LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012376.D
Acq On : 28 Oct 2022 14:05
Operator : CG/JU
Sample : N5232-16
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BHAB9

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-BP102522.M
Title : SVOA CALIBRATION

Signal : TIC: BP012376.D\data.ms

| peak # | R.T. min | first scan | max scan | last scan | PK TY | peak height | corr. area | corr. % max. | % of total |
|---|---|---|---|---|---|---|---|---|---|
| 1 | 3.034 | 6 | 8 | 27 | rVB | 701873 | 1057975 | 8.42% | 0.615% |
| 2 | 5.111 | 353 | 361 | 378 | rBV | 5310284 | 8707746 | 69.33% | 5.062% |
| 3 | 5.517 | 424 | 430 | 435 | rVV | 185628 | 285320 | 2.27% | 0.166% |
| 4 | 6.075 | 517 | 525 | 531 | rBV | 228415 | 355880 | 2.83% | 0.207% |
| 5 | 6.281 | 551 | 560 | 572 | rBV | 1753366 | 2830102 | 22.53% | 1.645% |
| 6 | 6.469 | 579 | 592 | 598 | rBV3 | 201354 | 361661 | 2.88% | 0.210% |
| 7 | 6.775 | 638 | 644 | 660 | rBV | 222037 | 394614 | 3.14% | 0.229% |
| 8 | 6.916 | 662 | 668 | 673 | rBV | 247127 | 343742 | 2.74% | 0.200% |
| 9 | 6.981 | 673 | 679 | 690 | rVB | 264468 | 478387 | 3.81% | 0.278% |
| 10 | 7.140 | 699 | 706 | 711 | rBV | 1077407 | 1746232 | 13.90% | 1.015% |
| 11 | 7.181 | 711 | 713 | 718 | rVV | 144339 | 211762 | 1.69% | 0.123% |
| 12 | 7.234 | 718 | 722 | 733 | rVB | 124362 | 221420 | 1.76% | 0.129% |
| 13 | 7.334 | 733 | 739 | 747 | rBV | 997513 | 1793223 | 14.28% | 1.042% |
| 14 | 7.505 | 762 | 768 | 774 | rBV | 552736 | 840752 | 6.69% | 0.489% |
| 15 | 7.805 | 814 | 819 | 824 | rBV | 799750 | 1233257 | 9.82% | 0.717% |
| 16 | 7.999 | 847 | 852 | 859 | rVB | 379279 | 577452 | 4.60% | 0.336% |
| 17 | 8.175 | 874 | 882 | 891 | rVB2 | 252751 | 514937 | 4.10% | 0.299% |
| 18 | 8.522 | 932 | 941 | 949 | rVB | 831056 | 1343097 | 10.69% | 0.781% |
| 19 | 8.611 | 949 | 956 | 968 | rVB5 | 87533 | 240486 | 1.91% | 0.140% |
| 20 | 8.710 | 968 | 973 | 978 | rBV3 | 140721 | 249891 | 1.99% | 0.145% |
| 21 | 8.799 | 983 | 988 | 998 | rVB3 | 197816 | 377240 | 3.00% | 0.219% |
| 22 | 8.905 | 998 | 1006 | 1012 | rBV | 323230 | 544981 | 4.34% | 0.317% |
| 23 | 8.975 | 1012 | 1018 | 1025 | rBV | 1253877 | 2184700 | 17.39% | 1.270% |
| 24 | 9.034 | 1025 | 1028 | 1033 | rVB2 | 137838 | 201788 | 1.61% | 0.117% |
| 25 | 9.428 | 1089 | 1095 | 1101 | rVB | 255745 | 414553 | 3.30% | 0.241% |
| 26 | 9.693 | 1134 | 1140 | 1154 | rVV3 | 1128480 | 2155877 | 17.16% | 1.253% |
| 27 | 9.852 | 1162 | 1167 | 1177 | rVB2 | 117822 | 245884 | 1.96% | 0.143% |
| 28 | 10.005 | 1182 | 1193 | 1208 | rBV4 | 207046 | 695551 | 5.54% | 0.404% |
| 29 | 10.234 | 1224 | 1232 | 1240 | rBV2 | 1716482 | 3318565 | 26.42% | 1.929% |
| 30 | 10.310 | 1240 | 1245 | 1248 | rVV3 | 188617 | 338157 | 2.69% | 0.197% |
| 31 | 10.352 | 1249 | 1252 | 1256 | rVV2 | 184560 | 341227 | 2.72% | 0.198% |
| 32 | 10.416 | 1261 | 1263 | 1269 | rVB | 170795 | 254170 | 2.02% | 0.148% |
| 33 | 10.605 | 1288 | 1295 | 1304 | rVV | 1206367 | 2183678 | 17.39% | 1.269% |
| 34 | 10.699 | 1304 | 1311 | 1314 | rVV | 1739141 | 2973705 | 23.68% | 1.729% |
| 35 | 10.746 | 1314 | 1319 | 1337 | rVV | 4784742 | 8662981 | 68.97% | 5.036% |
| 36 | 11.204 | 1392 | 1397 | 1406 | rVV2 | 142756 | 299311 | 2.38% | 0.174% |

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012376.D
Acq On : 28 Oct 2022 14:05
Operator : CG/JU
Sample : N5232-16
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BHAB9

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-BP102522.M
Title : SVOA CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 37 | 11.816 | 1495 | 1501 | 1507 | rVV3 | 115643 | 266075 | 2.12% | 0.155% |
| 38 | 11.957 | 1507 | 1525 | 1534 | rVV2 | 1075737 | 2768934 | 22.05% | 1.610% |
| 39 | 12.104 | 1541 | 1550 | 1561 | rVV | 98265 | 432049 | 3.44% | 0.251% |
| 40 | 12.428 | 1598 | 1605 | 1612 | rBV9 | 89365 | 238249 | 1.90% | 0.138% |
| 41 | 12.999 | 1696 | 1702 | 1706 | rVV | 401718 | 616586 | 4.91% | 0.358% |
| 42 | 13.046 | 1706 | 1710 | 1714 | rVV3 | 129229 | 215512 | 1.72% | 0.125% |
| 43 | 13.093 | 1714 | 1718 | 1728 | rVB | 312870 | 577170 | 4.60% | 0.336% |
| 44 | 13.510 | 1783 | 1789 | 1792 | rBV | 203319 | 299884 | 2.39% | 0.174% |
| 45 | 13.869 | 1840 | 1850 | 1863 | rBV | 3116049 | 4965440 | 39.53% | 2.886% |
| 46 | 14.146 | 1892 | 1897 | 1908 | rVV2 | 3449651 | 5233813 | 41.67% | 3.042% |
| 47 | 14.246 | 1909 | 1914 | 1925 | rVB | 296214 | 520683 | 4.15% | 0.303% |
| 48 | 14.457 | 1942 | 1950 | 1958 | rBV2 | 2194216 | 3845550 | 30.62% | 2.235% |
| 49 | 14.693 | 1985 | 1990 | 1999 | rVB2 | 238377 | 457713 | 3.64% | 0.266% |
| 50 | 15.204 | 2074 | 2077 | 2081 | rBV | 263109 | 337899 | 2.69% | 0.196% |
| 51 | 15.357 | 2099 | 2103 | 2112 | rVB5 | 149349 | 257204 | 2.05% | 0.150% |
| 52 | 15.451 | 2113 | 2119 | 2131 | rVB | 4603944 | 6949099 | 55.33% | 4.040% |
| 53 | 15.587 | 2137 | 2142 | 2148 | rVB | 1271888 | 1708704 | 13.60% | 0.993% |
| 54 | 15.687 | 2154 | 2159 | 2160 | rBV | 205766 | 259997 | 2.07% | 0.151% |
| 55 | 15.722 | 2160 | 2165 | 2180 | rVB | 5731394 | 8305157 | 66.12% | 4.828% |
| 56 | 15.881 | 2187 | 2192 | 2198 | rBV | 122195 | 200480 | 1.60% | 0.117% |
| 57 | 15.975 | 2201 | 2208 | 2215 | rVV3 | 743484 | 1488342 | 11.85% | 0.865% |
| 58 | 16.163 | 2235 | 2240 | 2246 | rVV | 950049 | 1505212 | 11.98% | 0.875% |
| 59 | 16.216 | 2247 | 2249 | 2258 | rVB2 | 221578 | 349040 | 2.78% | 0.203% |
| 60 | 16.492 | 2292 | 2296 | 2298 | rBV2 | 206519 | 293322 | 2.34% | 0.171% |
| 61 | 16.516 | 2298 | 2300 | 2305 | rVB2 | 227792 | 270723 | 2.16% | 0.157% |
| 62 | 16.681 | 2324 | 2328 | 2332 | rVV2 | 206254 | 238367 | 1.90% | 0.139% |
| 63 | 17.181 | 2408 | 2413 | 2415 | rBV | 804811 | 1060903 | 8.45% | 0.617% |
| 64 | 17.216 | 2415 | 2419 | 2428 | rVV | 2663421 | 4096838 | 32.62% | 2.382% |
| 65 | 17.316 | 2430 | 2436 | 2442 | rVV2 | 5523037 | 8022130 | 63.87% | 4.663% |
| 66 | 17.375 | 2442 | 2446 | 2452 | rVB | 211095 | 321535 | 2.56% | 0.187% |
| 67 | 17.581 | 2477 | 2481 | 2488 | rBV4 | 152801 | 303088 | 2.41% | 0.176% |
| 68 | 17.857 | 2523 | 2528 | 2531 | rBV | 473044 | 738847 | 5.88% | 0.429% |
| 69 | 17.910 | 2533 | 2537 | 2541 | rVV | 642736 | 978289 | 7.79% | 0.569% |
| 70 | 17.951 | 2541 | 2544 | 2555 | rVB | 216081 | 389116 | 3.10% | 0.226% |
| 71 | 18.104 | 2566 | 2570 | 2577 | rBV | 189596 | 304864 | 2.43% | 0.177% |
| 72 | 18.181 | 2578 | 2583 | 2588 | rVV | 1301991 | 1714317 | 13.65% | 0.997% |
| 73 | 18.257 | 2591 | 2596 | 2601 | rVB | 1040889 | 1356546 | 10.80% | 0.789% |
| 74 | 18.351 | 2608 | 2612 | 2615 | rBV | 816558 | 1165631 | 9.28% | 0.678% |
| 75 | 18.404 | 2615 | 2621 | 2626 | rVB | 1641947 | 2766080 | 22.02% | 1.608% |
| 76 | 18.557 | 2644 | 2647 | 2653 | rVB3 | 110291 | 219296 | 1.75% | 0.127% |

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012376.D
Acq On : 28 Oct 2022 14:05
Operator : CG/JU
Sample : N5232-16
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BHAB9

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-BP102522.M
Title : SVOA CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 77 | 18.775 | 2680 | 2684 | 2691 | rBV3 | 210856 | 426940 | 3.40% | 0.248% |
| 78 | 18.922 | 2705 | 2709 | 2719 | rVB2 | 335695 | 498499 | 3.97% | 0.290% |
| 79 | 19.004 | 2720 | 2723 | 2727 | rVV3 | 146104 | 218944 | 1.74% | 0.127% |
| 80 | 19.081 | 2729 | 2736 | 2741 | rVB2 | 508830 | 921231 | 7.33% | 0.536% |
| 81 | 19.133 | 2741 | 2745 | 2748 | rBV2 | 204593 | 273354 | 2.18% | 0.159% |
| 82 | 19.181 | 2748 | 2753 | 2765 | rVB2 | 1148724 | 1707253 | 13.59% | 0.992% |
| 83 | 19.281 | 2766 | 2770 | 2777 | rVV3 | 585435 | 870156 | 6.93% | 0.506% |
| 84 | 19.345 | 2777 | 2781 | 2791 | rVB4 | 285620 | 557171 | 4.44% | 0.324% |
| 85 | 19.557 | 2812 | 2817 | 2822 | rBV | 7550746 | 9690815 | 77.16% | 5.633% |
| 86 | 19.610 | 2822 | 2826 | 2835 | rVB | 3789290 | 5450576 | 43.40% | 3.168% |
| 87 | 19.781 | 2851 | 2855 | 2863 | rVV3 | 594086 | 1014380 | 8.08% | 0.590% |
| 88 | 19.857 | 2864 | 2868 | 2877 | rVB | 542477 | 816851 | 6.50% | 0.475% |
| 89 | 19.969 | 2883 | 2887 | 2890 | rVB | 554555 | 649969 | 5.17% | 0.378% |
| 90 | 20.010 | 2890 | 2894 | 2898 | rBV | 867648 | 1068605 | 8.51% | 0.621% |
| 91 | 20.139 | 2913 | 2916 | 2925 | rBV2 | 184425 | 346267 | 2.76% | 0.201% |
| 92 | 20.269 | 2934 | 2938 | 2943 | rVB | 770382 | 1007448 | 8.02% | 0.586% |
| 93 | 20.457 | 2965 | 2970 | 2975 | rBV3 | 493937 | 934061 | 7.44% | 0.543% |
| 94 | 20.551 | 2982 | 2986 | 2992 | rBV | 469799 | 596445 | 4.75% | 0.347% |
| 95 | 21.022 | 3063 | 3066 | 3075 | rVB2 | 331686 | 458216 | 3.65% | 0.266% |
| 96 | 21.310 | 3110 | 3115 | 3124 | rBV | 4011139 | 5157849 | 41.07% | 2.998% |
| 97 | 21.439 | 3132 | 3137 | 3141 | rBV | 3831616 | 5031823 | 40.06% | 2.925% |
| 98 | 21.475 | 3141 | 3143 | 3154 | rVB2 | 629721 | 877161 | 6.98% | 0.510% |
| 99 | 23.551 | 3489 | 3496 | 3505 | rVB2 | 6162701 | 12559806 | 100.00% | 7.301% |
| 100 | 23.704 | 3516 | 3522 | 3533 | rVB2 | 2866051 | 5873606 | 46.77% | 3.414% |

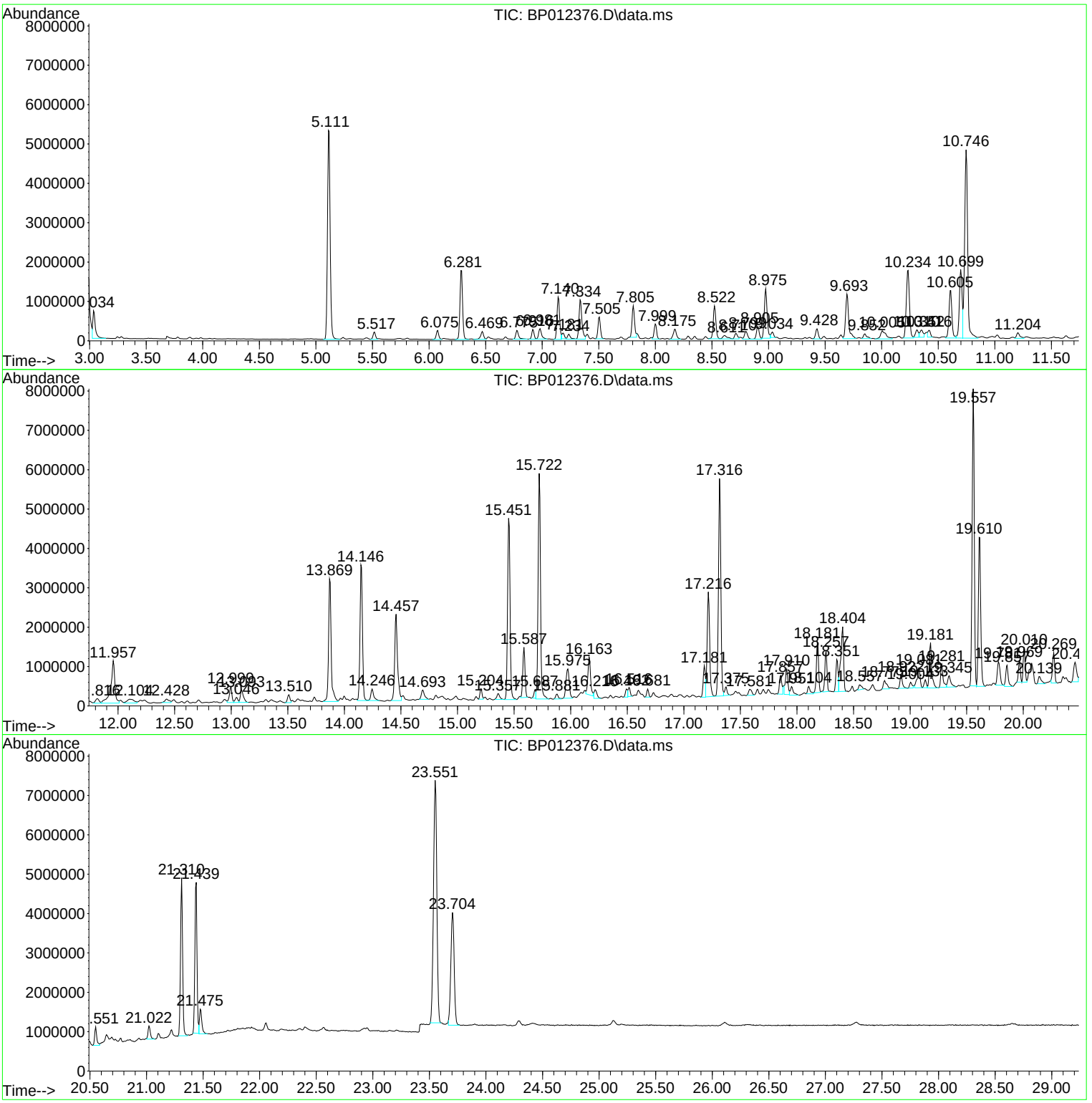
Sum of corrected areas: 172026414

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012376.D
Acq On : 28 Oct 2022 14:05
Operator : CG/JU
Sample : N5232-16
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BHAB9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012376.D
Acq On : 28 Oct 2022 14:05
Operator : CG/JU
Sample : N5232-16
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BHAB9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

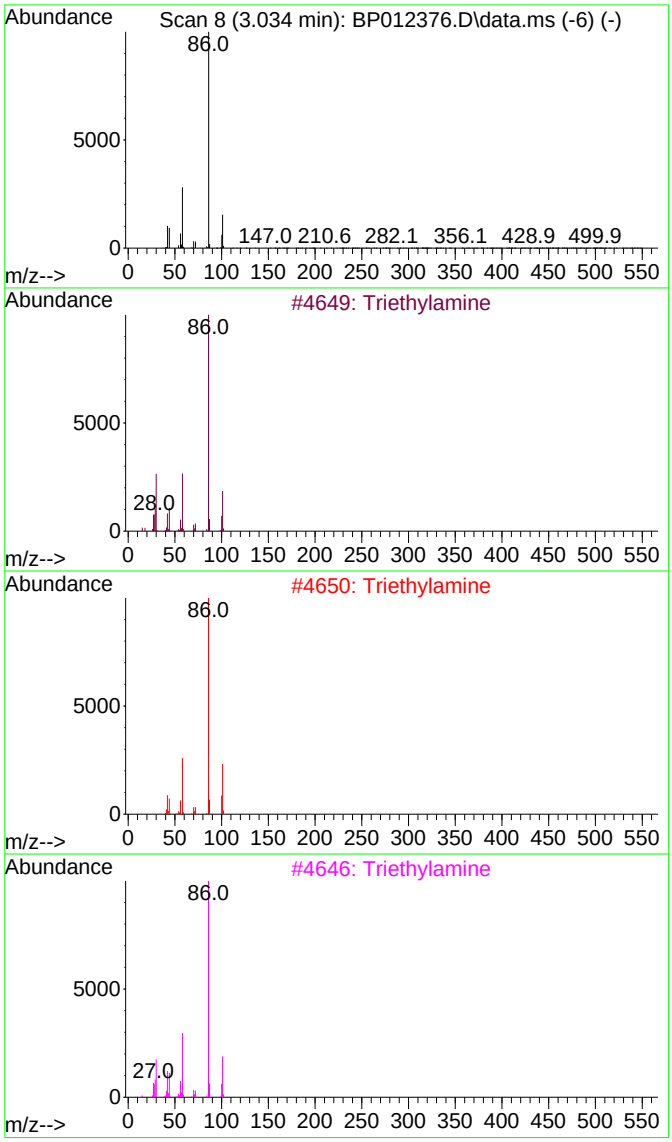
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

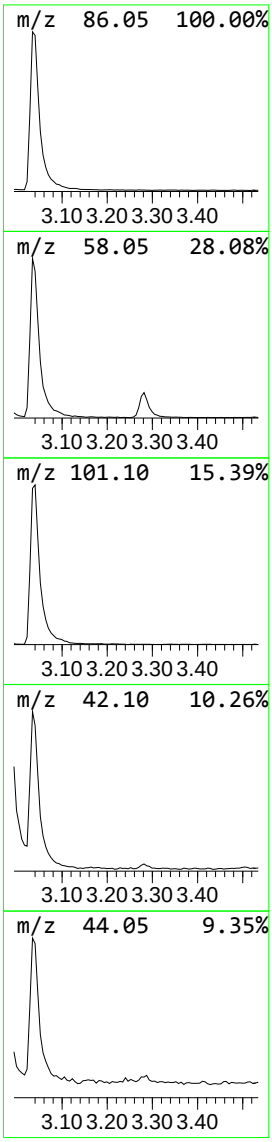
**********

Peak Number 1 Triethylamine Concentration Rank 7

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 3.034 | 17.16 ng/ul | 1057980 | 1,4-Dichlorobenzene-d4 | 7.805 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Triethylamine | 101 | C6H15N | 000121-44-8 | 91 |
| 2 | Triethylamine | 101 | C6H15N | 000121-44-8 | 91 |
| 3 | Triethylamine | 101 | C6H15N | 000121-44-8 | 91 |
| 4 | Triethylamine | 101 | C6H15N | 000121-44-8 | 86 |
| 5 | Tetraethyl ammonium fluoride | 149 | C8H20FN | 000665-46-3 | 78 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012376.D
Acq On : 28 Oct 2022 14:05
Operator : CG/JU
Sample : N5232-16
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BHAB9

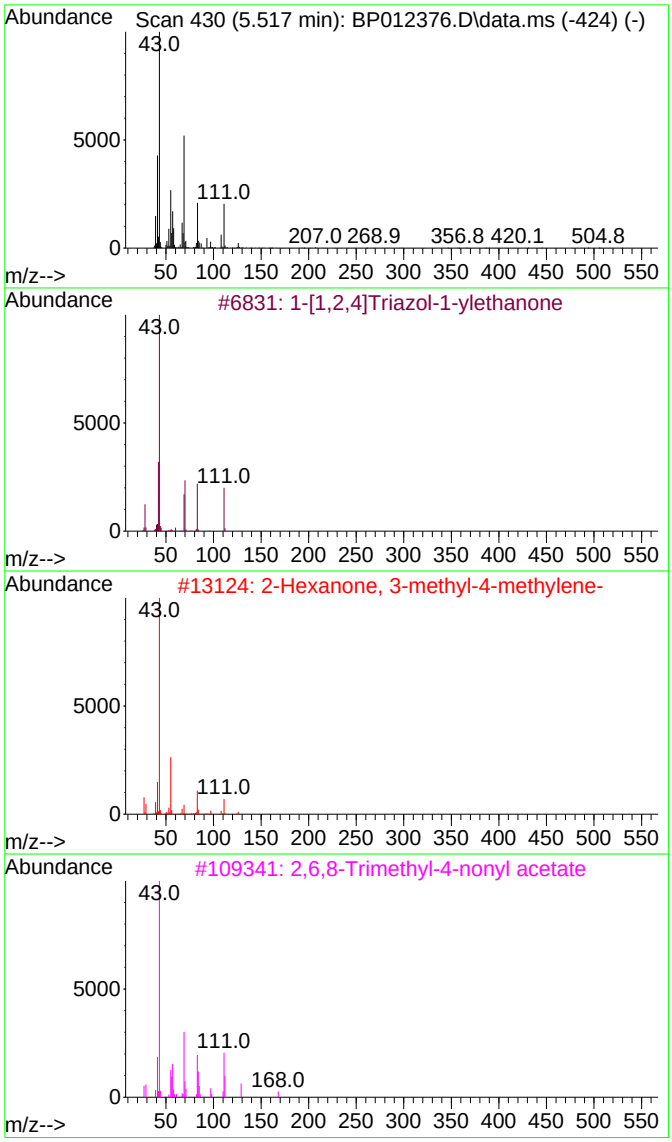
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

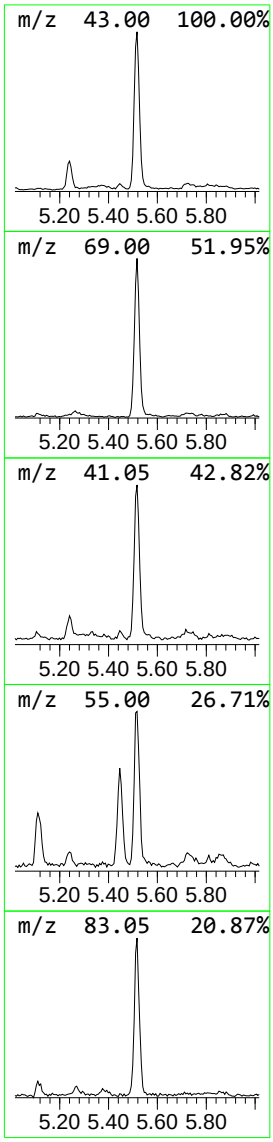
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 3 unknown-01 Concentration Rank 26

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 5.517 | 4.63 ng/ul | 285320 | 1,4-Dichlorobenzene-d4 | 7.805 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1-[1,2,4]Triazol-1-ylethanone | 111 | C4H5N3O | 015625-88-4 | 47 |
| 2 | 2-Hexanone, 3-methyl-4-methylene- | 126 | C8H14O | 020690-71-5 | 47 |
| 3 | 2,6,8-Trimethyl-4-nonyl acetate | 228 | C14H28O2 | 1000430-73-9 | 40 |
| 4 | Cyclopentane, (2-methylbutyl)- | 140 | C10H20 | 053366-38-4 | 25 |
| 5 | 2,2,4-Trimethyl-3-hexene | 126 | C9H18 | 059643-72-0 | 23 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012376.D
Acq On : 28 Oct 2022 14:05
Operator : CG/JU
Sample : N5232-16
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BHAB9

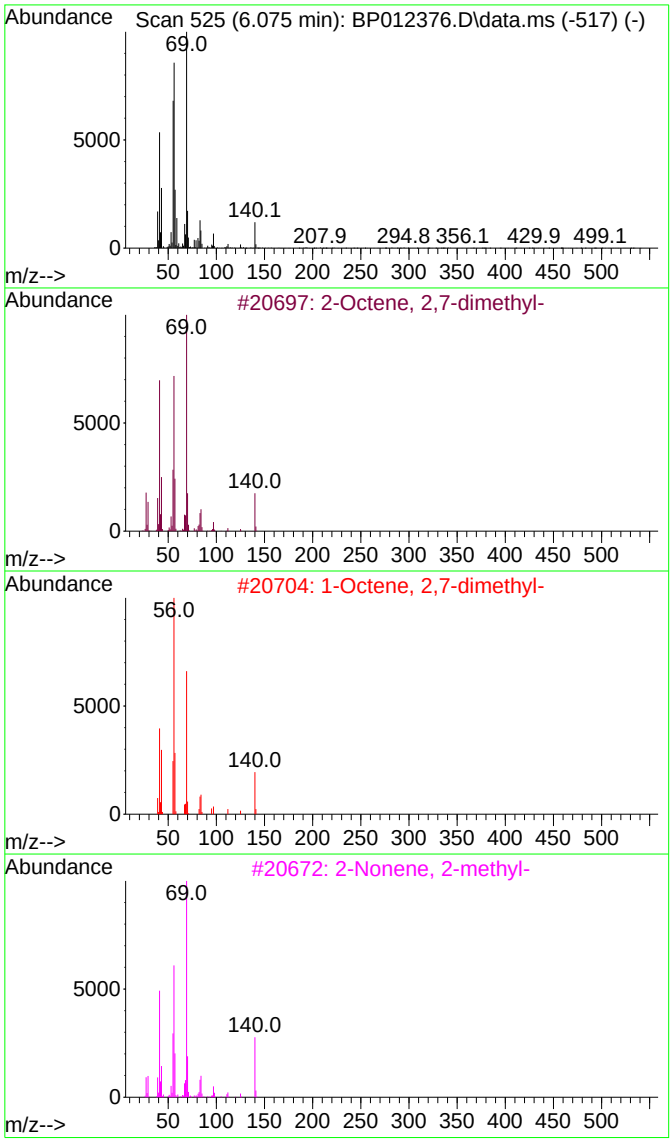
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

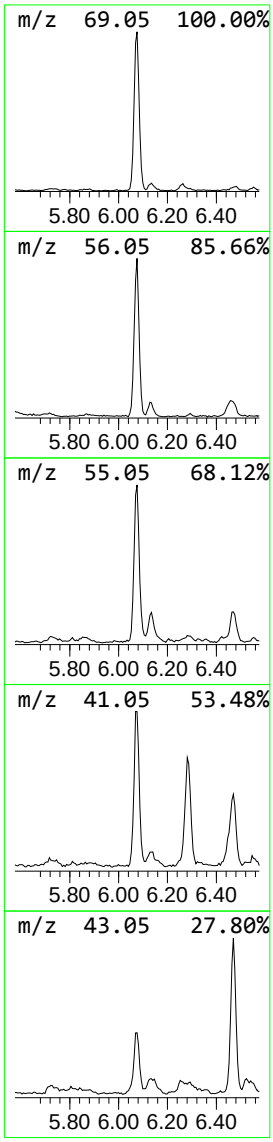
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 22

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 6.075 | 5.77 ng/ul | 355880 | 1,4-Dichlorobenzene-d4 | 7.805 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-Octene, 2,7-dimethyl- | 140 | C10H20 | 002050-81-9 | 90 |
| 2 | 1-Octene, 2,7-dimethyl- | 140 | C10H20 | 033718-03-5 | 80 |
| 3 | 2-Nonene, 2-methyl- | 140 | C10H20 | 002129-95-5 | 80 |
| 4 | 4-Nonene, 5-methyl- | 140 | C10H20 | 015918-07-7 | 72 |
| 5 | 3-Ethyl-4-octene | 140 | C10H20 | 019781-32-9 | 49 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012376.D
Acq On : 28 Oct 2022 14:05
Operator : CG/JU
Sample : N5232-16
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BHAB9

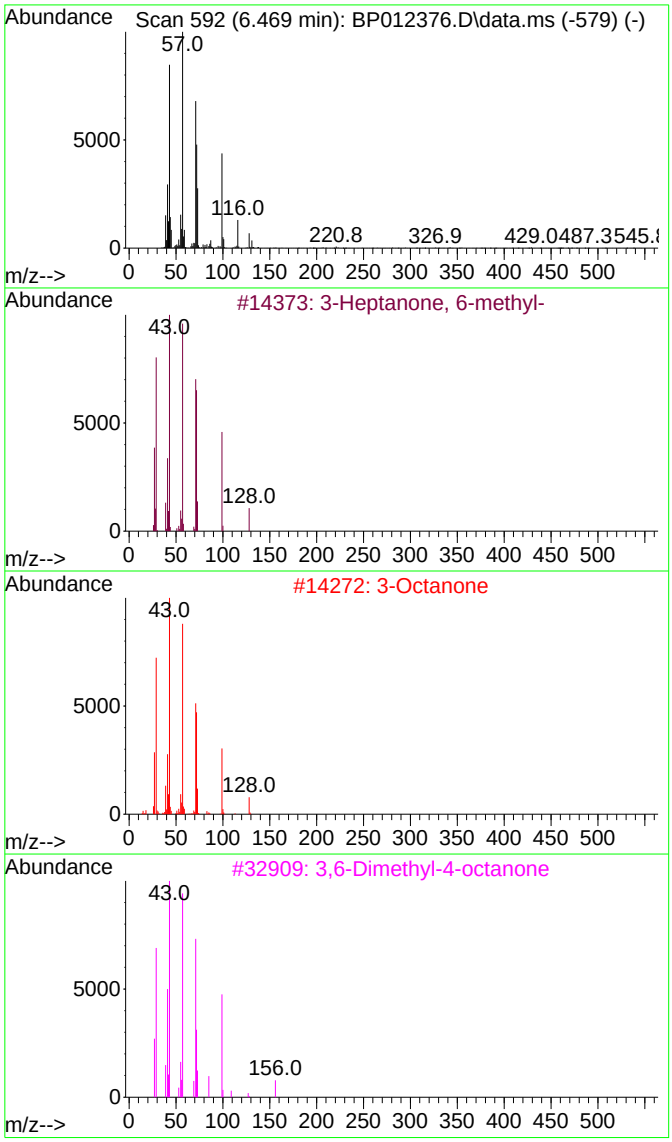
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

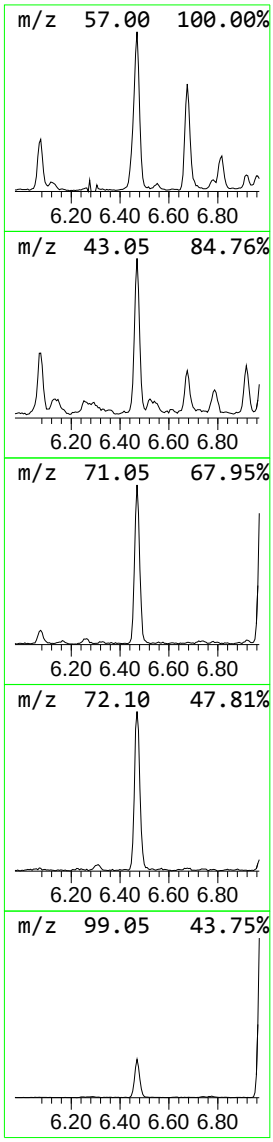
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 6 3-Heptanone, 6-methyl- Concentration Rank 21

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 6.469 | 5.87 ng/ul | 361661 | 1,4-Dichlorobenzene-d4 | 7.805 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 3-Heptanone, 6-methyl- | 128 | C8H16O | 000624-42-0 | 64 |
| 2 | 3-Octanone | 128 | C8H16O | 000106-68-3 | 58 |
| 3 | 3,6-Dimethyl-4-octanone | 156 | C10H20O | 034645-03-9 | 53 |
| 4 | 3-Octanone | 128 | C8H16O | 000106-68-3 | 50 |
| 5 | 3-Heptanone, 5-methyl- | 128 | C8H16O | 000541-85-5 | 50 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012376.D
Acq On : 28 Oct 2022 14:05
Operator : CG/JU
Sample : N5232-16
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BHAB9

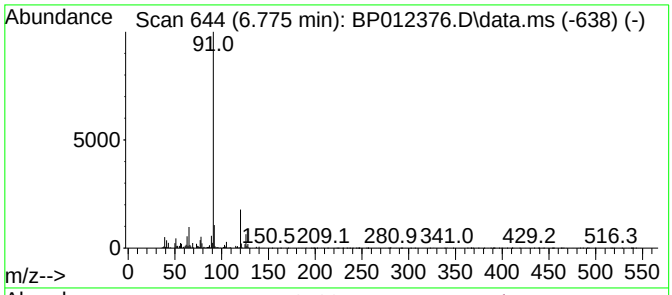
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

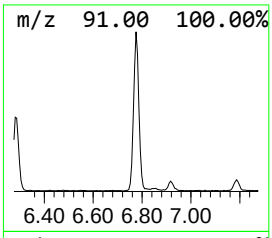
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

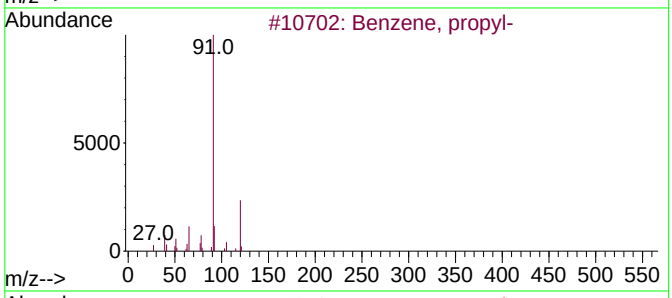
**********
Peak Number 7 Benzene, propyl- Concentration Rank 18

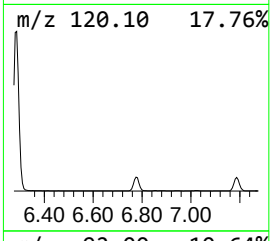
| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 6.775 | 6.40 ng/ul | 394614 | 1,4-Dichlorobenzene-d4 | 7.805 |

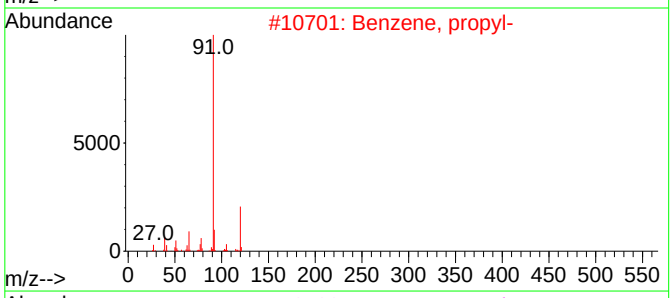
| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzene, propyl- | 120 | C9H12 | 000103-65-1 | 76 |
| 2 | Benzene, propyl- | 120 | C9H12 | 000103-65-1 | 76 |
| 3 | Benzene, propyl- | 120 | C9H12 | 000103-65-1 | 68 |
| 4 | Benzyl chloride | 126 | C7H7Cl | 000100-44-7 | 59 |
| 5 | Ethanol, 2-[(phenylmethyl)amino]- | 151 | C9H13NO | 000104-63-2 | 59 |

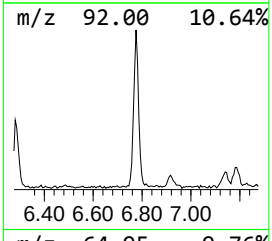


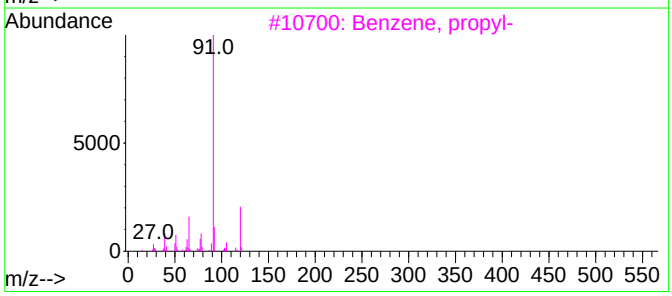



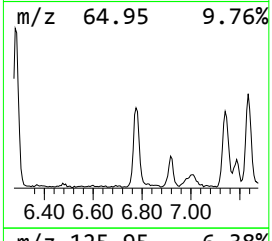



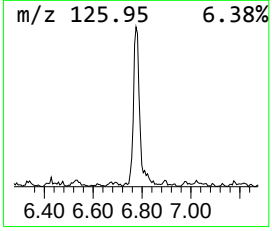



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012376.D
Acq On : 28 Oct 2022 14:05
Operator : CG/JU
Sample : N5232-16
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BHAB9

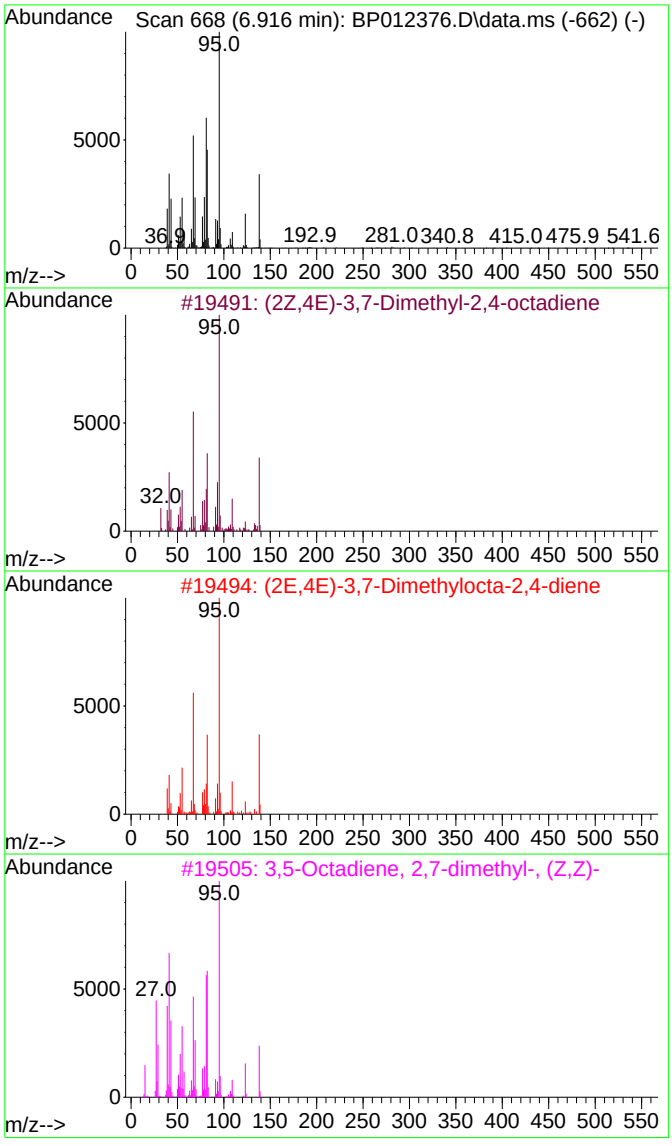
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

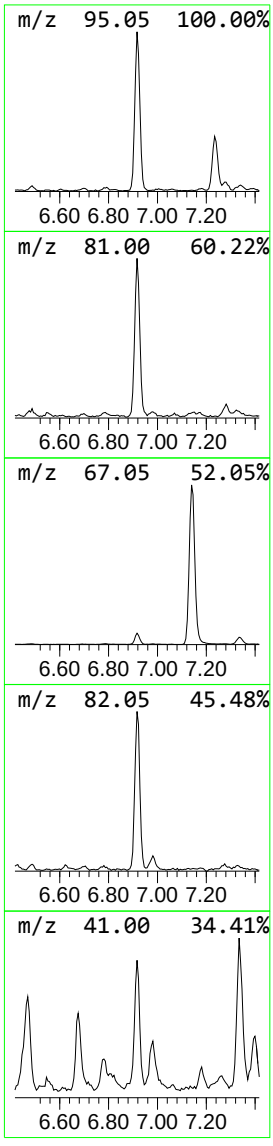
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 8 (2Z,4E)-3,7-Dimethyl-2,4-oc... Concentration Rank 24

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 6.916 | 5.57 ng/ul | 343742 | 1,4-Dichlorobenzene-d4 | 7.805 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | (2Z,4E)-3,7-Dimethyl-2,4-octadiene | 138 | C10H18 | 1000374-08-4 | 90 |
| 2 | (2E,4E)-3,7-Dimethylocta-2,4-diene | 138 | C10H18 | 006874-39-1 | 81 |
| 3 | 3,5-Octadiene, 2,7-dimethyl-, (Z... | 138 | C10H18 | 028980-73-6 | 64 |
| 4 | 3,5-Octadiene, 2,7-dimethyl-, (E... | 138 | C10H18 | 055682-64-9 | 64 |
| 5 | Cyclohexene, 1-methyl-4-(1-methy... | 138 | C10H18 | 005502-88-5 | 64 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012376.D
Acq On : 28 Oct 2022 14:05
Operator : CG/JU
Sample : N5232-16
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BHAB9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

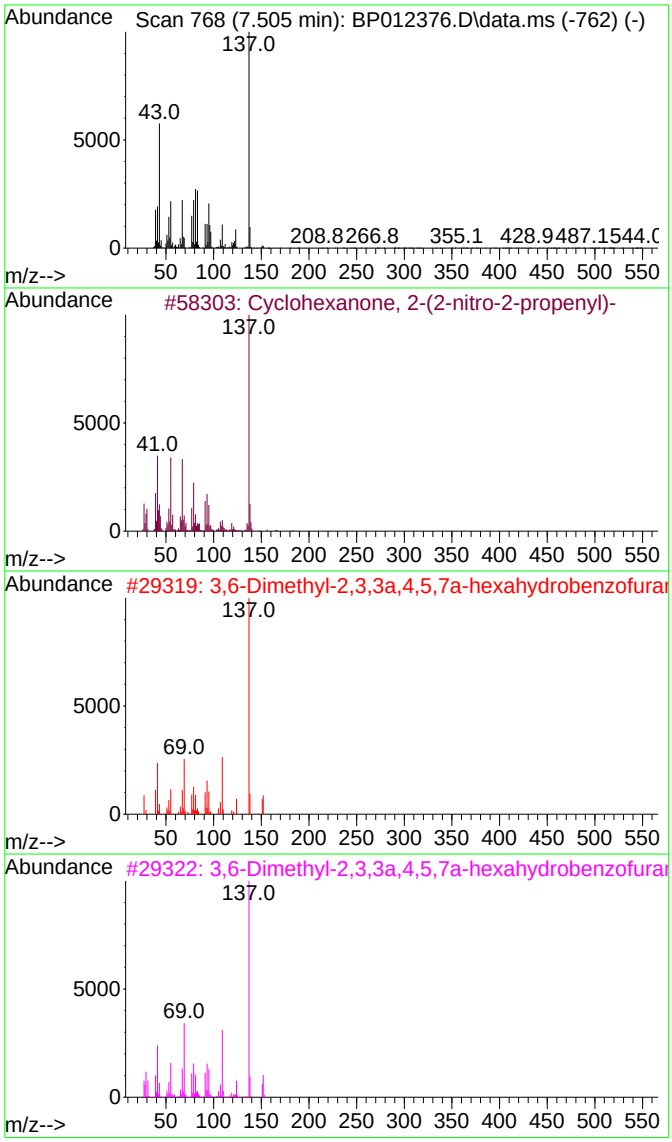
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

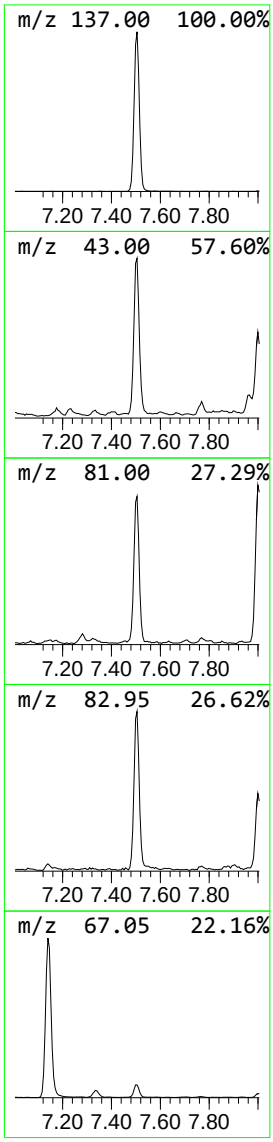
**********

Peak Number 9 unknown-02 Concentration Rank 8

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 7.505 | 13.63 ng/ul | 840752 | 1,4-Dichlorobenzene-d4 | 7.805 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Cyclohexanone, 2-(2-nitro-2-prop... | 183 | C9H13NO3 | 078551-08-3 | 45 |
| 2 | 3,6-Dimethyl-2,3,3a,4,5,7a-hexah... | 152 | C10H16O | 070786-44-6 | 45 |
| 3 | 3,6-Dimethyl-2,3,3a,4,5,7a-hexah... | 152 | C10H16O | 070786-44-6 | 42 |
| 4 | 4-Pyridinol, acetate (ester) | 137 | C7H7NO2 | 014210-20-9 | 40 |
| 5 | Dill ether | 152 | C10H16O | 074410-10-9 | 38 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012376.D
Acq On : 28 Oct 2022 14:05
Operator : CG/JU
Sample : N5232-16
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BHAB9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

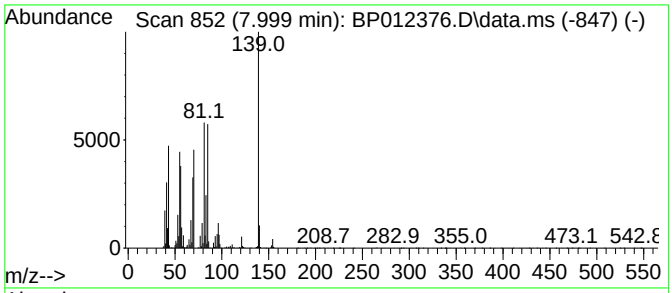
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

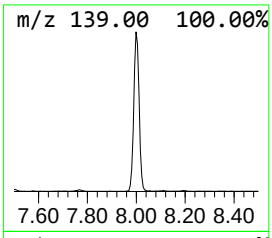
**********

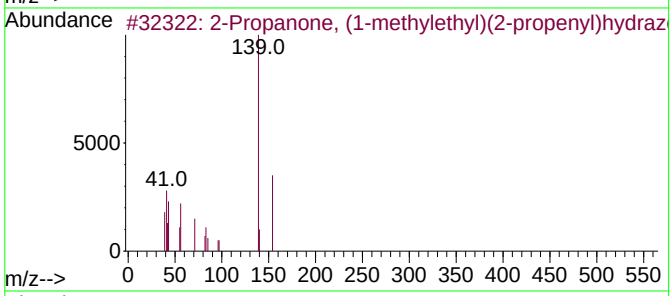
Peak Number 10 unknown-03 Concentration Rank 10

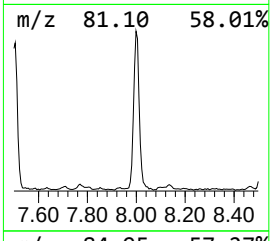
| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 7.999 | 9.36 ng/ul | 577452 | 1,4-Dichlorobenzene-d4 | 7.805 |

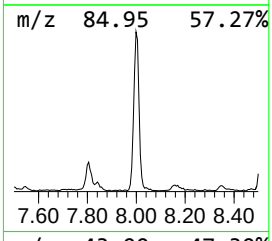
| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-Propanone, (1-methylethyl)(2-p... | 154 | C9H18N2 | 110213-09-7 | 47 |
| 2 | trans-Rose oxide | 154 | C10H18O | 000876-18-6 | 38 |
| 3 | (2S,4R)-4-Methyl-2-(2-methylprop... | 154 | C10H18O | 003033-23-6 | 37 |
| 4 | 2H-Pyran, tetrahydro-4-methyl-2-... | 154 | C10H18O | 016409-43-1 | 37 |
| 5 | 2H-Pyran, tetrahydro-4-methyl-2-... | 154 | C10H18O | 016409-43-1 | 37 |

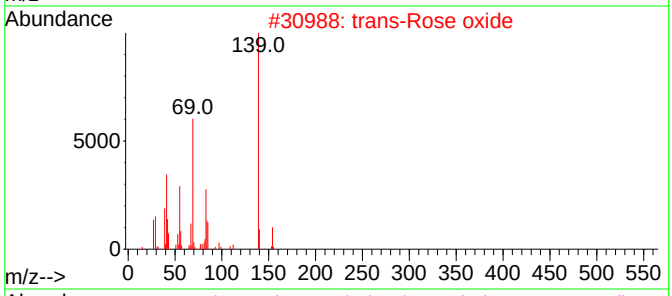


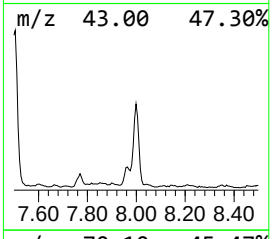



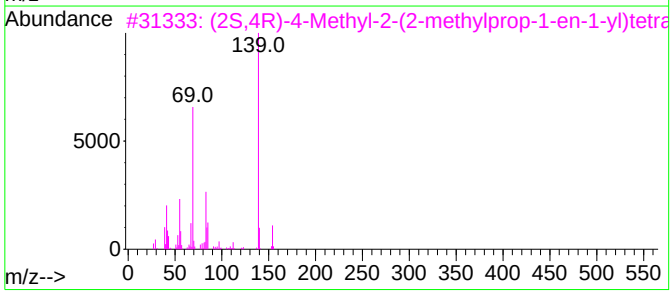



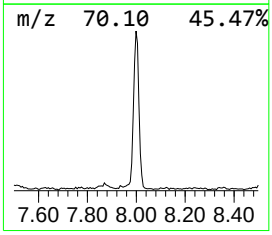



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012376.D
Acq On : 28 Oct 2022 14:05
Operator : CG/JU
Sample : N5232-16
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BHAB9

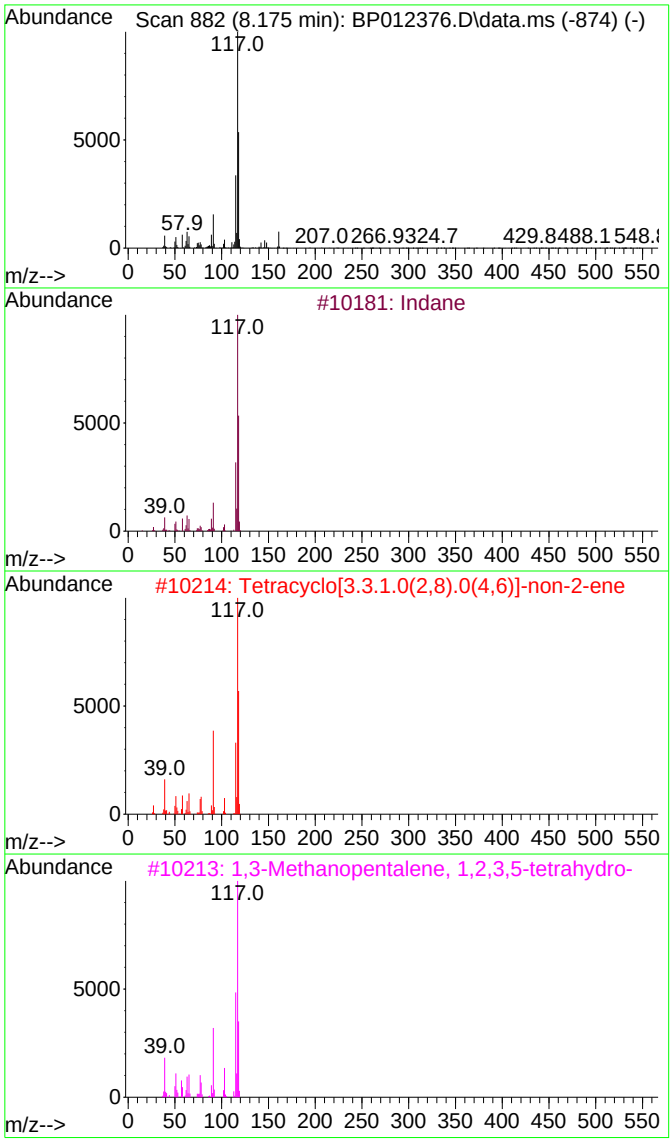
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

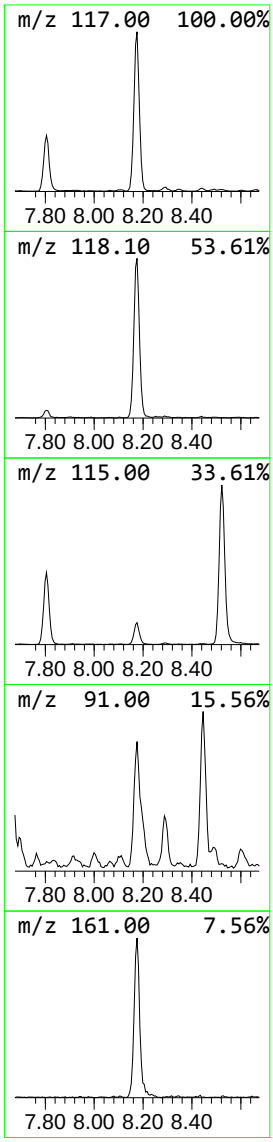
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 11 Indane Concentration Rank 13

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 8.175 | 8.35 ng/ul | 514937 | 1,4-Dichlorobenzene-d4 | 7.805 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Indane | 118 | C9H10 | 000496-11-7 | 76 |
| 2 | Tetracyclo[3.3.1.0(2,8).0(4,6)]-... | 118 | C9H10 | 1000191-13-7 | 74 |
| 3 | 1,3-Methanopentalene, 1,2,3,5-te... | 118 | C9H10 | 128600-88-4 | 64 |
| 4 | Benzene, cyclopropyl- | 118 | C9H10 | 000873-49-4 | 64 |
| 5 | Indane | 118 | C9H10 | 000496-11-7 | 60 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012376.D
Acq On : 28 Oct 2022 14:05
Operator : CG/JU
Sample : N5232-16
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BHAB9

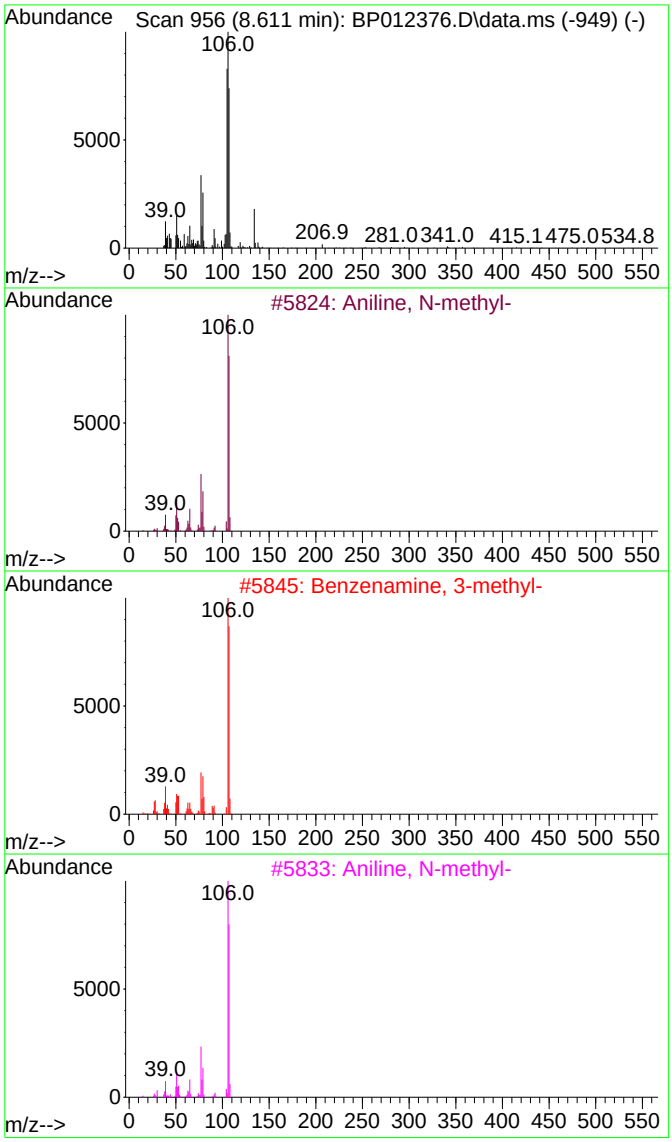
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

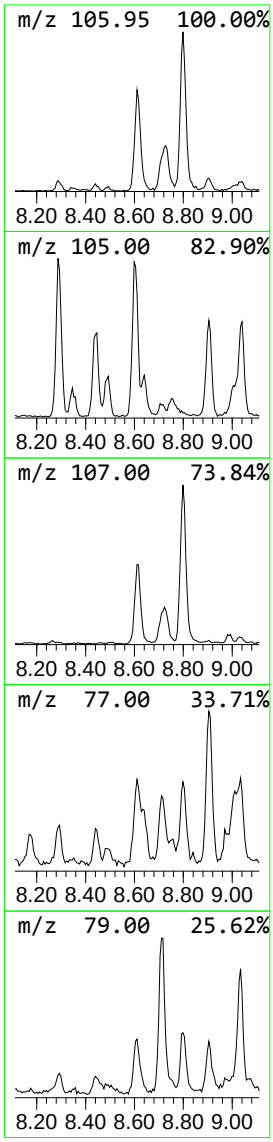
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 12 Aniline, N-methyl- Concentration Rank 33

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 8.611 | 3.90 ng/ul | 240486 | 1,4-Dichlorobenzene-d4 | 7.805 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Aniline, N-methyl- | 107 | C7H9N | 000100-61-8 | 90 |
| 2 | Benzenamine, 3-methyl- | 107 | C7H9N | 000108-44-1 | 87 |
| 3 | Aniline, N-methyl- | 107 | C7H9N | 000100-61-8 | 81 |
| 4 | p-Aminotoluene | 107 | C7H9N | 000106-49-0 | 80 |
| 5 | o-Toluidine | 107 | C7H9N | 000095-53-4 | 74 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012376.D
Acq On : 28 Oct 2022 14:05
Operator : CG/JU
Sample : N5232-16
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BHAB9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

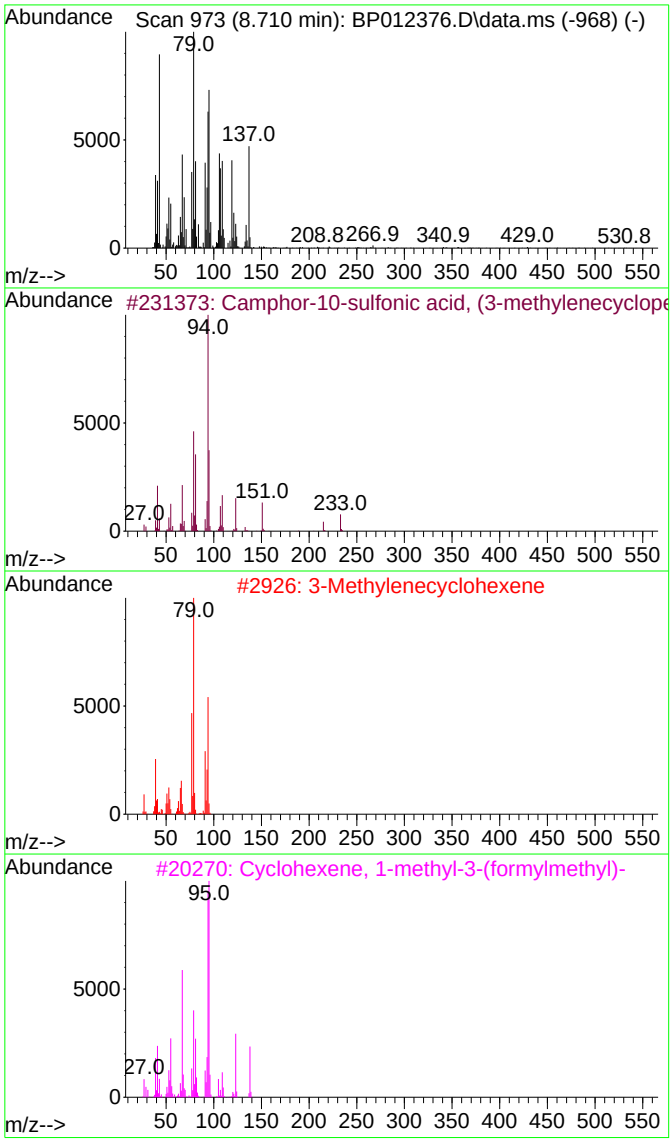
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

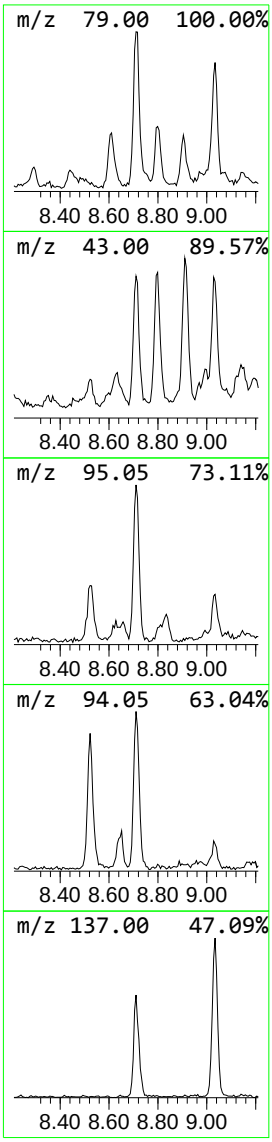
**********

Peak Number 13 unknown-04 Concentration Rank 29

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 8.710 | 4.05 ng/ul | 249891 | 1,4-Dichlorobenzene-d4 | 7.805 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Camphor-10-sulfonic acid, (3-met... | 326 | C17H26O4S | 1000156-37-8 | 43 |
| 2 | 3-Methylenecyclohexene | 94 | C7H10 | 001888-90-0 | 35 |
| 3 | Cyclohexene, 1-methyl-3-(formylm... | 138 | C9H14O | 129993-40-4 | 22 |
| 4 | Bicyclo[3.1.1]hept-3-en-2-ol, 4,... | 152 | C10H16O | 018881-04-4 | 22 |
| 5 | Bicyclo[4.1.0]hept-2-ene | 94 | C7H10 | 002566-57-6 | 20 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012376.D
Acq On : 28 Oct 2022 14:05
Operator : CG/JU
Sample : N5232-16
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BHAB9

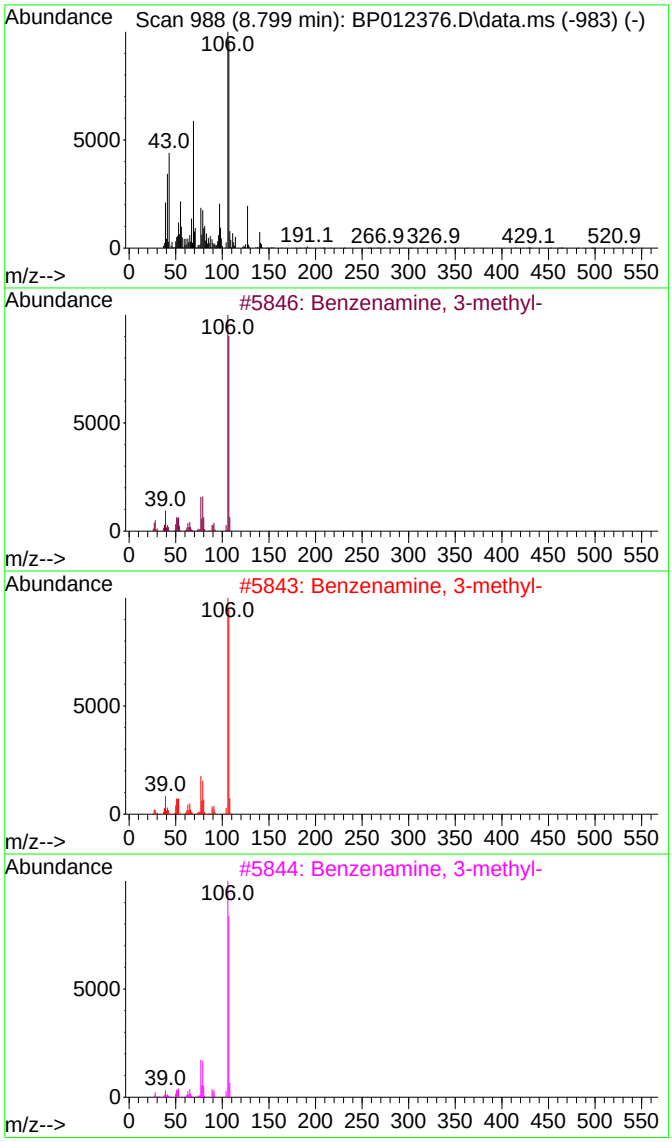
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

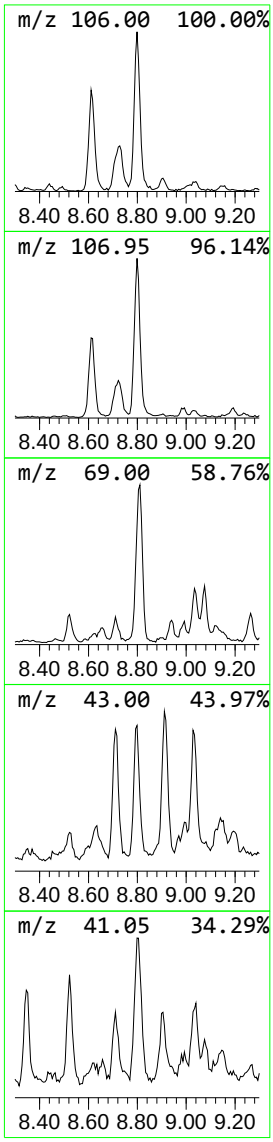
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 14 Benzenamine, 3-methyl- Concentration Rank 20

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 8.799 | 6.12 ng/ul | 377240 | 1,4-Dichlorobenzene-d4 | 7.805 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzenamine, 3-methyl- | 107 | C7H9N | 000108-44-1 | 93 |
| 2 | Benzenamine, 3-methyl- | 107 | C7H9N | 000108-44-1 | 93 |
| 3 | Benzenamine, 3-methyl- | 107 | C7H9N | 000108-44-1 | 92 |
| 4 | o-Toluidine | 107 | C7H9N | 000095-53-4 | 86 |
| 5 | o-Toluidine | 107 | C7H9N | 000095-53-4 | 70 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012376.D
Acq On : 28 Oct 2022 14:05
Operator : CG/JU
Sample : N5232-16
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BHAB9

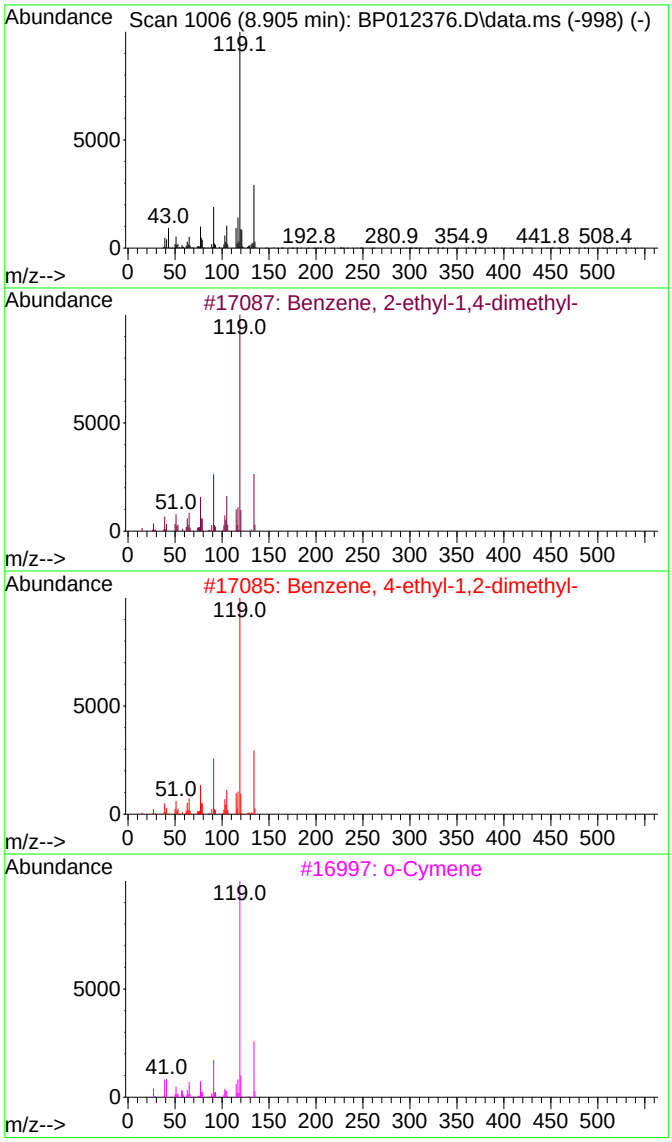
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

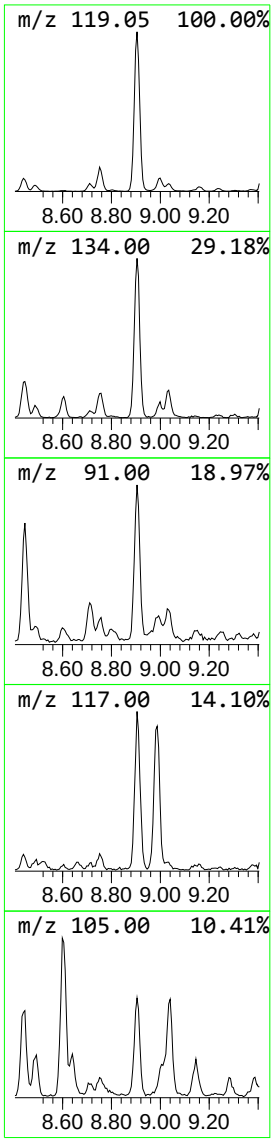
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 15 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 11

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 8.905 | 8.84 ng/ul | 544981 | 1,4-Dichlorobenzene-d4 | 7.805 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14 | 001758-88-9 | 96 |
| 2 | Benzene, 4-ethyl-1,2-dimethyl- | 134 | C10H14 | 000934-80-5 | 96 |
| 3 | o-Cymene | 134 | C10H14 | 000527-84-4 | 95 |
| 4 | p-Cymene | 134 | C10H14 | 000099-87-6 | 95 |
| 5 | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14 | 000933-98-2 | 95 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012376.D
Acq On : 28 Oct 2022 14:05
Operator : CG/JU
Sample : N5232-16
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BHAB9

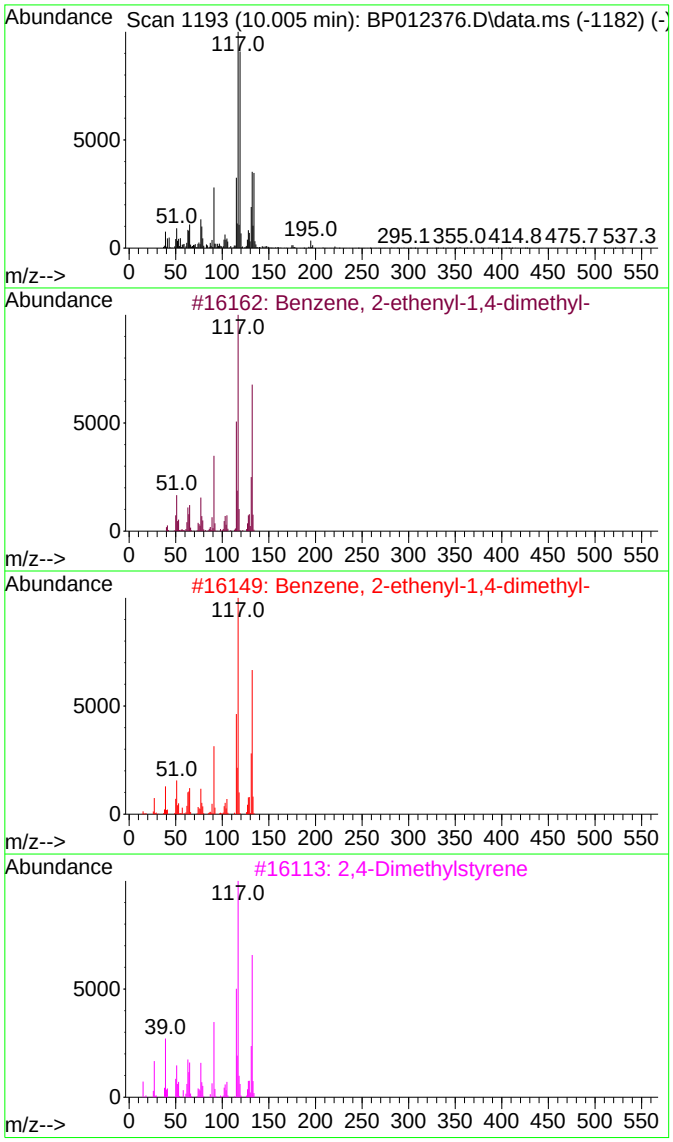
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

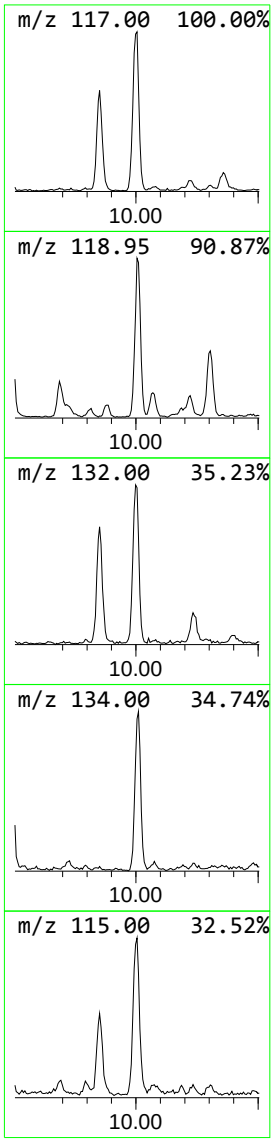
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 19 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 19

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.005 | 6.37 ng/ul | 695551 | Naphthalene-d8 | 10.610 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzene, 2-ethenyl-1,4-dimethyl- | 132 | C10H12 | 002039-89-6 | 86 |
| 2 | Benzene, 2-ethenyl-1,4-dimethyl- | 132 | C10H12 | 002039-89-6 | 83 |
| 3 | 2,4-Dimethylstyrene | 132 | C10H12 | 002234-20-0 | 80 |
| 4 | Benzene, (1-methyl-1-propenyl)-,... | 132 | C10H12 | 000768-00-3 | 55 |
| 5 | Benzene, 1-methyl-2-(2-propenyl)- | 132 | C10H12 | 001587-04-8 | 55 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012376.D
Acq On : 28 Oct 2022 14:05
Operator : CG/JU
Sample : N5232-16
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BHAB9

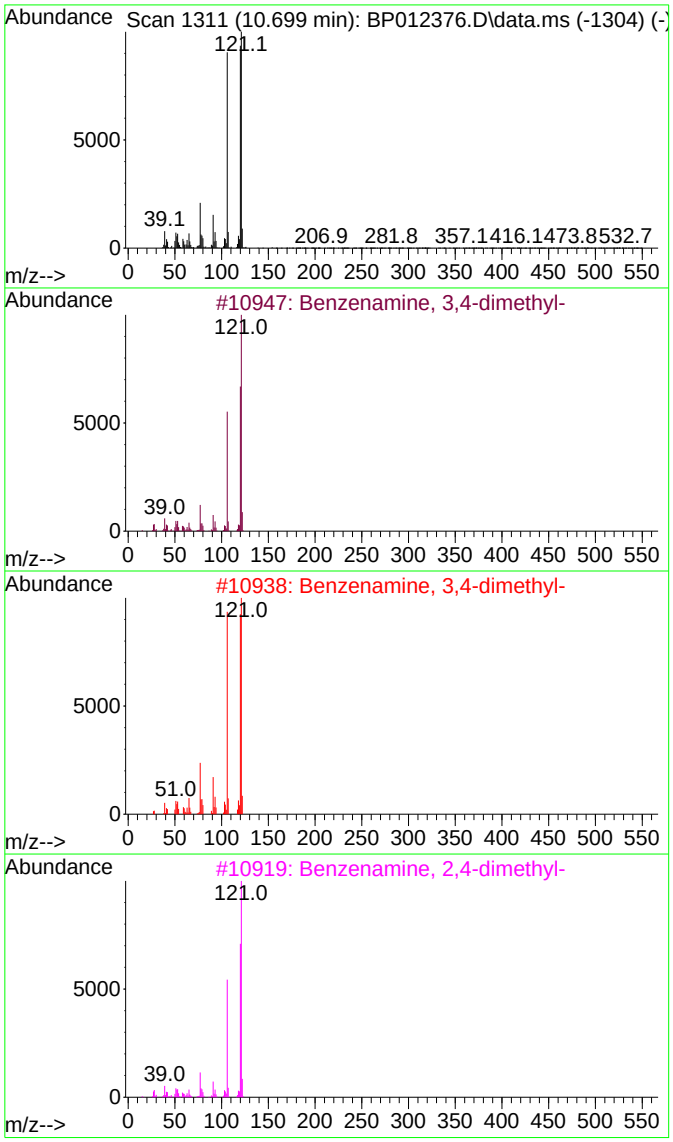
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

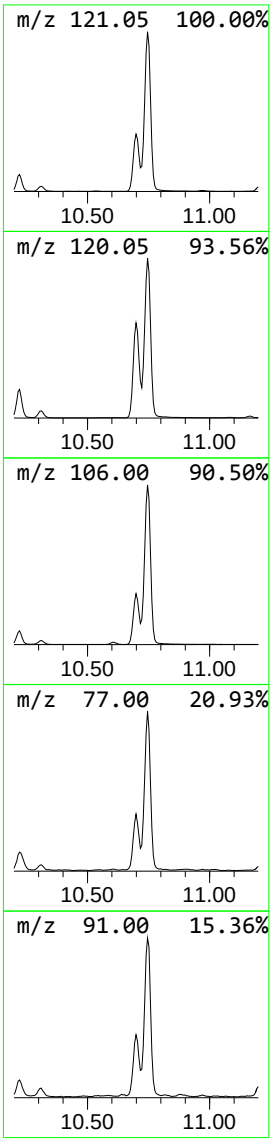
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 23 Benzenamine, 3,4-dimethyl- Concentration Rank 3

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.699 | 27.24 ng/ul | 2973710 | Naphthalene-d8 | 10.610 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzenamine, 3,4-dimethyl- | 121 | C8H11N | 000095-64-7 | 97 |
| 2 | Benzenamine, 3,4-dimethyl- | 121 | C8H11N | 000095-64-7 | 96 |
| 3 | Benzenamine, 2,4-dimethyl- | 121 | C8H11N | 000095-68-1 | 96 |
| 4 | Benzenamine, 3,4-dimethyl- | 121 | C8H11N | 000095-64-7 | 96 |
| 5 | Benzenamine, 2,4-dimethyl- | 121 | C8H11N | 000095-68-1 | 96 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012376.D
Acq On : 28 Oct 2022 14:05
Operator : CG/JU
Sample : N5232-16
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BHAB9

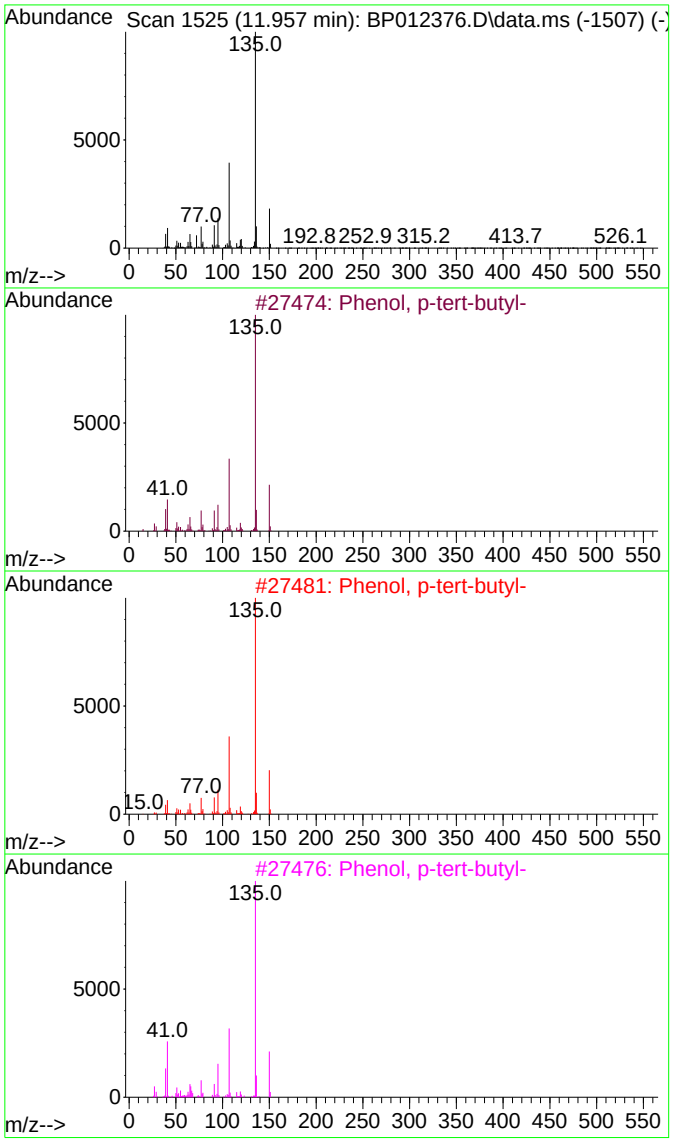
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

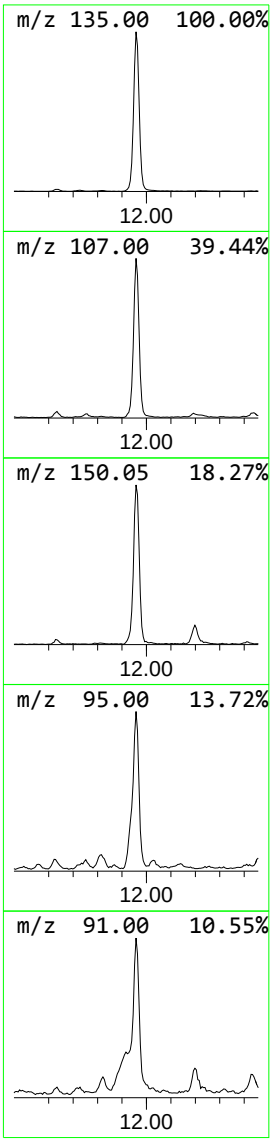
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 26 Phenol, p-tert-butyl- Concentration Rank 4

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.957 | 25.36 ng/ul | 2768930 | Naphthalene-d8 | 10.610 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenol, p-tert-butyl- | 150 | C10H14O | 000098-54-4 | 97 |
| 2 | Phenol, p-tert-butyl- | 150 | C10H14O | 000098-54-4 | 95 |
| 3 | Phenol, p-tert-butyl- | 150 | C10H14O | 000098-54-4 | 91 |
| 4 | Phenol, m-tert-butyl- | 150 | C10H14O | 000585-34-2 | 87 |
| 5 | Phenol, m-tert-butyl- | 150 | C10H14O | 000585-34-2 | 81 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012376.D
Acq On : 28 Oct 2022 14:05
Operator : CG/JU
Sample : N5232-16
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BHAB9

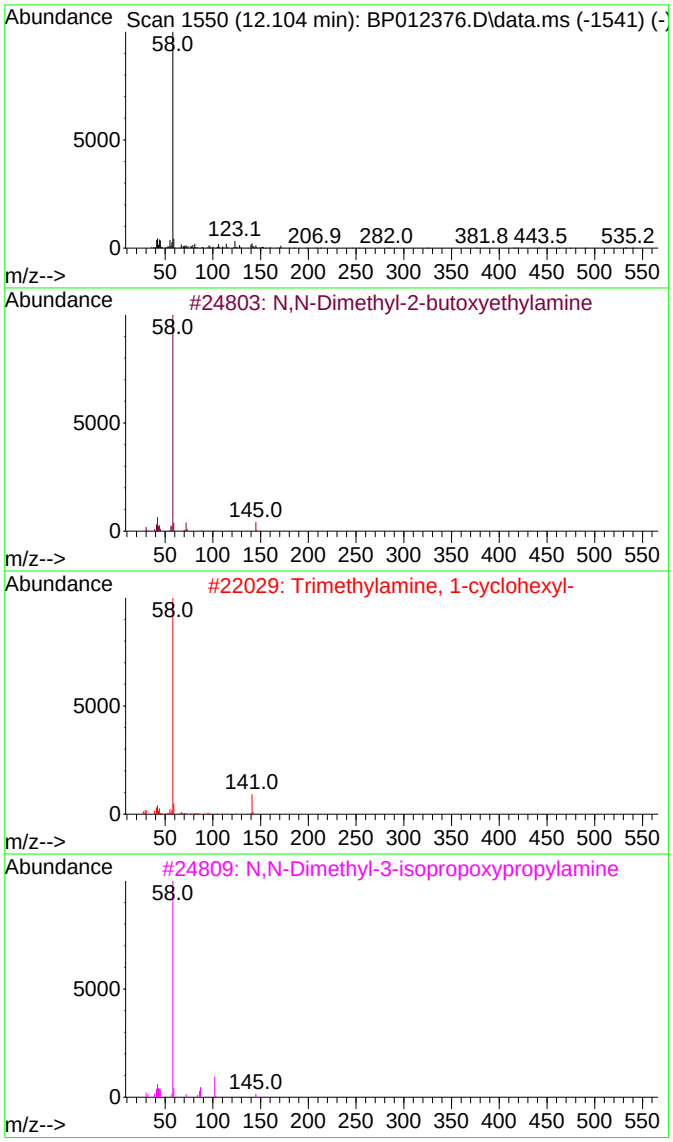
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

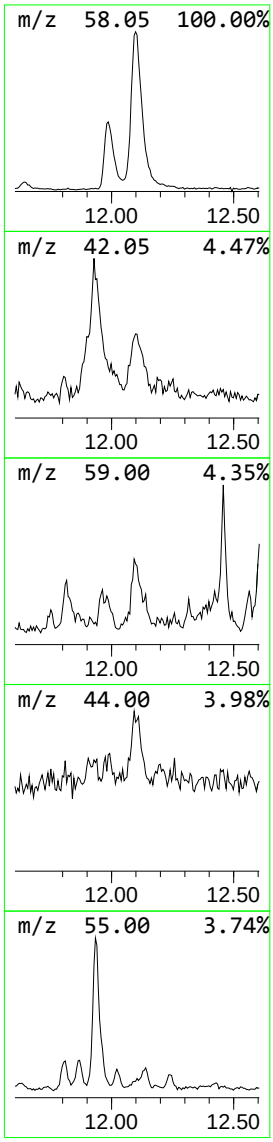
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 27 N,N-Dimethyl-2-butoxyethyla... Concentration Rank 30

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.104 | 3.96 ng/ul | 432049 | Naphthalene-d8 | 10.610 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | N,N-Dimethyl-2-butoxyethylamine | 145 | C8H19NO | 071126-58-4 | 72 |
| 2 | Trimethylamine, 1-cyclohexyl- | 141 | C9H19N | 016607-80-0 | 72 |
| 3 | N,N-Dimethyl-3-isopropoxypropyla... | 145 | C8H19NO | 071126-66-4 | 64 |
| 4 | 2-[(dimethylamino)methyl]cyclope... | 141 | C8H15NO | 006947-99-5 | 64 |
| 5 | Benzedrex | 155 | C10H21N | 000101-40-6 | 64 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012376.D
Acq On : 28 Oct 2022 14:05
Operator : CG/JU
Sample : N5232-16
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BHAB9

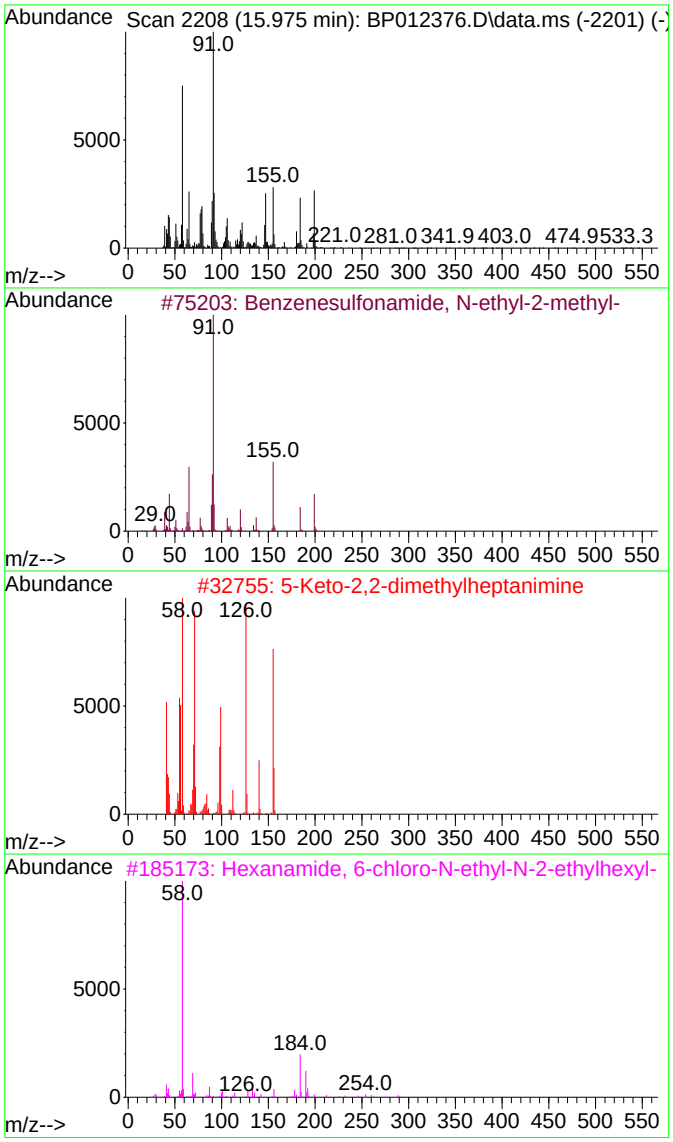
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

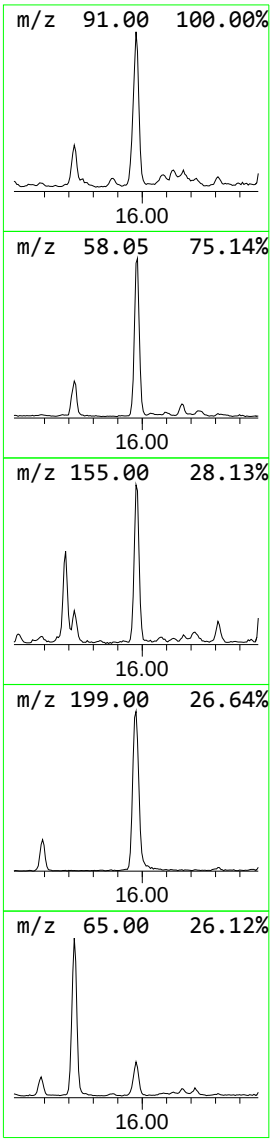
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 32 Benzenesulfonamide, N-ethyl... Concentration Rank 16

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.975 | 7.27 ng/ul | 1488340 | Phenanthrene-d10 | 17.216 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzenesulfonamide, N-ethyl-2-me... | 199 | C9H13NO2S | 001077-56-1 | 72 |
| 2 | 5-Keto-2,2-dimethylheptanimine | 155 | C9H17NO | 1010129-16-7 | 38 |
| 3 | Hexanamide, 6-chloro-N-ethyl-N-2... | 289 | C16H32ClNO | 1000415-59-3 | 35 |
| 4 | 3-Hexyne, 2,2-dimethyl-1-dimethy... | 153 | C10H19N | 1000150-17-9 | 25 |
| 5 | N-Methyl-bk-MMDA-2 | 237 | C12H15NO4 | 1010475-85-3 | 25 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012376.D
Acq On : 28 Oct 2022 14:05
Operator : CG/JU
Sample : N5232-16
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BHAB9

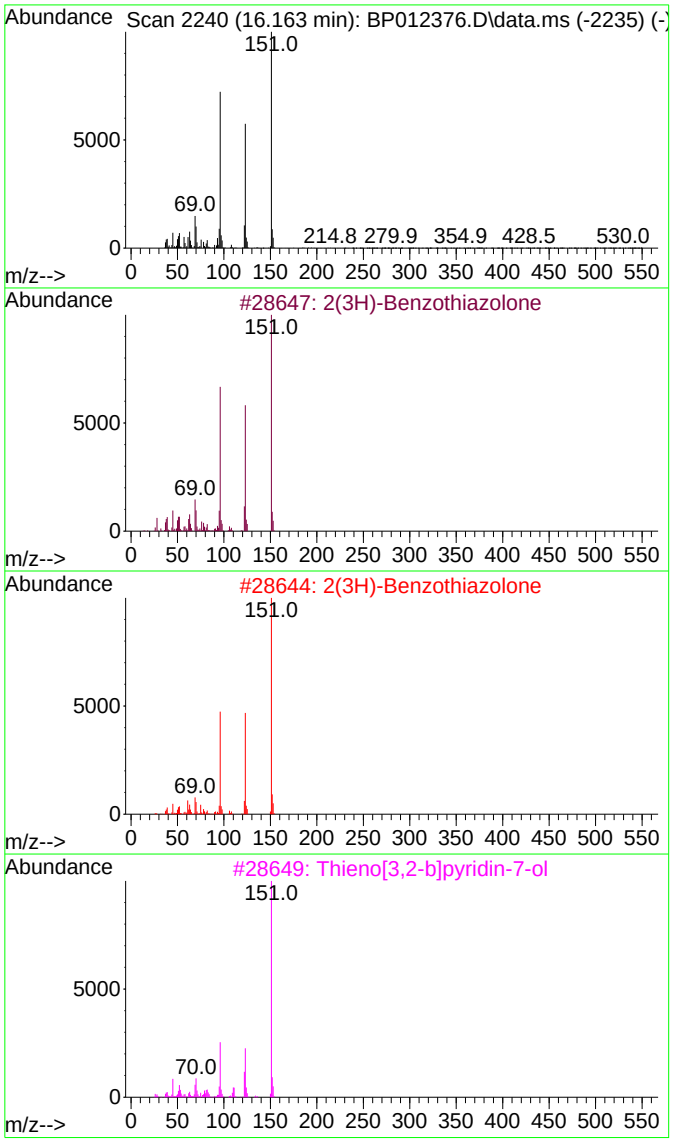
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

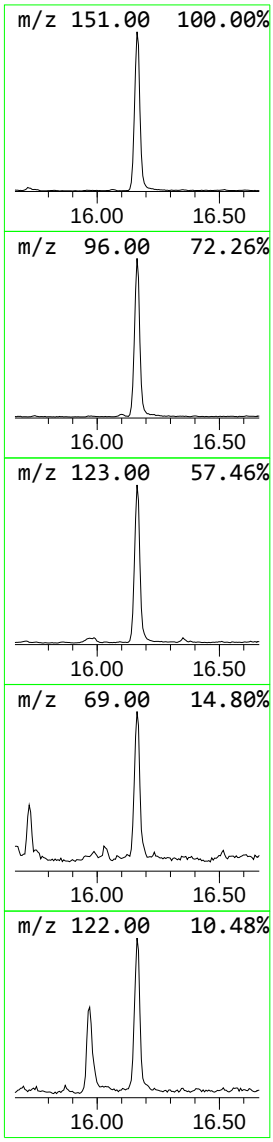
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 33 2(3H)-Benzothiazolone Concentration Rank 15

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.163 | 7.35 ng/ul | 1505210 | Phenanthrene-d10 | 17.216 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2(3H)-Benzothiazolone | 151 | C7H5NOS | 000934-34-9 | 97 |
| 2 | 2(3H)-Benzothiazolone | 151 | C7H5NOS | 000934-34-9 | 91 |
| 3 | Thieno[3,2-b]pyridin-7-ol | 151 | C7H5NOS | 107818-20-2 | 64 |
| 4 | 6-Methyl-7-oxo-4,7-dihydro-triaz... | 151 | C5H5N5O | 057250-39-2 | 50 |
| 5 | 4-Mercaptoimidazo[4,5-c]pyridine | 151 | C6H5N3S | 034550-51-1 | 38 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012376.D
Acq On : 28 Oct 2022 14:05
Operator : CG/JU
Sample : N5232-16
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BHAB9

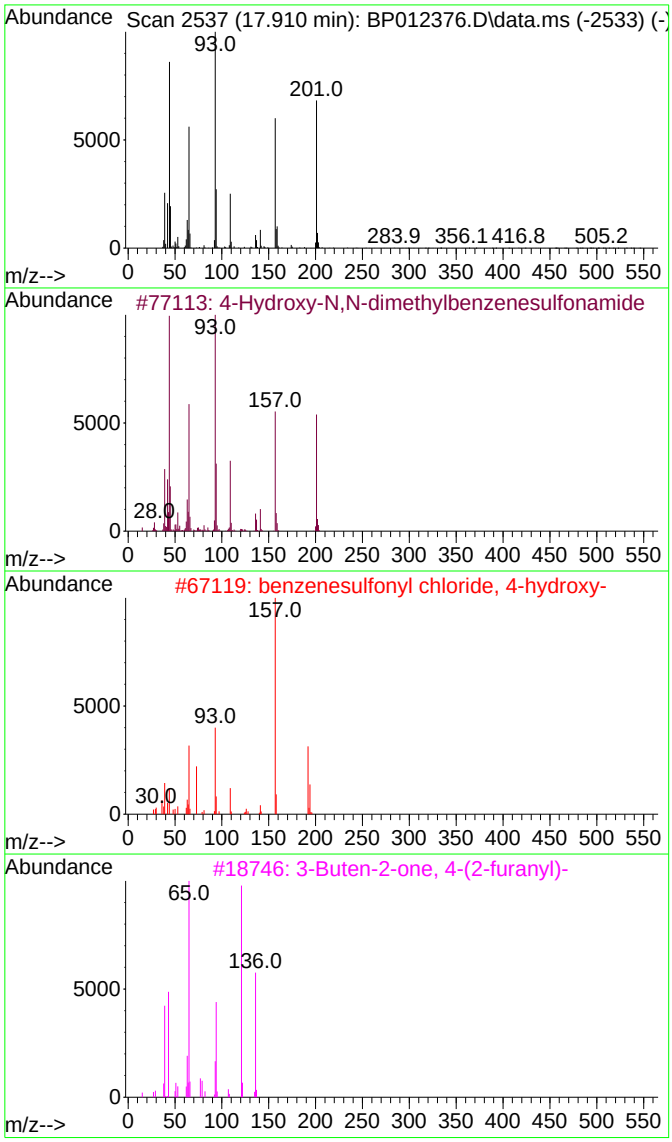
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

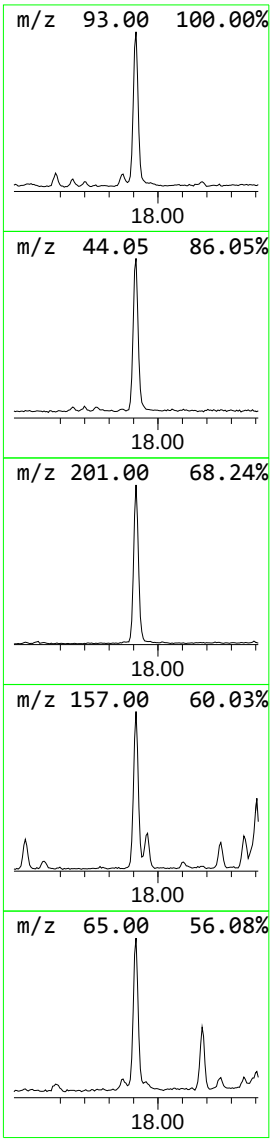
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 35 4-Hydroxy-N,N-dimethylbenze... Concentration Rank 25

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 17.910 | 4.78 ng/ul | 978289 | Phenanthrene-d10 | 17.216 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 4-Hydroxy-N,N-dimethylbenzenesul... | 201 | C8H11NO3S | 015020-57-2 | 90 |
| 2 | benzenesulfonyl chloride, 4-hydr... | 192 | C6H5ClO3S | 1000400-26-6 | 28 |
| 3 | 3-Buten-2-one, 4-(2-furanyl)- | 136 | C8H8O2 | 000623-15-4 | 16 |
| 4 | Phenoxyacetonitrile | 133 | C8H7NO | 003598-14-9 | 12 |
| 5 | Benzene, 1-methyl-2-nitroso- | 121 | C7H7NO | 000611-23-4 | 10 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012376.D
Acq On : 28 Oct 2022 14:05
Operator : CG/JU
Sample : N5232-16
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BHAB9

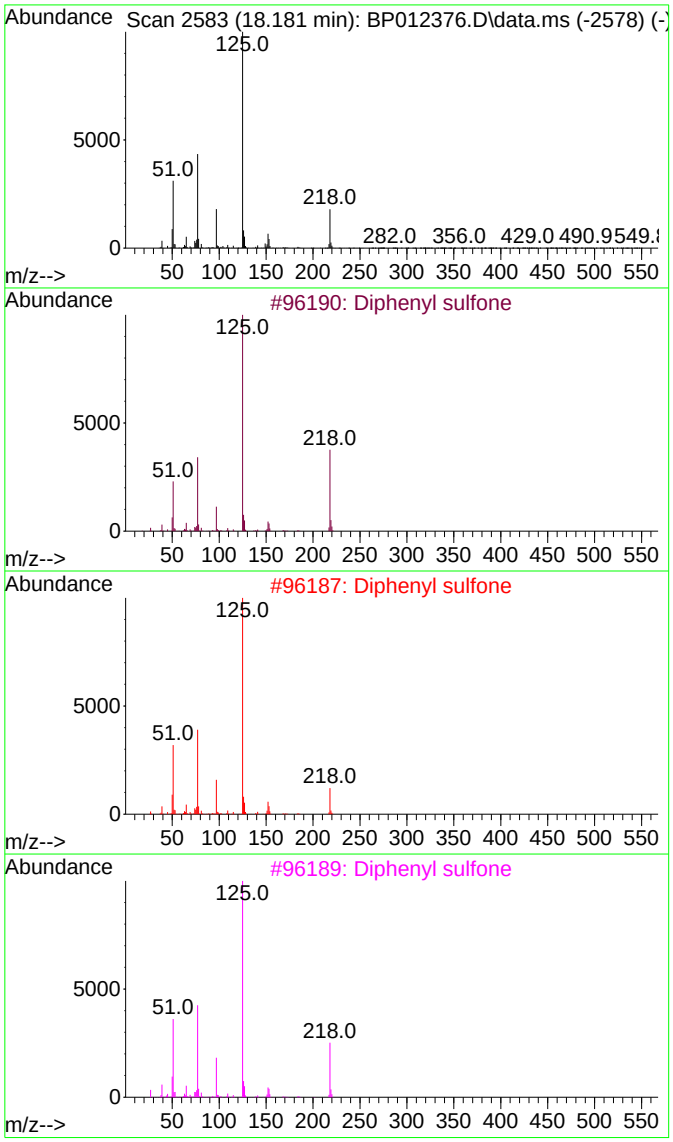
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

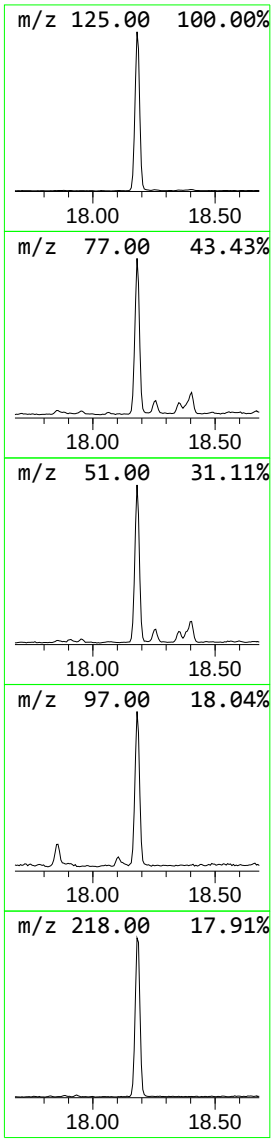
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 36 Diphenyl sulfone Concentration Rank 12

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.181 | 8.37 ng/ul | 1714320 | Phenanthrene-d10 | 17.216 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Diphenyl sulfone | 218 | C12H10O2S | 000127-63-9 | 96 |
| 2 | Diphenyl sulfone | 218 | C12H10O2S | 000127-63-9 | 96 |
| 3 | Diphenyl sulfone | 218 | C12H10O2S | 000127-63-9 | 95 |
| 4 | 2-Propenenitrile, 3-(phenylsulfo... | 193 | C9H7NO2S | 001424-51-7 | 72 |
| 5 | Diphenyl sulfone | 218 | C12H10O2S | 000127-63-9 | 70 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012376.D
Acq On : 28 Oct 2022 14:05
Operator : CG/JU
Sample : N5232-16
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BHAB9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

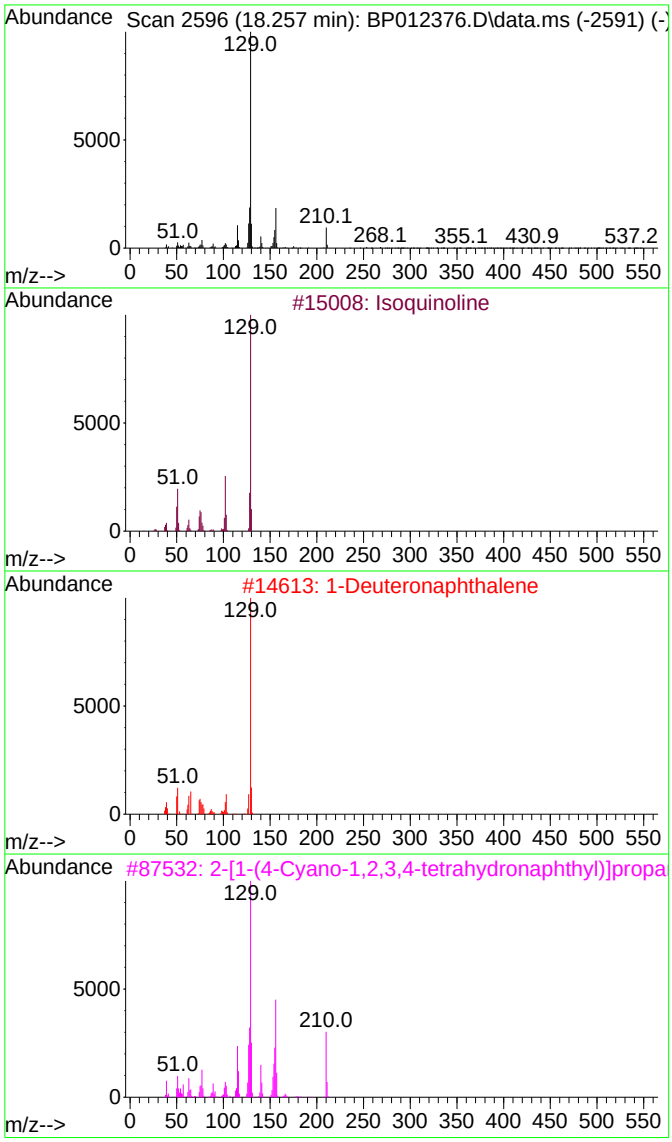
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

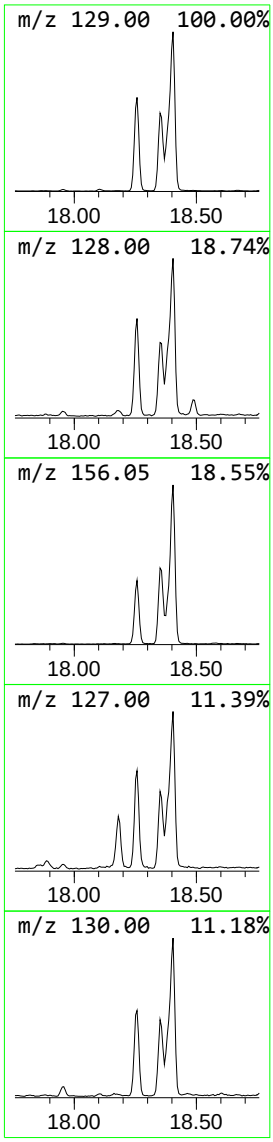
**********

Peak Number 37 Isoquinoline Concentration Rank 17

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.257 | 6.62 ng/ul | 1356550 | Phenanthrene-d10 | 17.216 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Isoquinoline | 129 | C9H7N | 000119-65-3 | 52 |
| 2 | 1-Deuteronaphthalene | 129 | C10H7D | 000875-62-7 | 50 |
| 3 | 2-[1-(4-Cyano-1,2,3,4-tetrahydro... | 210 | C14H14N2 | 057964-39-3 | 43 |
| 4 | Benzeneacetonitrile, .alpha.-met... | 129 | C9H7N | 000495-10-3 | 40 |
| 5 | Isoquinoline | 129 | C9H7N | 000119-65-3 | 40 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012376.D
Acq On : 28 Oct 2022 14:05
Operator : CG/JU
Sample : N5232-16
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BHAB9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

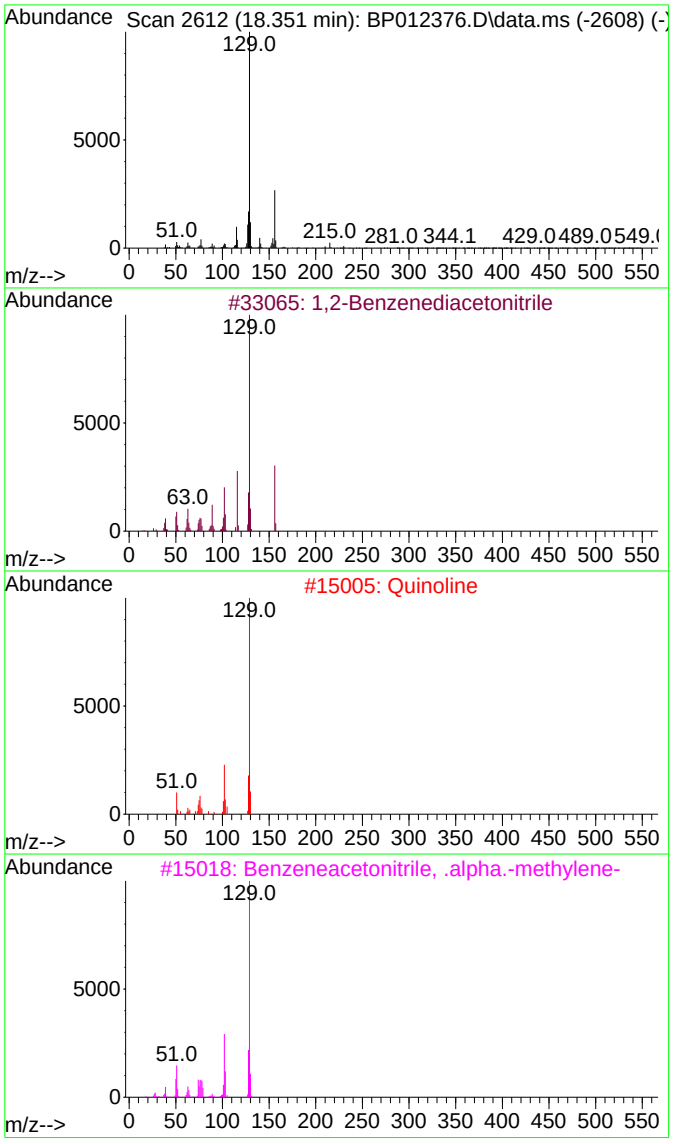
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

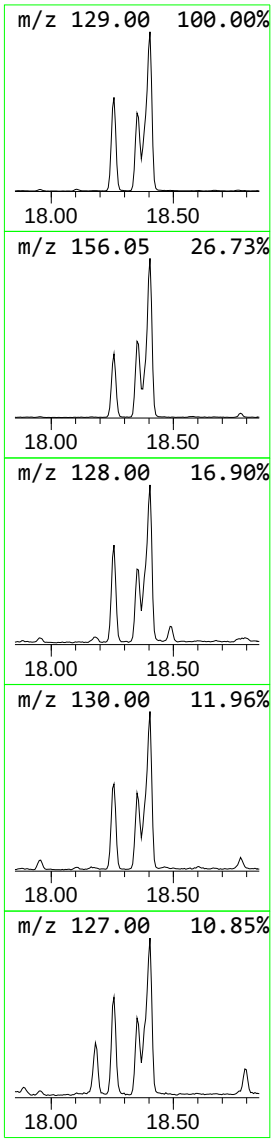
**********

Peak Number 38 1,2-Benzenediacetonitrile Concentration Rank 23

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.351 | 5.69 ng/ul | 1165630 | Phenanthrene-d10 | 17.216 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1,2-Benzenediacetonitrile | 156 | C10H8N2 | 000613-73-0 | 59 |
| 2 | Quinoline | 129 | C9H7N | 000091-22-5 | 50 |
| 3 | Benzeneacetonitrile, .alpha.-met... | 129 | C9H7N | 000495-10-3 | 47 |
| 4 | 2-[1-(4-Cyano-1,2,3,4-tetrahydro... | 210 | C14H14N2 | 057964-39-3 | 46 |
| 5 | 3-[1-(4-Cyano-1,2,3,4-tetrahydro... | 210 | C14H14N2 | 057964-40-6 | 45 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012376.D
Acq On : 28 Oct 2022 14:05
Operator : CG/JU
Sample : N5232-16
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BHAB9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

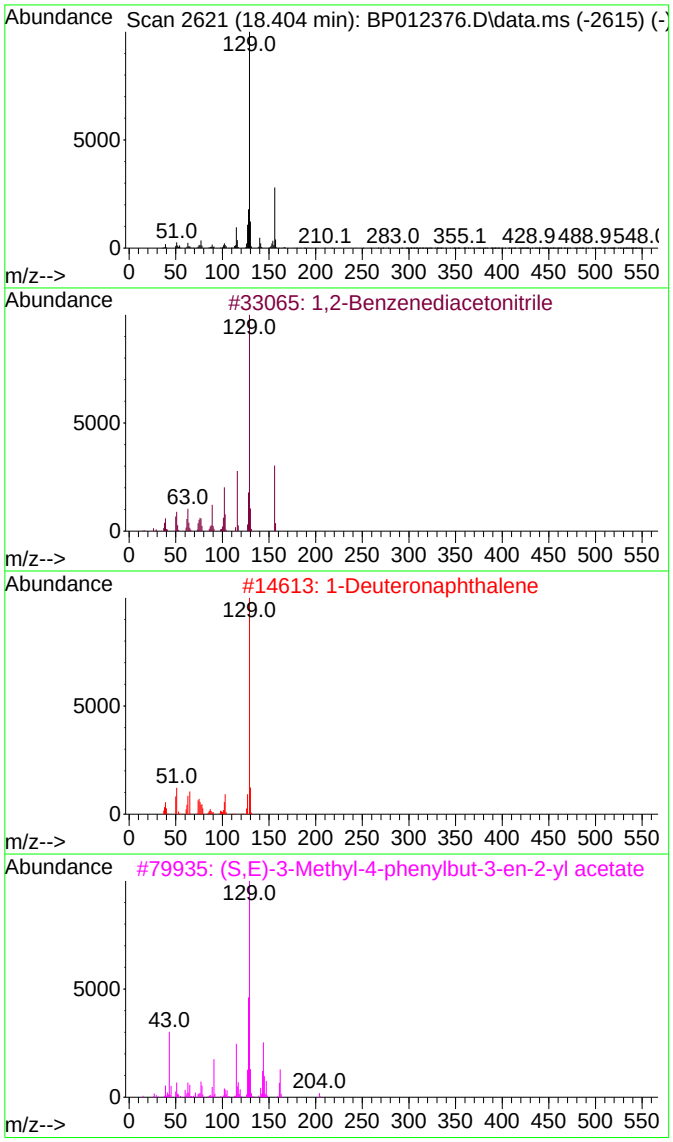
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

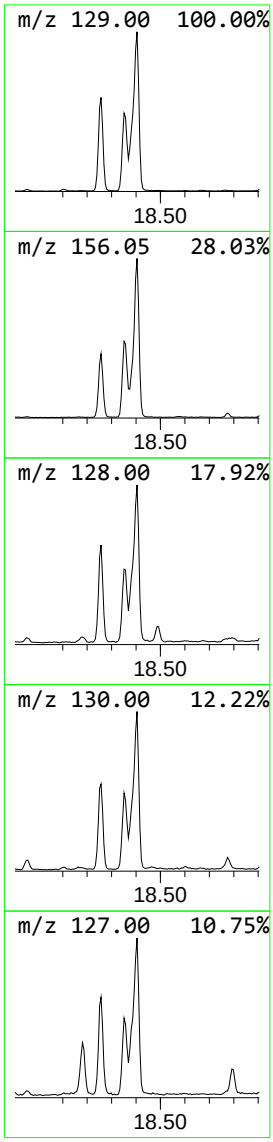
**********

Peak Number 39 unknown-05 Concentration Rank 9

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.404 | 13.50 ng/ul | 2766080 | Phenanthrene-d10 | 17.216 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1,2-Benzenediacetonitrile | 156 | C10H8N2 | 000613-73-0 | 50 |
| 2 | 1-Deuteronaphthalene | 129 | C10H7D | 000875-62-7 | 47 |
| 3 | (S,E)-3-Methyl-4-phenylbut-3-en-... | 204 | C13H16O2 | 187736-05-6 | 40 |
| 4 | Isoquinoline | 129 | C9H7N | 000119-65-3 | 40 |
| 5 | (R,E)-3-Methyl-4-phenylbut-3-en-... | 204 | C13H16O2 | 187735-90-6 | 40 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012376.D
Acq On : 28 Oct 2022 14:05
Operator : CG/JU
Sample : N5232-16
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BHAB9

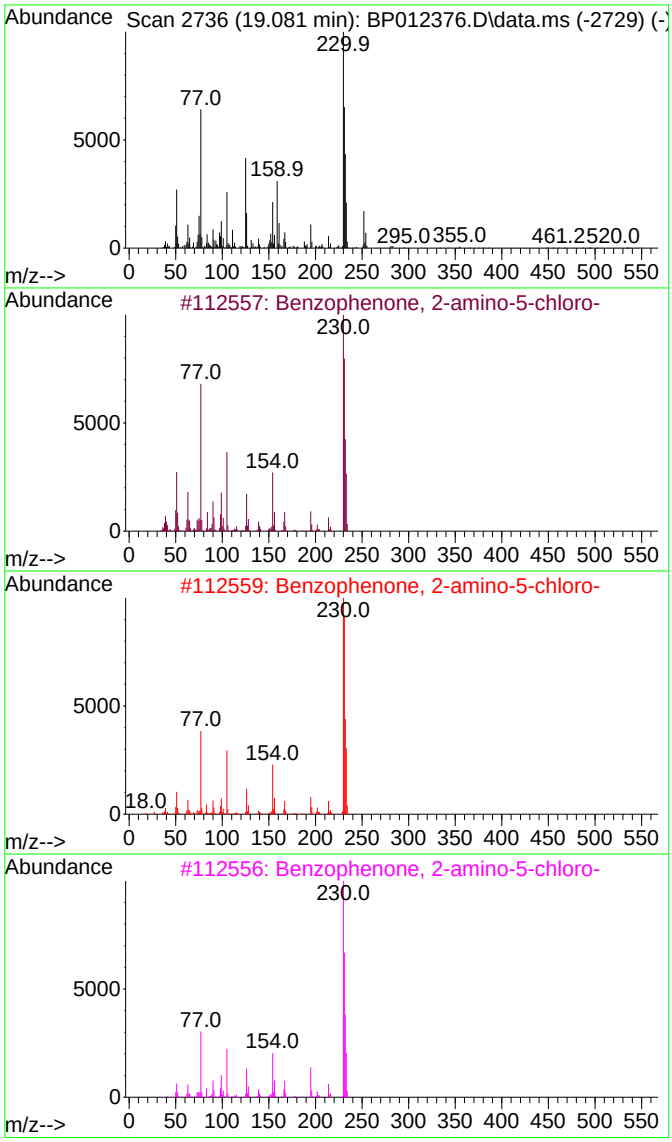
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

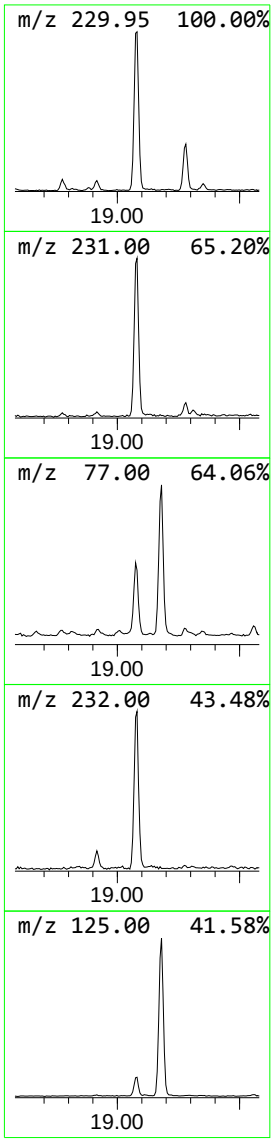
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 42 Benzophenone, 2-amino-5-chl... Concentration Rank 27

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 19.081 | 4.50 ng/ul | 921231 | Phenanthrene-d10 | 17.216 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzophenone, 2-amino-5-chloro- | 231 | C13H10ClNO | 000719-59-5 | 98 |
| 2 | Benzophenone, 2-amino-5-chloro- | 231 | C13H10ClNO | 000719-59-5 | 98 |
| 3 | Benzophenone, 2-amino-5-chloro- | 231 | C13H10ClNO | 000719-59-5 | 98 |
| 4 | Benz[c]isoxazole, 1,3-dihydro-5-... | 231 | C13H10ClNO | 1010266-69-2 | 98 |
| 5 | Benzophenone, 2-amino-5-chloro- | 231 | C13H10ClNO | 000719-59-5 | 98 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012376.D
Acq On : 28 Oct 2022 14:05
Operator : CG/JU
Sample : N5232-16
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BHAB9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

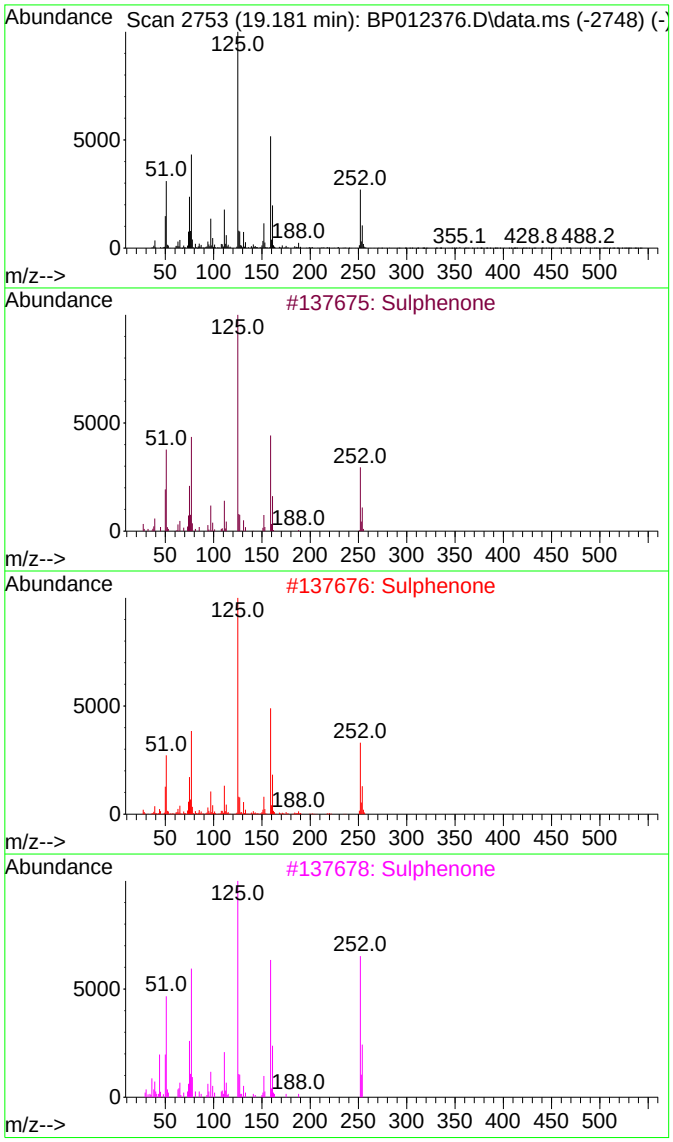
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

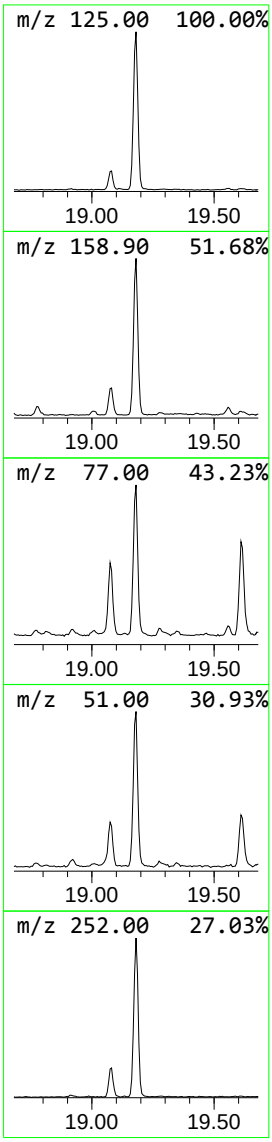
**********

Peak Number 43 Sulphenone Concentration Rank 14

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 19.180 | 8.33 ng/ul | 1707250 | Phenanthrene-d10 | 17.216 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Sulphenone | 252 | C12H9ClO2S | 000080-00-2 | 96 |
| 2 | Sulphenone | 252 | C12H9ClO2S | 000080-00-2 | 95 |
| 3 | Sulphenone | 252 | C12H9ClO2S | 000080-00-2 | 94 |
| 4 | Sulphenone | 252 | C12H9ClO2S | 000080-00-2 | 94 |
| 5 | Diphenyl sulfone | 218 | C12H10O2S | 000127-63-9 | 35 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012376.D
Acq On : 28 Oct 2022 14:05
Operator : CG/JU
Sample : N5232-16
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BHAB9

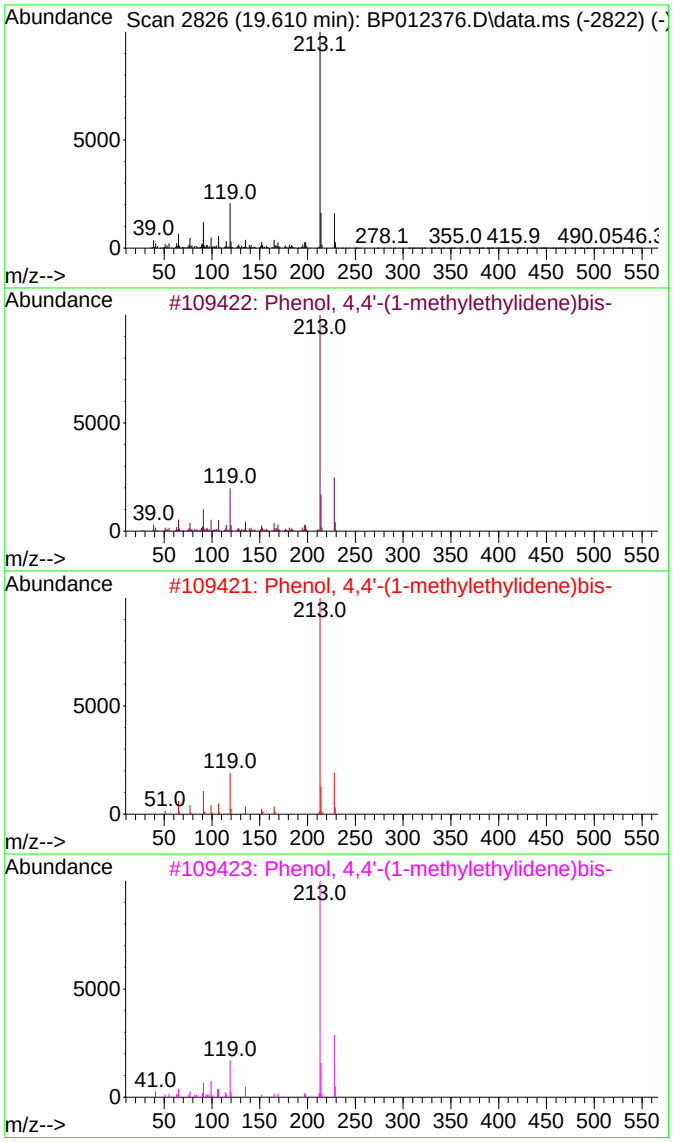
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

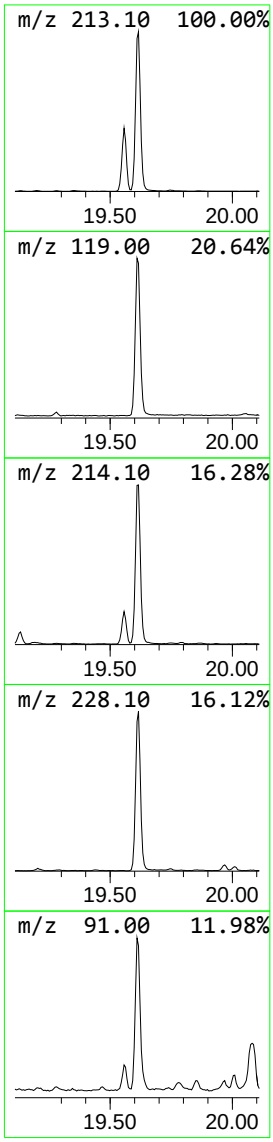
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 46 Phenol, 4,4'-(1-methylethyl... Concentration Rank 5

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 19.610 | 21.14 ng/ul | 5450580 | Chrysene-d12 | 21.310 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenol, 4,4'-(1-methylethylidene... | 228 | C15H16O2 | 000080-05-7 | 96 |
| 2 | Phenol, 4,4'-(1-methylethylidene... | 228 | C15H16O2 | 000080-05-7 | 95 |
| 3 | Phenol, 4,4'-(1-methylethylidene... | 228 | C15H16O2 | 000080-05-7 | 90 |
| 4 | Phenol, 4,4'-(1-methylethylidene... | 228 | C15H16O2 | 000080-05-7 | 83 |
| 5 | 3,4'-Isopropylidenediphenol | 228 | C15H16O2 | 046765-25-7 | 70 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012376.D
Acq On : 28 Oct 2022 14:05
Operator : CG/JU
Sample : N5232-16
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BHAB9

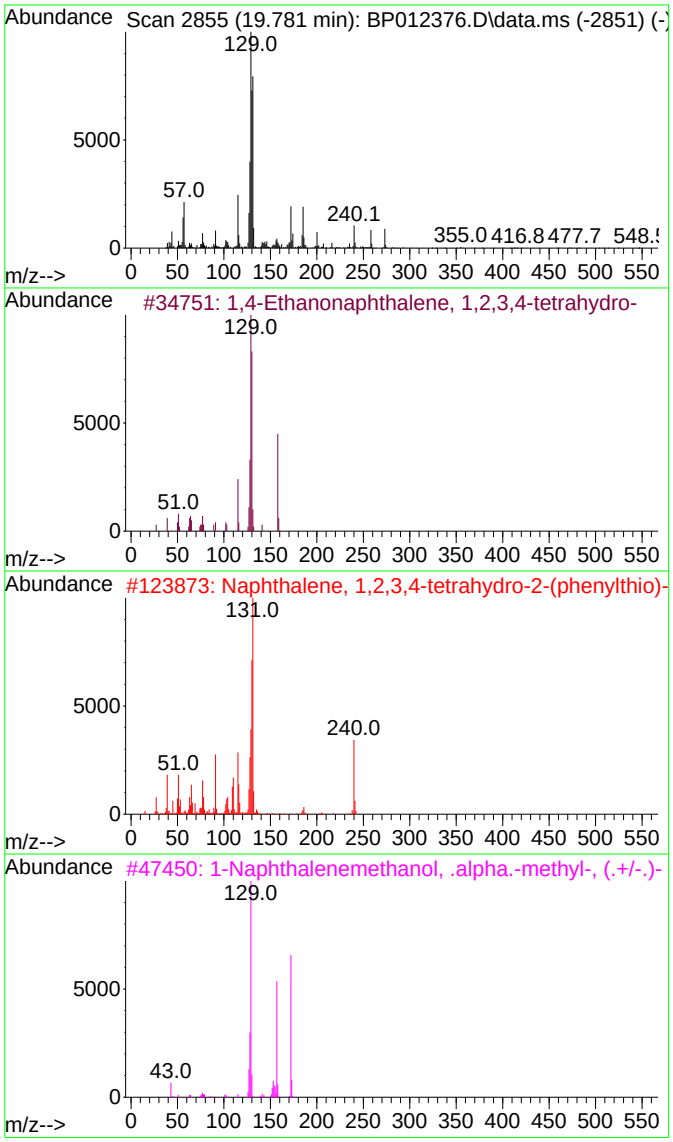
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

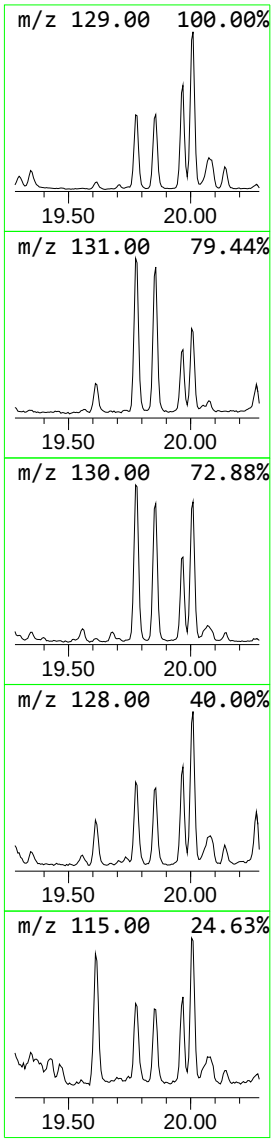
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 47 1,4-Ethanonaphthalene, 1,2,... Concentration Rank 31

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 19.781 | 3.93 ng/ul | 1014380 | Chrysene-d12 | 21.310 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1,4-Ethanonaphthalene, 1,2,3,4-t... | 158 | C12H14 | 004175-52-4 | 55 |
| 2 | Naphthalene, 1,2,3,4-tetrahydro-... | 240 | C16H16S | 023853-68-1 | 46 |
| 3 | 1-Naphthalenemethanol, .alpha.-m... | 172 | C12H12O | 057605-95-5 | 43 |
| 4 | 1,4-Ethanonaphthalen-2(1H)-one, ... | 172 | C12H12O | 013153-76-9 | 38 |
| 5 | (1-Ethylbuta-1,3-dienyl)benzene | 158 | C12H14 | 1000210-05-6 | 38 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012376.D
Acq On : 28 Oct 2022 14:05
Operator : CG/JU
Sample : N5232-16
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BHAB9

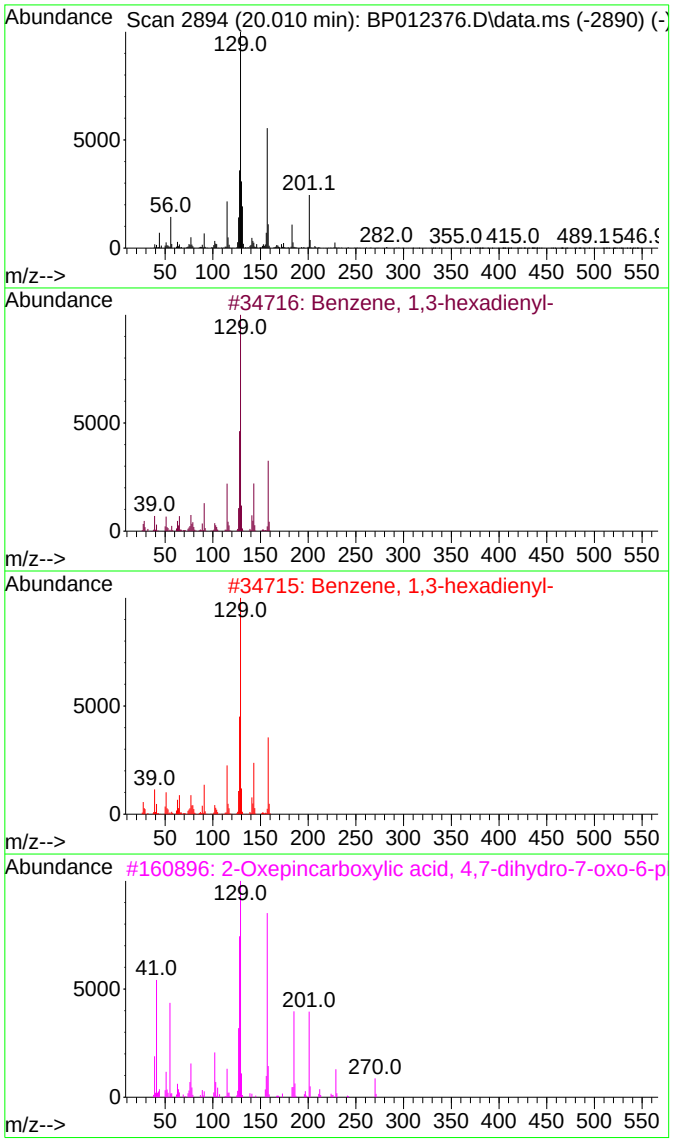
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

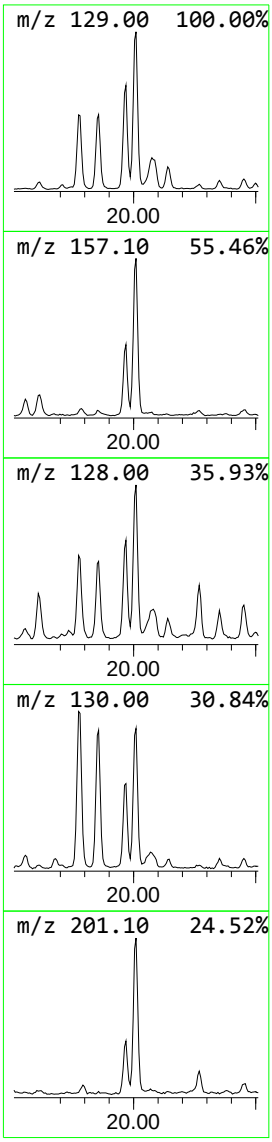
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 50 Benzene, 1,3-hexadienyl- Concentration Rank 28

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 20.010 | 4.14 ng/ul | 1068610 | Chrysene-d12 | 21.310 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzene, 1,3-hexadienyl- | 158 | C12H14 | 041635-77-2 | 50 |
| 2 | Benzene, 1,3-hexadienyl- | 158 | C12H14 | 041635-77-2 | 45 |
| 3 | 2-Oxepincarboxylic acid, 4,7-dih... | 270 | C16H14O4 | 1000142-99-5 | 40 |
| 4 | 2-Quinolinecarboxylic acid, meth... | 187 | C11H9NO2 | 019575-07-6 | 40 |
| 5 | Naphthalene, 1,2-dichloro-1,2,3,... | 200 | C10H10Cl2 | 031448-63-2 | 37 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012376.D
Acq On : 28 Oct 2022 14:05
Operator : CG/JU
Sample : N5232-16
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BHAB9

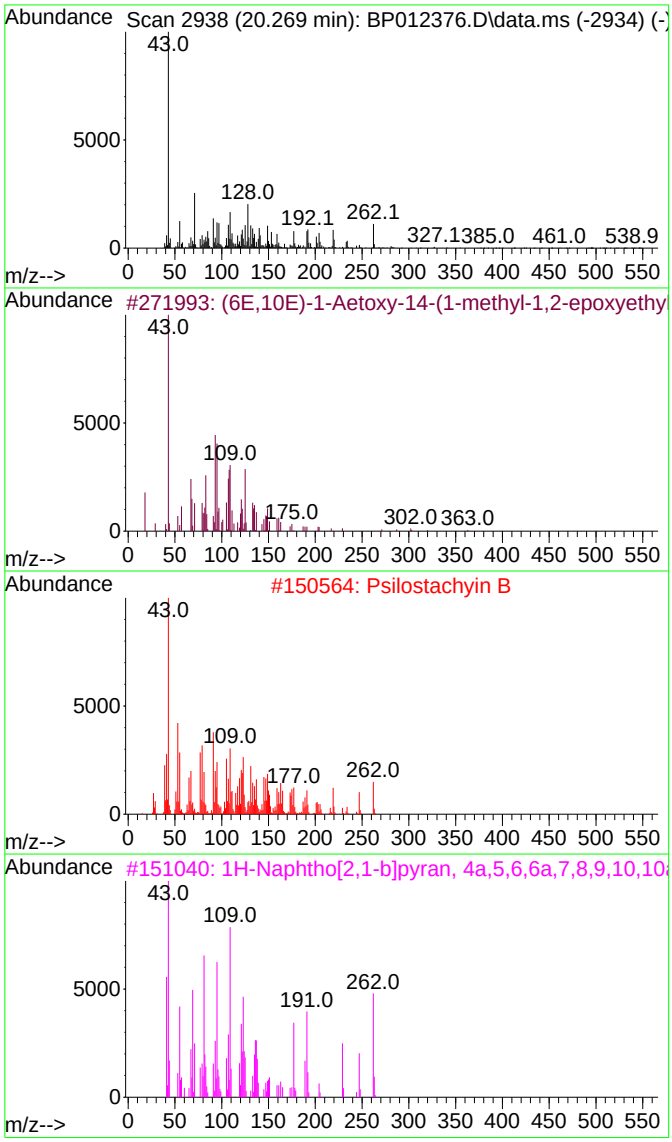
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

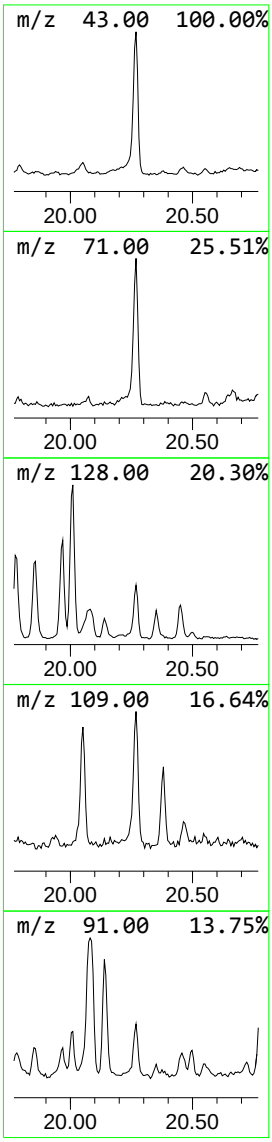
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 51 unknown-06 Concentration Rank 32

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 20.269 | 3.91 ng/ul | 1007450 | Chrysene-d12 | 21.310 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | (6E,10E)-1-Aetoxy-14-(1-methyl-1... | 362 | C22H34O4 | 1000426-95-6 | 27 |
| 2 | Psilostachyin B | 262 | C15H18O4 | 006995-02-4 | 25 |
| 3 | 1H-Naphtho[2,1-b]pyran, 4a,5,6,6... | 262 | C18H30O | 005153-92-4 | 15 |
| 4 | 3-[(E)-2-Chloro-1-methyl-1-buten... | 192 | C8H13ClO3 | 105949-78-8 | 14 |
| 5 | 2H-Cyclohepta[b]furan-2-one, 6-[... | 306 | C17H22O5 | 000580-49-4 | 12 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012376.D
Acq On : 28 Oct 2022 14:05
Operator : CG/JU
Sample : N5232-16
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BHAB9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

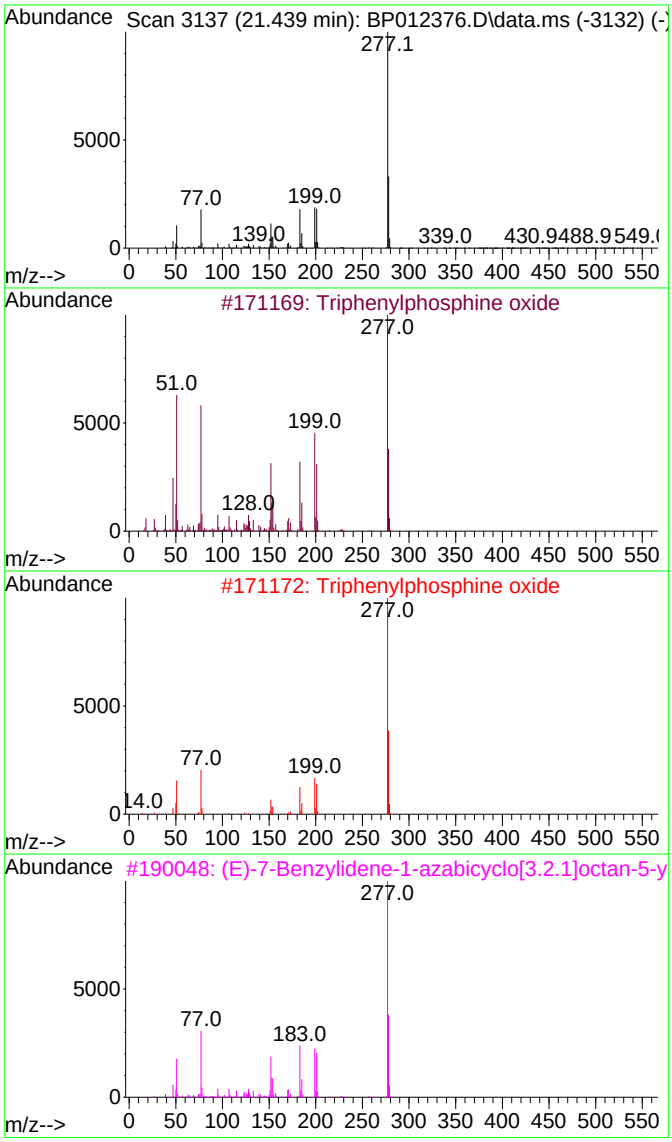
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

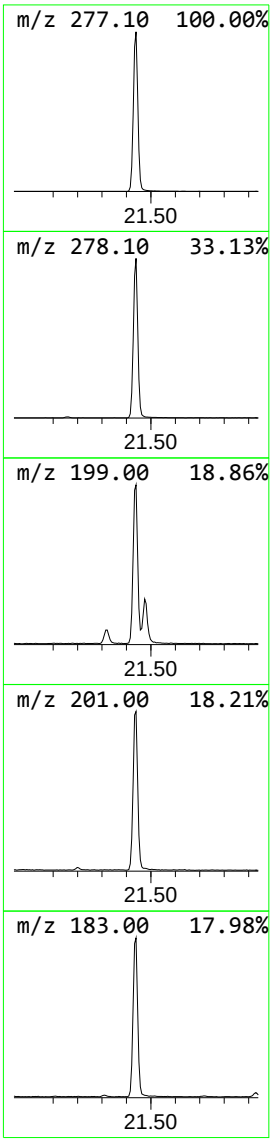
**********

Peak Number 54 Triphenylphosphine oxide Concentration Rank 6

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 21.439 | 19.51 ng/ul | 5031820 | Chrysene-d12 | 21.310 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Triphenylphosphine oxide | 278 | C18H15OP | 000791-28-6 | 94 |
| 2 | Triphenylphosphine oxide | 278 | C18H15OP | 000791-28-6 | 94 |
| 3 | (E)-7-Benzylidene-1-azabicyclo[3... | 293 | C15H19NO3S | 1000483-83-4 | 91 |
| 4 | Triphenylphosphine oxide | 278 | C18H15OP | 000791-28-6 | 50 |
| 5 | Triphenylphosphine oxide | 278 | C18H15OP | 000791-28-6 | 50 |



Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012376.D
Acq On : 28 Oct 2022 14:05
Operator : CG/JU
Sample : N5232-16
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BHAB9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.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 |
| Triethylamine | 3.034 | 17.2 | ng/ul | 1057980 | 1 | 7.805 | 1233260 | 20.0 |
| unknown-01 | 5.517 | 4.6 | ng/ul | 285320 | 1 | 7.805 | 1233260 | 20.0 |
| (DEL) Alkane: S... | 6.075 | 5.8 | ng/ul | 355880 | 1 | 7.805 | 1233260 | 20.0 |
| 3-Heptanone, 6-... | 6.469 | 5.9 | ng/ul | 361661 | 1 | 7.805 | 1233260 | 20.0 |
| Benzene, propyl- | 6.775 | 6.4 | ng/ul | 394614 | 1 | 7.805 | 1233260 | 20.0 |
| (2Z,4E)-3,7-Dim... | 6.916 | 5.6 | ng/ul | 343742 | 1 | 7.805 | 1233260 | 20.0 |
| unknown-02 | 7.505 | 13.6 | ng/ul | 840752 | 1 | 7.805 | 1233260 | 20.0 |
| unknown-03 | 7.999 | 9.4 | ng/ul | 577452 | 1 | 7.805 | 1233260 | 20.0 |
| Indane | 8.175 | 8.3 | ng/ul | 514937 | 1 | 7.805 | 1233260 | 20.0 |
| Aniline, N-methyl- | 8.611 | 3.9 | ng/ul | 240486 | 1 | 7.805 | 1233260 | 20.0 |
| unknown-04 | 8.710 | 4.0 | ng/ul | 249891 | 1 | 7.805 | 1233260 | 20.0 |
| Benzenamine, 3-... | 8.799 | 6.1 | ng/ul | 377240 | 1 | 7.805 | 1233260 | 20.0 |
| Benzene, 2-ethy... | 8.905 | 8.8 | ng/ul | 544981 | 1 | 7.805 | 1233260 | 20.0 |
| Benzene, 2-ethe... | 10.005 | 6.4 | ng/ul | 695551 | 2 | 10.610 | 2183680 | 20.0 |
| Benzenamine, 3,... | 10.699 | 27.2 | ng/ul | 2973710 | 2 | 10.610 | 2183680 | 20.0 |
| Phenol, p-tert-... | 11.957 | 25.4 | ng/ul | 2768930 | 2 | 10.610 | 2183680 | 20.0 |
| N,N-Dimethyl-2-... | 12.104 | 4.0 | ng/ul | 432049 | 2 | 10.610 | 2183680 | 20.0 |
| Benzenesulfonam... | 15.975 | 7.3 | ng/ul | 1488340 | 4 | 17.216 | 4096840 | 20.0 |
| 2(3H)-Benzothia... | 16.163 | 7.3 | ng/ul | 1505210 | 4 | 17.216 | 4096840 | 20.0 |
| 4-Hydroxy-N,N-d... | 17.910 | 4.8 | ng/ul | 978289 | 4 | 17.216 | 4096840 | 20.0 |
| Diphenyl sulfone | 18.181 | 8.4 | ng/ul | 1714320 | 4 | 17.216 | 4096840 | 20.0 |
| Isoquinoline | 18.257 | 6.6 | ng/ul | 1356550 | 4 | 17.216 | 4096840 | 20.0 |
| 1,2-Benzenediac... | 18.351 | 5.7 | ng/ul | 1165630 | 4 | 17.216 | 4096840 | 20.0 |
| unknown-05 | 18.404 | 13.5 | ng/ul | 2766080 | 4 | 17.216 | 4096840 | 20.0 |
| Benzophenone, 2... | 19.081 | 4.5 | ng/ul | 921231 | 4 | 17.216 | 4096840 | 20.0 |
| Sulphenone | 19.180 | 8.3 | ng/ul | 1707250 | 4 | 17.216 | 4096840 | 20.0 |
| Phenol, 4,4'-(1... | 19.610 | 21.1 | ng/ul | 5450580 | 5 | 21.310 | 5157850 | 20.0 |
| 1,4-Ethanonapht... | 19.781 | 3.9 | ng/ul | 1014380 | 5 | 21.310 | 5157850 | 20.0 |
| Benzene, 1,3-he... | 20.010 | 4.1 | ng/ul | 1068610 | 5 | 21.310 | 5157850 | 20.0 |
| unknown-06 | 20.269 | 3.9 | ng/ul | 1007450 | 5 | 21.310 | 5157850 | 20.0 |
| Triphenylphosph... | 21.439 | 19.5 | ng/ul | 5031820 | 5 | 21.310 | 5157850 | 20.0 |