

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110421\
 Data File : BM032868.D
 Acq On : 04 Nov 2021 10:45
 Operator : CG/JU
 Sample : PB140471BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB140471BL

Quant Time: Nov 04 14:28:53 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM110221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 02 16:07:18 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.489	152	105283	20.000	ng	0.00
21) Naphthalene-d8	10.272	136	461354	20.000	ng	# 0.00
39) Acenaphthene-d10	14.148	164	319031	20.000	ng	0.00
64) Phenanthrene-d10	16.883	188	674975	20.000	ng	#-0.01
76) Chrysene-d12	21.065	240	660867	20.000	ng	#-0.01
86) Perylene-d12	23.183	264	682782	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.078	112	955793	158.271	ng	0.00
7) Phenol-d6	6.695	99	1244430	138.374	ng	0.00
23) Nitrobenzene-d5	8.642	82	763634	90.594	ng	0.00
42) 2,4,6-Tribromophenol	15.642	330	442508	123.850	ng	0.00
45) 2-Fluorobiphenyl	12.760	172	1790999	87.070	ng	0.00
79) Terphenyl-d14	19.518	244	2602196	85.967	ng	0.00

Target Compounds	Qvalue

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

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