

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041125\
 Data File : BP024271.D
 Acq On : 11 Apr 2025 18:23
 Operator : RC/JU
 Sample : Q1752-02
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TP-1

Quant Time: Apr 11 22:52:51 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP031225.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 10 11:43:37 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.739	152	305111	20.000	ng	0.00
21) Naphthalene-d8	10.516	136	1196968	20.000	ng	-0.01
39) Acenaphthene-d10	14.369	164	711365	20.000	ng	-0.02
64) Phenanthrene-d10	17.174	188	1453687	20.000	ng	-0.02
76) Chrysene-d12	21.609	240	1552111	20.000	ng	-0.04
86) Perylene-d12	24.986	264	1910036	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.357	112	2823134	147.701	ng	0.00
7) Phenol-d6	6.922	99	3266307	127.789	ng	0.00
23) Nitrobenzene-d5	8.886	82	2284018	107.661	ng	0.00
42) 2,4,6-Tribromophenol	15.892	330	1685107	188.058	ng	0.00
45) 2-Fluorobiphenyl	12.986	172	4798705	99.468	ng	-0.01
79) Terphenyl-d14	19.898	244	8324288	110.709	ng	-0.02

Target Compounds	Qvalue

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

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