

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG100719\
 Data File : BG043076.D
 Acq On : 7 Oct 2019 20:27
 Operator : JU
 Sample : K5154-24
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MW-101S

Quant Time: Oct 08 02:58:51 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	56297	20.00	ng	-0.02
21) Naphthalene-d8	10.86	136	213482	20.00	ng	-0.02
39) Acenaphthene-d10	14.68	164	166218	20.00	ng	-0.02
64) Phenanthrene-d10	17.42	188	406046	20.00	ng	-0.02
76) Chrysene-d12	21.69	240	417644	20.00	ng	-0.03
87) Perylene-d12	24.91	264	488321	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.64	112	251392	79.69	ng	-0.02
7) Phenol-d6	7.23	99	240007	52.83	ng	-0.02
23) Nitrobenzene-d5	9.22	82	368711	83.95	ng	-0.02
42) 2,4,6-Tribromophenol	16.16	330	423220	148.07	ng	-0.02
45) 2-Fluorobiphenyl	13.30	172	974089	84.96	ng	-0.02
79) Terphenyl-d14	20.03	244	2014391	102.78	ng	-0.02
Target Compounds						
2) 1,4-Dioxane	3.54	88	4919	3.740	ng	Qvalue # 81
10) Phenol	7.26	94	15056	3.612	ng	99
50) Dimethylphthalate	14.13	163	100744	8.089	ng	96

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

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