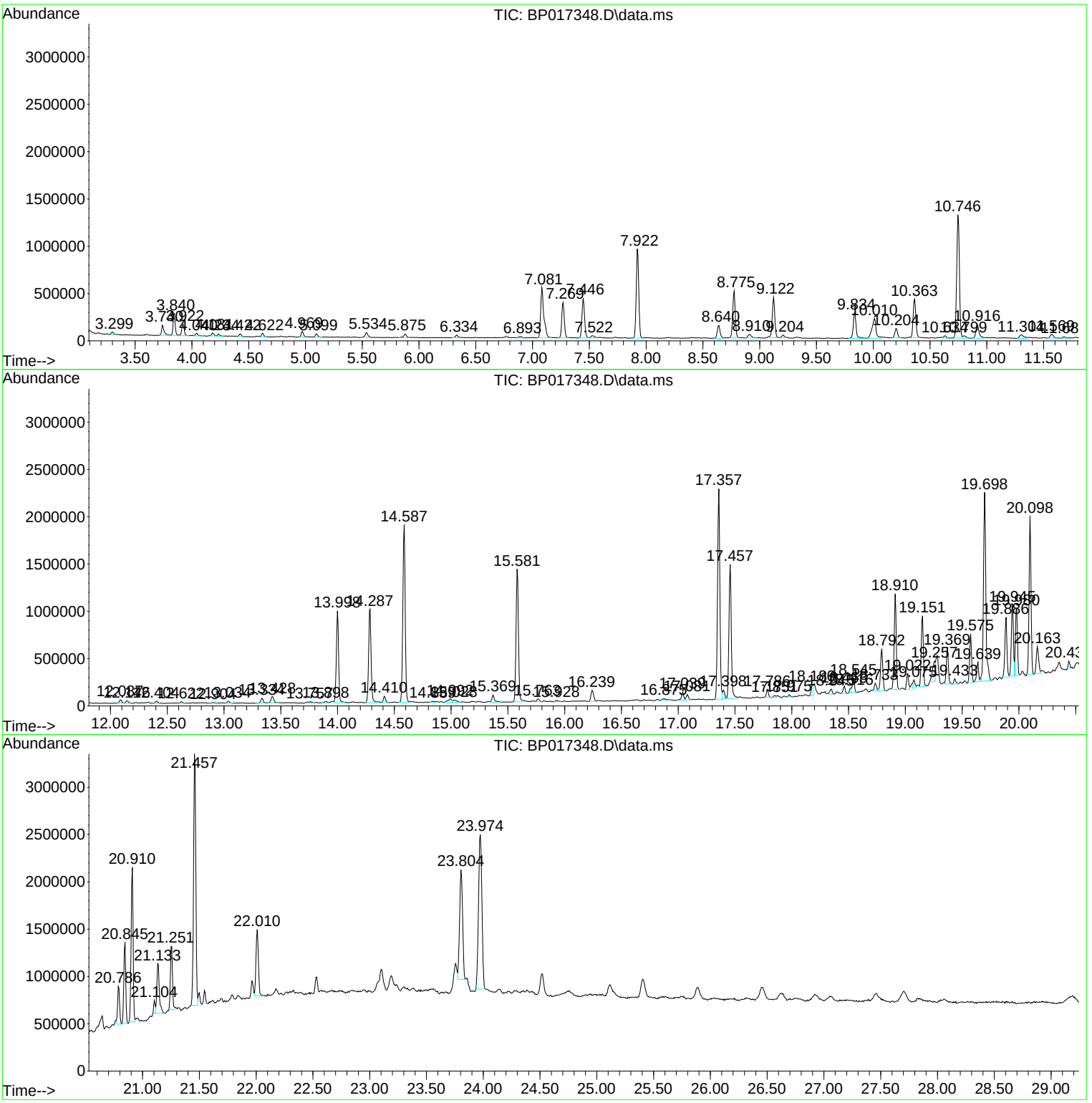
Misc :
ALS Vial : 40 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EZYE8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
Data File : BP017348.D
Acq On : 26 Sep 2023 07:11
Operator : MA/JU
Sample : 04472-08
Misc :
ALS Vial : 40 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EZYE8

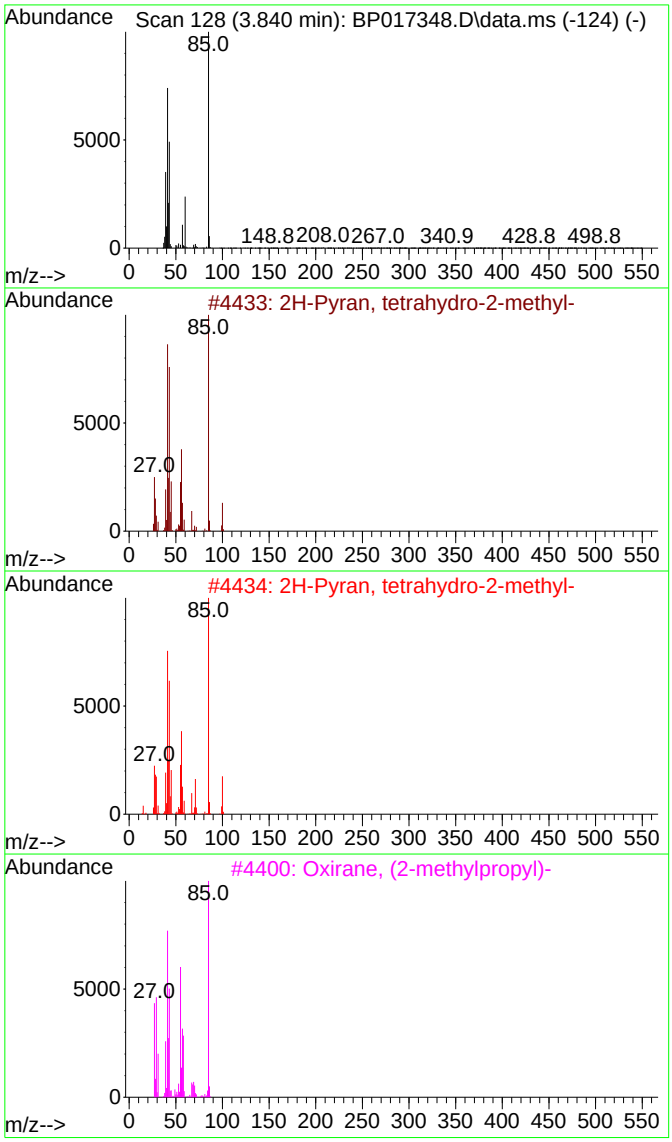
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

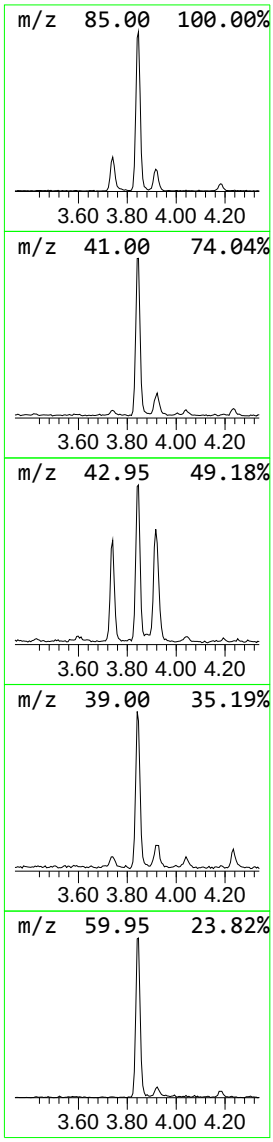
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 1 unknown-01 Concentration Rank 13

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 3.840 | 4.16 ng/ul | 303158 | 1,4-Dichlorobenzene-d4 | 7.922 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2H-Pyran, tetrahydro-2-methyl- | 100 | C6H12O | 010141-72-7 | 38 |
| 2 | 2H-Pyran, tetrahydro-2-methyl- | 100 | C6H12O | 010141-72-7 | 38 |
| 3 | Oxirane, (2-methylpropyl)- | 100 | C6H12O | 023850-78-4 | 36 |
| 4 | Methacrylamide | 85 | C4H7NO | 000079-39-0 | 33 |
| 5 | 2-Pyrrolidinone | 85 | C4H7NO | 000616-45-5 | 9 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
Data File : BP017348.D
Acq On : 26 Sep 2023 07:11
Operator : MA/JU
Sample : 04472-08
Misc :
ALS Vial : 40 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EZYE8

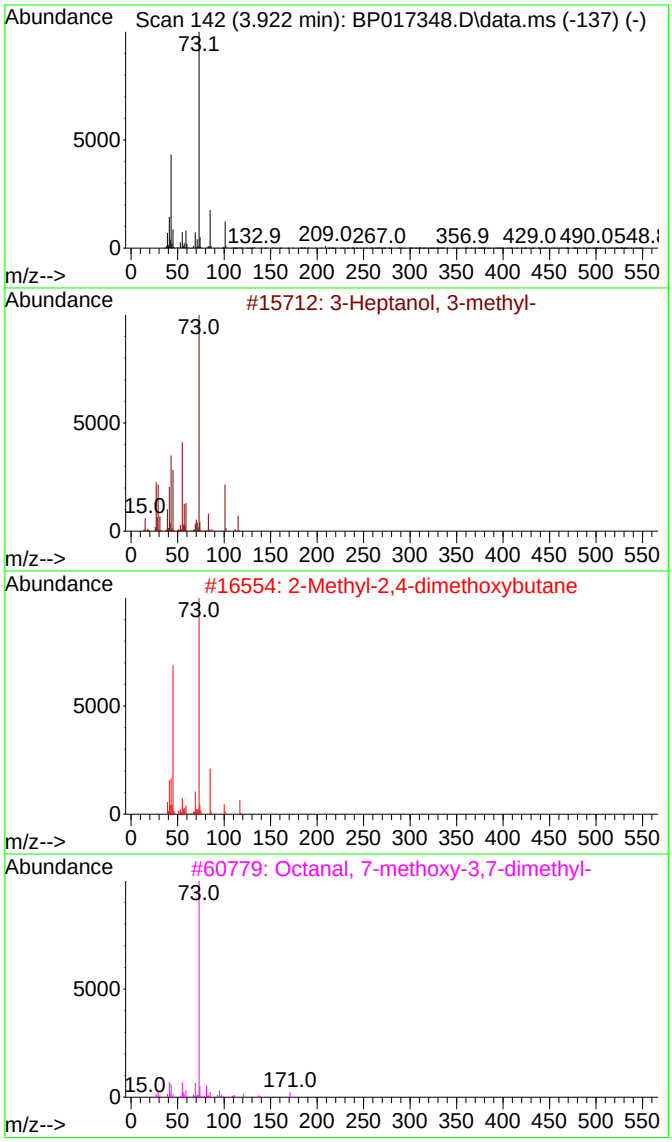
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

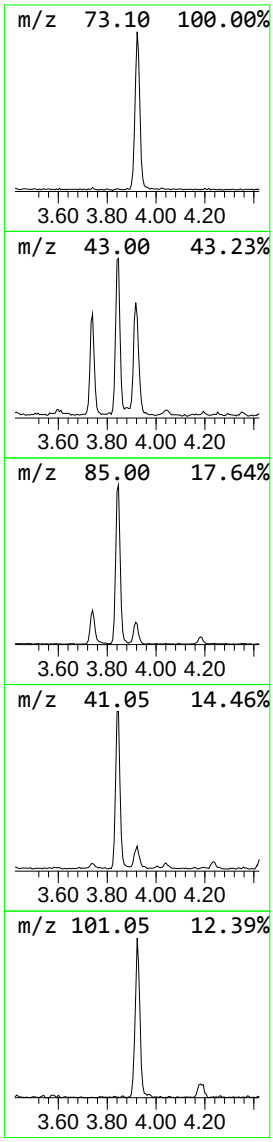
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 2 unknown-02 Concentration Rank 17

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 3.922 | 2.50 ng/ul | 182013 | 1,4-Dichlorobenzene-d4 | 7.922 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 3-Heptanol, 3-methyl- | 130 | C8H18O | 005582-82-1 | 37 |
| 2 | 2-Methyl-2,4-dimethoxybutane | 132 | C7H16O2 | 039836-89-0 | 33 |
| 3 | Octanal, 7-methoxy-3,7-dimethyl- | 186 | C11H22O2 | 003613-30-7 | 23 |
| 4 | Pentane, 3-methoxy- | 102 | C6H14O | 036839-67-5 | 17 |
| 5 | Silane, trimethyl(3-methylbutyl)- | 144 | C8H20Si | 018291-15-1 | 12 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
Data File : BP017348.D
Acq On : 26 Sep 2023 07:11
Operator : MA/JU
Sample : 04472-08
Misc :
ALS Vial : 40 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EZYE8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

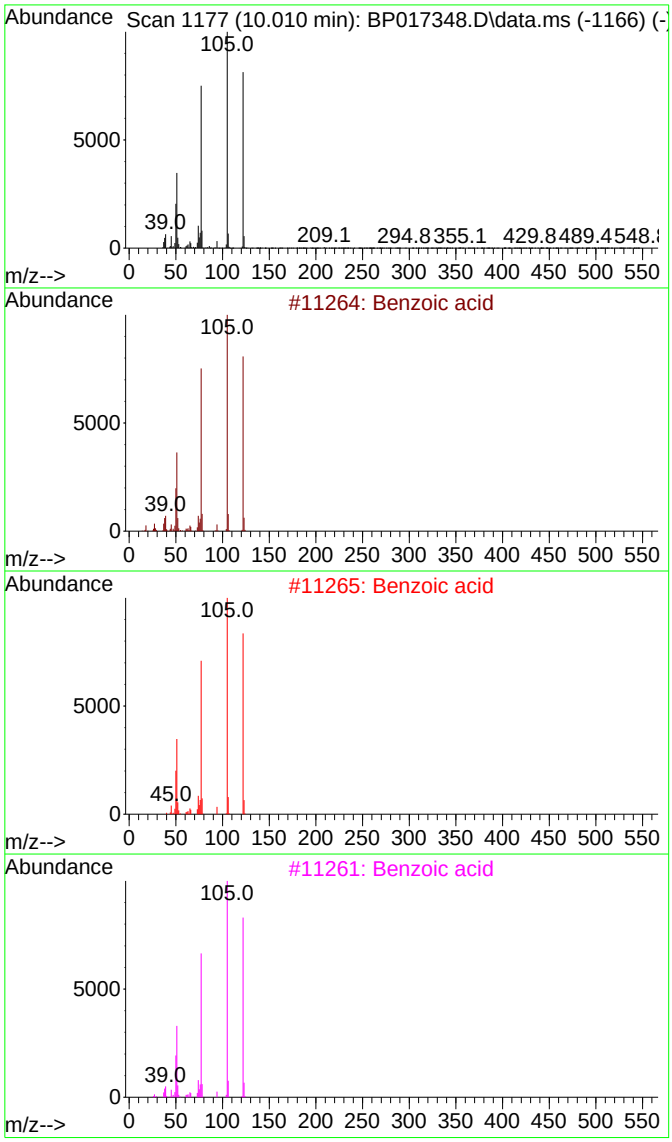
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

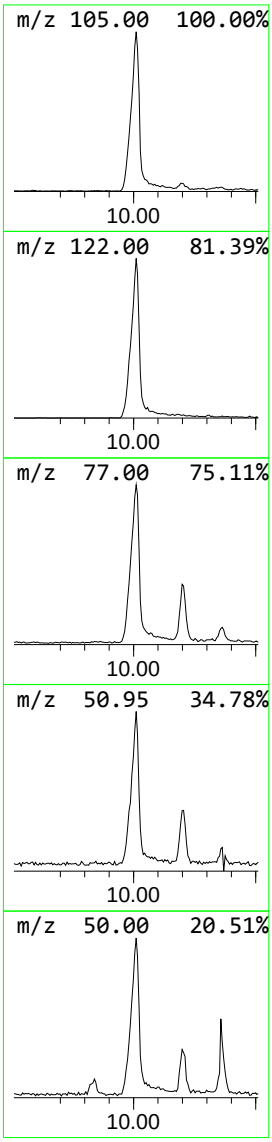
**********

Peak Number 3 Benzoic acid Concentration Rank 8

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.010 | 4.77 ng/ul | 510828 | Naphthalene-d8 | 10.746 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzoic acid | 122 | C7H6O2 | 000065-85-0 | 96 |
| 2 | Benzoic acid | 122 | C7H6O2 | 000065-85-0 | 96 |
| 3 | Benzoic acid | 122 | C7H6O2 | 000065-85-0 | 96 |
| 4 | Benzoic acid | 122 | C7H6O2 | 000065-85-0 | 94 |
| 5 | Cyclobutane-1,1-dicarboxamide, N... | 382 | C20H18N2O6 | 1000253-25-3 | 91 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
Data File : BP017348.D
Acq On : 26 Sep 2023 07:11
Operator : MA/JU
Sample : 04472-08
Misc :
ALS Vial : 40 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
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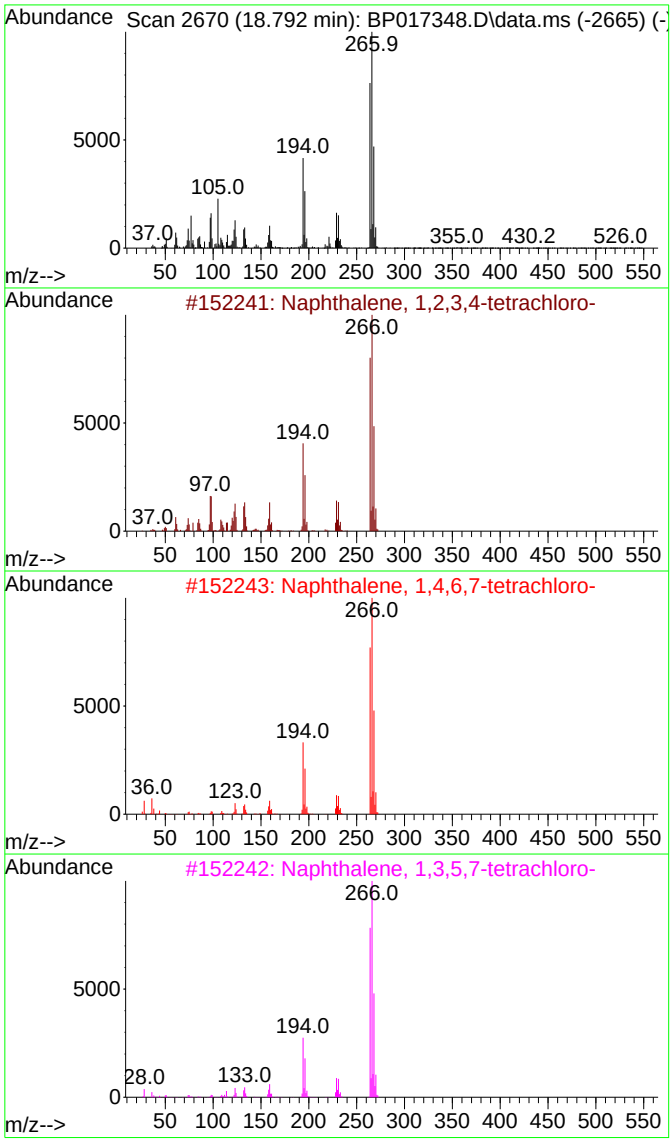
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

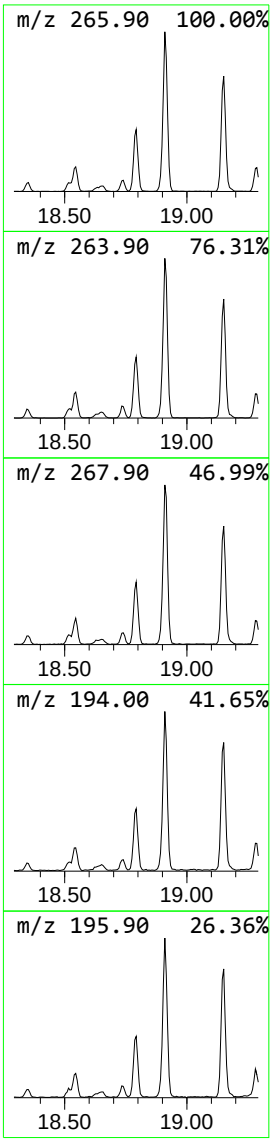
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 4 Naphthalene, 1,2,3,4-tetrac... Concentration Rank 14

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.792 | 3.86 ng/ul | 608442 | Phenanthrene-d10 | 17.357 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Naphthalene, 1,2,3,4-tetrachloro- | 264 | C10H4Cl4 | 020020-02-4 | 99 |
| 2 | Naphthalene, 1,4,6,7-tetrachloro- | 264 | C10H4Cl4 | 055720-43-9 | 99 |
| 3 | Naphthalene, 1,3,5,7-tetrachloro- | 264 | C10H4Cl4 | 053555-64-9 | 99 |
| 4 | Naphthalene, 1,2,3,4-tetrachloro- | 264 | C10H4Cl4 | 020020-02-4 | 95 |
| 5 | Naphthalene, 1,2,3,4-tetrachloro- | 264 | C10H4Cl4 | 020020-02-4 | 94 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
Data File : BP017348.D
Acq On : 26 Sep 2023 07:11
Operator : MA/JU
Sample : 04472-08
Misc :
ALS Vial : 40 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EZYE8

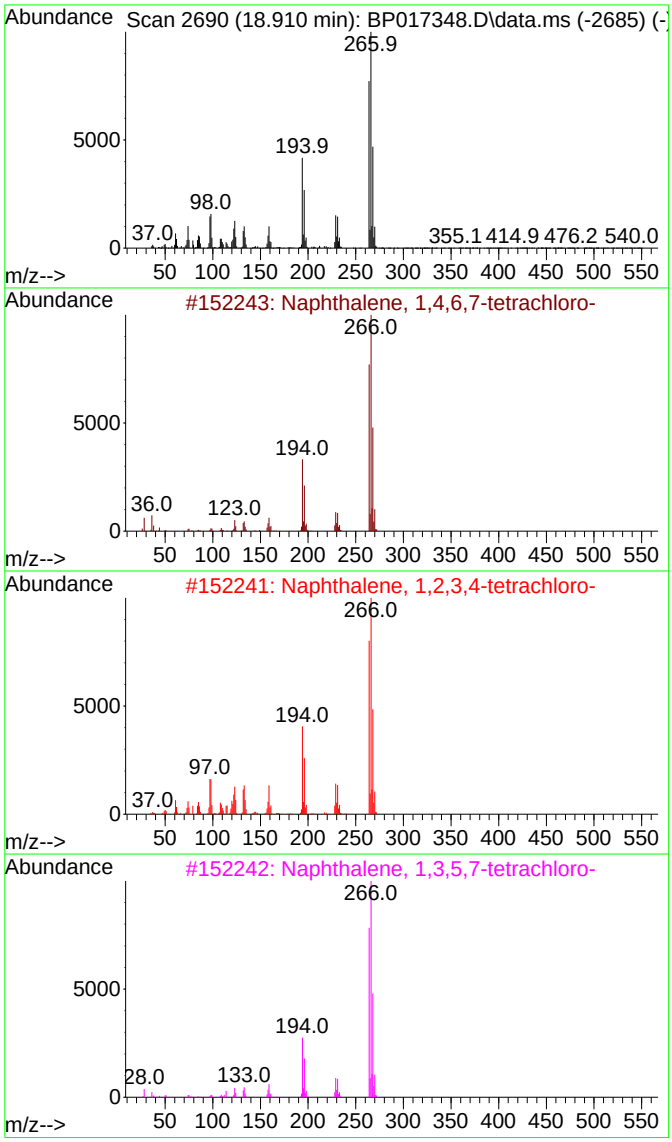
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

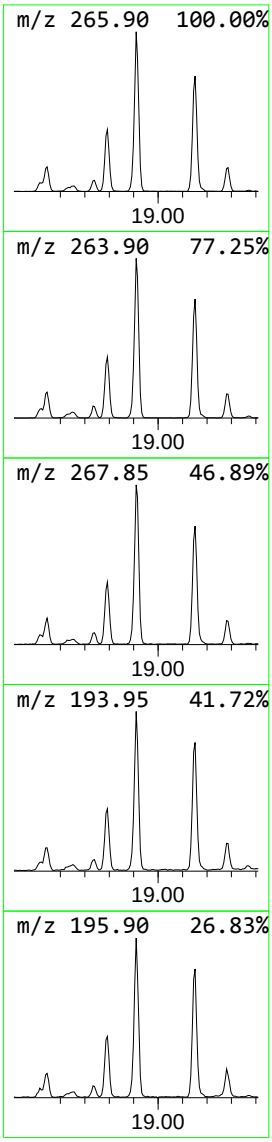
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 5 Naphthalene, 1,4,6,7-tetrac... Concentration Rank 3

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.910 | 8.07 ng/ul | 1273240 | Phenanthrene-d10 | 17.357 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Naphthalene, 1,4,6,7-tetrachloro- | 264 | C10H4Cl4 | 055720-43-9 | 99 |
| 2 | Naphthalene, 1,2,3,4-tetrachloro- | 264 | C10H4Cl4 | 020020-02-4 | 99 |
| 3 | Naphthalene, 1,3,5,7-tetrachloro- | 264 | C10H4Cl4 | 053555-64-9 | 99 |
| 4 | Naphthalene, 1,2,3,4-tetrachloro- | 264 | C10H4Cl4 | 020020-02-4 | 98 |
| 5 | Naphthalene, 1,2,3,4-tetrachloro- | 264 | C10H4Cl4 | 020020-02-4 | 95 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
Data File : BP017348.D
Acq On : 26 Sep 2023 07:11
Operator : MA/JU
Sample : 04472-08
Misc :
ALS Vial : 40 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EZYE8

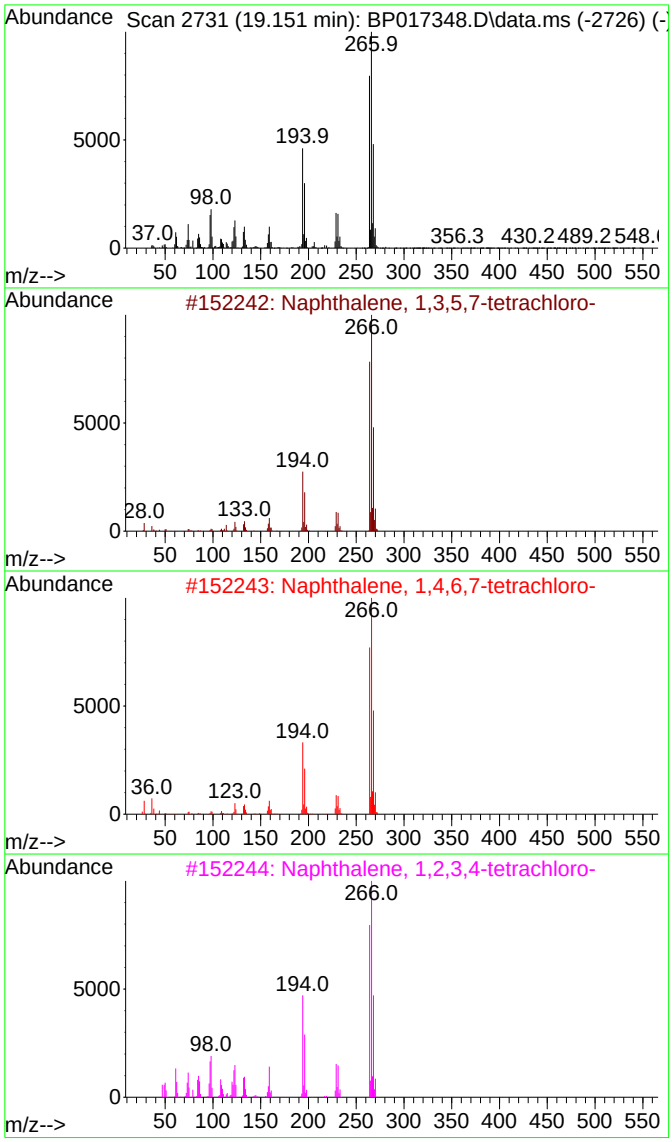
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

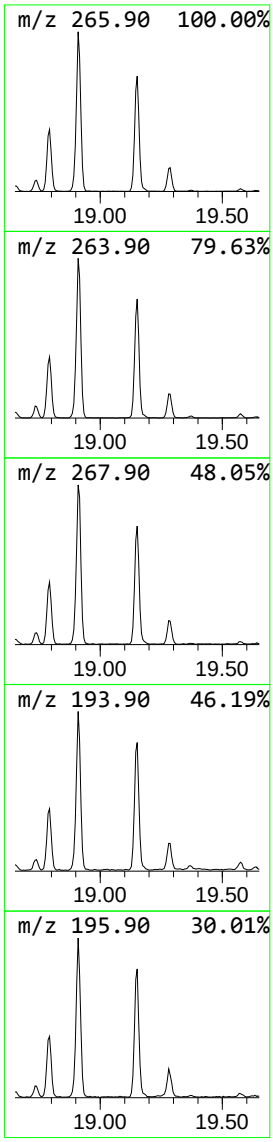
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 6 Naphthalene, 1,3,5,7-tetrac... Concentration Rank 4

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 19.151 | 6.00 ng/ul | 945948 | Phenanthrene-d10 | 17.357 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Naphthalene, 1,3,5,7-tetrachloro- | 264 | C10H4Cl4 | 053555-64-9 | 99 |
| 2 | Naphthalene, 1,4,6,7-tetrachloro- | 264 | C10H4Cl4 | 055720-43-9 | 99 |
| 3 | Naphthalene, 1,2,3,4-tetrachloro- | 264 | C10H4Cl4 | 020020-02-4 | 95 |
| 4 | Tetrachloroterephthalonitrile | 264 | C8Cl4N2 | 001897-41-2 | 95 |
| 5 | Naphthalene, 1,2,3,4-tetrachloro- | 264 | C10H4Cl4 | 020020-02-4 | 94 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
Data File : BP017348.D
Acq On : 26 Sep 2023 07:11
Operator : MA/JU
Sample : 04472-08
Misc :
ALS Vial : 40 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EZYE8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

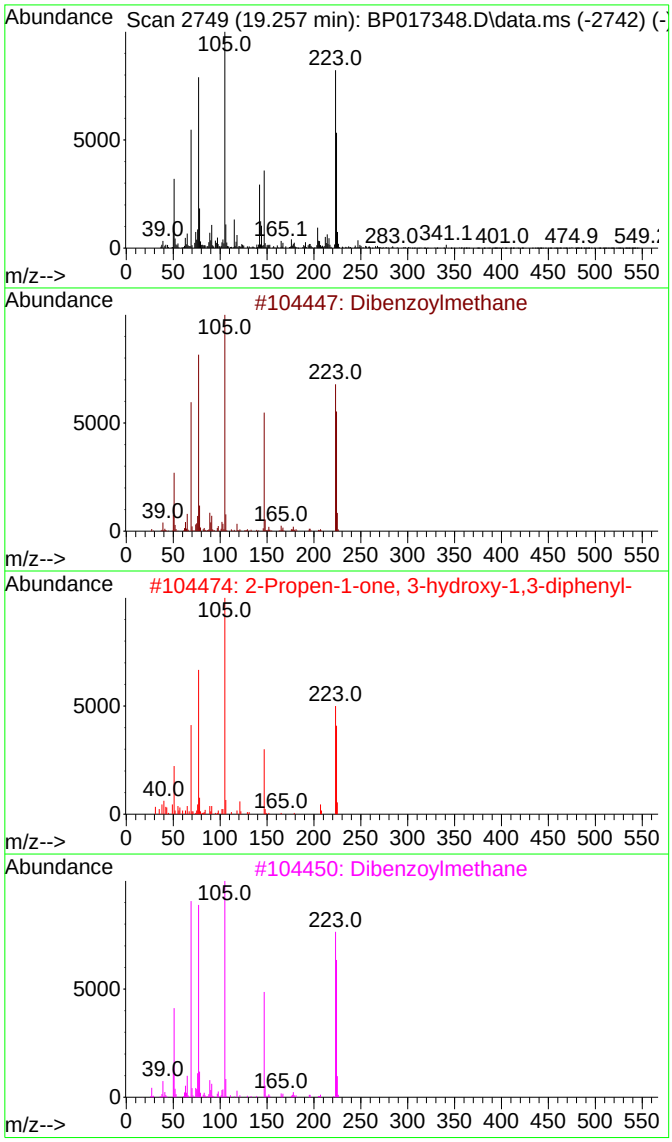
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

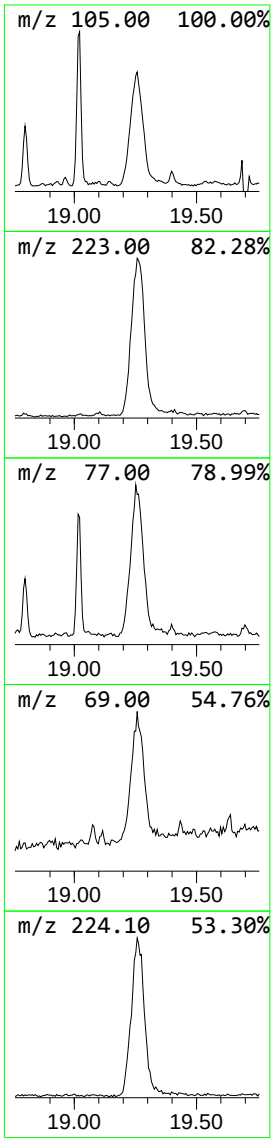
**********

Peak Number 7 Dibenzoylmethane Concentration Rank 18

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 19.257 | 2.34 ng/ul | 368425 | Phenanthrene-d10 | 17.357 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Dibenzoylmethane | 224 | C15H12O2 | 000120-46-7 | 93 |
| 2 | 2-Propen-1-one, 3-hydroxy-1,3-di... | 224 | C15H12O2 | 001704-15-0 | 90 |
| 3 | Dibenzoylmethane | 224 | C15H12O2 | 000120-46-7 | 64 |
| 4 | Dibenzoylmethane | 224 | C15H12O2 | 000120-46-7 | 64 |
| 5 | 1,3-Propanedione, 2-bromo-1,3-di... | 302 | C15H11BrO2 | 000728-84-7 | 52 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
Data File : BP017348.D
Acq On : 26 Sep 2023 07:11
Operator : MA/JU
Sample : 04472-08
Misc :
ALS Vial : 40 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EZYE8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

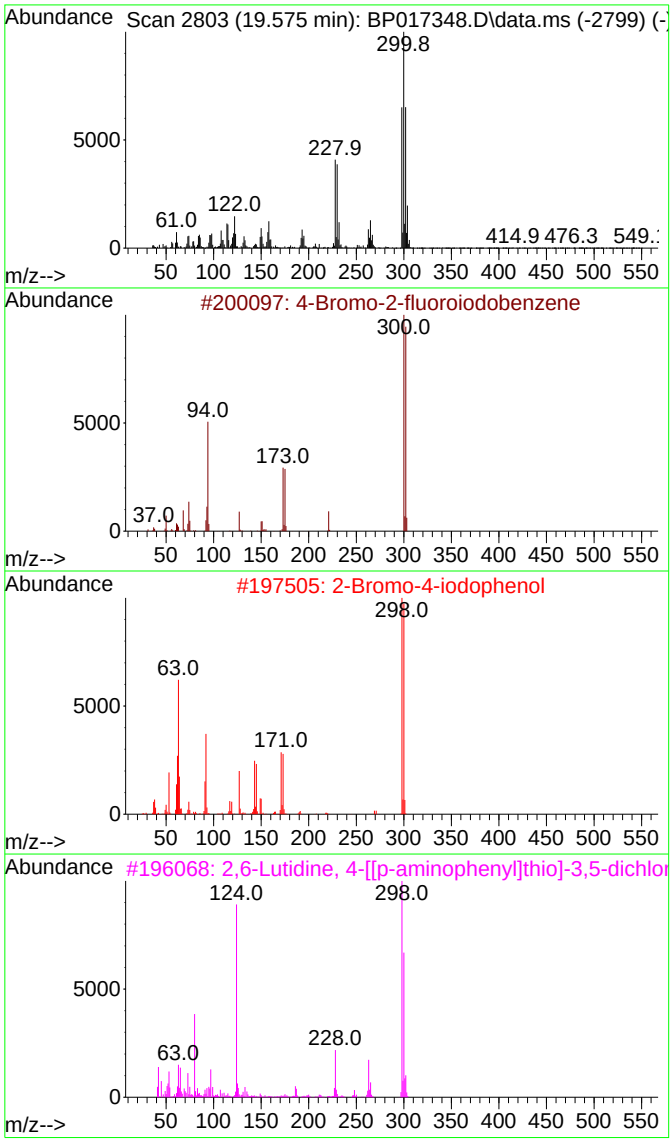
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

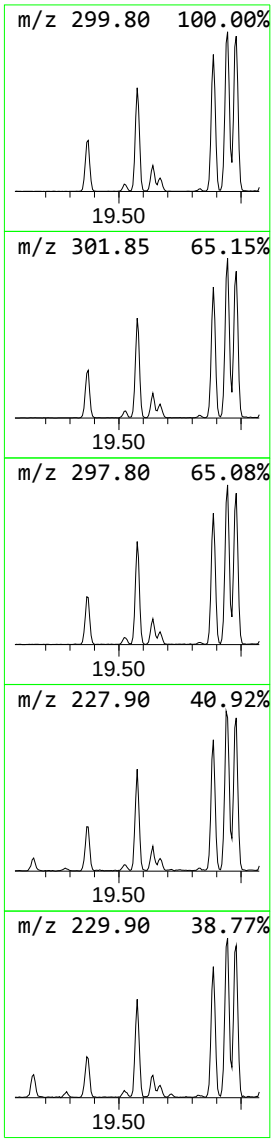
**********

Peak Number 8 unknown-03 Concentration Rank 15

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 19.575 | 3.32 ng/ul | 606187 | Chrysene-d12 | 21.457 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 4-Bromo-2-fluoroiodobenzene | 300 | C6H3BrFI | 105931-73-5 | 27 |
| 2 | 2-Bromo-4-iodophenol | 298 | C6H4BrIO | 133430-98-5 | 10 |
| 3 | 2,6-Lutidine, 4-[[p-aminophenyl]... | 298 | C13H12Cl2N2S | 1000252-27-9 | 10 |
| 4 | (E,E)-2,5-Bis[2-(2-thienyl)ethen... | 300 | C16H12S3 | 026263-68-3 | 10 |
| 5 | 3-Cyclopropyl-5-[3-(trifluoromet... | 300 | C10H7F3N6S | 1000435-68-4 | 10 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
Data File : BP017348.D
Acq On : 26 Sep 2023 07:11
Operator : MA/JU
Sample : 04472-08
Misc :
ALS Vial : 40 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EZYE8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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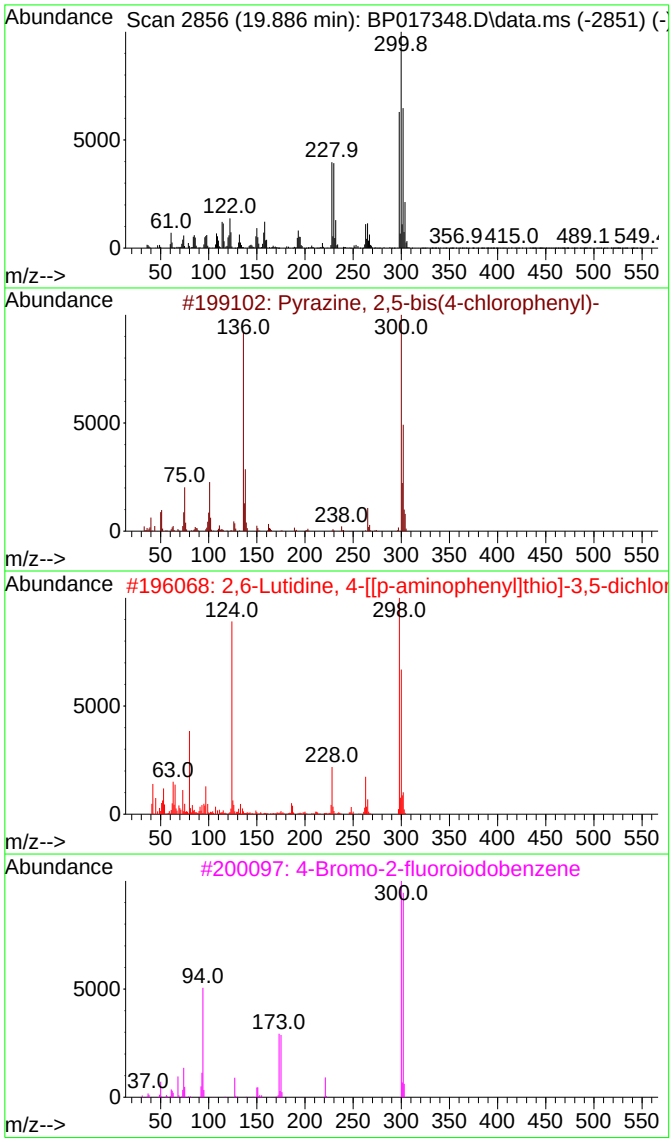
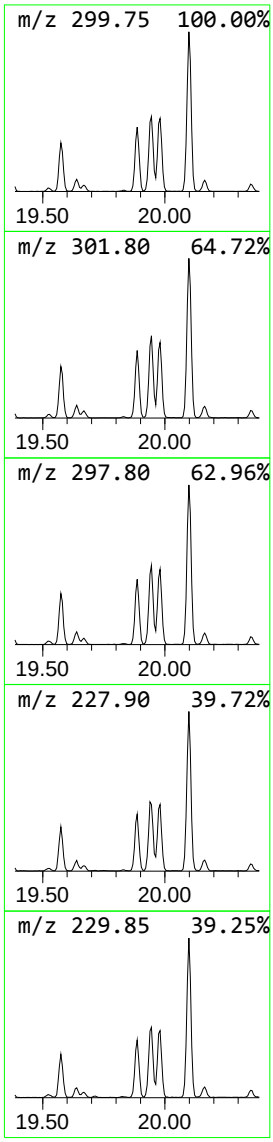
Peak Number 9 unknown-04 Concentration Rank 11

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 19.886 | 4.36 ng/ul | 794290 | Chrysene-d12 | 21.457 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Pyrazine, 2,5-bis(4-chlorophenyl)- | 300 | C16H10Cl2N2 | 074376-33-3 | 35 |
| 2 | 2,6-Lutidine, 4-[[p-aminophenyl]... | 298 | C13H12Cl2N2S | 1000252-27-9 | 22 |
| 3 | 4-Bromo-2-fluoroiodobenzene | 300 | C6H3BrFI | 105931-73-5 | 16 |
| 4 | 3-Chloro-6-methyl-1-benzothiophe... | 298 | C13H15ClO2SSi | 1000485-00-5 | 14 |
| 5 | 3-Cyclopropyl-5-[3-(trifluoromet... | 300 | C10H7F3N6S | 1000435-68-4 | 14 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
Data File : BP017348.D
Acq On : 26 Sep 2023 07:11
Operator : MA/JU
Sample : 04472-08
Misc :
ALS Vial : 40 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EZYE8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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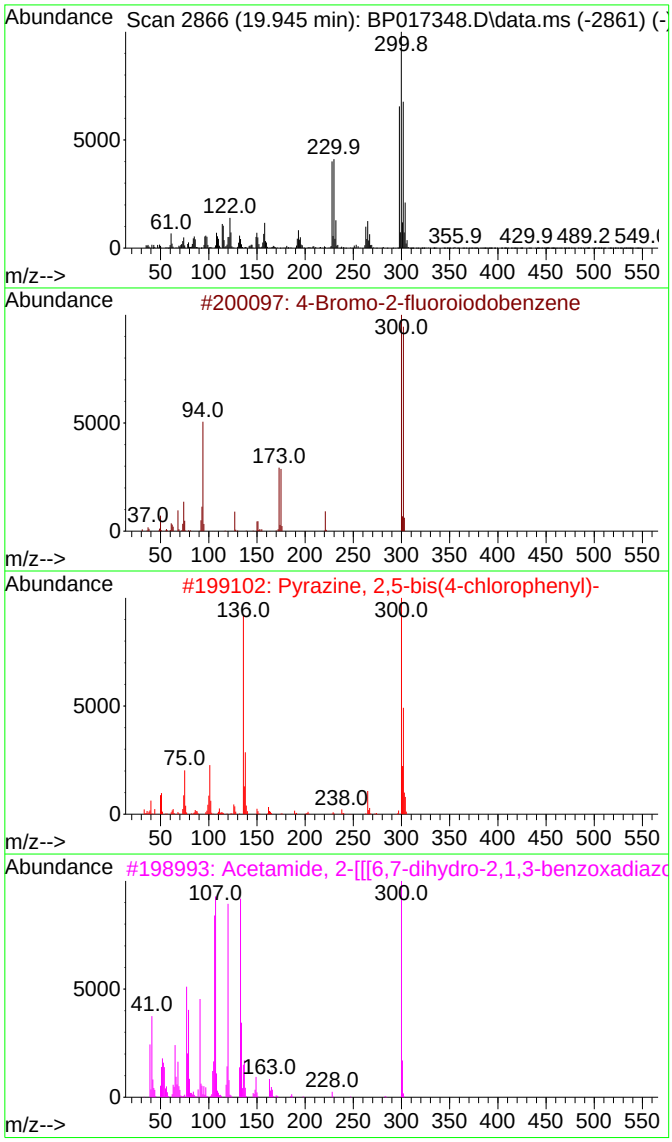
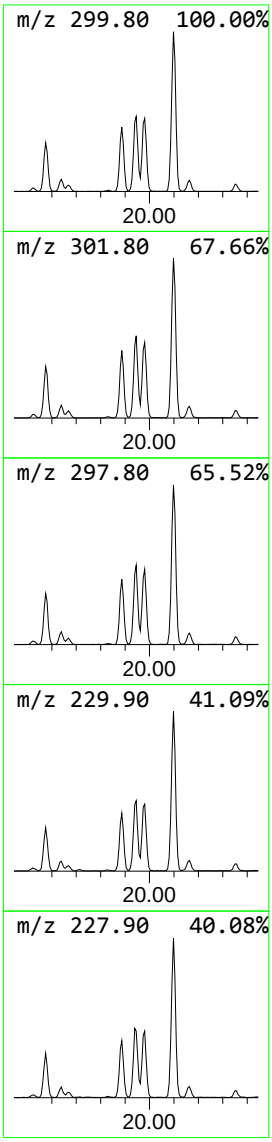
Peak Number 10 unknown-05 Concentration Rank 6

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 19.945 | 5.50 ng/ul | 1002140 | Chrysene-d12 | 21.457 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 4-Bromo-2-fluoroiodobenzene | 300 | C6H3BrFI | 105931-73-5 | 27 |
| 2 | Pyrazine, 2,5-bis(4-chlorophenyl)- | 300 | C16H10Cl2N2 | 074376-33-3 | 17 |
| 3 | Acetamide, 2-[[[6,7-dihydro-2,1,... | 300 | C15H16N4O3 | 1000319-88-7 | 14 |
| 4 | 2-Naphthalenol, 1,6-dibromo- | 300 | C10H6Br2O | 016239-18-2 | 10 |
| 5 | Cobalt, cyclopentadienyl(1-hexyl... | 298 | C13H19CoS2 | 1000062-71-0 | 10 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
Data File : BP017348.D
Acq On : 26 Sep 2023 07:11
Operator : MA/JU
Sample : 04472-08
Misc :
ALS Vial : 40 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EZYE8

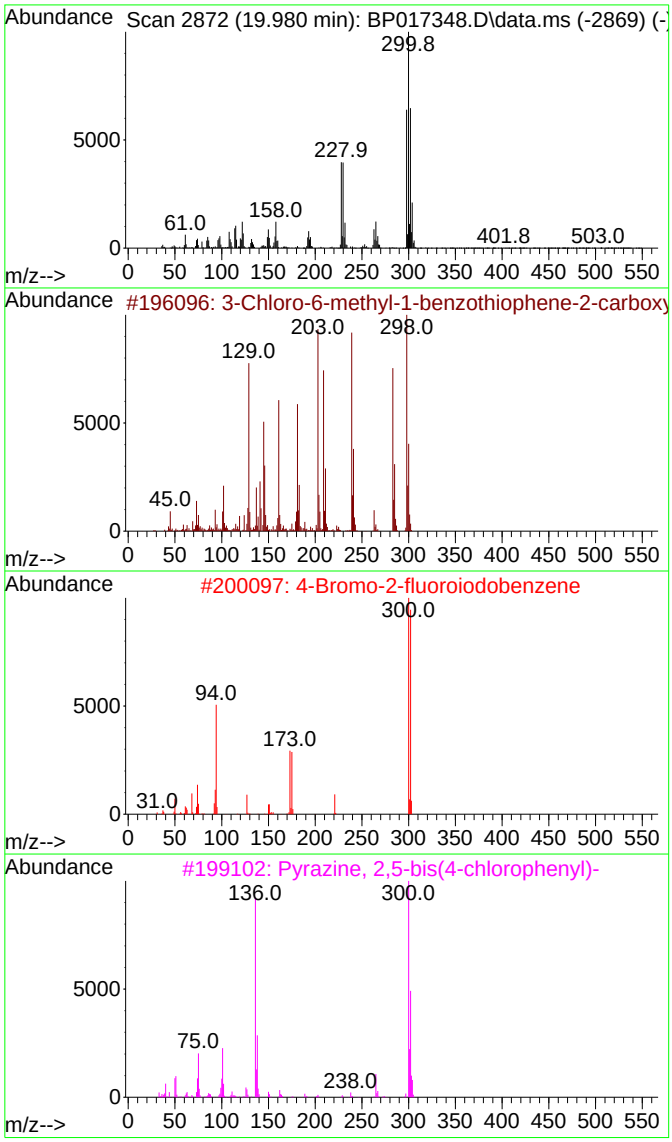
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

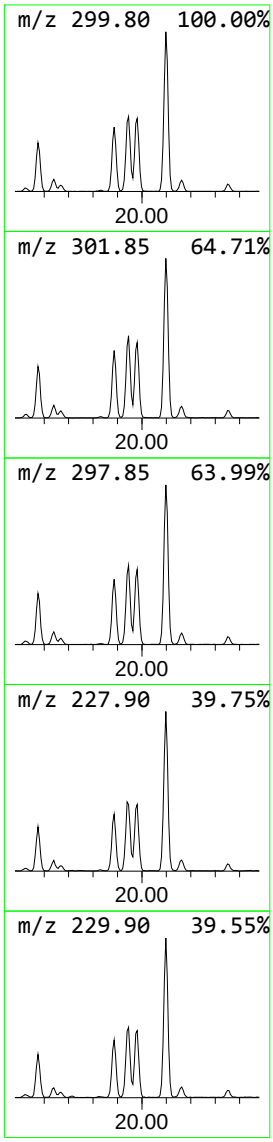
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 11 3-Chloro-6-methyl-1-benzoth... Concentration Rank 9

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 19.980 | 4.62 ng/ul | 842502 | Chrysene-d12 | 21.457 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 3-Chloro-6-methyl-1-benzothiophe... | 298 | C13H15ClO2SSi | 1000485-00-5 | 53 |
| 2 | 4-Bromo-2-fluoroiodobenzene | 300 | C6H3BrFI | 105931-73-5 | 27 |
| 3 | Pyrazine, 2,5-bis(4-chlorophenyl)- | 300 | C16H10Cl2N2 | 074376-33-3 | 25 |
| 4 | 1-(4-Methoxyphenyl)-4-cyanopyrim... | 300 | C18H12N4O | 1000286-51-5 | 10 |
| 5 | 3-[5-(Cyclohex-2-en-1-yl)-3H-1,2... | 302 | C17H14N6 | 1000443-34-3 | 9 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
Data File : BP017348.D
Acq On : 26 Sep 2023 07:11
Operator : MA/JU
Sample : 04472-08
Misc :
ALS Vial : 40 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EZYE8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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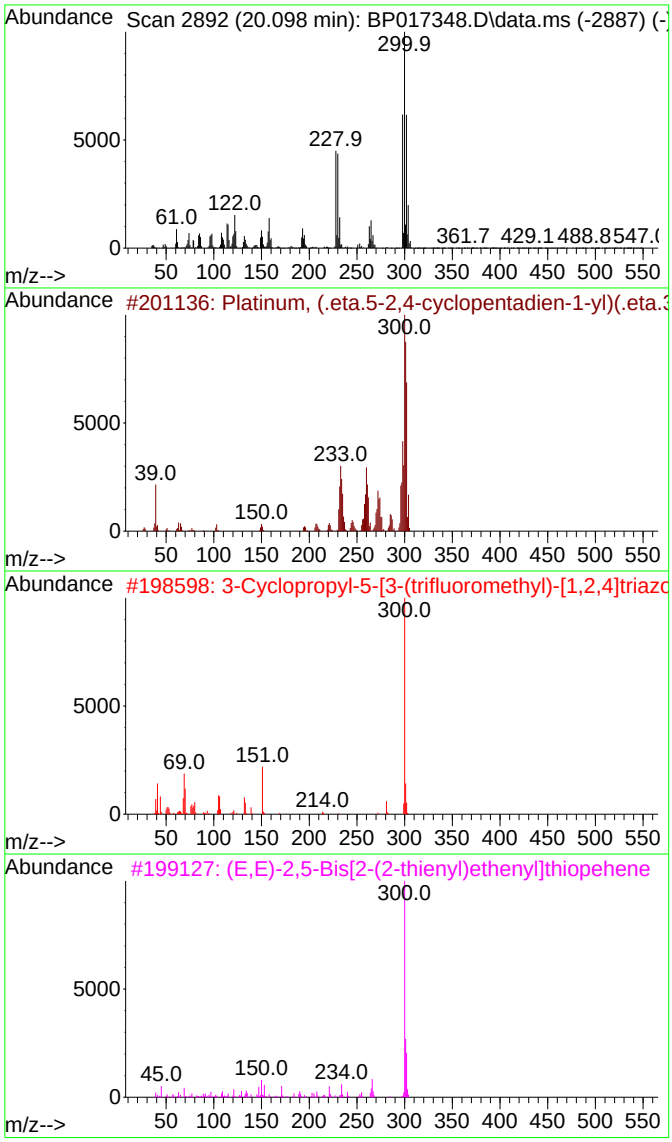
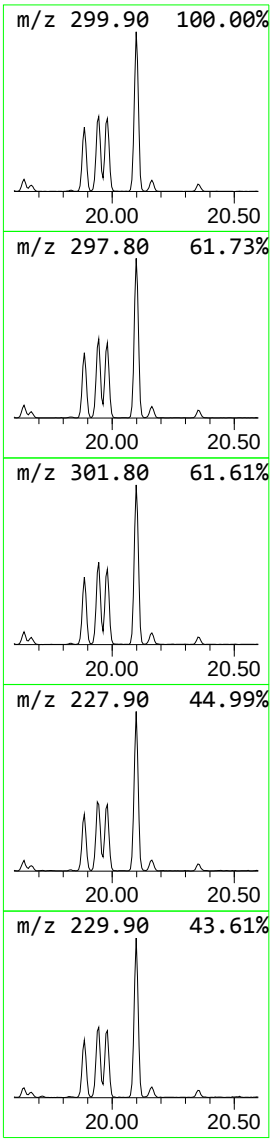
Peak Number 12 unknown-06 Concentration Rank 1

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 20.098 | 11.11 ng/ul | 2025570 | Chrysene-d12 | 21.457 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Platinum, (.eta.5-2,4-cyclopenta... | 301 | C8H10Pt | 035770-29-7 | 17 |
| 2 | 3-Cyclopropyl-5-[3-(trifluoromet... | 300 | C10H7F3N6S | 1000435-68-4 | 14 |
| 3 | (E,E)-2,5-Bis[2-(2-thienyl)ethen... | 300 | C16H12S3 | 026263-68-3 | 14 |
| 4 | Dimethyl 2,2'-diaminobiphenyl-4,... | 300 | C16H16N2O4 | 023933-57-5 | 9 |
| 5 | Acridin-9-yl-(3-methoxy-phenyl)-... | 300 | C20H16N2O | 075775-67-6 | 9 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
Data File : BP017348.D
Acq On : 26 Sep 2023 07:11
Operator : MA/JU
Sample : 04472-08
Misc :
ALS Vial : 40 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EZYE8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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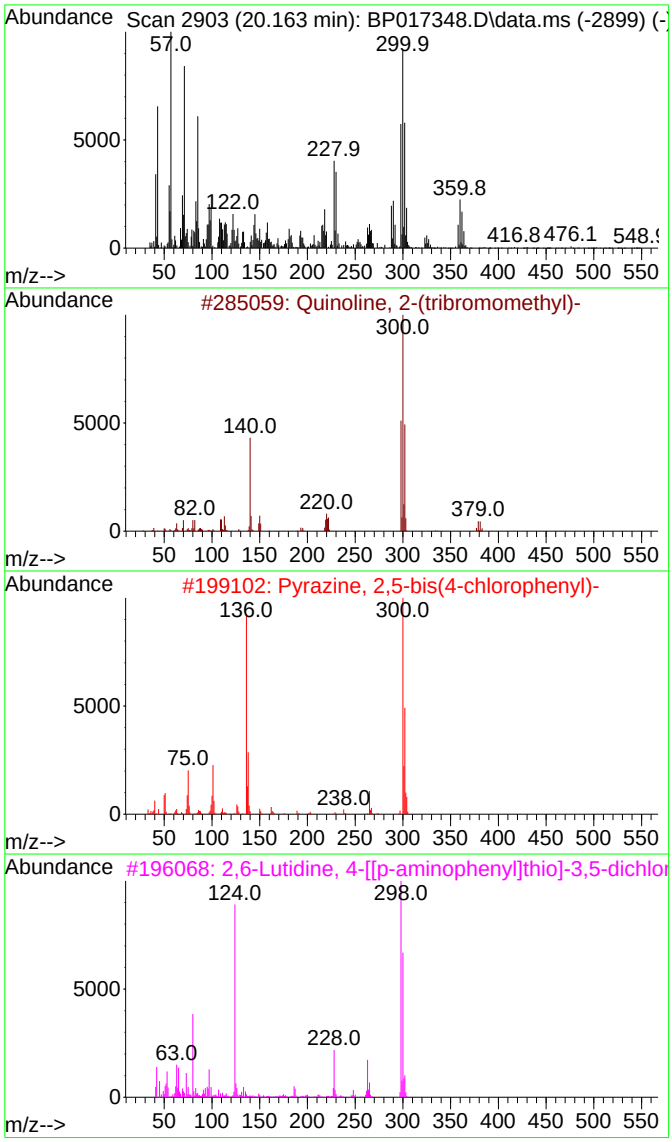
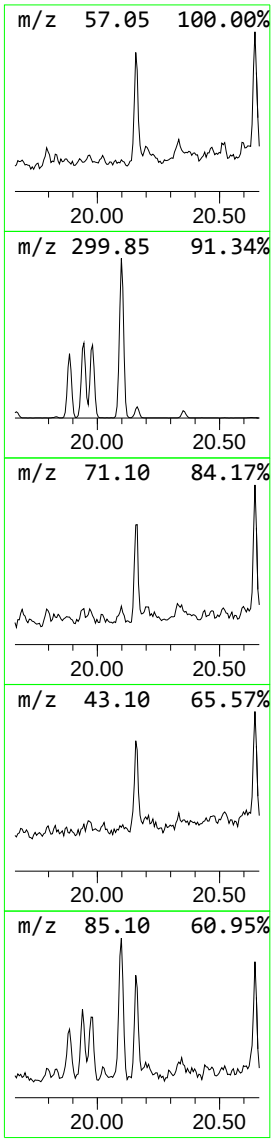
Peak Number 13 unknown-07 Concentration Rank 19

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 20.163 | 2.29 ng/ul | 418048 | Chrysene-d12 | 21.457 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Quinoline, 2-(tribromomethyl)- | 377 | C10H6Br3N | 000613-53-6 | 41 |
| 2 | Pyrazine, 2,5-bis(4-chlorophenyl)- | 300 | C16H10Cl2N2 | 074376-33-3 | 22 |
| 3 | 2,6-Lutidine, 4-[[p-aminophenyl]... | 298 | C13H12Cl2N2S | 1000252-27-9 | 11 |
| 4 | Dimethyl 2,2'-diaminobiphenyl-4,... | 300 | C16H16N2O4 | 023933-57-5 | 10 |
| 5 | Benzothiophene-3-carbonitrile, 4... | 300 | C16H13ClN2S | 069438-07-9 | 10 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
Data File : BP017348.D
Acq On : 26 Sep 2023 07:11
Operator : MA/JU
Sample : 04472-08
Misc :
ALS Vial : 40 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EZYE8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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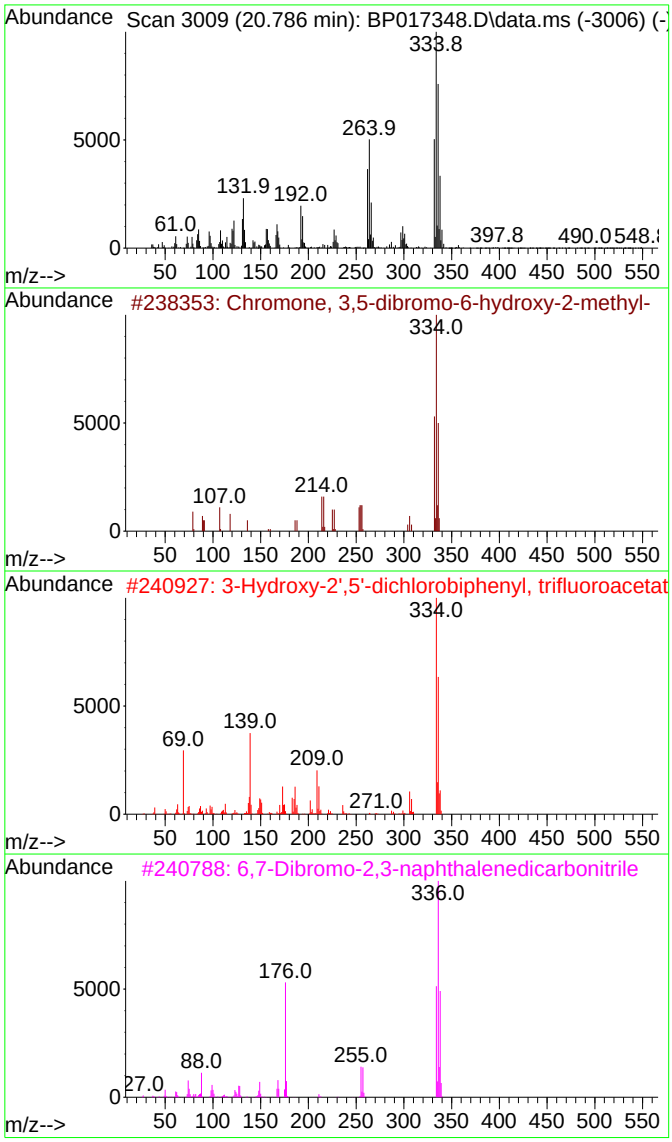
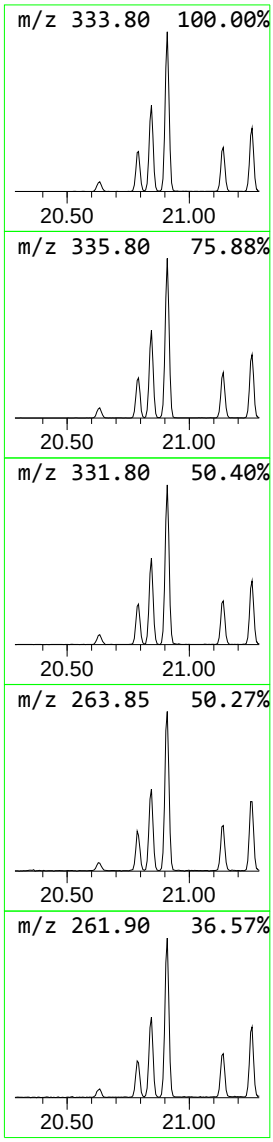
Peak Number 14 unknown-08 Concentration Rank 16

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 20.786 | 2.70 ng/ul | 491876 | Chrysene-d12 | 21.457 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Chromone, 3,5-dibromo-6-hydroxy-... | 332 | C10H6Br2O3 | 030095-77-3 | 43 |
| 2 | 3-Hydroxy-2',5'-dichlorobiphenyl... | 334 | C14H7Cl2F3O2 | 1010447-64-2 | 38 |
| 3 | 6,7-Dibromo-2,3-naphthalenedicar... | 334 | C12H4Br2N2 | 074815-81-9 | 22 |
| 4 | Anthracene, 9,10-dibromo- | 334 | C14H8Br2 | 000523-27-3 | 12 |
| 5 | Ambroxol | 376 | C13H18Br2N2O | 018683-91-5 | 11 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
Data File : BP017348.D
Acq On : 26 Sep 2023 07:11
Operator : MA/JU
Sample : 04472-08
Misc :
ALS Vial : 40 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EZYE8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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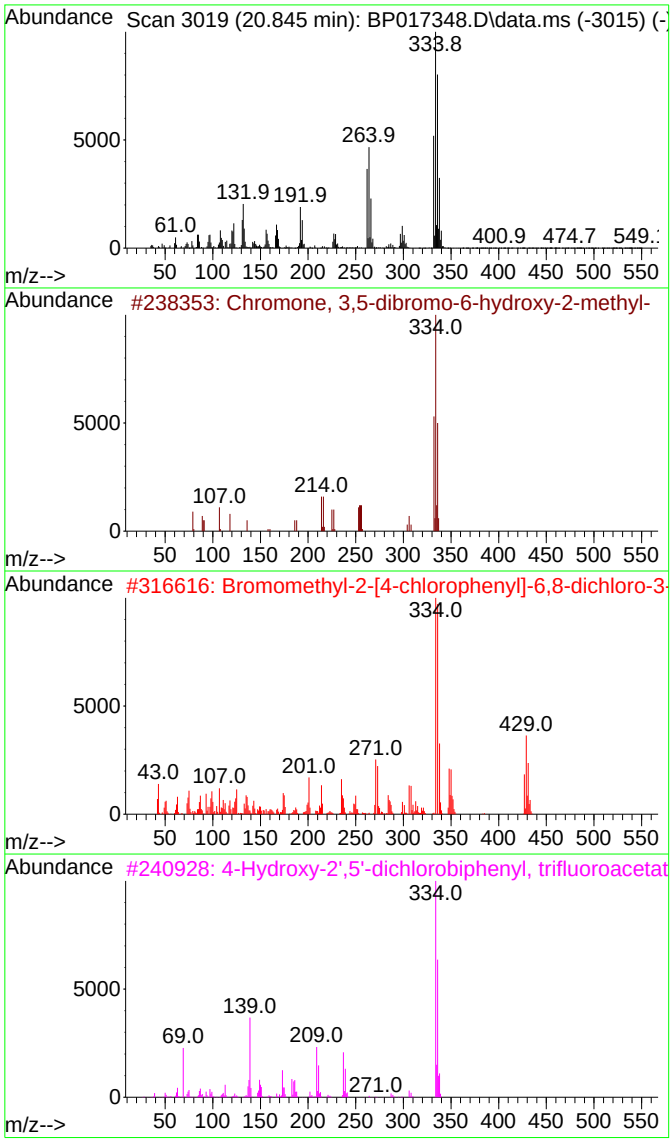
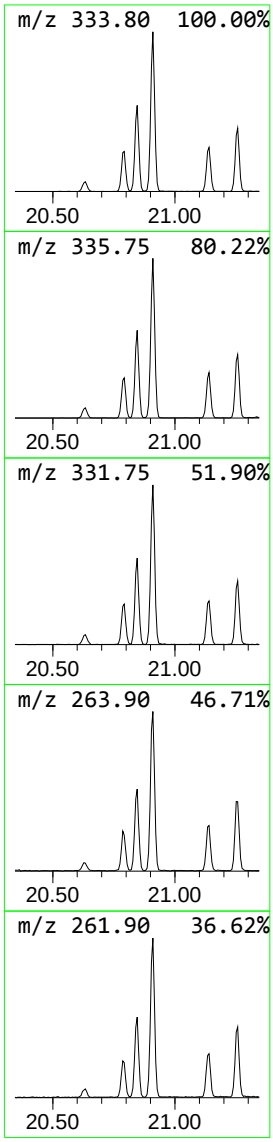
Peak Number 15 unknown-09 Concentration Rank 7

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 20.845 | 5.47 ng/ul | 997241 | Chrysene-d12 | 21.457 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Chromone, 3,5-dibromo-6-hydroxy-... | 332 | C10H6Br2O3 | 030095-77-3 | 38 |
| 2 | Bromomethyl-2-[4-chlorophenyl]-6... | 427 | C17H9BrCl3NO | 1000254-06-5 | 37 |
| 3 | 4-Hydroxy-2',5'-dichlorobiphenyl... | 334 | C14H7Cl2F3O2 | 1000447-64-8 | 32 |
| 4 | Anthracene, 1,8-dibromo- | 334 | C14H8Br2 | 131276-24-9 | 25 |
| 5 | 6,7-Dibromo-2,3-naphthalenedicar... | 334 | C12H4Br2N2 | 074815-81-9 | 22 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
Data File : BP017348.D
Acq On : 26 Sep 2023 07:11
Operator : MA/JU
Sample : 04472-08
Misc :
ALS Vial : 40 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EZYE8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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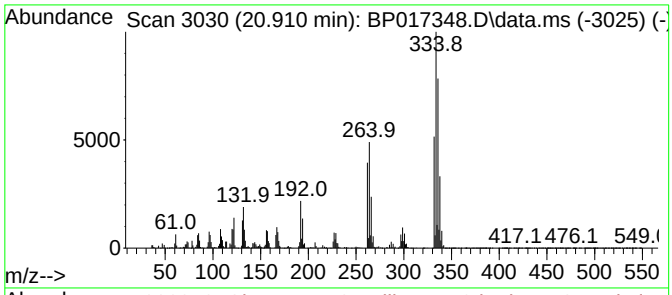
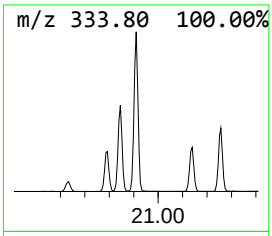
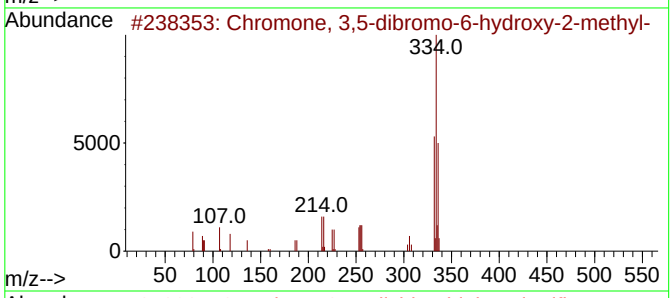
Peak Number 16 unknown-10 Concentration Rank 2

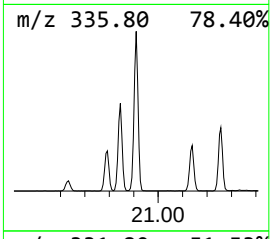
| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 20.910 | 10.87 ng/ul | 1981490 | Chrysene-d12 | 21.457 |

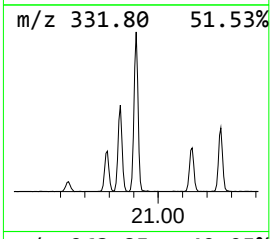
| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Chromone, 3,5-dibromo-6-hydroxy-... | 332 | C10H6Br2O3 | 030095-77-3 | 30 |
| 2 | 3-Hydroxy-2',5'-dichlorobiphenyl... | 334 | C14H7Cl2F3O2 | 1010447-64-2 | 25 |
| 3 | 6,7-Dibromo-2,3-naphthalenedicar... | 334 | C12H4Br2N2 | 074815-81-9 | 22 |
| 4 | Imidazole, 4,5-diiodo-2-methyl- | 334 | C4H4I2N2 | 073746-44-8 | 10 |
| 5 | Anthracene, 9,10-dibromo- | 334 | C14H8Br2 | 000523-27-3 | 10 |

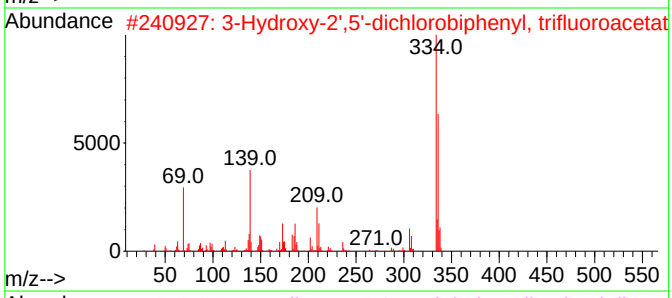


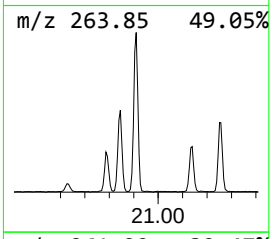



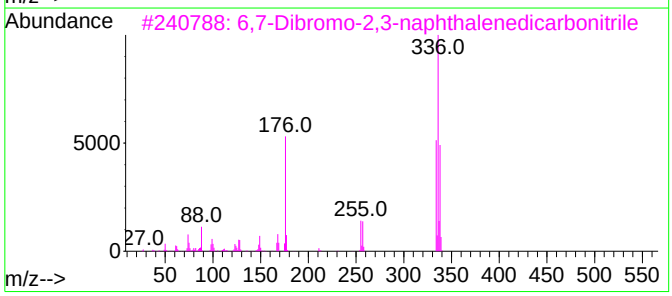



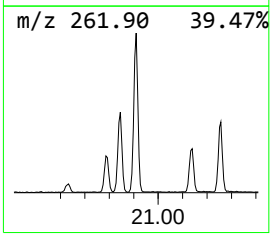



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
Data File : BP017348.D
Acq On : 26 Sep 2023 07:11
Operator : MA/JU
Sample : 04472-08
Misc :
ALS Vial : 40 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EZYE8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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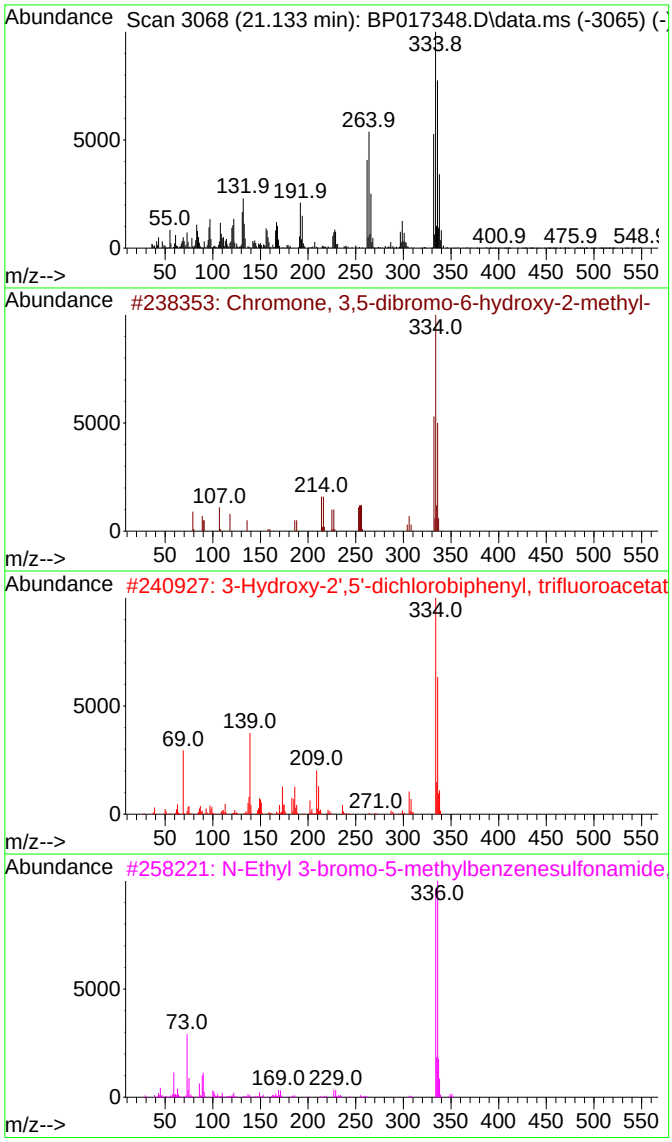
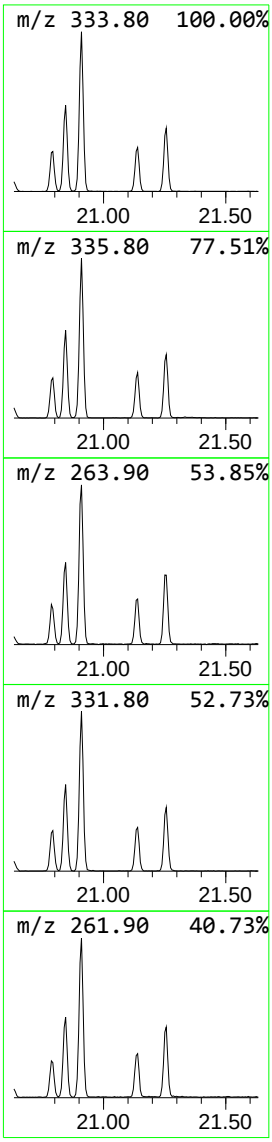
Peak Number 17 unknown-11 Concentration Rank 12

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 21.133 | 4.33 ng/ul | 788698 | Chrysene-d12 | 21.457 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Chromone, 3,5-dibromo-6-hydroxy-... | 332 | C10H6Br2O3 | 030095-77-3 | 35 |
| 2 | 3-Hydroxy-2',5'-dichlorobiphenyl... | 334 | C14H7Cl2F3O2 | 1010447-64-2 | 32 |
| 3 | N-Ethyl 3-bromo-5-methylbenzenes... | 349 | C12H20BrNO2SSi | 1010505-08-7 | 25 |
| 4 | 6,7-Dibromo-2,3-naphthalenedicar... | 334 | C12H4Br2N2 | 074815-81-9 | 22 |
| 5 | Ambroxol | 376 | C13H18Br2N2O | 018683-91-5 | 15 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
Data File : BP017348.D
Acq On : 26 Sep 2023 07:11
Operator : MA/JU
Sample : 04472-08
Misc :
ALS Vial : 40 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EZYE8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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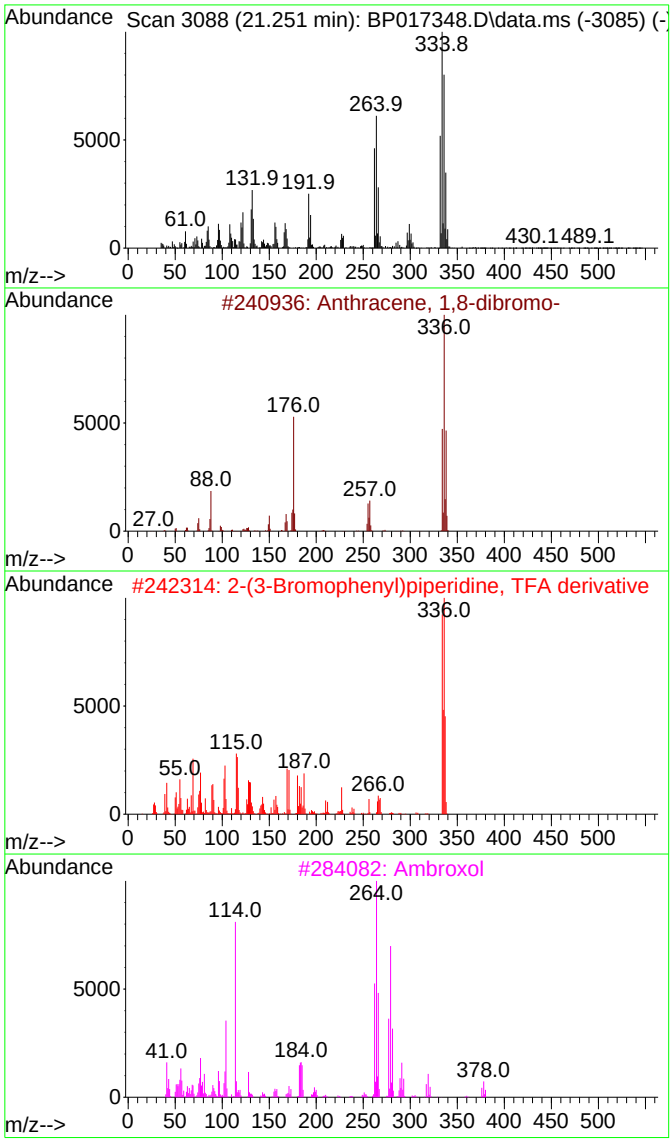
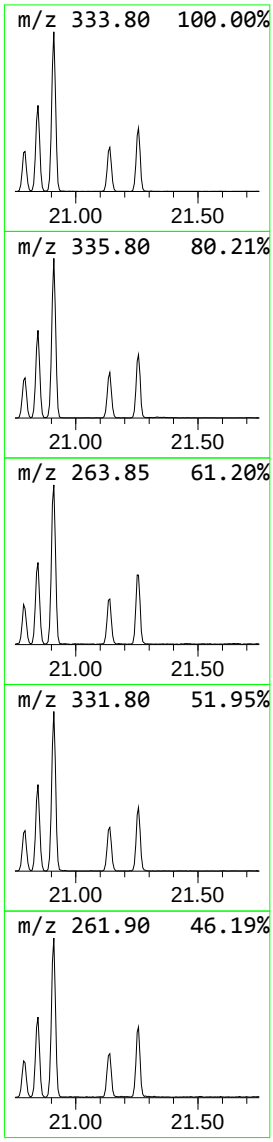
Peak Number 18 unknown-12 Concentration Rank 10

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 21.251 | 4.58 ng/ul | 834479 | Chrysene-d12 | 21.457 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Anthracene, 1,8-dibromo- | 334 | C14H8Br2 | 131276-24-9 | 25 |
| 2 | 2-(3-Bromophenyl)piperidine, TFA... | 335 | C13H13BrF3NO | 1000511-00-8 | 25 |
| 3 | Ambroxol | 376 | C13H18Br2N2O | 018683-91-5 | 25 |
| 4 | 6,7-Dibromo-2,3-naphthalenedicar... | 334 | C12H4Br2N2 | 074815-81-9 | 22 |
| 5 | (2,3,4,5-Tetrachloro-2,4-cyclope... | 334 | C13H6Cl4O2 | 1000243-99-3 | 22 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
Data File : BP017348.D
Acq On : 26 Sep 2023 07:11
Operator : MA/JU
Sample : 04472-08
Misc :
ALS Vial : 40 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EZYE8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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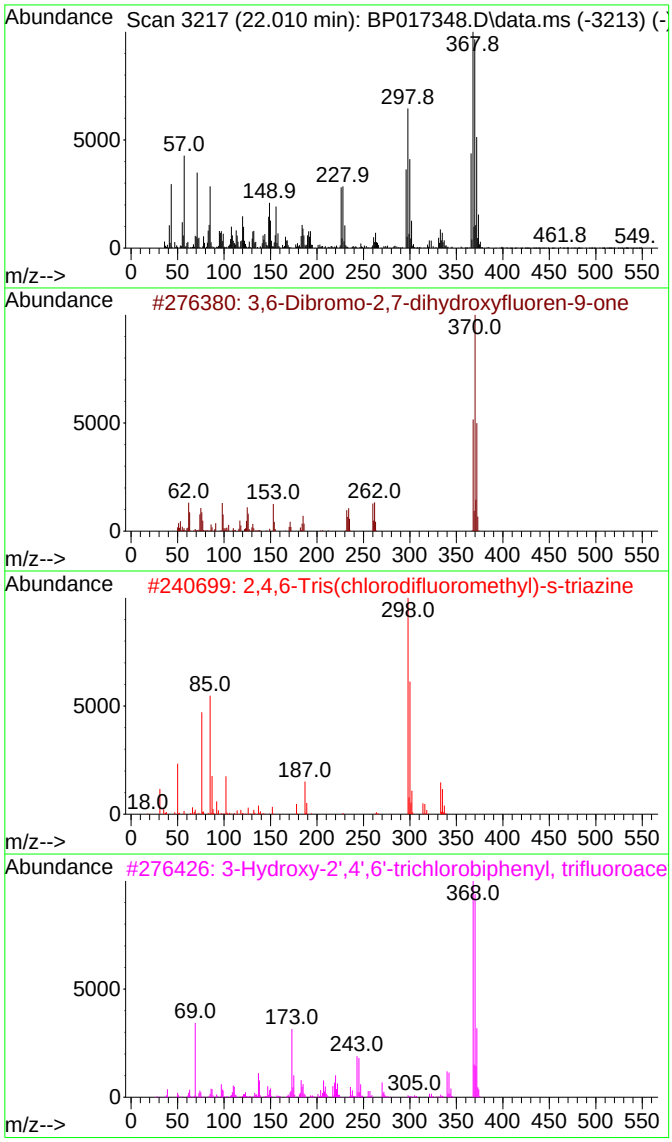
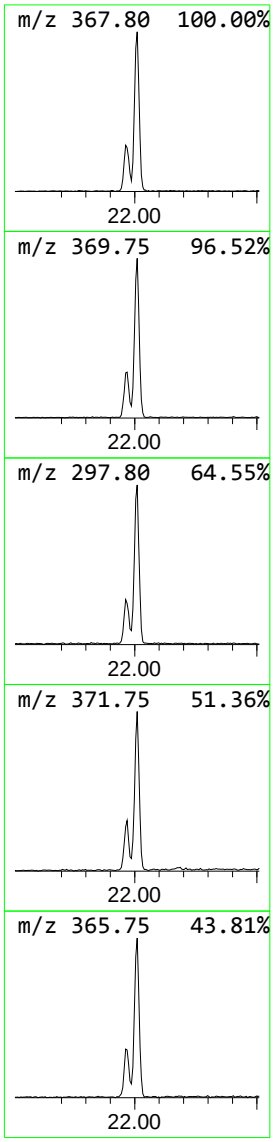
Peak Number 19 unknown-13 Concentration Rank 5

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 22.010 | 5.53 ng/ul | 1008140 | Chrysene-d12 | 21.457 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 3,6-Dibromo-2,7-dihydroxyfluoren... | 368 | C13H6Br2O3 | 084487-72-9 | 27 |
| 2 | 2,4,6-Tris(chlorodifluoromethyl)... | 333 | C6Cl3F6N3 | 1000343-67-7 | 25 |
| 3 | 3-Hydroxy-2',4',6'-trichlorobiph... | 368 | C14H6Cl3F3O2 | 1010447-69-7 | 17 |
| 4 | 3-Amino-4-(4-bromophenyl)-2-thio... | 383 | C15H18BrNO2SSi | 1010471-74-8 | 12 |
| 5 | 3-[2,4-Dichlorophenyl]-1-hydroxy... | 368 | C20H10Cl2O3 | 1000212-49-5 | 12 |



Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
Data File : BP017348.D
Acq On : 26 Sep 2023 07:11
Operator : MA/JU
Sample : 04472-08
Misc :
ALS Vial : 40 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EZYE8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

| | | | | | --Internal Standard--- | | | |
|---|---|---|---|---|---|---|---|---|
| TIC Top Hit name | RT | EstConc | Units | Response | # | RT | Resp | Conc |
| unknown-01 | 3.840 | 4.2 | ng/ul | 303158 | 1 | 7.922 | 1455930 | 20.0 |
| unknown-02 | 3.922 | 2.5 | ng/ul | 182013 | 1 | 7.922 | 1455930 | 20.0 |
| Benzoic acid | 10.010 | 4.8 | ng/ul | 510828 | 2 | 10.746 | 2144070 | 20.0 |
| Naphthalene, 1,... | 18.792 | 3.9 | ng/ul | 608442 | 4 | 17.357 | 3155320 | 20.0 |
| Naphthalene, 1,... | 18.910 | 8.1 | ng/ul | 1273240 | 4 | 17.357 | 3155320 | 20.0 |
| Naphthalene, 1,... | 19.151 | 6.0 | ng/ul | 945948 | 4 | 17.357 | 3155320 | 20.0 |
| Dibenzoylmethane | 19.257 | 2.3 | ng/ul | 368425 | 4 | 17.357 | 3155320 | 20.0 |
| unknown-03 | 19.575 | 3.3 | ng/ul | 606187 | 5 | 21.457 | 3646950 | 20.0 |
| unknown-04 | 19.886 | 4.4 | ng/ul | 794290 | 5 | 21.457 | 3646950 | 20.0 |
| unknown-05 | 19.945 | 5.5 | ng/ul | 1002140 | 5 | 21.457 | 3646950 | 20.0 |
| 3-Chloro-6-meth... | 19.980 | 4.6 | ng/ul | 842502 | 5 | 21.457 | 3646950 | 20.0 |
| unknown-06 | 20.098 | 11.1 | ng/ul | 2025570 | 5 | 21.457 | 3646950 | 20.0 |
| unknown-07 | 20.163 | 2.3 | ng/ul | 418048 | 5 | 21.457 | 3646950 | 20.0 |
| unknown-08 | 20.786 | 2.7 | ng/ul | 491876 | 5 | 21.457 | 3646950 | 20.0 |
| unknown-09 | 20.845 | 5.5 | ng/ul | 997241 | 5 | 21.457 | 3646950 | 20.0 |
| unknown-10 | 20.910 | 10.9 | ng/ul | 1981490 | 5 | 21.457 | 3646950 | 20.0 |
| unknown-11 | 21.133 | 4.3 | ng/ul | 788698 | 5 | 21.457 | 3646950 | 20.0 |
| unknown-12 | 21.251 | 4.6 | ng/ul | 834479 | 5 | 21.457 | 3646950 | 20.0 |
| unknown-13 | 22.010 | 5.5 | ng/ul | 1008140 | 5 | 21.457 | 3646950 | 20.0 |