

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100719\
 Data File : BG043073.D
 Acq On : 7 Oct 2019 18:32
 Operator : JU
 Sample : K5154-21
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MW-102S

Quant Time: Oct 22 05:53:41 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG092519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 25 15:35:52 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.06	152	47610	20.00	ng	-0.02
21) Naphthalene-d8	10.86	136	189366	20.00	ng	-0.02
39) Acenaphthene-d10	14.68	164	153609	20.00	ng	-0.02
64) Phenanthrene-d10	17.42	188	400517	20.00	ng	-0.02
76) Chrysene-d12	21.69	240	434587	20.00	ng	-0.03
87) Perylene-d12	24.91	264	505185	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.64	112	213948	80.20	ng	-0.02
7) Phenol-d6	7.23	99	205649	53.53	ng	-0.02
23) Nitrobenzene-d5	9.22	82	301299	77.34	ng	-0.02
42) 2,4,6-Tribromophenol	16.16	330	365830	138.50	ng	-0.02
45) 2-Fluorobiphenyl	13.30	172	814237	76.84	ng	-0.02
79) Terphenyl-d14	20.03	244	1908492	93.58	ng	-0.01
Target Compounds						
10) Phenol	7.25	94	26582	7.542	ng	94
50) Dimethylphthalate	14.13	163	133820	11.627	ng	96

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

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