

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM031622\
 Data File : BM034257.D
 Acq On : 16 Mar 2022 13:43
 Operator : CG/JU
 Sample : PB143171BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB143171BL

Quant Time: Mar 16 14:25:30 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM031522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 16 02:38:26 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.948	152	143145	20.000	ng	0.00
21) Naphthalene-d8	10.760	136	524287	20.000	ng	0.00
39) Acenaphthene-d10	14.583	164	281008	20.000	ng	0.00
64) Phenanthrene-d10	17.324	188	486997	20.000	ng	0.00
76) Chrysene-d12	21.494	240	438676	20.000	ng	0.00
86) Perylene-d12	23.906	264	479829	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.501	112	1250969	156.778	ng	0.00
7) Phenol-d6	7.107	99	1514364	132.555	ng	0.00
23) Nitrobenzene-d5	9.107	82	1012710	94.043	ng	0.00
42) 2,4,6-Tribromophenol	16.071	330	457254	120.723	ng	0.00
45) 2-Fluorobiphenyl	13.219	172	2049449	99.661	ng	0.00
79) Terphenyl-d14	19.942	244	2458965	97.400	ng	0.00

Target Compounds	Qvalue

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

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