

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP010924\  
 Data File : BP019313.D  
 Acq On : 09 Jan 2024 13:11  
 Operator : MA/JU  
 Sample : 05953-07ME  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 GCFA9ME

Manual Integrations  
 APPROVED

Reviewed By :Yogesh Patel 01/10/2024  
 Supervised By :mohammad ahmed 01/10/2024

Quant Time: Jan 10 00:04:06 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP010824.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Jan 09 03:05:24 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc    | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|---------|--------|----------|
| -----                         |        |      |          |         |        |          |
| Internal Standards            |        |      |          |         |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.781  | 152  | 142203   | 20.000  | ng/u1  | 0.00     |
| 20) Naphthalene-d8            | 10.575 | 136  | 558525   | 20.000  | ng/u1  | 0.00     |
| 38) Acenaphthene-d10          | 14.428 | 164  | 351653   | 20.000  | ng/u1  | 0.00     |
| 64) Phenanthrene-d10          | 17.186 | 188  | 707597   | 20.000  | ng/u1  | 0.00     |
| 79) Chrysene-d12              | 21.286 | 240  | 620581   | 20.000  | ng/u1  | 0.00     |
| 88) Perylene-d12              | 23.645 | 264  | 769721   | 20.000  | ng/u1  | 0.00     |
|                               |        |      |          |         |        |          |
| System Monitoring Compounds   |        |      |          |         |        |          |
| 3) 1,4-Dioxane-d8             | 3.222  | 96   | 14662    | 3.899   | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.646  | 84   | 57888    | 5.323   | ng/u1  | 0.00     |
| 7) Phenol-d5                  | 6.969  | 99   | 287064   | 21.121  | ng/u1  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.122  | 67   | 176388   | 21.697  | ng/u1  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.316  | 132  | 211820   | 22.688  | ng/u1  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.510  | 113  | 228882   | 21.227  | ng/u1  | 0.00     |
| 21) Nitrobenzene-d5           | 8.952  | 128  | 105320   | 22.414  | ng/u1  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.669  | 143  | 125072   | 22.908  | ng/u1  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.210 | 165  | 242310   | 21.968  | ng/u1  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.728 | 131  | 167349   | 12.011  | ng/u1  | 0.00     |
| 46) Dimethylphthalate-d6      | 13.845 | 166  | 685343   | 22.060  | ng/u1  | 0.00     |
| 49) Acenaphthylene-d8         | 14.116 | 160  | 740290   | 22.229  | ng/u1  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.663 | 143  | 96483    | 20.224  | ng/u1  | 0.00     |
| 60) Fluorene-d10              | 15.422 | 176  | 581680   | 22.869  | ng/u1  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.563 | 200  | 121179   | 23.196  | ng/u1  | 0.00     |
| 73) Anthracene-d10            | 17.281 | 188  | 871436   | 24.693  | ng/u1  | 0.00     |
| 81) Pyrene-d10                | 19.527 | 212  | 1022071  | 25.257  | ng/u1  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.492 | 264  | 1091054  | 26.068  | ng/u1  | 0.00     |
|                               |        |      |          |         |        |          |
| Target Compounds              |        |      |          |         | Qvalue |          |
| 58) 2,3,4,6-Tetrachlorophenol | 15.063 | 232  | 54828m   | 6.145   | ng/u1  |          |
| 71) Pentachlorophenol         | 16.845 | 266  | 3020360  | 436.582 | ng/u1  | 98       |
| 82) Pyrene                    | 19.557 | 202  | 73256    | 1.509   | ng/u1  | 98       |
| -----                         |        |      |          |         |        |          |

(#) = qualifier out of range (m) = manual integration (+) = signals summed

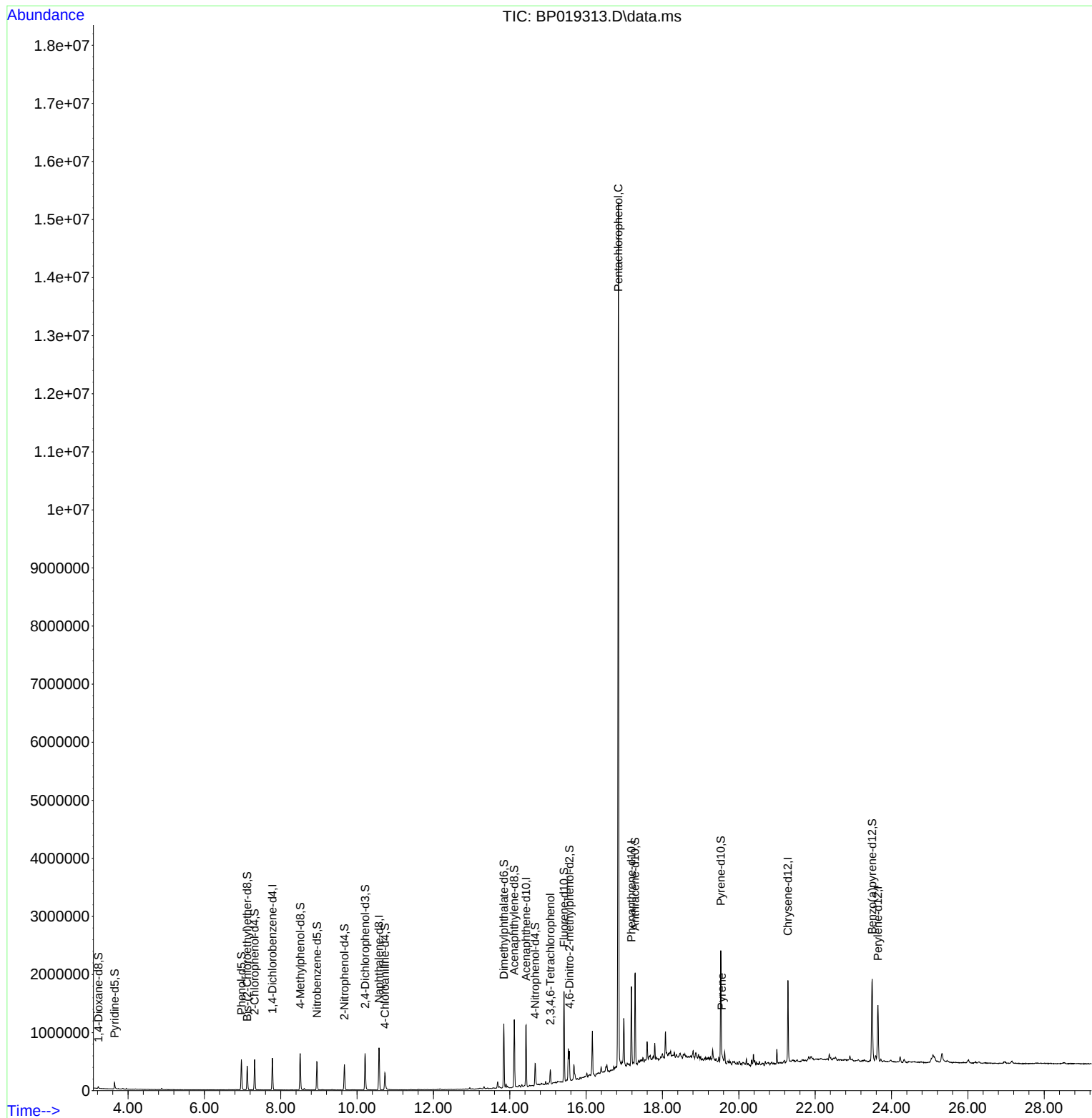
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP010924\  
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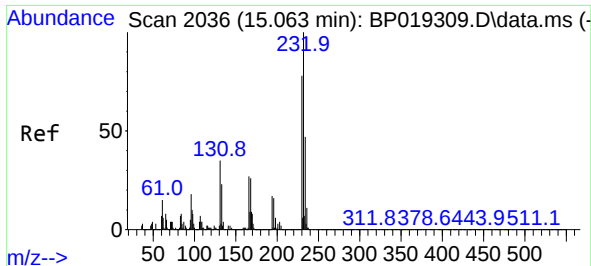
Instrument :  
BNA\_P  
ClientSampleId :  
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Quant Time: Jan 10 00:04:06 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP010824.MA.M  
Quant Title : SVOA CALIBRATION  
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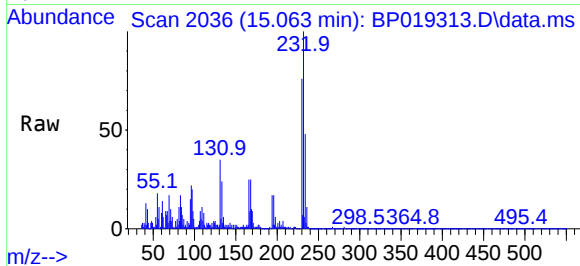
Reviewed By :Yogesh Patel 01/10/2024  
Supervised By :mohammad ahmed 01/10/2024





#58  
2,3,4,6-Tetrachlorophenol  
Concen: 6.145 ng/ul m  
RT: 15.063 min Scan# 2036  
Delta R.T. 0.000 min  
Lab File: BP019313.D  
Acq: 09 Jan 2024 13:11

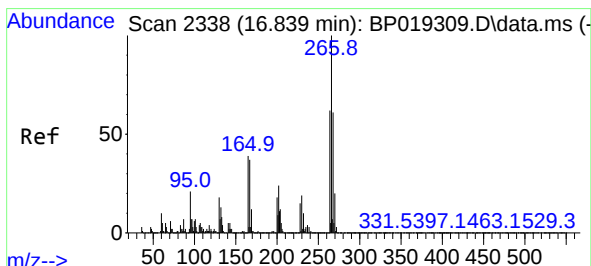
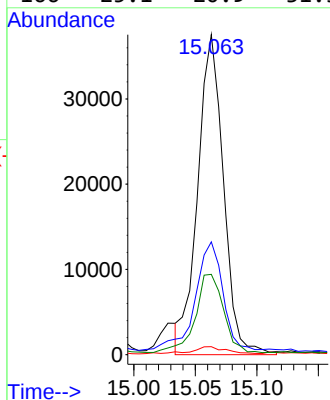
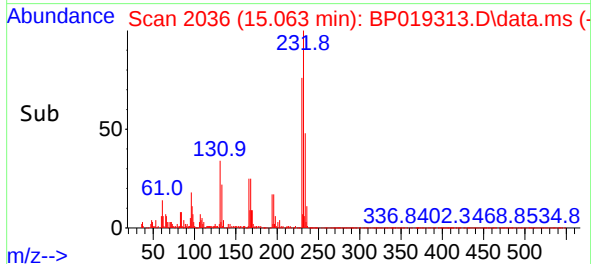
Instrument :  
BNA\_P  
ClientSampleId :  
GCFA9ME



Tgt Ion:232 Resp: 54828  
Ion Ratio Lower Upper  
232 100  
131 35.3 27.9 41.9  
130 2.5 1.8 2.8  
166 25.1 20.9 31.3

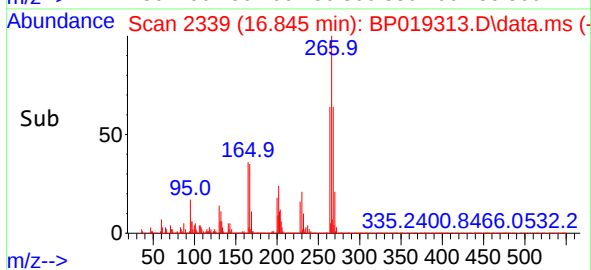
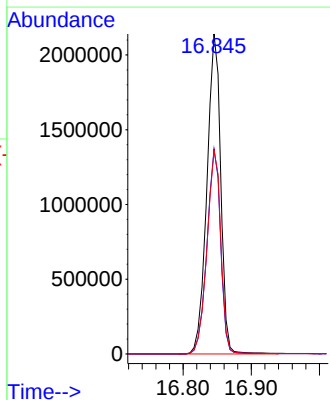
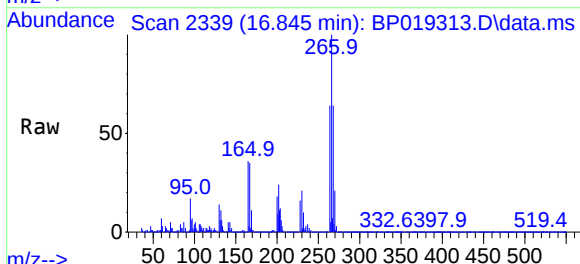
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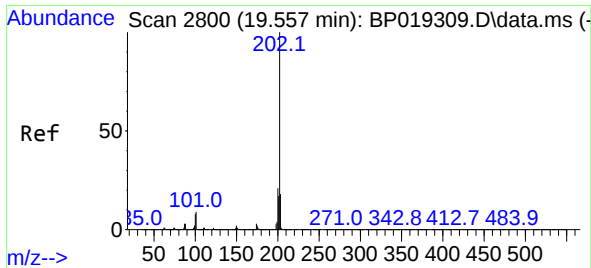
Reviewed By :Yogesh Patel 01/10/2024  
Supervised By :mohammad ahmed 01/10/2024



#71  
Pentachlorophenol  
Concen: 436.582 ng/ul  
RT: 16.845 min Scan# 2339  
Delta R.T. 0.006 min  
Lab File: BP019313.D  
Acq: 09 Jan 2024 13:11

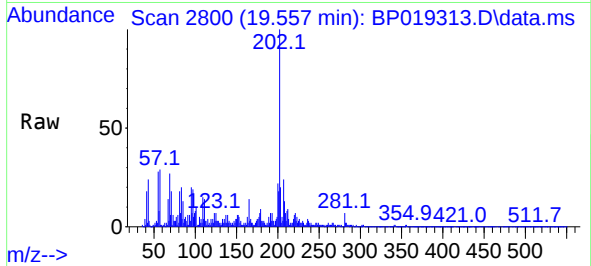
Tgt Ion:266 Resp: 3020360  
Ion Ratio Lower Upper  
266 100  
264 64.0 49.4 74.2  
268 63.6 51.6 77.4





#82  
 Pyrene  
 Concen: 1.509 ng/ul  
 RT: 19.557 min Scan# 21  
 Delta R.T. -0.000 min  
 Lab File: BP019313.D  
 Acq: 09 Jan 2024 13:11

Instrument :  
 BNA\_P  
 ClientSampleId :  
 GCFA9ME



Tgt Ion: 202 Resp: 73250

| Ion | Ratio | Lower | Upper |
|-----|-------|-------|-------|
| 202 | 100   |       |       |
| 101 | 10.0  | 7.2   | 10.8  |
| 100 | 7.8   | 5.8   | 8.6   |

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