

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP112420\  
 Data File : BP004082.D  
 Acq On : 24 Nov 2020 18:43  
 Operator : CG/JU  
 Sample : L4829-19DL 2X  
 Misc :  
 ALS Vial : 13 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 MW-6DL

Manual Integrations  
 APPROVED

mohammad  
 11/25/2020 1:22:03 PM

Quant Time: Nov 24 23:58:40 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\8270E-BP110320.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Nov 23 16:42:41 2020  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc  | Units | Dev(Min) |
|-----------------------------|--------|------|----------|-------|-------|----------|
| -----                       |        |      |          |       |       |          |
| Internal Standards          |        |      |          |       |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 8.269  | 152  | 121595   | 20.00 | ng    | 0.00     |
| 21) Naphthalene-d8          | 11.098 | 136  | 475245   | 20.00 | ng    | # 0.00   |
| 39) Acenaphthene-d10        | 14.892 | 164  | 299345   | 20.00 | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.622 | 188  | 626598   | 20.00 | ng    | # 0.00   |
| 76) Chrysene-d12            | 21.657 | 240  | 547996   | 20.00 | ng    | 0.00     |
| 86) Perylene-d12            | 24.133 | 264  | 576662   | 20.00 | ng    | 0.00     |
|                             |        |      |          |       |       |          |
| System Monitoring Compounds |        |      |          |       |       |          |
| 5) 2-Fluorophenol           | 5.805  | 112  | 244957   | 33.62 | ng    | 0.00     |
| 7) Phenol-d6                | 7.457  | 99   | 207735   | 19.86 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.434  | 82   | 421302   | 43.42 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 16.375 | 330  | 267150   | 71.35 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.510 | 172  | 985987   | 46.05 | ng    | 0.00     |
| 79) Terphenyl-d14           | 20.121 | 244  | 1110408  | 39.15 | ng    | 0.00     |
|                             |        |      |          |       |       |          |
| Target Compounds            |        |      |          |       |       | Qvalue   |
| 20) 3+4-Methylphenols       | 9.075  | 107  | 29363    | 3.32  | ng    | # 77     |
| 27) 2,4-Dimethylphenol      | 10.275 | 122  | 52826m   | 8.41  | ng    |          |
| 31) Naphthalene             | 11.151 | 128  | 1189781  | 46.65 | ng    | 99       |
| 37) 2-Methylnaphthalene     | 12.734 | 142  | 392667   | 21.52 | ng    | 98       |
| 38) 1-Methylnaphthalene     | 12.951 | 142  | 240014   | 13.79 | ng    | 98       |
| 50) Dimethylphthalate       | 14.316 | 163  | 104521   | 4.47  | ng    | 99       |
| -----                       |        |      |          |       |       |          |

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

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP112420\  
Data File : BP004082.D  
Acq On : 24 Nov 2020 18:43  
Operator : CG/JU  
Sample : L4829-19DL 2X  
Misc :  
ALS Vial : 13 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
MW-6DL

Manual Integrations  
APPROVED

mohammad  
11/25/2020 1:22:03 PM

Quant Time: Nov 24 23:58:40 2020  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\8270E-BP110320.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
QLast Update : Mon Nov 23 16:42:41 2020  
Response via : Initial Calibration

