

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092021\
 Data File : BP007050.D
 Acq On : 20 Sep 2021 13:57
 Operator : CG/JU
 Sample : M3687-28
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A004015

Quant Time: Sep 20 14:26:41 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	7.910	152	280749	20.000	ng	0.00
21) Naphthalene-d8	10.710	136	1064084	20.000	ng	# 0.00
39) Acenaphthene-d10	14.534	164	662195	20.000	ng	0.00
64) Phenanthrene-d10	17.286	188	1402379	20.000	ng	# 0.00
76) Chrysene-d12	21.363	240	1474742	20.000	ng	# 0.00
86) Perylene-d12	23.786	264	1619303	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.481	112	1633169	94.990	ng	0.00
7) Phenol-d6	7.075	99	2050489	87.378	ng	0.00
23) Nitrobenzene-d5	9.069	82	1207211	63.719	ng	0.00
42) 2,4,6-Tribromophenol	16.022	330	862398	108.292	ng	0.00
45) 2-Fluorobiphenyl	13.163	172	2805012	58.477	ng	0.00
79) Terphenyl-d14	19.839	244	4275649	55.626	ng	0.00

Target Compounds	Qvalue

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

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