

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072222\
 Data File : BM035903.D
 Acq On : 22 Jul 2022 12:47
 Operator : CG/JU
 Sample : PB146339BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB146339BL

Quant Time: Jul 22 14:28:54 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM071822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 22 14:25:12 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.887	152	303450	20.000	ng	0.00
21) Naphthalene-d8	10.692	136	1205957	20.000	ng	0.00
39) Acenaphthene-d10	14.521	164	627142	20.000	ng	0.00
64) Phenanthrene-d10	17.262	188	1119627	20.000	ng	0.00
76) Chrysene-d12	21.433	240	983981	20.000	ng	0.00
86) Perylene-d12	23.809	264	1063366	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.451	112	2672820	136.552	ng	0.00
7) Phenol-d6	7.057	99	3165806	128.620	ng	0.00
23) Nitrobenzene-d5	9.057	82	1929984	85.594	ng	0.00
42) 2,4,6-Tribromophenol	16.004	330	873480	124.517	ng	0.00
45) 2-Fluorobiphenyl	13.157	172	3974490	89.581	ng	0.00
79) Terphenyl-d14	19.886	244	4902867	96.946	ng	0.00

Target Compounds	Qvalue

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

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