

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM031822\
 Data File : BM034277.D
 Acq On : 18 Mar 2022 10:23
 Operator : CG/JU
 Sample : PB143392BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB143392BL

Quant Time: Mar 18 18:01:53 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	7.943	152	202357	20.000	ng	0.00
21) Naphthalene-d8	10.760	136	813092	20.000	ng	0.00
39) Acenaphthene-d10	14.583	164	526734	20.000	ng	0.00
64) Phenanthrene-d10	17.324	188	1049833	20.000	ng	0.00
76) Chrysene-d12	21.495	240	932401	20.000	ng	0.00
86) Perylene-d12	23.912	264	883861	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.496	112	1662978	147.429	ng	0.00
7) Phenol-d6	7.107	99	2135168	132.207	ng	0.00
23) Nitrobenzene-d5	9.107	82	1399925	83.825	ng	0.00
42) 2,4,6-Tribromophenol	16.071	330	866103	121.991	ng	0.00
45) 2-Fluorobiphenyl	13.213	172	3169478	82.225	ng	0.00
79) Terphenyl-d14	19.942	244	4800743	89.466	ng	0.00

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

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