

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011719\
 Data File : BG039192.D
 Acq On : 17 Jan 2019 18:40
 Operator : JU/SJ
 Sample : K1149-01
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SW001-190114-01

Quant Time: Jan 18 04:45:24 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG010919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 18 04:20:14 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.01	152	23595	20.00	ng	0.00
21) Naphthalene-d8	10.83	136	90193	20.00	ng	0.00
39) Acenaphthene-d10	14.66	164	69558	20.00	ng	0.00
64) Phenanthrene-d10	17.41	188	187591	20.00	ng	0.00
76) Chrysene-d12	21.68	240	210172	20.00	ng	0.00
87) Perylene-d12	24.94	264	205468	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.57	112	92191	74.12	ng	0.00
7) Phenol-d6	7.17	99	75256	42.04	ng	0.00
23) Nitrobenzene-d5	9.19	82	145715	88.40	ng	0.00
42) 2,4,6-Tribromophenol	16.15	330	155108	133.03	ng	0.00
45) 2-Fluorobiphenyl	13.28	172	449810	88.50	ng	0.00
79) Terphenyl-d14	20.01	244	962042	84.68	ng	0.00
Target Compounds						
10) Phenol	7.20	94	11205	6.244	ng	91
17) 2-Methylphenol	8.46	107	4952	3.974	ng	94
20) 3+4-Methylphenols	8.79	107	9819	5.527	ng	90
31) Naphthalene	10.88	128	67802	14.989	ng	99

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

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