

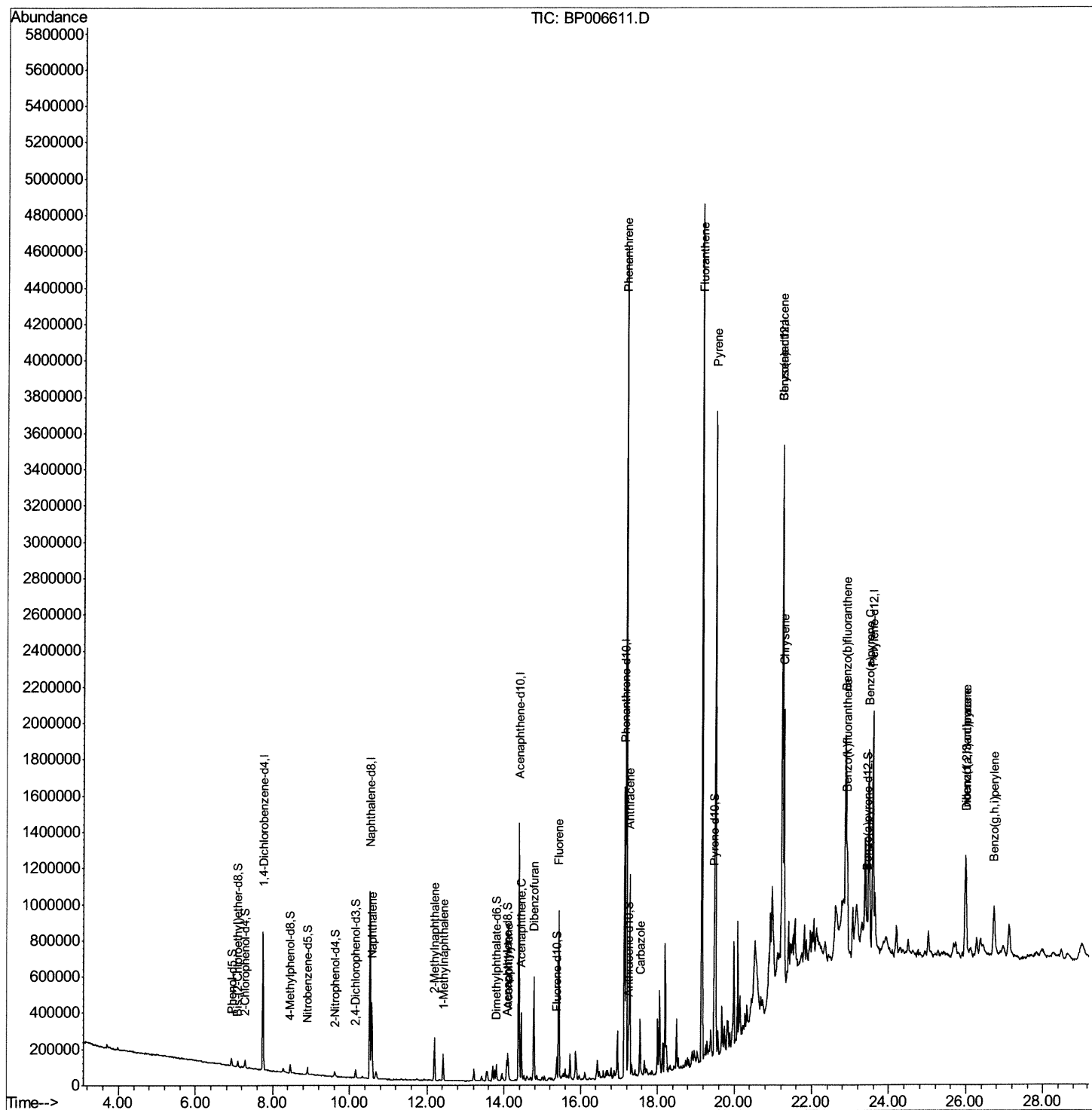
Data Path : Z:\svoasrv\HPCHEM1\BNA P\Data\BP080921\  
Data File : BP006611.D  
Acq On : 09 Aug 2021 13:30  
Operator : CG/JU  
Sample : M3169-04DL 30X  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
C00A3DL

Manual Integrations  
APPROVED

mohammad  
8/11/2021 9:24:04 PM

Quant Time: Aug 09 14:21:35 2021  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SFAM-EPA-BP072421.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Sat Aug 07 13:59:25 2021  
Response via : Initial Calibration



# Quantitation Report (Qedit)

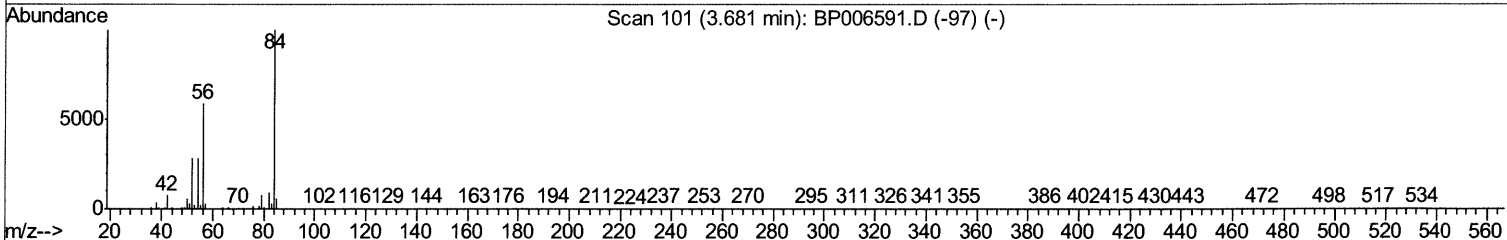
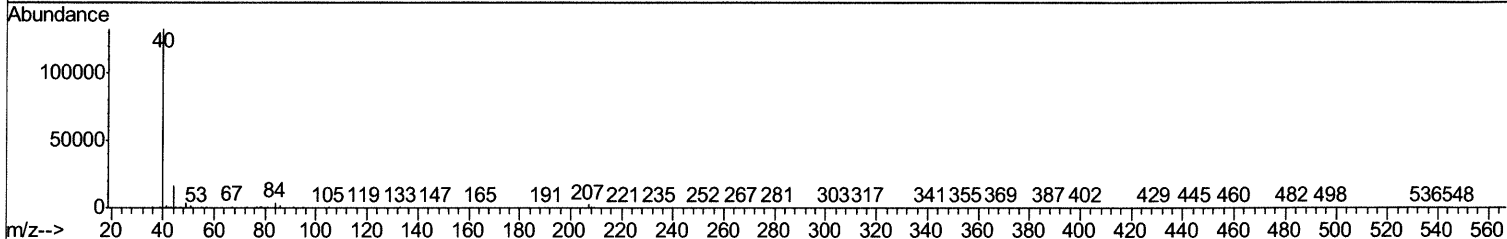
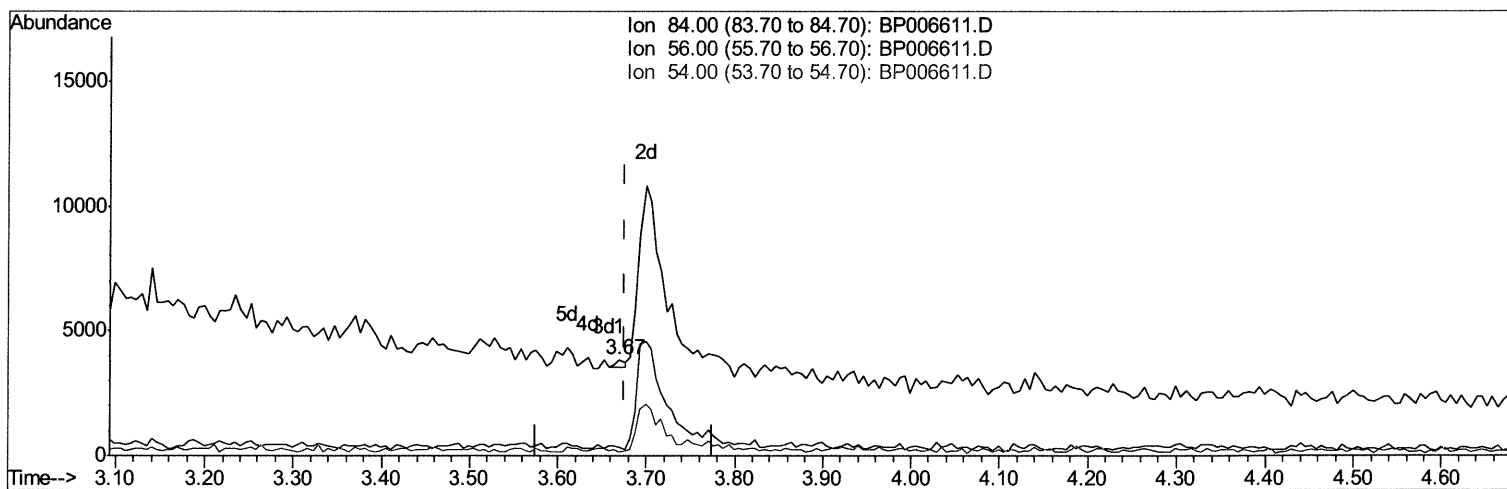
Data Path : Z:\svoasrv\HPCHEM1\BNA P\Data\BP080921\  
 Data File : BP006611.D  
 Acq On : 09 Aug 2021 13:30  
 Operator : CG/JU  
 Sample : M3169-04DL 30X  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 C00A3DL

Manual Integrations  
 APPROVED

mohammad  
 8/11/2021 9:24:04 PM

Quant Time: Aug 09 14:08:16 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SFAM-EPA-BP072421.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat Aug 07 13:59:25 2021  
 Response via : Initial Calibration



TIC: BP006611.D

(4) Pyridine-d5 (S)  
 3.670min (-0.006) 0.01ng/ul  
 response 217

| Ion   | Exp%  | Act%  |
|-------|-------|-------|
| 84.00 | 100   | 100   |
| 56.00 | 58.40 | 8.71# |
| 54.00 | 27.10 | 3.64# |
| 0.00  | 0.00  | 0.00  |

## Quantitation Report (Qedit)

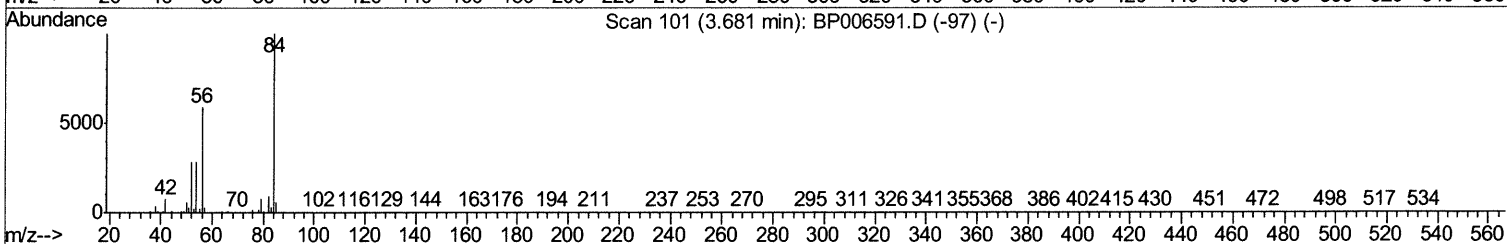
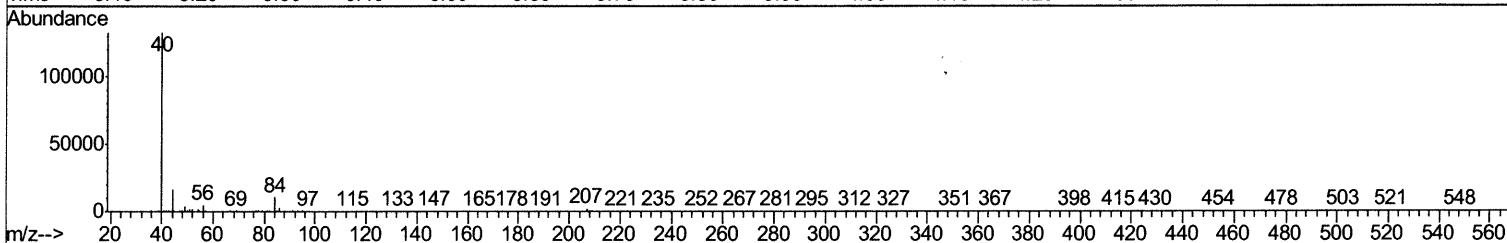
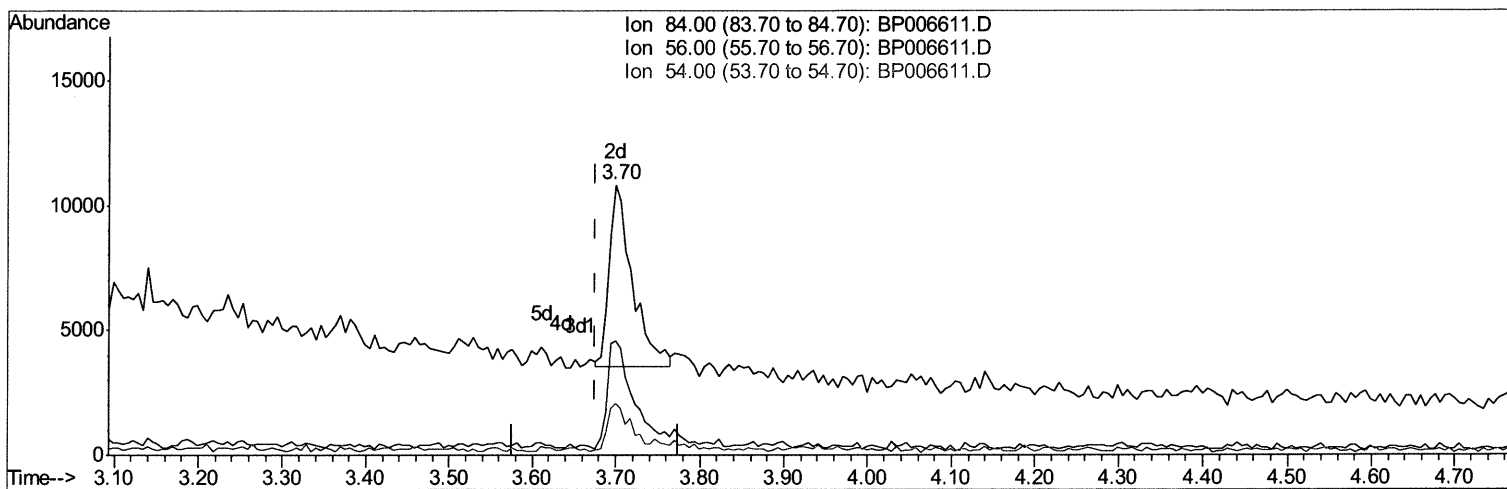
Data Path : Z:\svoasrv\HPCHEM1\BNA P\Data\BP080921\  
Data File : BP006611.D  
Acq On : 09 Aug 2021 13:30  
Operator : CG/JU  
Sample : M3169-04DL 30X  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
C00A3DL

Manual Integrations  
APPROVED

mohammad  
8/11/2021 9:24:04 PM

Quant Time: Aug 09 14:08:16 2021  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SFAM-EPA-BP072421.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Sat Aug 07 13:59:25 2021  
Response via : Initial Calibration



TIC: BP006611.D

(4) Pyridine-d5 (S)

3.699min (+0.024) 0.92ng/ul m 08/12/21 JU

response 14005

| Ion   | Exp%  | Act%   |
|-------|-------|--------|
| 84.00 | 100   | 100    |
| 56.00 | 58.40 | 42.34# |
| 54.00 | 27.10 | 18.90# |
| 0.00  | 0.00  | 0.00   |

# Quantitation Report (Qedit)

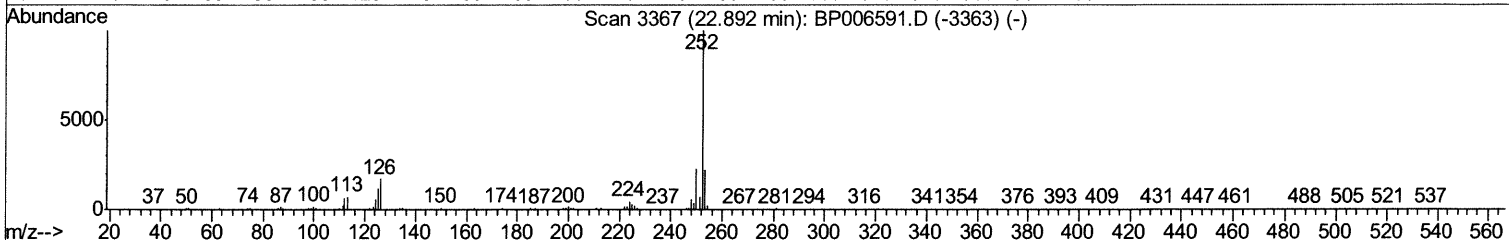
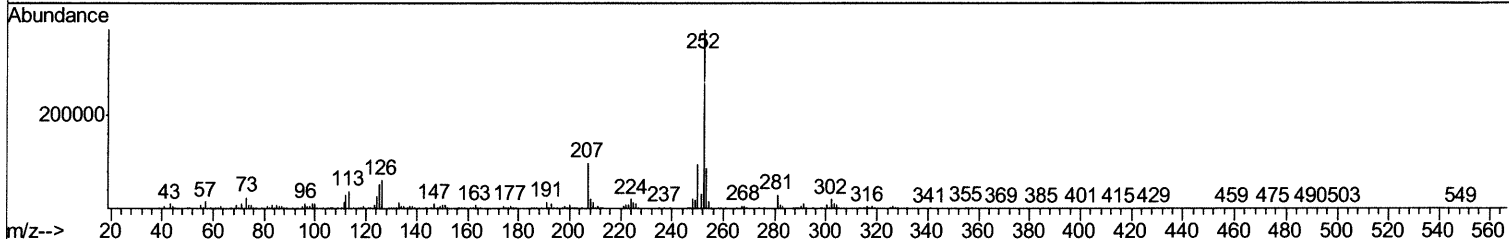
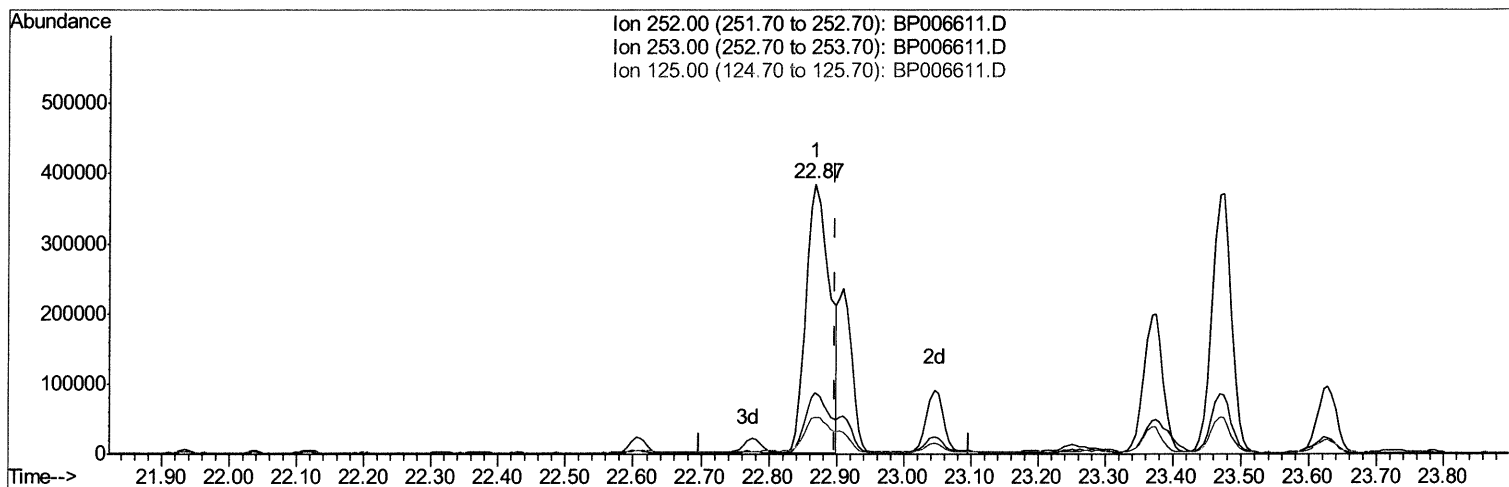
Data Path : Z:\svoasrv\HPCHEM1\BNA P\Data\BP080921\  
 Data File : BP006611.D  
 Acq On : 09 Aug 2021 13:30  
 Operator : CG/JU  
 Sample : M3169-04DL 30X  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 C00A3DL

Manual Integrations  
 APPROVED

mohammad  
 8/11/2021 9:24:04 PM

Quant Time: Aug 09 14:08:16 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SFAM-EPA-BP072421.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat Aug 07 13:59:25 2021  
 Response via : Initial Calibration



TIC: BP006611.D

(91) Benzo(k)fluoranthene

22.869min (-0.029) 17.95ng/ul

response 935341

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 252.00 | 100   | 100   |
| 253.00 | 22.10 | 22.80 |
| 125.00 | 12.00 | 13.63 |
| 0.00   | 0.00  | 0.00  |

## Quantitation Report (Qedit)

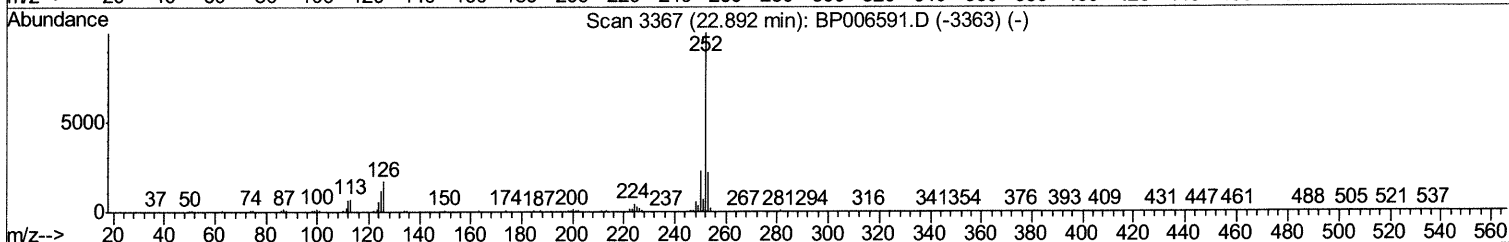
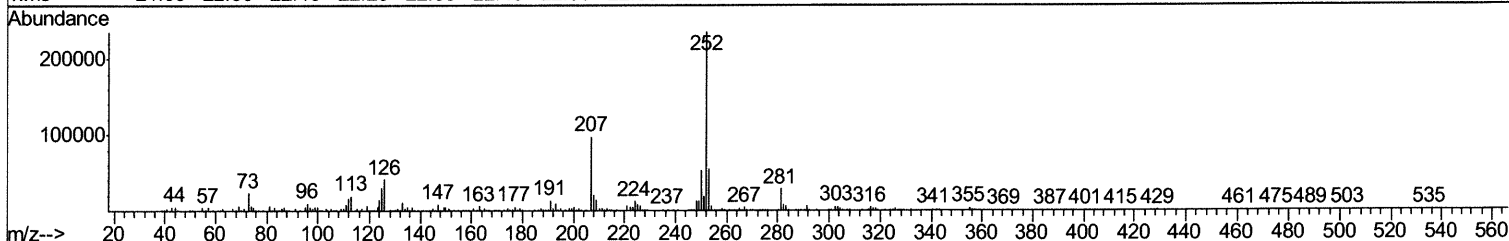
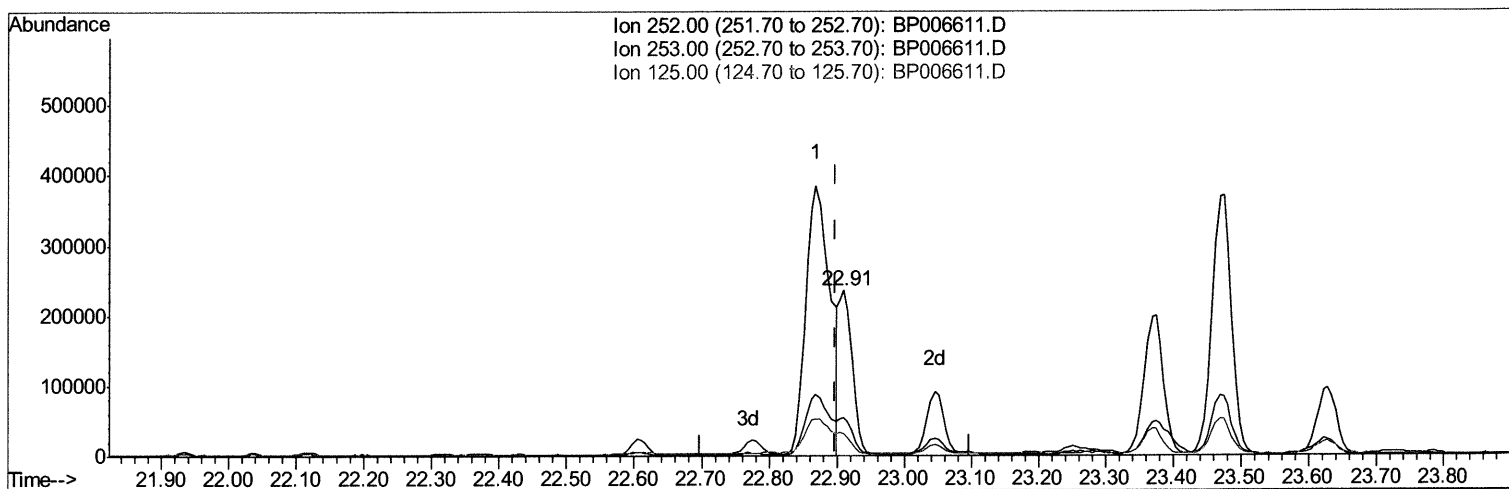
Data Path : Z:\svoasrv\HPCHEM1\BNA P\Data\BP080921\  
Data File : BP006611.D  
Acq On : 09 Aug 2021 13:30  
Operator : CG/JU  
Sample : M3169-04DL 30X  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
C00A3DL

Manual Integrations  
APPROVED

mohammad  
8/11/2021 9:24:04 PM

Quant Time: Aug 09 14:08:16 2021  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SFAM-EPA-BP072421.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Sat Aug 07 13:59:25 2021  
Response via : Initial Calibration



TIC: BP006611.D

(91) Benzo(k)fluoranthene

22.910min (+0.012) 6.38ng/ul m 08/12/21 JU

response 332708

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 252.00 | 100   | 100   |
| 253.00 | 22.10 | 23.27 |
| 125.00 | 12.00 | 13.31 |
| 0.00   | 0.00  | 0.00  |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080921\  
 Data File : BP006611.D  
 Acq On : 09 Aug 2021 13:30  
 Operator : CG/JU  
 Sample : M3169-04DL 30X  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 C00A3DL

Manual Integrations  
 APPROVED

mohammad  
 8/11/2021 9:24:04 PM

Quant Time: Aug 09 14:21:35 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SFAM-EPA-BP072421.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat Aug 07 13:59:25 2021  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.75  | 152  | 190617   | 20.00 | ng/ul | 0.00     |
| 20) Naphthalene-d8        | 10.53 | 136  | 811058   | 20.00 | ng/ul | 0.00     |
| 38) Acenaphthene-d10      | 14.38 | 164  | 448683   | 20.00 | ng/ul | 0.00     |
| 64) Phenanthrene-d10      | 17.14 | 188  | 853507   | 20.00 | ng/ul | 0.00     |
| 79) Chrysene-d12          | 21.24 | 240  | 807603   | 20.00 | ng/ul | 0.00     |
| 88) Perylene-d12          | 23.57 | 264  | 864438   | 20.00 | ng/ul | 0.02     |

## System Monitoring Compounds

|                                |       |     |          |      |         |      |
|--------------------------------|-------|-----|----------|------|---------|------|
| 3) 1,4-Dioxane-d8              | 3.28  | 96  | 1336     | 0.26 | ng/uL   | 0.01 |
| 4) Pyridine-d5                 | 3.70  | 84  | 14005m > | 0.92 | ng/ul > | 0.02 |
| 7) Phenol-d5                   | 6.93  | 99  | 23182    | 1.29 | ng/ul   | 0.00 |
| 9) Bis-(2-Chloroethyl)ether-d  | 7.09  | 67  | 15410    | 1.38 | ng/ul   | 0.00 |
| 11) 2-Chlorophenol-d4          | 7.28  | 132 | 18619    | 1.40 | ng/ul   | 0.00 |
| 15) 4-Methylphenol-d8          | 8.46  | 113 | 18526    | 1.32 | ng/ul   | 0.00 |
| 21) Nitrobenzene-d5            | 8.90  | 128 | 10056    | 1.67 | ng/ul   | 0.00 |
| 24) 2-Nitrophenol-d4           | 9.62  | 143 | 9166     | 1.59 | ng/ul   | 0.00 |
| 28) 2,4-Dichlorophenol-d3      | 10.16 | 165 | 14712    | 1.30 | ng/ul   | 0.00 |
| 31) 4-Chloroaniline-d4         | 10.68 | 131 | 17035    | 0.92 | ng/ul   | 0.00 |
| 46) Dimethylphthalate-d6       | 13.80 | 166 | 52051    | 1.63 | ng/ul   | 0.00 |
| 49) Acenaphthylene-d8          | 14.07 | 160 | 62468    | 1.59 | ng/ul   | 0.00 |
| 54) 4-Nitrophenol-d4           | 0.00  | 143 | 0d       | 0.00 | ng/ul   |      |
| 60) Fluorene-d10               | 15.38 | 176 | 43449    | 1.62 | ng/ul   | 0.00 |
| 65) 4,6-Dinitro-2-methylphenol | 15.51 | 200 | 1886     | 0.40 | ng/ul   | 0.01 |
| 73) Anthracene-d10             | 17.24 | 188 | 67062    | 1.74 | ng/ul   | 0.00 |
| 81) Pyrene-d10                 | 19.48 | 212 | 73519    | 1.77 | ng/ul   | 0.00 |
| 92) Benzo(a)pyrene-d12         | 23.42 | 264 | 72849    | 1.66 | ng/ul   | 0.01 |

## Target Compounds

|                            |       |     |           |        |         | Ovalue      |
|----------------------------|-------|-----|-----------|--------|---------|-------------|
| 30) Naphthalene            | 10.58 | 128 | 327849    | 7.791  | ng/ul   | 100         |
| 36) 2-Methylnaphthalene    | 12.20 | 142 | 104655    | 3.630  | ng/ul   | 100         |
| 37) 1-Methylnaphthalene    | 12.42 | 142 | 63906     | 2.202  | ng/ul   | 100         |
| 50) Acenaphthylene         | 14.10 | 152 | 118368    | 3.098  | ng/ul   | 97          |
| 52) Acenaphthene           | 14.45 | 153 | 140303    | 5.005  | ng/ul   | 98          |
| 56) Dibenzofuran           | 14.78 | 168 | 320349    | 8.521  | ng/ul   | 100         |
| 61) Fluorene               | 15.43 | 166 | 377880    | 12.220 | ng/ul   | 100         |
| 72) Phenanthrene           | 17.18 | 178 | 2516603   | 54.911 | ng/ul   | 99          |
| 74) Anthracene             | 17.27 | 178 | 643417    | 13.846 | ng/ul   | 100         |
| 77) Carbazole              | 17.55 | 167 | 186420    | 4.385  | ng/ul   | 97          |
| 80) Fluoranthene           | 19.16 | 202 | 2503335   | 48.908 | ng/ul   | 99          |
| 82) Pyrene                 | 19.51 | 202 | 1925235   | 36.248 | ng/ul   | 99          |
| 85) Benzo(a)anthracene     | 21.23 | 228 | 922138    | 18.427 | ng/ul   | 97          |
| 87) Chrysene               | 21.27 | 228 | 716690    | 14.941 | ng/ul   | 97          |
| 90) Benzo(b)fluoranthene   | 22.87 | 252 | 935341    | 17.035 | ng/ul   | 98          |
| 91) Benzo(k)fluoranthene   | 22.91 | 252 | 332708m > | 6.385  | ng/ul > | 08/12/21 34 |
| 93) Benzo(a)pyrene         | 23.47 | 252 | 729782    | 15.206 | ng/ul   | 99          |
| 94) Indeno(1,2,3-cd)pyrene | 25.99 | 276 | 446343    | 7.265  | ng/ul   | 97          |
| 95) Dibenzo(a,h)anthracene | 25.99 | 278 | 112651    | 2.165  | ng/ul#  | 76          |
| 96) Benzo(a,h,i)perylene   | 26.72 | 276 | 347993    | 6.856  | ng/ul   | 100         |

Data Path : Z:\svoasrv\HPCHEM1\BNA P\Data\BP080921\  
Data File : BP006611.D  
Acq On : 09 Aug 2021 13:30  
Operator : CG/JU  
Sample : M3169-04DL 30X  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
C00A3DL

Manual Integrations  
APPROVED

mohammad  
8/11/2021 9:24:04 PM

Quant Time: Aug 09 14:21:35 2021  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SFAM-EPA-BP072421.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Sat Aug 07 13:59:25 2021  
Response via : Initial Calibration

| Internal Standards                                                          | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|-----------------------------------------------------------------------------|------|------|----------|------|-------|----------|
| -----                                                                       |      |      |          |      |       |          |
| (# ) = qualifier out of range (m) = manual integration (+) = signals summed |      |      |          |      |       |          |