

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092722\
 Data File : BP011810.D
 Acq On : 27 Sep 2022 11:40
 Operator : CG/JU
 Sample : PB147841TB
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB147841TB

Quant Time: Sep 28 02:06:05 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 28 02:03:33 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.910	152	120847	20.000	ng	0.00
21) Naphthalene-d8	10.722	136	529750	20.000	ng	0.00
39) Acenaphthene-d10	14.551	164	353384	20.000	ng	0.00
64) Phenanthrene-d10	17.310	188	883591	20.000	ng	0.00
76) Chrysene-d12	21.392	240	1018815	20.000	ng	0.00
86) Perylene-d12	23.839	264	1082050	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.469	112	1159489	168.348	ng	0.00
7) Phenol-d6	7.069	99	1591651	156.292	ng	0.00
23) Nitrobenzene-d5	9.075	82	788108	77.867	ng	0.00
42) 2,4,6-Tribromophenol	16.045	330	412319	173.280	ng	0.00
45) 2-Fluorobiphenyl	13.175	172	2033685	91.007	ng	0.00
79) Terphenyl-d14	19.863	244	4416660	99.678	ng	0.00

Target Compounds	Qvalue

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

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