

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100223\  
Data File : BP017512.D  
Acq On : 03 Oct 2023 12:07  
Operator : MA/JU  
Sample : 04417-30  
Misc :  
ALS Vial : 36 Sample Multiplier: 1

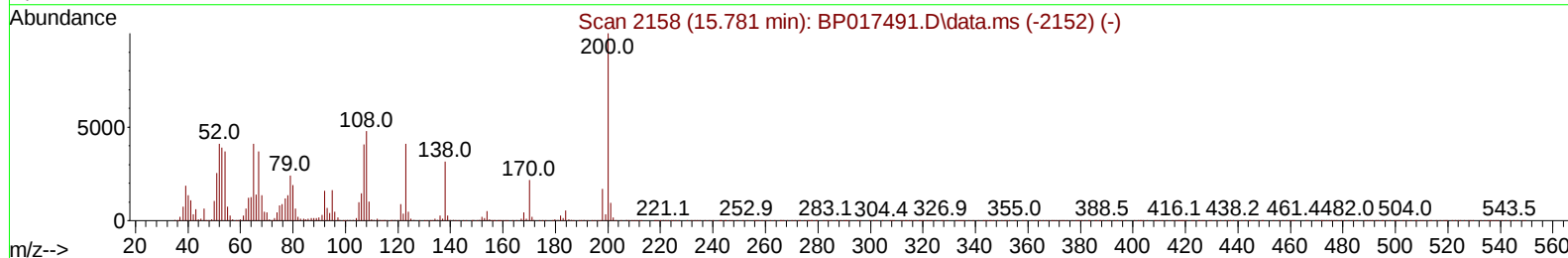
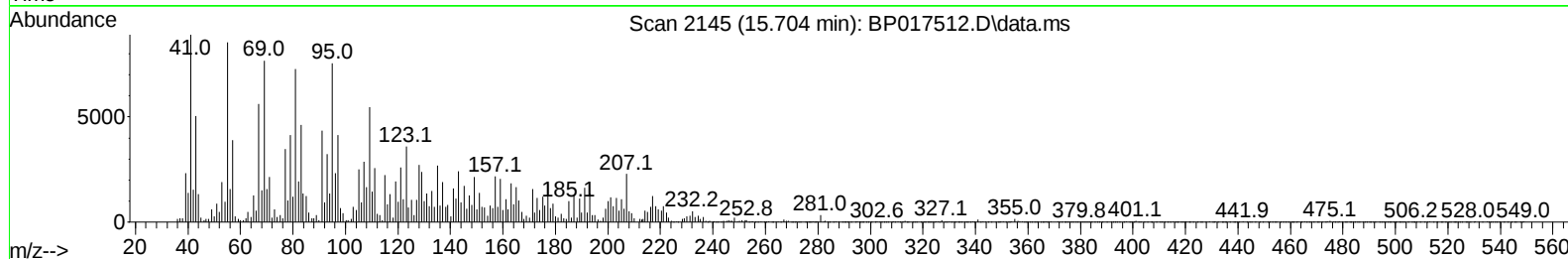
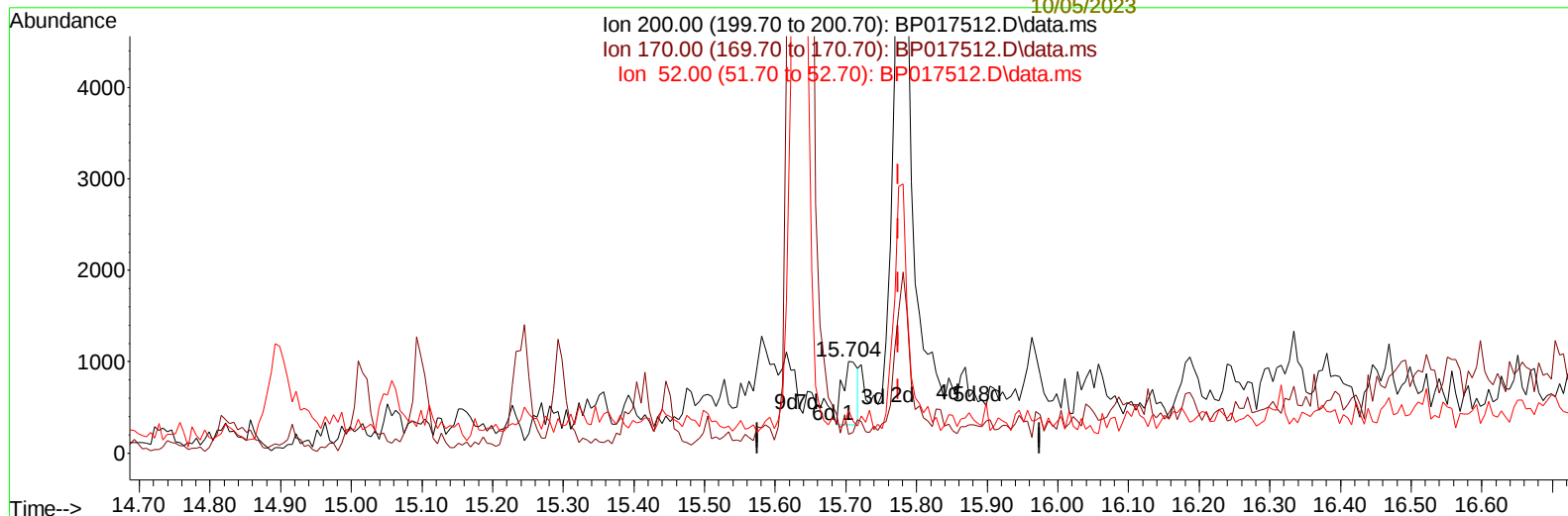
Instrument :  
BNA\_P  
ClientSampleId :  
EZYM9

Manual IntegrationsAPPROVED

Quant Time: Oct 04 00:38:53 2023  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092023.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Tue Oct 03 22:59:42 2023  
Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/04/2023  
Supervised By :mohammad ahmed 10/05/2023

Supervised By :mohammad  
ahmed  
10/05/2023



TIC: BP017512.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.704min (-0.070) 0.31 ng/u1

response 1003

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 200.00 | 100.00 | 100.00 |
| 170.00 | 18.80  | 20.22  |
| 52.00  | 41.50  | 46.91  |
| 0.00   | 0.00   | 0.00   |

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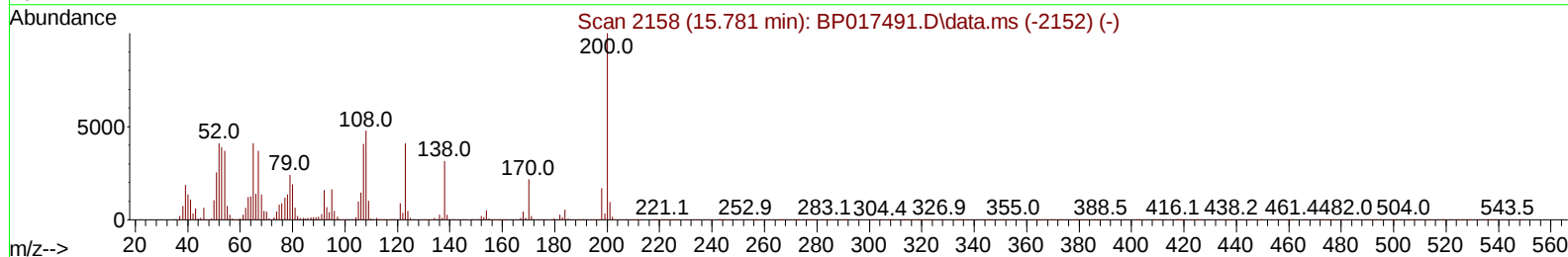
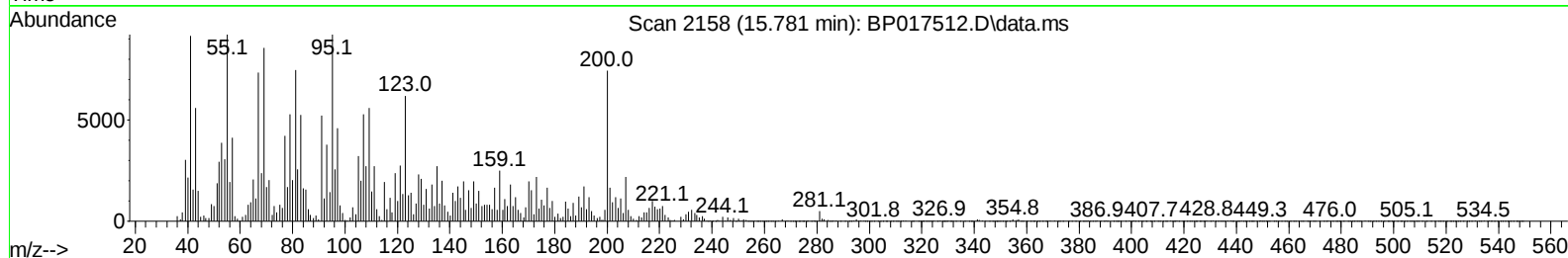
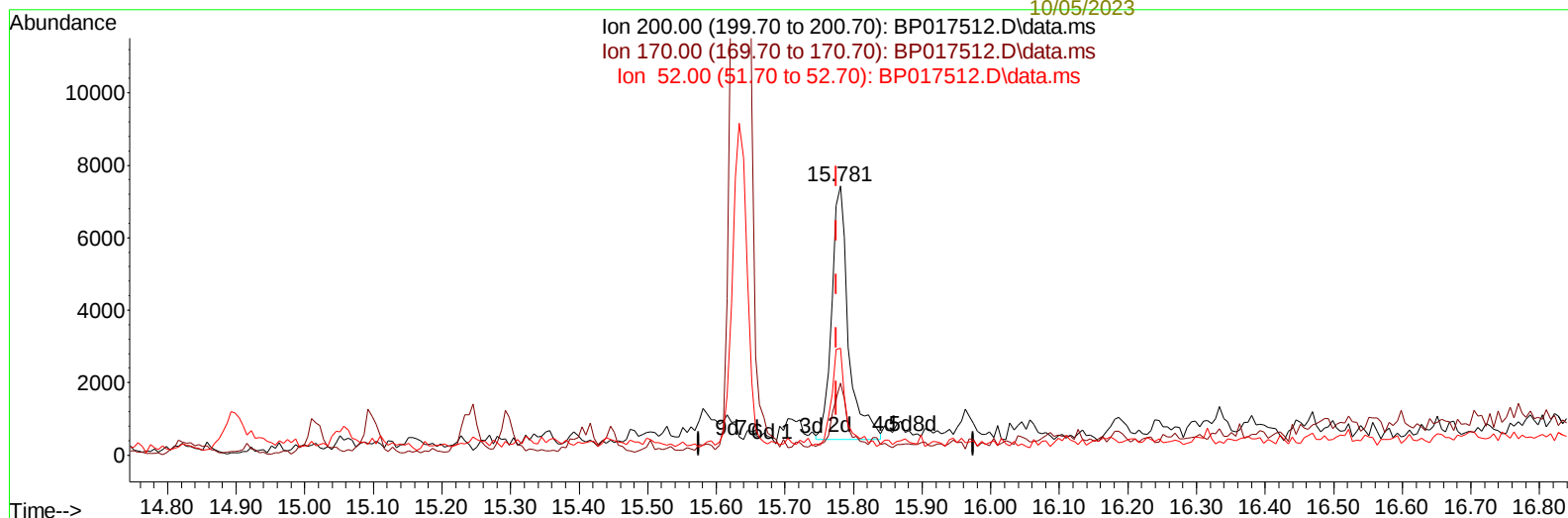
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(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.781min (+ 0.006) 3.71 ng/ul m

response 11938

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 200.00 | 100.00 | 100.00 |
| 170.00 | 18.80  | 26.69# |
| 52.00  | 41.50  | 39.64  |
| 0.00   | 0.00   | 0.00   |

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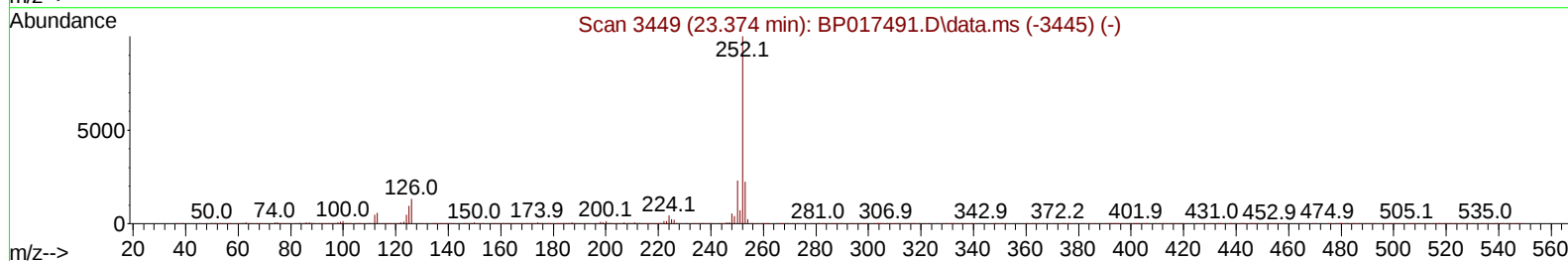
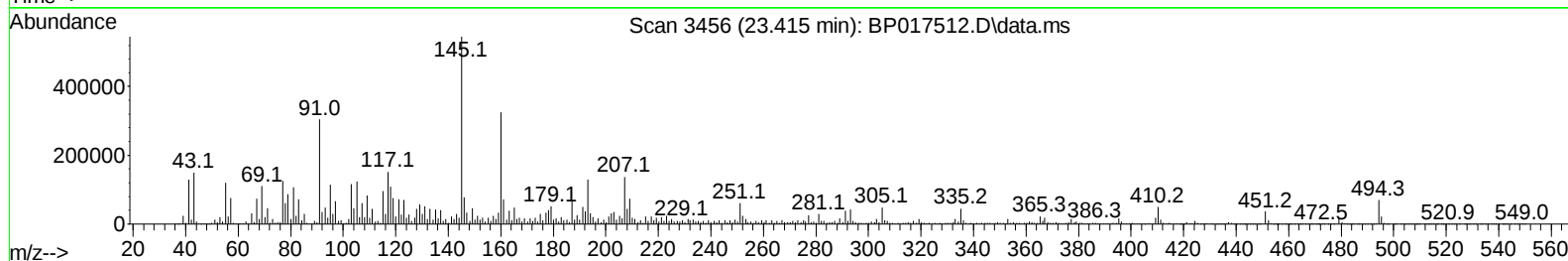
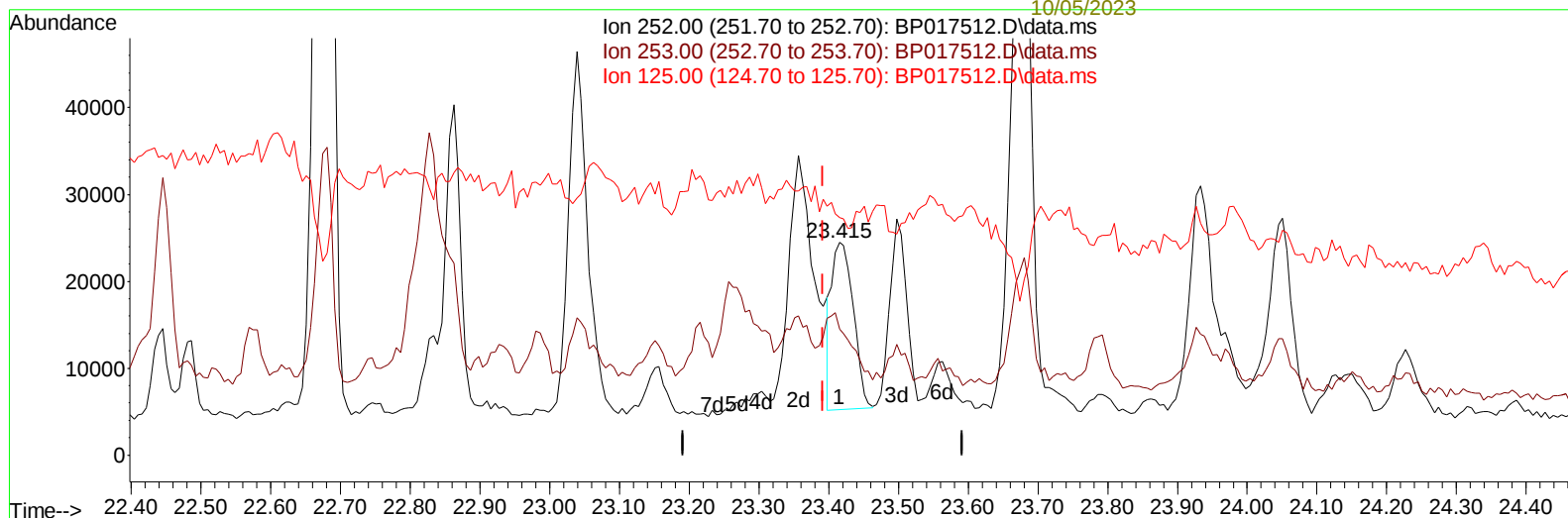
Instrument :  
 BNA\_P  
 ClientSampleId :  
 EZYM9

Manual IntegrationsAPPROVED

Quant Time: Oct 04 00:41:22 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092023.MA.M  
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**(91) Benzo(k)fluoranthene**

**23.415min (+ 0.024) 1.06 ng/u1**

**response 40913**

| Ion    | Exp%   | Act%    |
|--------|--------|---------|
| 252.00 | 100.00 | 100.00  |
| 253.00 | 22.40  | 59.36#  |
| 125.00 | 9.70   | 111.65# |
| 0.00   | 0.00   | 0.00    |

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 ALS Vial : 36 Sample Multiplier: 1

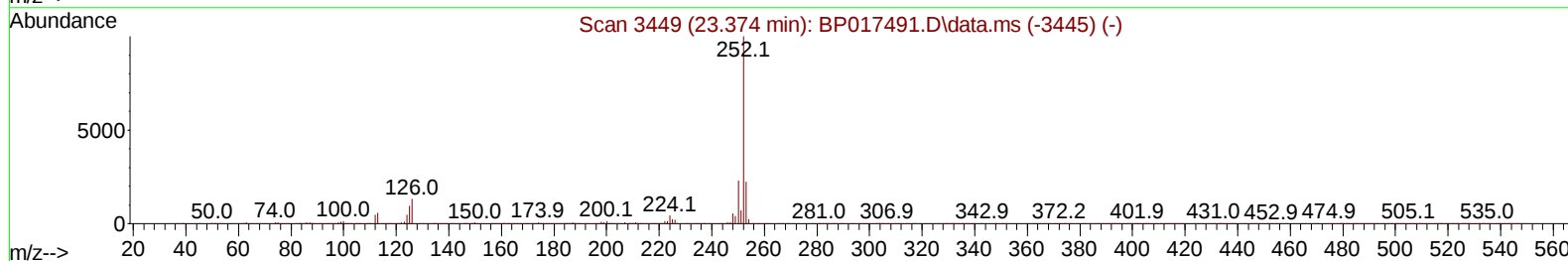
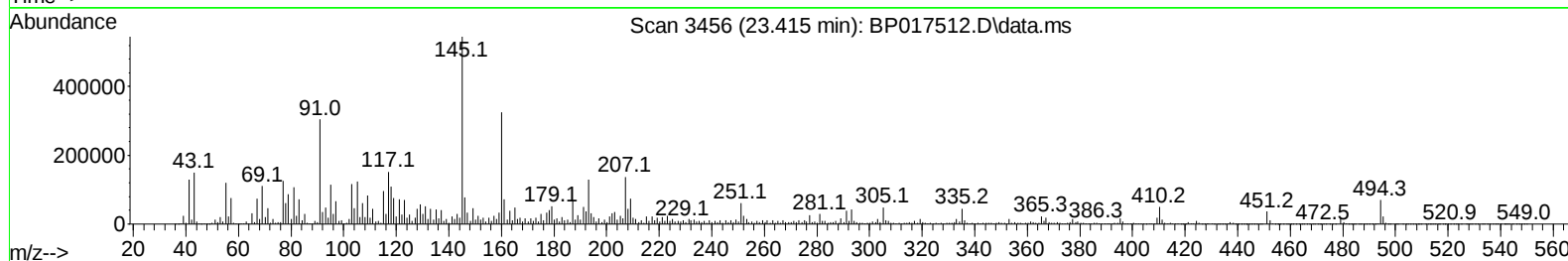
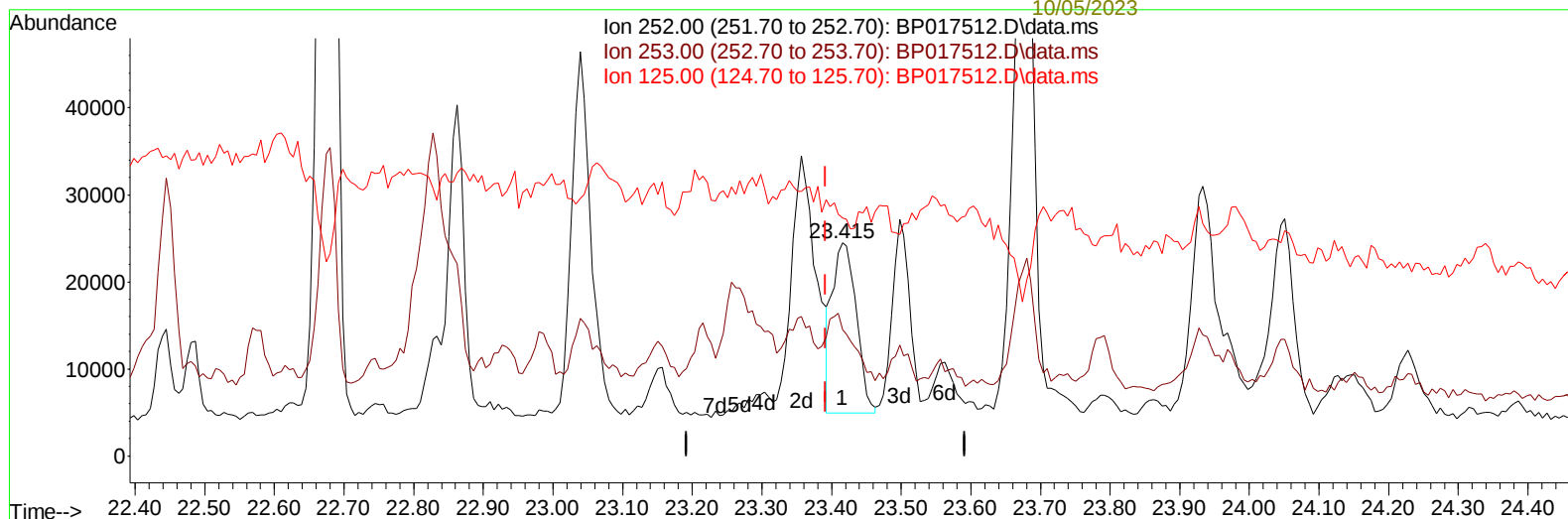
Instrument :  
 BNA\_P  
 ClientSampleId :  
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Manual IntegrationsAPPROVED

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TIC: BP017512.D\data.ms

**(91) Benzo(k)fluoranthene**

**23.415min (+ 0.024) 1.23 ng/ul m**

**response 47193**

| Ion    | Exp%   | Act%    |
|--------|--------|---------|
| 252.00 | 100.00 | 100.00  |
| 253.00 | 22.40  | 59.36#  |
| 125.00 | 9.70   | 111.65# |
| 0.00   | 0.00   | 0.00    |

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 Acq On : 03 Oct 2023 12:07  
 Operator : MA/JU  
 Sample : 04417-30  
 Misc :  
 ALS Vial : 36 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 EZYM9

## Manual IntegrationsAPPROVED

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Quant Time: Oct 04 00:41:50 2023  
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| Compound                      | R.T.   | Q   | Ion | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|-----|-----|----------|--------|--------|----------|
| -----                         |        |     |     |          |        |        |          |
| Internal Standards            |        |     |     |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.934  | 152 |     | 117622   | 20.000 | ng/ul  | 0.00     |
| 20) Naphthalene-d8            | 10.775 | 136 |     | 444015   | 20.000 | ng/ul  | 0.00     |
| 38) Acenaphthene-d10          | 14.634 | 164 |     | 258497   | 20.000 | ng/ul  | 0.00     |
| 64) Phenanthrene-d10          | 17.422 | 188 |     | 571659   | 20.000 | ng/ul  | 0.00     |
| 79) Chrysene-d12              | 21.569 | 240 |     | 562669   | 20.000 | ng/ul  | # 0.00   |
| 88) Perylene-d12              | 24.168 | 264 |     | 647047   | 20.000 | ng/ul  | 0.02     |
| System Monitoring Compounds   |        |     |     |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.287  | 96  |     | 7958     | 2.779  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.734  | 84  |     | 18200    | 2.273  | ng/ul  | 0.02     |
| 7) Phenol-d5                  | 7.093  | 99  |     | 196655   | 20.357 | ng/ul  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.275  | 67  |     | 140541   | 24.107 | ng/ul  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.457  | 132 |     | 158330   | 20.665 | ng/ul  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.657  | 113 |     | 103417   | 13.770 | ng/ul  | 0.00     |
| 21) Nitrobenzene-d5           | 9.146  | 128 |     | 87494    | 27.257 | ng/ul  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.857  | 143 |     | 62564    | 17.950 | ng/ul  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.393 | 165 |     | 129235   | 19.626 | ng/ul  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.946 | 131 |     | 112442   | 11.788 | ng/ul  | 0.00     |
| 46) Dimethylphthalate-d6      | 14.045 | 166 |     | 506759   | 27.366 | ng/ul  | 0.00     |
| 49) Acenaphthylene-d8         | 14.328 | 160 |     | 568595   | 26.697 | ng/ul  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.892 | 143 |     | 16890    | 5.235  | ng/ul  | 0.02     |
| 60) Fluorene-d10              | 15.634 | 176 |     | 470474   | 29.511 | ng/ul  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.781 | 200 |     | 11938m   | 3.705  | ng/ul  | 0.00     |
| 73) Anthracene-d10            | 17.522 | 188 |     | 708887   | 27.960 | ng/ul  | 0.00     |
| 81) Pyrene-d10                | 19.780 | 212 |     | 900971   | 29.610 | ng/ul  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.998 | 264 |     | 870975   | 27.958 | ng/ul  | 0.03     |
| Target Compounds              |        |     |     |          |        |        |          |
|                               |        |     |     |          |        | Qvalue |          |
| 8) Phenol                     | 7.122  | 94  |     | 82937    | 8.536  | ng/ul  | 95       |
| 16) Acetophenone              | 8.793  | 105 |     | 60993    | 5.255  | ng/ul  | 98       |
| 80) Fluoranthene              | 19.445 | 202 |     | 47552    | 1.340  | ng/ul# | 95       |
| 82) Pyrene                    | 19.810 | 202 |     | 67799    | 1.797  | ng/ul  | 95       |
| 90) Benzo(b)fluoranthene      | 23.357 | 252 |     | 69813    | 1.830  | ng/ul# | 1        |
| 91) Benzo(k)fluoranthene      | 23.415 | 252 |     | 47193m   | 1.226  | ng/ul  |          |
| 93) Benzo(a)pyrene            | 24.051 | 252 |     | 53997    | 1.479  | ng/ul# | 1        |
| 94) Indeno(1,2,3-cd)pyrene    | 26.927 | 276 |     | 70268    | 1.536  | ng/ul# | 80       |
| 95) Dibenzo(a,h)anthracene    | 26.951 | 278 |     | 44340    | 1.164  | ng/ul# | 49       |
| 96) Benzo(g,h,i)perylene      | 27.792 | 276 |     | 84372    | 2.287  | ng/ul# | 86       |
| -----                         |        |     |     |          |        |        |          |

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

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