

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG100419\
 Data File : BG043038.D
 Acq On : 4 Oct 2019 21:28
 Operator : JU
 Sample : K5154-11
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MW-103I

Quant Time: Oct 05 06:35:24 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	74034	20.00	ng	-0.01
21) Naphthalene-d8	10.86	136	285954	20.00	ng	-0.02
39) Acenaphthene-d10	14.68	164	203015	20.00	ng	-0.01
64) Phenanthrene-d10	17.42	188	480380	20.00	ng	-0.01
76) Chrysene-d12	21.69	240	513886	20.00	ng	-0.03
87) Perylene-d12	24.91	264	601085	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.64	112	303061	73.05	ng	-0.01
7) Phenol-d6	7.23	99	272852	45.67	ng	-0.01
23) Nitrobenzene-d5	9.22	82	425047	72.25	ng	-0.01
42) 2,4,6-Tribromophenol	16.17	330	402179	115.21	ng	-0.01
45) 2-Fluorobiphenyl	13.31	172	1025703	73.24	ng	-0.02
79) Terphenyl-d14	20.03	244	1886053	78.21	ng	-0.01
Target Compounds						
2) 1,4-Dioxane	3.55	88	18989	10.979	ng	Qvalue 87
10) Phenol	7.26	94	37819	6.900	ng	98
50) Dimethylphthalate	14.13	163	277788	18.262	ng	98

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

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