

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111721\
 Data File : BP008002.D
 Acq On : 17 Nov 2021 10:20
 Operator : CG/JU
 Sample : PB140347BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB140347BL

Quant Time: Nov 17 11:09:22 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	130707	20.000	ng	0.00
21) Naphthalene-d8	10.640	136	506750	20.000	ng	0.00
39) Acenaphthene-d10	14.481	164	344268	20.000	ng	0.00
64) Phenanthrene-d10	17.234	188	781163	20.000	ng	# 0.00
76) Chrysene-d12	21.327	240	941888	20.000	ng	# 0.00
86) Perylene-d12	23.733	264	1064192	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.428	112	1216809	151.020	ng	0.00
7) Phenol-d6	7.016	99	1347111	125.676	ng	0.00
23) Nitrobenzene-d5	8.999	82	835582	88.239	ng	0.00
42) 2,4,6-Tribromophenol	15.975	330	742050	136.995	ng	0.00
45) 2-Fluorobiphenyl	13.104	172	2085859	84.988	ng	0.00
79) Terphenyl-d14	19.804	244	3977449	85.627	ng	0.00

Target Compounds	Qvalue

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

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