

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112421\
 Data File : BM033249.D
 Acq On : 24 Nov 2021 16:24
 Operator : CG/JU
 Sample : PB140809BL
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB140809BL

Quant Time: Nov 26 04:08:51 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM112421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 24 15:16:33 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.937	152	111290	20.000	ng	0.00
21) Naphthalene-d8	10.731	136	416941	20.000	ng	# 0.00
39) Acenaphthene-d10	14.560	164	242339	20.000	ng	0.00
64) Phenanthrene-d10	17.295	188	482064	20.000	ng	# 0.00
76) Chrysene-d12	21.453	240	473986	20.000	ng	0.00
86) Perylene-d12	23.789	264	489389	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.507	112	589159	87.422	ng	0.00
7) Phenol-d6	7.096	99	756857	83.856	ng	0.00
23) Nitrobenzene-d5	9.095	82	759587	84.574	ng	0.00
42) 2,4,6-Tribromophenol	16.042	330	204382	77.367	ng	0.00
45) 2-Fluorobiphenyl	13.183	172	1466859	85.365	ng	0.00
79) Terphenyl-d14	19.901	244	2190406	91.380	ng	0.00

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

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