

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091621\
 Data File : BP007013.D
 Acq On : 17 Sep 2021 01:08
 Operator : CG/JU
 Sample : M3687-31
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A003145

Quant Time: Sep 17 02:41:07 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	7.911	152	319962	20.000	ng	-0.01
21) Naphthalene-d8	10.710	136	1248602	20.000	ng	#-0.01
39) Acenaphthene-d10	14.534	164	759401	20.000	ng	-0.01
64) Phenanthrene-d10	17.281	188	1602946	20.000	ng	#-0.01
76) Chrysene-d12	21.357	240	1579830	20.000	ng	#-0.02
86) Perylene-d12	23.769	264	1588692	20.000	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.481	112	2079555	106.130	ng	0.00
7) Phenol-d6	7.075	99	2635755	98.553	ng	0.00
23) Nitrobenzene-d5	9.069	82	1478870	66.523	ng	-0.01
42) 2,4,6-Tribromophenol	16.022	330	929260	101.751	ng	-0.01
45) 2-Fluorobiphenyl	13.163	172	3662039	66.572	ng	-0.01
79) Terphenyl-d14	19.839	244	5754711	69.888	ng	-0.01

Target Compounds	Qvalue

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

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