

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111521\
 Data File : BP007988.D
 Acq On : 15 Nov 2021 18:45
 Operator : CG/JU
 Sample : PB140347BL
 Misc : LOQ-WATER 10ppm
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB140347BL

Quant Time: Nov 16 03:16:30 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	135526	20.000	ng	0.00
21) Naphthalene-d8	10.634	136	494884	20.000	ng	# 0.00
39) Acenaphthene-d10	14.474	164	343143	20.000	ng	0.00
64) Phenanthrene-d10	17.233	188	780100	20.000	ng	# 0.00
76) Chrysene-d12	21.327	240	945043	20.000	ng	# 0.00
86) Perylene-d12	23.733	264	1074314	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.428	112	1160087	138.861	ng	0.00
7) Phenol-d6	7.016	99	1432182	128.861	ng	0.00
23) Nitrobenzene-d5	8.998	82	807959	87.368	ng	0.00
42) 2,4,6-Tribromophenol	15.969	330	734477	136.041	ng	0.00
45) 2-Fluorobiphenyl	13.104	172	2051410	83.859	ng	0.00
79) Terphenyl-d14	19.804	244	3887376	83.409	ng	0.00

Target Compounds	Qvalue

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

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