

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020922\
 Data File : BM034032.D
 Acq On : 09 Feb 2022 17:44
 Operator : CG/JU
 Sample : PB142558BL
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB142558BL

Quant Time: Feb 10 04:06:23 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM020922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 09 17:52:36 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.900	152	138166	20.000	ng	0.00
21) Naphthalene-d8	10.715	136	529969	20.000	ng	# 0.00
39) Acenaphthene-d10	14.544	164	313172	20.000	ng	0.00
64) Phenanthrene-d10	17.270	188	679156	20.000	ng	0.00
76) Chrysene-d12	21.437	240	606748	20.000	ng	# 0.00
86) Perylene-d12	23.825	264	601663	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.444	112	1111088	130.211	ng	0.00
7) Phenol-d6	7.066	99	1418473	136.361	ng	0.00
23) Nitrobenzene-d5	9.071	82	924133	81.907	ng	0.00
42) 2,4,6-Tribromophenol	16.031	330	498813	143.886	ng	0.00
45) 2-Fluorobiphenyl	13.170	172	2239521	93.889	ng	0.00
79) Terphenyl-d14	19.883	244	2570049	83.773	ng	0.00

Target Compounds	Qvalue

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

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