

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM022322\
 Data File : BM034047.D
 Acq On : 23 Feb 2022 10:58
 Operator : CG/JU
 Sample : PB142857BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB142857BL

Quant Time: Feb 23 11:38:47 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM022222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 22 17:55:56 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.019	152	226442	20.000	ng	0.00
21) Naphthalene-d8	10.836	136	863546	20.000	ng	# 0.00
39) Acenaphthene-d10	14.654	164	577338	20.000	ng	0.00
64) Phenanthrene-d10	17.389	188	1247006	20.000	ng	# 0.00
76) Chrysene-d12	21.565	240	1122420	20.000	ng	0.00
86) Perylene-d12	24.024	264	982184	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.566	112	1793334	141.177	ng	0.00
7) Phenol-d6	7.172	99	2239980	134.287	ng	0.00
23) Nitrobenzene-d5	9.178	82	1423280	88.945	ng	0.00
42) 2,4,6-Tribromophenol	16.136	330	989859	141.267	ng	0.00
45) 2-Fluorobiphenyl	13.283	172	3842277	88.667	ng	0.00
79) Terphenyl-d14	20.012	244	6464092	100.285	ng	0.00

Target Compounds	Qvalue

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

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