

# Quantitation Report (QT Reviewed)

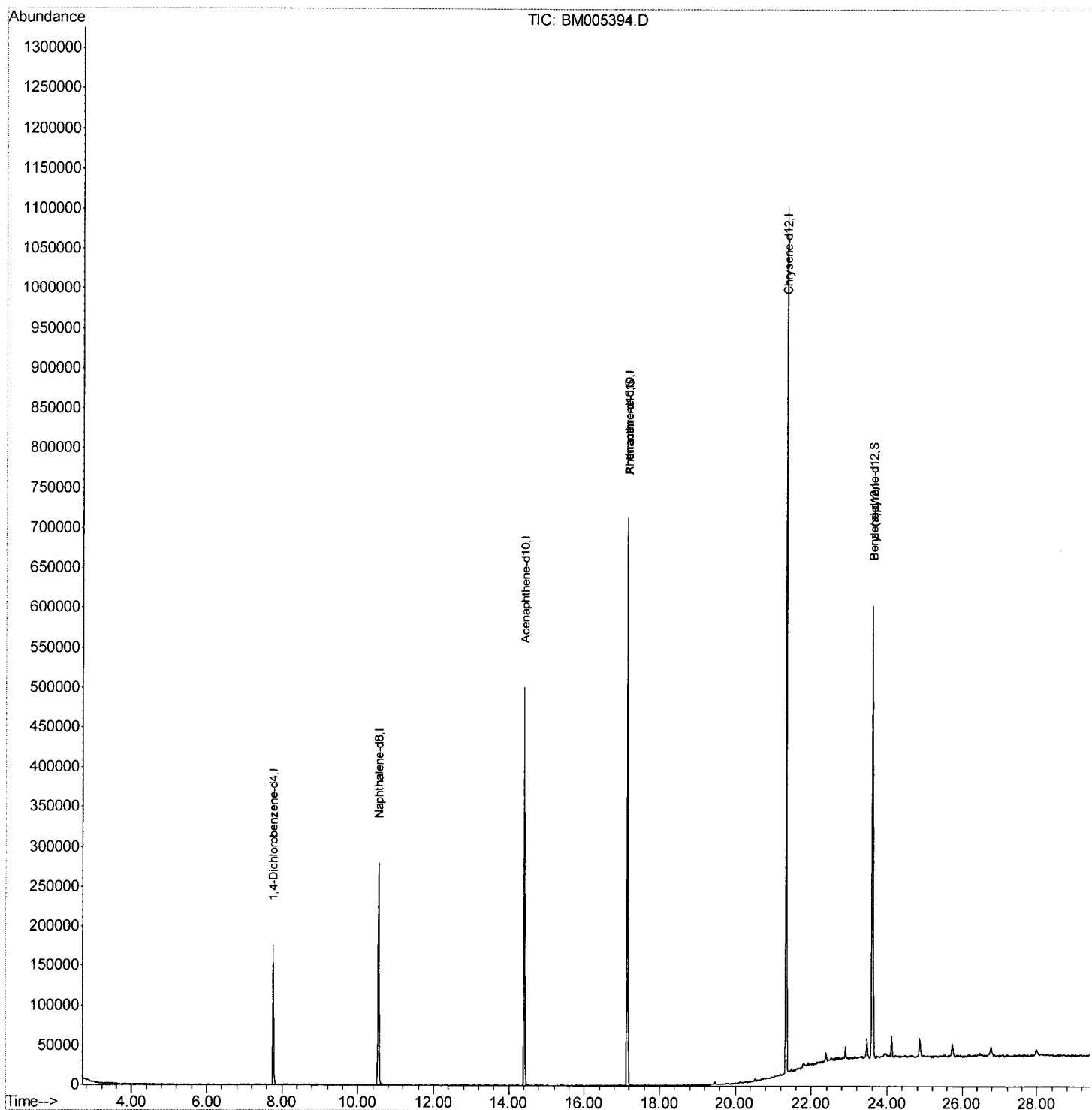
Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM051116\  
 Data File : BM005394.D  
 Acq On : 11 May 2016 19:28  
 Operator : UM/SJ  
 Sample : H2918-05  
 Misc :  
 ALS Vial : 34 Sample Multiplier: 1

Instrument :  
 BNA M  
 ClientSampleId :  
 SVOC-GPC-BLANK

Manual Integrations  
 APPROVED

sohil  
 5/12/2016 7:08:04 PM

Quant Time: May 12 04:29:34 2016  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM050516.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu May 12 02:20:05 2016  
 Response via : Initial Calibration



# Quantitation Report (Qedit)

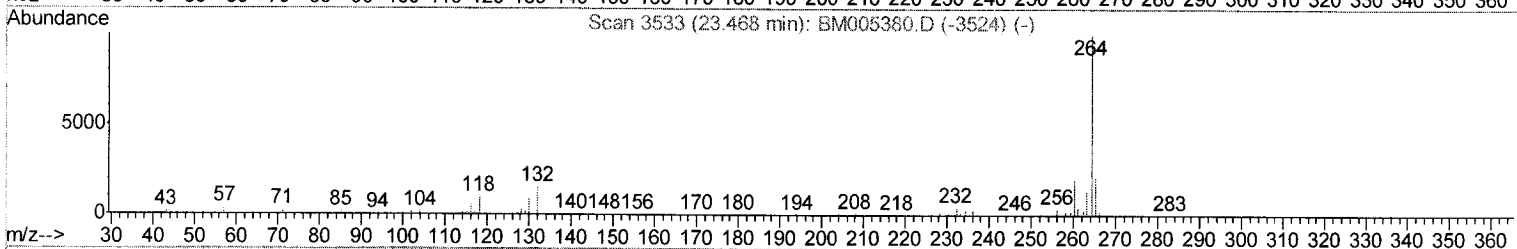
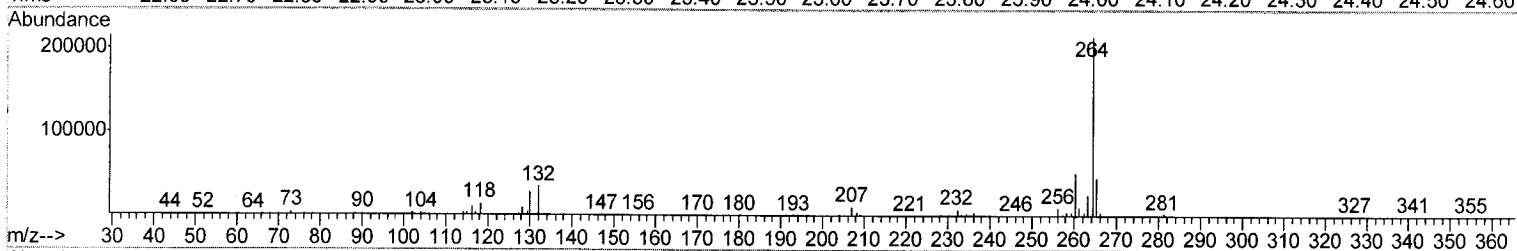
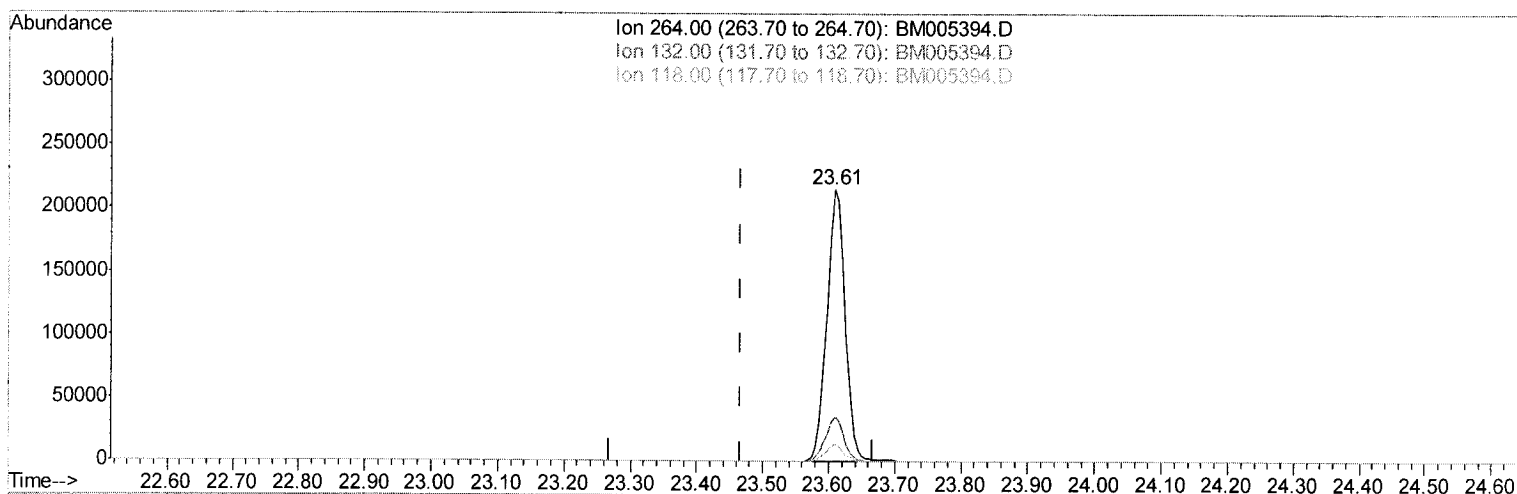
Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM051116\  
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Manual Integrations  
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 5/12/2016 7:08:04 PM

Quant Time: May 12 03:33:39 2016  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM050516.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu May 12 02:20:05 2016  
 Response via : Initial Calibration



TIC: BM005394.D

(87) Benzo(a)pyrene-d12 (S)

23.609min (+0.141) 22.60ng/ul m

response 419756

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 264.00 | 100   | 100   |
| 132.00 | 17.20 | 16.55 |
| 118.00 | 10.70 | 6.34# |
| 0.00   | 0.00  | 0.00  |

> U. M  
 05/16/16

# Quantitation Report (Qedit)

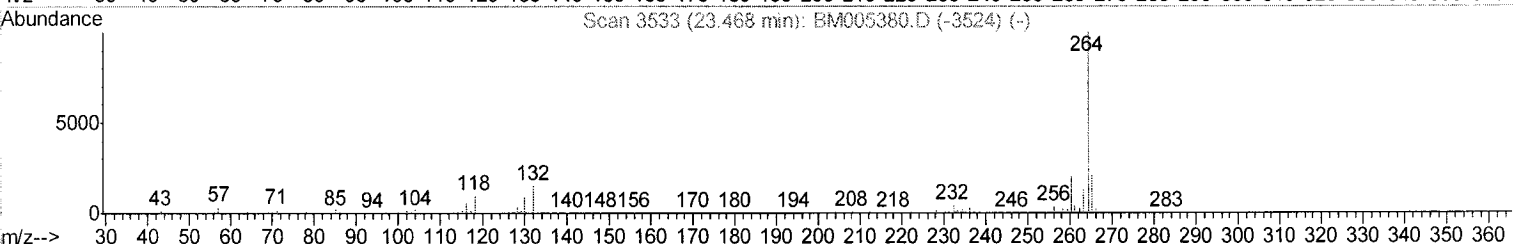
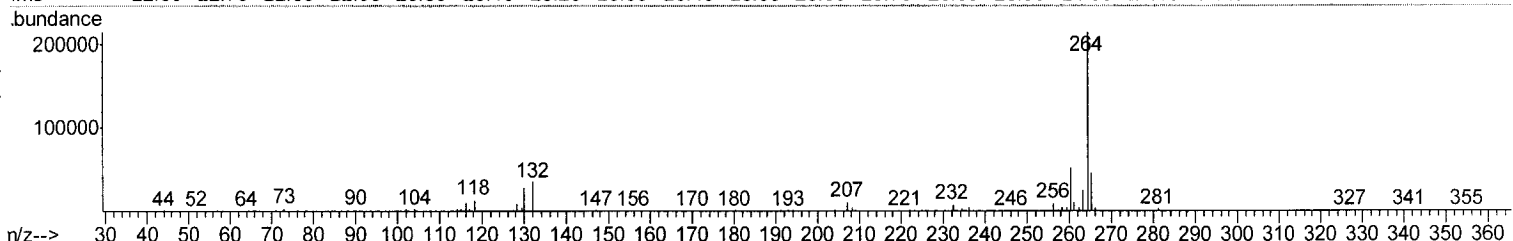
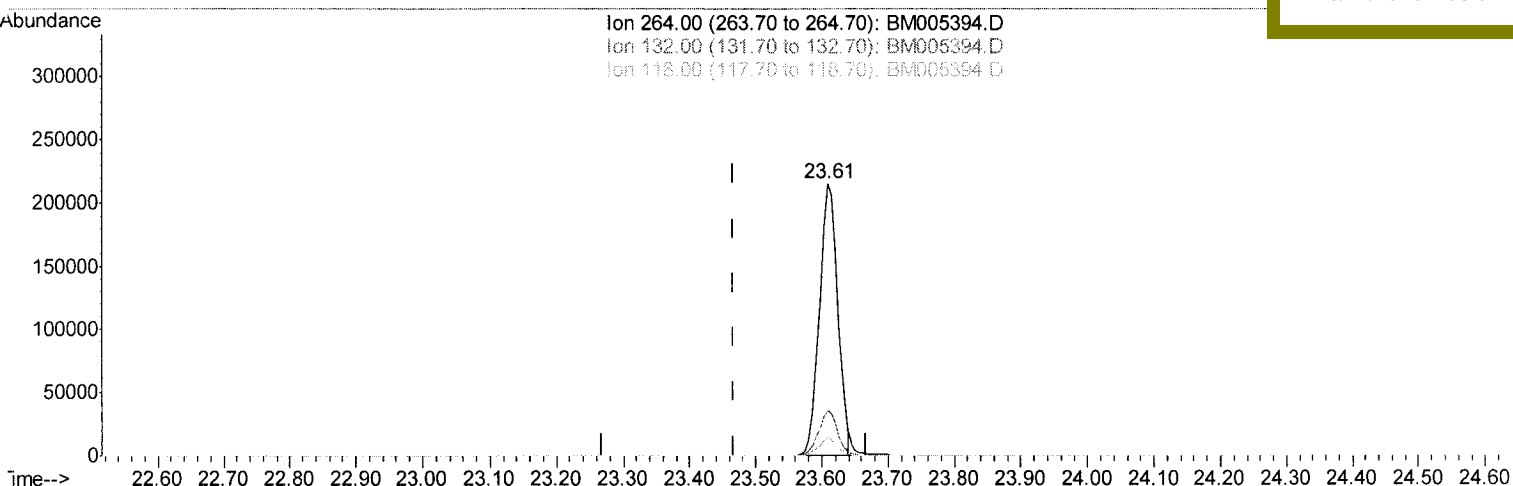
Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM0511116\  
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 Acq On : 11 May 2016 19:28  
 Operator : UM/SJ  
 Sample : H2918-05  
 Misc :  
 ALS Vial : 34 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SVOC-GPC-BLANK

Quant Time: May 12 03:33:39 2016  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM050516.M  
 Quant Title : SVOA CALIBRATION  
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 Response via : Initial Calibration

Manual Integrations  
 APPROVED

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TIC: BM005394.D

(87) Benzo(a)pyrene-d12 (S)

23.609min (+0.141) 22.15ng/ul

response 411524

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 264.00 | 100   | 100   |
| 132.00 | 17.20 | 16.55 |
| 118.00 | 10.70 | 6.34# |
| 0.00   | 0.00  | 0.00  |

Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM051116\  
 Data File : BM005394.D  
 Acq On : 11 May 2016 19:28  
 Operator : UM/SJ  
 Sample : H2918-05  
 Misc :  
 ALS Vial : 34 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SVOC-GPC-BLANK

Quant Time: May 12 04:29:34 2016  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM050516.M  
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Manual Integrations  
 APPROVED

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| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev (Min) |
|---------------------------|-------|------|----------|-------|-------|-----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.76  | 152  | 50478    | 20.00 | ng/ul | 0.00      |
| 18) Naphthalene-d8        | 10.55 | 136  | 241643   | 20.00 | ng/ul | 0.00      |
| 35) Acenaphthene-d10      | 14.40 | 164  | 171555   | 20.00 | ng/ul | 0.00      |
| 61) Phenanthrene-d10      | 17.15 | 188  | 436327   | 20.00 | ng/ul | 0.00      |
| 75) Chrysene-d12          | 21.34 | 240  | 527219   | 20.00 | ng/ul | 0.00      |
| 83) Perylene-d12          | 23.61 | 264  | 419644   | 20.00 | ng/ul | 0.00      |

## System Monitoring Compounds

|                                 |       |     |           |       |       |       |
|---------------------------------|-------|-----|-----------|-------|-------|-------|
| 3) 1,4-Dioxane-d8               | 0.00  | 96  | 0         | 0.00  | ng/uL |       |
| 5) Phenol-d5                    | 0.00  | 99  | 0         | 0.00  | ng/ul |       |
| 7) Bis- (2-Chloroethyl) ether-d | 0.00  | 67  | 0         | 0.00  | ng/ul |       |
| 9) 2-Chlorophenol-d4            | 0.00  | 132 | 0         | 0.00  | ng/ul |       |
| 13) 4-Methylphenol-d8           | 0.00  | 113 | 0         | 0.00  | ng/ul |       |
| 19) Nitrobenzene-d5             | 0.00  | 128 | 0         | 0.00  | ng/ul |       |
| 22) 2-Nitrophenol-d4            | 0.00  | 143 | 0         | 0.00  | ng/ul |       |
| 26) 2,4-Dichlorophenol-d3       | 0.00  | 165 | 0         | 0.00  | ng/ul |       |
| 29) 4-Chloroaniline-d4          | 0.00  | 131 | 0         | 0.00  | ng/ul |       |
| 43) Dimethylphthalate-d6        | 0.00  | 166 | 0         | 0.00  | ng/ul |       |
| 46) Acenaphthylene-d8           | 0.00  | 160 | 0         | 0.00  | ng/ul |       |
| 51) 4-Nitrophenol-d4            | 0.00  | 143 | 0         | 0.00  | ng/ul |       |
| 57) Fluorene-d10                | 15.41 | 176 | 631       | 0.05  | ng/ul | 0.01  |
| 62) 4,6-Dinitro-2-methylphenol  | 0.00  | 200 | 0         | 0.00  | ng/ul |       |
| 70) Anthracene-d10              | 17.15 | 188 | 436327    | 22.62 | ng/ul | -0.10 |
| 76) Pyrene-d10                  | 0.00  | 212 | 0         | 0.00  | ng/ul |       |
| 87) Benzo(a)pyrene-d12          | 23.61 | 264 | 419756m > | 22.60 | ng/ul | 0.14  |

## Target Compounds

Qvalue

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

U.M  
 05/16/16