

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032322\
 Data File : BM034333.D
 Acq On : 23 Mar 2022 10:58
 Operator : CG/JU
 Sample : PB143395BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB143395BL

Quant Time: Mar 24 02:28:05 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	7.936	152	153115	20.000	ng	-0.01
21) Naphthalene-d8	10.748	136	614472	20.000	ng	-0.02
39) Acenaphthene-d10	14.577	164	407670	20.000	ng	-0.01
64) Phenanthrene-d10	17.318	188	875303	20.000	ng	-0.01
76) Chrysene-d12	21.489	240	889439	20.000	ng	-0.01
86) Perylene-d12	23.900	264	946671	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.495	112	831385	97.409	ng	0.00
7) Phenol-d6	7.101	99	1060211	86.759	ng	0.00
23) Nitrobenzene-d5	9.101	82	1019192	80.754	ng	-0.01
42) 2,4,6-Tribromophenol	16.065	330	450059	81.905	ng	0.00
45) 2-Fluorobiphenyl	13.207	172	2380059	79.778	ng	-0.01
79) Terphenyl-d14	19.936	244	4353913	85.058	ng	-0.01

Target Compounds	Qvalue

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

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