

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110821\
 Data File : BP007870.D
 Acq On : 08 Nov 2021 20:23
 Operator : CG/JU
 Sample : M4579-06
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SB-12-4.5-5.0

Quant Time: Nov 09 08:45:14 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP110221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 08 16:36:40 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.863	152	109054	20.000	ng	0.00
21) Naphthalene-d8	10.663	136	414994	20.000	ng	0.00
39) Acenaphthene-d10	14.498	164	265169	20.000	ng	0.00
64) Phenanthrene-d10	17.251	188	598208	20.000	ng	-0.01
76) Chrysene-d12	21.345	240	717936	20.000	ng	0.00
86) Perylene-d12	23.751	264	821271	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.452	112	1133845	176.328	ng	0.00
7) Phenol-d6	7.040	99	1386231	148.942	ng	0.00
23) Nitrobenzene-d5	9.022	82	799804	103.527	ng	-0.01
42) 2,4,6-Tribromophenol	15.992	330	583302	154.136	ng	0.00
45) 2-Fluorobiphenyl	13.122	172	1898470	104.262	ng	-0.01
79) Terphenyl-d14	19.816	244	3285778	93.195	ng	0.00

Target Compounds	Qvalue

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

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