

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092022\
 Data File : BP011760.D
 Acq On : 20 Sep 2022 13:26
 Operator : CG/JU
 Sample : PB147737BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB147737BL

Quant Time: Sep 21 06:45:11 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 20 04:42:52 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	556038	20.000	ng	0.00
21) Naphthalene-d8	10.746	136	2251575	20.000	ng	0.00
39) Acenaphthene-d10	14.575	164	1231070	20.000	ng	0.00
64) Phenanthrene-d10	17.328	188	2329494	20.000	ng	# 0.00
76) Chrysene-d12	21.404	240	1641980	20.000	ng	# 0.00
86) Perylene-d12	23.857	264	1326546	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.493	112	5681536	185.987	ng	0.00
7) Phenol-d6	7.099	99	7082751	173.934	ng	0.00
23) Nitrobenzene-d5	9.099	82	4389891	115.631	ng	0.00
42) 2,4,6-Tribromophenol	16.063	330	1577062	163.956	ng	0.00
45) 2-Fluorobiphenyl	13.198	172	8169519	103.154	ng	0.00
79) Terphenyl-d14	19.880	244	10340751	136.949	ng	0.00

Target Compounds Qvalue

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

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