

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091621\
 Data File : BP007004.D
 Acq On : 16 Sep 2021 20:02
 Operator : CG/JU
 Sample : M3709-05
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 250EM-DISPOSAL-2

Quant Time: Sep 17 02:39:01 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 16 15:16:52 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.910	152	304618	20.000	ng	-0.01
21) Naphthalene-d8	10.710	136	1149082	20.000	ng	#-0.01
39) Acenaphthene-d10	14.534	164	695697	20.000	ng	-0.01
64) Phenanthrene-d10	17.280	188	1514884	20.000	ng	#-0.01
76) Chrysene-d12	21.363	240	1524864	20.000	ng	#-0.02
86) Perylene-d12	23.774	264	1536602	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.487	112	2464921	132.134	ng	0.00
7) Phenol-d6	7.075	99	2722313	106.917	ng	0.00
23) Nitrobenzene-d5	9.069	82	1876465	91.718	ng	-0.01
42) 2,4,6-Tribromophenol	16.022	330	1308425	156.388	ng	-0.01
45) 2-Fluorobiphenyl	13.163	172	4560156	90.489	ng	-0.01
79) Terphenyl-d14	19.839	244	7964948	100.217	ng	-0.01

Target Compounds	Qvalue

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

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