

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092322\
 Data File : BP011780.D
 Acq On : 23 Sep 2022 11:23
 Operator : CG/JU
 Sample : PB147850BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB147850BL

Quant Time: Sep 24 01:33:22 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	205283	20.000	ng	0.00
21) Naphthalene-d8	10.728	136	895912	20.000	ng	0.00
39) Acenaphthene-d10	14.563	164	589917	20.000	ng	0.00
64) Phenanthrene-d10	17.316	188	1321683	20.000	ng	0.00
76) Chrysene-d12	21.398	240	1696839	20.000	ng	0.00
86) Perylene-d12	23.851	264	1848819	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.475	112	1688690	144.336	ng	0.00
7) Phenol-d6	7.075	99	2370133	137.008	ng	0.00
23) Nitrobenzene-d5	9.081	82	1352134	78.994	ng	-0.01
42) 2,4,6-Tribromophenol	16.051	330	528385	133.021	ng	0.00
45) 2-Fluorobiphenyl	13.187	172	3016565	80.865	ng	0.00
79) Terphenyl-d14	19.875	244	7281426	98.668	ng	0.00

Target Compounds	Qvalue

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

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