

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092022\
 Data File : BP011767.D
 Acq On : 20 Sep 2022 18:41
 Operator : CG/JU
 Sample : PB147694TB
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB147694TB

Quant Time: Sep 21 06:51:08 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	7.928	152	535002	20.000	ng	0.00
21) Naphthalene-d8	10.740	136	2233432	20.000	ng	0.00
39) Acenaphthene-d10	14.569	164	1246295	20.000	ng	0.00
64) Phenanthrene-d10	17.322	188	2299617	20.000	ng	# 0.00
76) Chrysene-d12	21.404	240	1446589	20.000	ng	# 0.00
86) Perylene-d12	23.851	264	1148886	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.493	112	5401027	183.756	ng	0.00
7) Phenol-d6	7.093	99	6775675	172.936	ng	0.00
23) Nitrobenzene-d5	9.099	82	4146237	110.100	ng	0.00
42) 2,4,6-Tribromophenol	16.063	330	1427774	146.622	ng	0.00
45) 2-Fluorobiphenyl	13.198	172	7777197	97.001	ng	0.00
79) Terphenyl-d14	19.875	244	9091313	136.665	ng	0.00

Target Compounds Qvalue

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

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