

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092622\
 Data File : BP011800.D
 Acq On : 26 Sep 2022 17:02
 Operator : CG/JU
 Sample : N4790-02
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 250FERRYBLVD-SOIL-1

Quant Time: Sep 27 03:20:01 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 22 02:45:14 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.916	152	359514	20.000	ng	0.00
21) Naphthalene-d8	10.728	136	1586807	20.000	ng	0.00
39) Acenaphthene-d10	14.557	164	937217	20.000	ng	-0.01
64) Phenanthrene-d10	17.316	188	1918284	20.000	ng	0.00
76) Chrysene-d12	21.398	240	1721090	20.000	ng	0.00
86) Perylene-d12	23.845	264	1639653	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.475	112	3301820	161.144	ng	0.00
7) Phenol-d6	7.081	99	4256154	140.484	ng	0.00
23) Nitrobenzene-d5	9.087	82	2666500	87.954	ng	0.00
42) 2,4,6-Tribromophenol	16.051	330	1200380	190.213	ng	0.00
45) 2-Fluorobiphenyl	13.187	172	5545304	93.567	ng	0.00
79) Terphenyl-d14	19.869	244	7995146	106.813	ng	0.00
Target Compounds						
27) 2,4-Dimethylphenol	9.899	122	65631	3.267	ng	Qvalue 99

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

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