

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112621\
 Data File : BM033266.D
 Acq On : 26 Nov 2021 15:53
 Operator : CG/JU
 Sample : PB141003BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB141003BL

Quant Time: Nov 26 17:28:01 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	131428	20.000	ng	0.00
21) Naphthalene-d8	10.731	136	502542	20.000	ng	# 0.00
39) Acenaphthene-d10	14.554	164	286103	20.000	ng	0.00
64) Phenanthrene-d10	17.295	188	522148	20.000	ng	0.00
76) Chrysene-d12	21.453	240	489768	20.000	ng	0.00
86) Perylene-d12	23.783	264	552235	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.513	112	1031193	129.568	ng	0.00
7) Phenol-d6	7.101	99	1329125	124.696	ng	0.00
23) Nitrobenzene-d5	9.095	82	920155	85.001	ng	0.00
42) 2,4,6-Tribromophenol	16.042	330	336657	107.945	ng	0.00
45) 2-Fluorobiphenyl	13.183	172	1727678	85.163	ng	0.00
79) Terphenyl-d14	19.901	244	2040247	82.373	ng	0.00

Target Compounds	Qvalue

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

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