

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111521\
 Data File : BP007989.D
 Acq On : 15 Nov 2021 19:19
 Operator : CG/JU
 Sample : PB140390BL
 Misc : LOQ-WATER 10ppm
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB140390BL

Quant Time: Nov 16 03:16:59 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 15 05:36:49 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.840	152	136027	20.000	ng	0.00
21) Naphthalene-d8	10.634	136	486628	20.000	ng	0.00
39) Acenaphthene-d10	14.475	164	340680	20.000	ng	0.00
64) Phenanthrene-d10	17.234	188	800885	20.000	ng	# 0.00
76) Chrysene-d12	21.327	240	944612	20.000	ng	# 0.00
86) Perylene-d12	23.733	264	1041640	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.428	112	1151022	137.268	ng	0.00
7) Phenol-d6	7.016	99	1410328	126.427	ng	0.00
23) Nitrobenzene-d5	8.999	82	800410	88.020	ng	0.00
42) 2,4,6-Tribromophenol	15.975	330	758610	141.527	ng	0.00
45) 2-Fluorobiphenyl	13.104	172	2067915	85.145	ng	0.00
79) Terphenyl-d14	19.804	244	3982782	85.495	ng	0.00

Target Compounds	Qvalue

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

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