

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM011022\
 Data File : BM033687.D
 Acq On : 10 Jan 2022 15:26
 Operator : CG/JU
 Sample : PB141902BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB141902BL

Quant Time: Jan 11 02:25:30 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM010522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 05 17:07:26 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.978	152	171460	20.000	ng	0.00
21) Naphthalene-d8	10.789	136	702190	20.000	ng	-0.02
39) Acenaphthene-d10	14.613	164	402179	20.000	ng	-0.01
64) Phenanthrene-d10	17.348	188	639835	20.000	ng	-0.01
76) Chrysene-d12	21.512	240	625489	20.000	ng	-0.01
86) Perylene-d12	23.935	264	683239	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.525	112	1482067	136.325	ng	0.00
7) Phenol-d6	7.137	99	1854189	126.546	ng	0.00
23) Nitrobenzene-d5	9.142	82	1268076	83.401	ng	-0.01
42) 2,4,6-Tribromophenol	16.095	330	541636	128.941	ng	-0.01
45) 2-Fluorobiphenyl	13.248	172	2670619	86.975	ng	-0.01
79) Terphenyl-d14	19.959	244	2551631	73.429	ng	-0.01

Target Compounds	Qvalue

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

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