

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092021\
 Data File : BP007045.D
 Acq On : 20 Sep 2021 10:53
 Operator : CG/JU
 Sample : PB139134BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB139134BL

Quant Time: Sep 20 11:32:01 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 17 10:50:16 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.911	152	296622	20.000	ng	0.00
21) Naphthalene-d8	10.710	136	1156890	20.000	ng	# 0.00
39) Acenaphthene-d10	14.540	164	727252	20.000	ng	0.00
64) Phenanthrene-d10	17.287	188	1552788	20.000	ng	# 0.00
76) Chrysene-d12	21.363	240	1596109	20.000	ng	# 0.00
86) Perylene-d12	23.786	264	1727906	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.481	112	2279709	125.500	ng	0.00
7) Phenol-d6	7.075	99	2785767	112.358	ng	0.00
23) Nitrobenzene-d5	9.069	82	1583425	76.872	ng	0.00
42) 2,4,6-Tribromophenol	16.028	330	1198101	136.988	ng	0.00
45) 2-Fluorobiphenyl	13.163	172	4079325	77.436	ng	0.00
79) Terphenyl-d14	19.839	244	6795372	81.684	ng	0.00

Target Compounds	Qvalue

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

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