

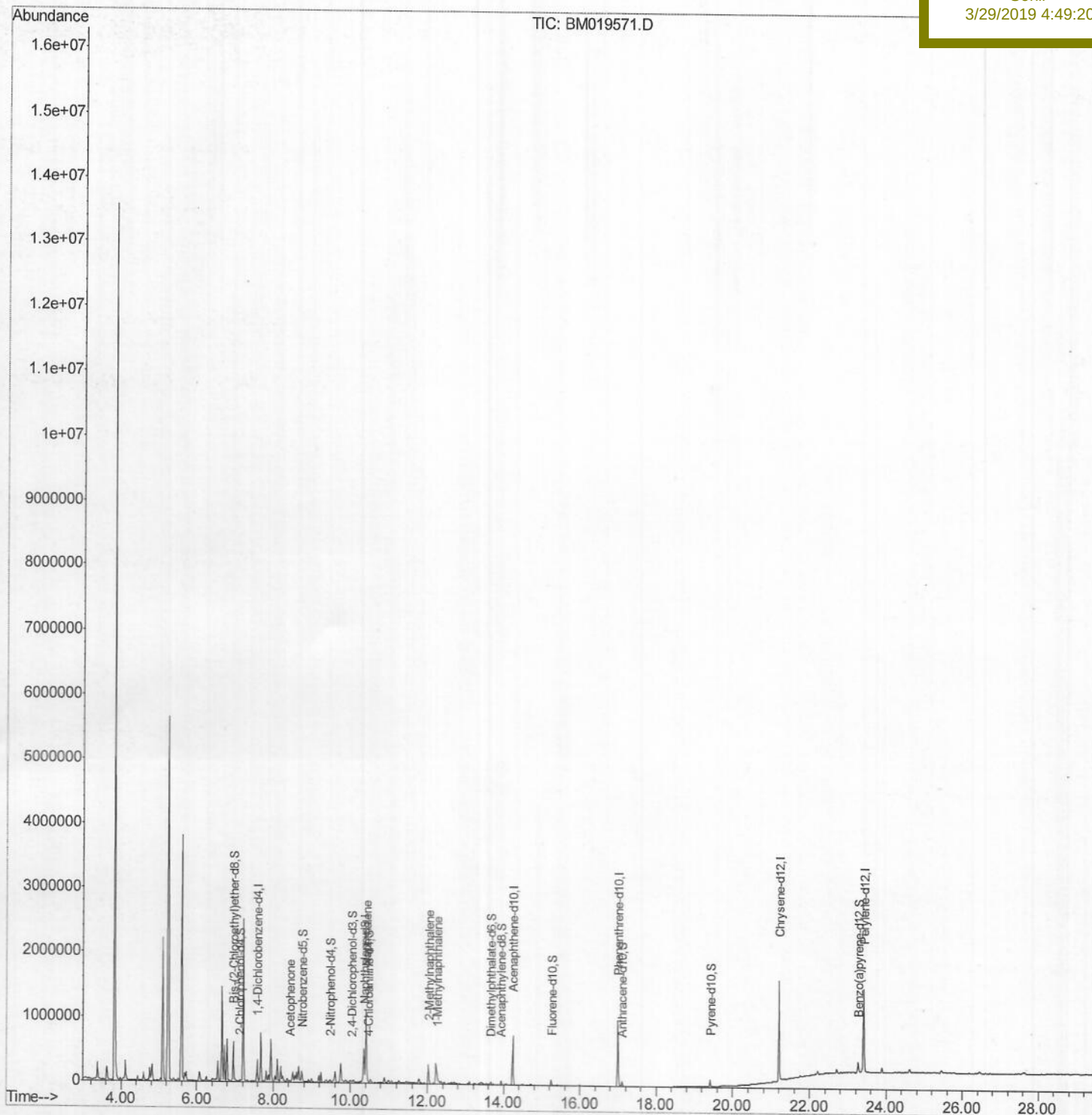
Data Path : Z:\SVOASRV\HPCHEM1\BNA\_M\DATA\BM032819\  
Data File : BM019571.D  
Acq On : 29 Mar 2019 09:28  
Operator : JU/SJ  
Sample : K2063-05DL2 20X  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Quant Time: Mar 29 12:30:05 2019  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Fri Mar 29 09:53:26 2019  
Response via : Initial Calibration

Instrument :  
BNA\_M  
ClientSampleId :  
DB6R7DL2

Manual Integrations  
APPROVED

Sohil  
3/29/2019 4:49:20 PM



# Quantitation Report (Qedit)

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032819\

Data File : BM019571.D

Acq On : 29 Mar 2019 09:28

Operator : JU/SJ

Sample : K2063-05DL2 20X

Misc :

ALS Vial : 33 Sample Multiplier: 1

Quant Time: Mar 29 11:25:49 2019

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M

Quant Title : SVOA CALIBRATION

QLast Update : Fri Mar 29 09:53:26 2019

Response via : Initial Calibration

Instrument :

BNA\_M

ClientSampled :

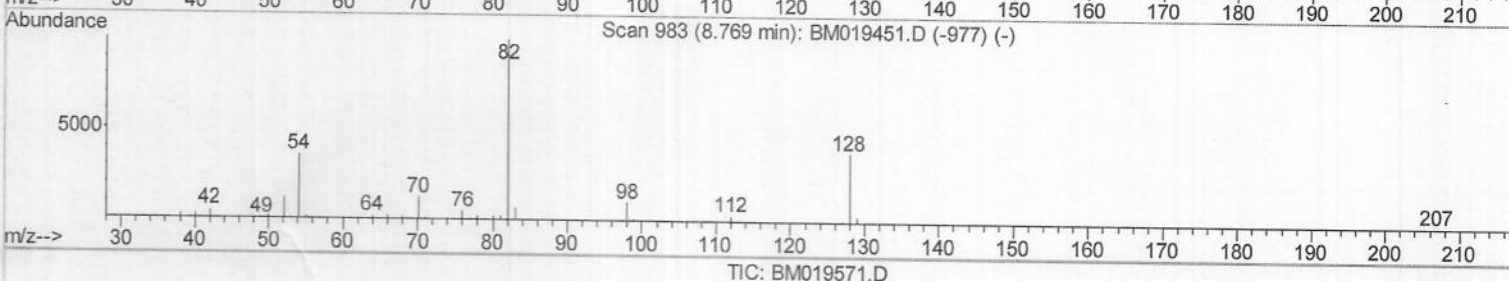
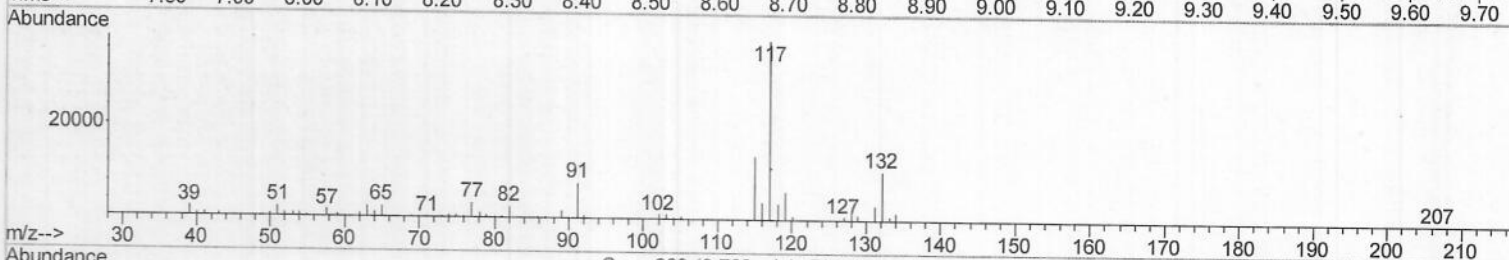
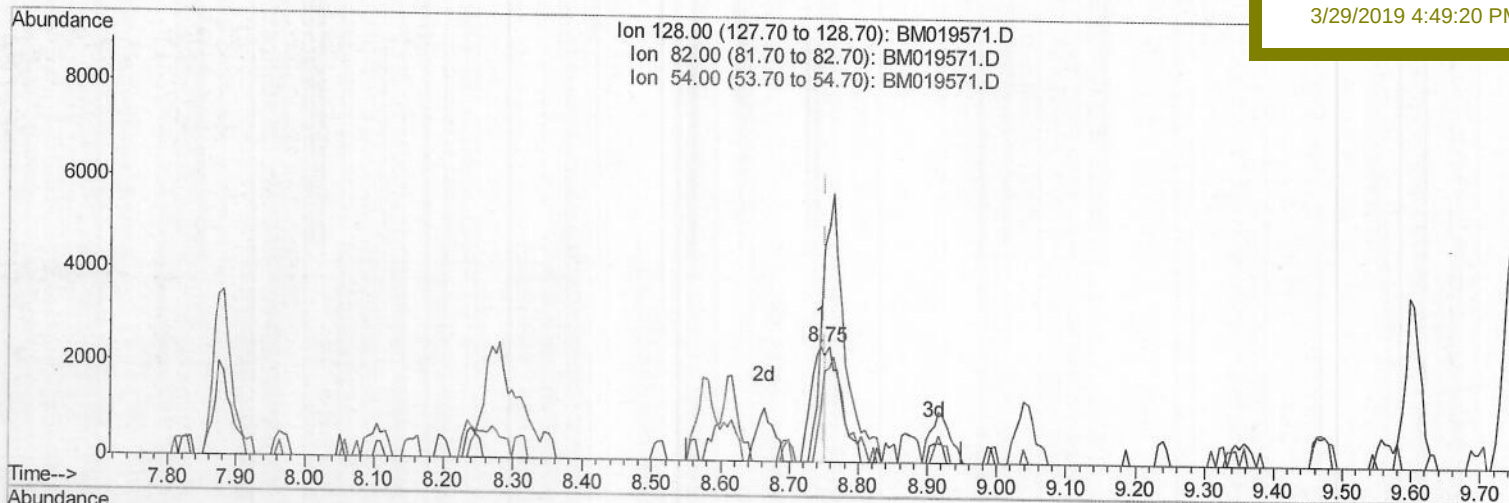
DB6R7DL2

Manual Integrations

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(19) Nitrobenzene-d5 (S)

8.745min (-0.006) 1.24ng/ul

response 3666

| Ion    | Exp%   | Act%    |
|--------|--------|---------|
| 128.00 | 100    | 100     |
| 82.00  | 244.60 | 113.98# |
| 54.00  | 108.20 | 42.43#  |
| 0.00   | 0.00   | 0.00    |

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Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032819\

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Acq On : 29 Mar 2019 09:28

Operator : JU/SJ

Sample : K2063-05DL2 20X

Misc :

ALS Vial : 33 Sample Multiplier: 1

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Quant Title : SVOA CALIBRATION

QLast Update : Fri Mar 29 09:53:26 2019

Response via : Initial Calibration

Instrument :

BNA\_M

ClientSampleId :

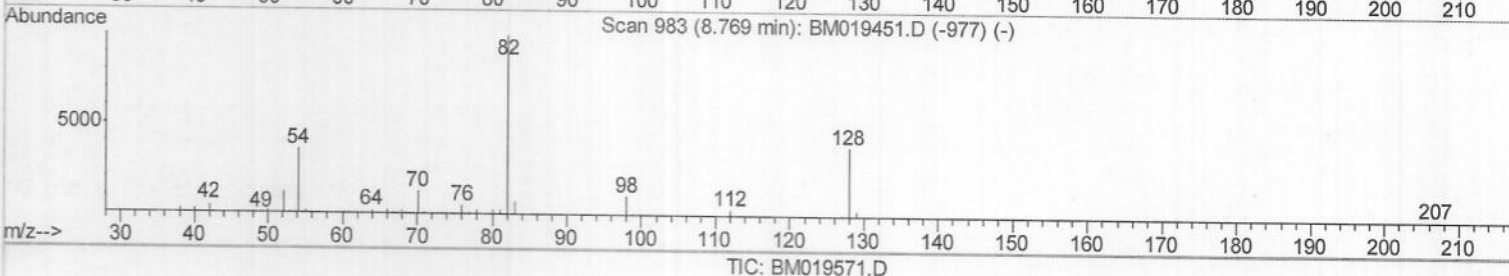
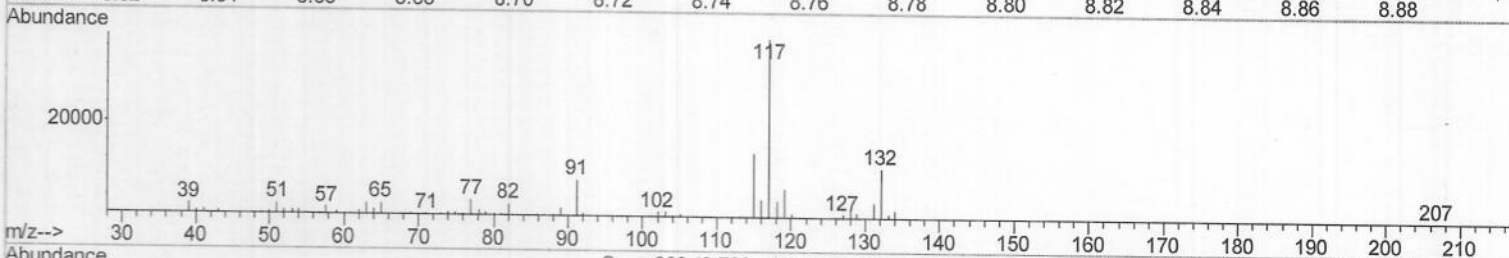
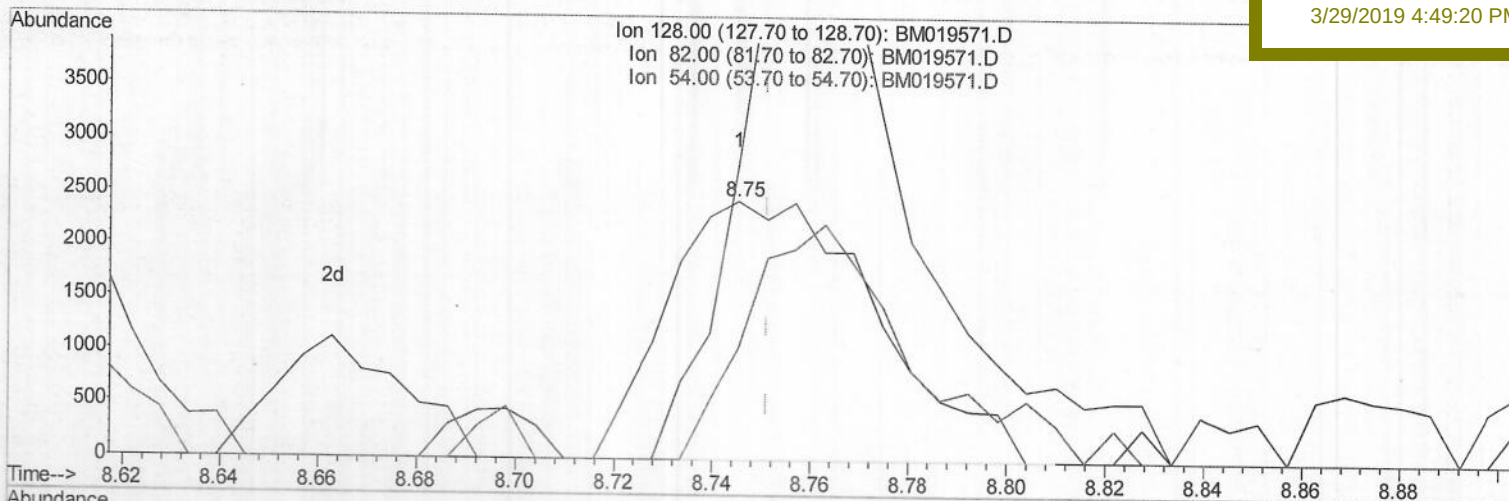
DB6R7DL2

Manual Integrations

APPROVED

Sohil

3/29/2019 4:49:20 PM



(19) Nitrobenzene-d5 (S)

8.745min (-0.006) 2.41ng/ul m

response 7141

Ion Exp% Act%

128.00 100 100

82.00 244.60 113.98#

54.00 108.20 42.43#

0.00 0.00 0.00

27  
04101119



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_M\DATA\BM032819\  
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 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Mar 29 09:53:26 2019  
 Response via : Initial Calibration

Instrument :  
 BNA\_M  
 ClientSampleId :  
 DB6R7DL2

Manual Integrations  
 APPROVED

Sohil  
 3/29/2019 4:49:20 PM

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.58  | 152  | 87621    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.36 | 136  | 415066   | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.24 | 164  | 231369   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 17.00 | 188  | 558206   | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.20 | 240  | 701850   | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 23.41 | 264  | 888807   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |       |      |       |       |
|--------------------------------|-------|-----|-------|------|-------|-------|
| 3) 1,4-Dioxane-d8              | 3.09  | 96  | 1202  | 0.54 | ng/uL | -0.03 |
| 5) Phenol-d5                   | 6.79  | 99  | 3475  | 0.45 | ng/ul | 0.03  |
| 7) Bis-(2-Chloroethyl)ether-d  | 6.95  | 67  | 7653  | 1.47 | ng/ul | 0.01  |
| 9) 2-Chlorophenol-d4           | 7.12  | 132 | 8674  | 1.47 | ng/ul | 0.01  |
| 13) 4-Methylphenol-d8          | 8.32  | 113 | 4926  | 0.84 | ng/ul | 0.02  |
| 19) Nitrobenzene-d5            | 8.75  | 128 | 7141m | 2.41 | ng/ul | 0.00  |
| 22) 2-Nitrophenol-d4           | 9.47  | 143 | 4334  | 1.32 | ng/ul | 0.00  |
| 26) 2,4-Dichlorophenol-d3      | 10.02 | 165 | 7571  | 1.22 | ng/ul | 0.02  |
| 29) 4-Chloroaniline-d4         | 10.46 | 131 | 15594 | 2.09 | ng/ul | -0.06 |
| 44) Dimethylphthalate-d6       | 13.66 | 166 | 35275 | 1.89 | ng/ul | 0.00  |
| 47) Acenaphthylene-d8          | 13.93 | 160 | 38457 | 1.65 | ng/ul | 0.00  |
| 52) 4-Nitrophenol-d4           | 0.00  | 143 | 0     | 0.00 | ng/ul |       |
| 58) Fluorene-d10               | 15.24 | 176 | 32570 | 2.02 | ng/ul | 0.00  |
| 63) 4,6-Dinitro-2-methylphenol | 15.42 | 200 | 1120  | 0.30 | ng/ul | 0.04  |
| 71) Anthracene-d10             | 17.10 | 188 | 47296 | 1.77 | ng/ul | 0.00  |
| 79) Pyrene-d10                 | 19.41 | 212 | 55718 | 1.72 | ng/ul | 0.00  |
| 90) Benzo(a)pyrene-d12         | 23.27 | 264 | 77718 | 1.65 | ng/ul | 0.00  |

## Target Compounds

|                         |       |     |        |        |       | Qvalue |
|-------------------------|-------|-----|--------|--------|-------|--------|
| 14) Acetophenone        | 8.43  | 105 | 12948  | 1.357  | ng/ul | 97     |
| 28) Naphthalene         | 10.41 | 128 | 815548 | 37.929 | ng/ul | 99     |
| 34) 2-Methylnaphthalene | 12.03 | 142 | 141230 | 9.374  | ng/ul | 99     |
| 35) 1-Methylnaphthalene | 12.25 | 142 | 69264  | 4.869  | ng/ul | 98     |

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