

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092021\
 Data File : BP007063.D
 Acq On : 20 Sep 2021 22:33
 Operator : CG/JU
 Sample : M3817-01
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SW-RR-02-(1.0)

Quant Time: Sep 21 01:04:12 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	7.905	152	349528	20.000	ng	0.00
21) Naphthalene-d8	10.704	136	1352751	20.000	ng	# 0.00
39) Acenaphthene-d10	14.534	164	829976	20.000	ng	0.00
64) Phenanthrene-d10	17.281	188	1698734	20.000	ng	# 0.00
76) Chrysene-d12	21.363	240	1663578	20.000	ng	# 0.00
86) Perylene-d12	23.774	264	1702117	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.481	112	1179575	55.107	ng	0.00
7) Phenol-d6	7.069	99	870413	29.792	ng	0.00
23) Nitrobenzene-d5	9.069	82	1981716	82.279	ng	0.00
42) 2,4,6-Tribromophenol	16.022	330	1328934	133.141	ng	0.00
45) 2-Fluorobiphenyl	13.163	172	4622337	76.884	ng	0.00
79) Terphenyl-d14	19.839	244	6649165	76.685	ng	0.00
Target Compounds						
						Qvalue
20) 3+4-Methylphenols	8.669	107	390231	15.239	ng	87
31) Naphthalene	10.757	128	284017	3.789	ng	99
32) Benzoic acid	9.969	122	63725	9.994	ng	95

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

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