

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121621\
 Data File : BM033489.D
 Acq On : 16 Dec 2021 17:59
 Operator : CG/JU
 Sample : PB141339BL
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB141339BL

Quant Time: Dec 17 08:00:00 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM121621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 16 17:10:15 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.943	152	31539	20.000	ng	0.00
21) Naphthalene-d8	10.754	136	122727	20.000	ng	0.00
39) Acenaphthene-d10	14.584	164	78635	20.000	ng	0.00
64) Phenanthrene-d10	17.336	188	150130	20.000	ng	# 0.00
76) Chrysene-d12	21.513	240	85727	20.000	ng	0.00
86) Perylene-d12	23.936	264	59144	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.502	112	190490	104.220	ng	0.00
7) Phenol-d6	7.108	99	242785	98.980	ng	0.00
23) Nitrobenzene-d5	9.113	82	282172	98.722	ng	0.00
42) 2,4,6-Tribromophenol	16.078	330	74573	84.305	ng	0.00
45) 2-Fluorobiphenyl	13.213	172	551098	96.946	ng	0.00
79) Terphenyl-d14	19.954	244	635813	116.764	ng	0.00

Target Compounds	Qvalue

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

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