

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092021\
 Data File : BP007061.D
 Acq On : 20 Sep 2021 21:25
 Operator : CG/JU
 Sample : M3817-07
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SW-CB-03-(1.0)

Quant Time: Sep 21 01:03:42 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	313395	20.000	ng	0.00
21) Naphthalene-d8	10.705	136	1227086	20.000	ng	# 0.00
39) Acenaphthene-d10	14.534	164	774991	20.000	ng	0.00
64) Phenanthrene-d10	17.281	188	1626683	20.000	ng	# 0.00
76) Chrysene-d12	21.357	240	1603501	20.000	ng	#-0.01
86) Perylene-d12	23.774	264	1634828	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.481	112	1062665	55.370	ng	0.00
7) Phenol-d6	7.069	99	759477	28.993	ng	0.00
23) Nitrobenzene-d5	9.069	82	1891410	86.571	ng	0.00
42) 2,4,6-Tribromophenol	16.022	330	1383939	148.489	ng	0.00
45) 2-Fluorobiphenyl	13.163	172	4757308	84.743	ng	0.00
79) Terphenyl-d14	19.839	244	6707189	80.253	ng	0.00

Target Compounds	Qvalue

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

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