Quantitation Report (QT/LSC Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111225\
Data File : BP026115.D
Acq On : 12 Nov 2025 16:36
Operator : RC/JU
Sample : Q3584-07
Misc :
ALS Vial : 9 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
SEEP-1

Quant Time: Nov 12 17:20:17 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP102925.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Oct 30 11:51:34 2025
Response via : Initial Calibration

| Compound | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|---|---|---|---|---|---|---|
| Internal Standards | | | | | | |
| 1) 1,4-Dichlorobenzene-d4 | 7.672 | 152 | 324415 | 20.000 | ng | -0.02 |
| 21) Naphthalene-d8 | 10.437 | 136 | 1294546 | 20.000 | ng | -0.02 |
| 39) Acenaphthene-d10 | 14.313 | 164 | 810714 | 20.000 | ng | 0.00 |
| 64) Phenanthrene-d10 | 17.125 | 188 | 1608201 | 20.000 | ng | 0.00 |
| 76) Chrysene-d12 | 21.583 | 240 | 1784622 | 20.000 | ng | 0.02 |
| 86) Perylene-d12 | 24.924 | 264 | 1987960 | 20.000 | ng | 0.04 |
| System Monitoring Compounds | | | | | | |
| 5) 2-Fluorophenol | 5.290 | 112 | 3195 | 0.163 | ng | 0.00 |
| 7) Phenol-d6 | 6.831 | 99 | 18786 | 0.751 | ng | -0.02 |
| 23) Nitrobenzene-d5 | 8.808 | 82 | 6007 | 0.289 | ng | -0.02 |
| 42) 2,4,6-Tribromophenol | 15.831 | 330 | 3446 | 0.393 | ng | 0.00 |
| 45) 2-Fluorobiphenyl | 12.925 | 172 | 15193 | 0.269 | ng | 0.00 |
| 79) Terphenyl-d14 | 19.878 | 244 | 31631 | 0.360 | ng | 0.01 |
| Target Compounds | | | | | | Qvalue |
| 2) 1,4-Dioxane | 3.231 | 88 | 82925 | 10.005 | ng | # 84 |
| 6) Aniline | 7.002 | 93 | 90162 | 2.700 | ng | 100 |
| 10) Phenol | 6.872 | 94 | 127525 | 4.590 | ng | 99 |
| 20) 3+4-Methylphenols | 8.425 | 107 | 397532 | 15.708 | ng | 91 |
| 32) Benzoic acid | 9.684 | 122 | 22202 | 5.619 | ng | 97 |

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

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