

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101819\
 Data File : BG043288.D
 Acq On : 18 Oct 2019 23:13
 Operator : JU
 Sample : K5411-04
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 T1-658

Quant Time: Oct 19 01:25:21 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG092519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 09 01:11:45 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.03	152	48301	20.00	ng	-0.02
21) Naphthalene-d8	10.84	136	178447	20.00	ng	-0.02
39) Acenaphthene-d10	14.66	164	125440	20.00	ng	-0.02
64) Phenanthrene-d10	17.41	188	281851	20.00	ng	-0.01
76) Chrysene-d12	21.68	240	320506	20.00	ng	-0.01
87) Perylene-d12	24.89	264	375083	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.62	112	127651	47.16	ng	-0.02
7) Phenol-d6	7.21	99	114716	29.43	ng	-0.02
23) Nitrobenzene-d5	9.20	82	263788	71.85	ng	-0.02
42) 2,4,6-Tribromophenol	16.15	330	183101	84.89	ng	-0.01
45) 2-Fluorobiphenyl	13.29	172	696631	80.51	ng	-0.02
79) Terphenyl-d14	20.02	244	1266575	84.21	ng	0.00
Target Compounds						
10) Phenol	7.24	94	14818	4.144	ng	93
15) Benzyl Alcohol	8.28	79	9113	3.110	ng	89
20) 3+4-Methylphenols	8.83	107	7777	2.268	ng	83
50) Dimethylphthalate	14.11	163	65615	6.981	ng	99
71) Phenanthrene	17.45	178	36551	2.656	ng	98
74) Di-n-butylphthalate	18.36	149	86864	5.715	ng	100

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

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