

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110419\
 Data File : BG043496.D
 Acq On : 5 Nov 2019 00:53
 Operator : JU
 Sample : K5662-01 5X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 S-1

Quant Time: Nov 05 17:33:10 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG102119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 04 15:22:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.01	152	45341	20.00	ng	0.00
21) Naphthalene-d8	10.82	136	182678	20.00	ng	0.00
39) Acenaphthene-d10	14.64	164	124837	20.00	ng	0.00
64) Phenanthrene-d10	17.39	188	289106	20.00	ng	0.00
76) Chrysene-d12	21.66	240	297044	20.00	ng	0.00
87) Perylene-d12	24.85	264	355496	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	8297	3.35	ng	0.02
7) Phenol-d6	7.20	99	43841	12.31	ng	0.00
23) Nitrobenzene-d5	9.18	82	30138	8.51	ng	0.00
42) 2,4,6-Tribromophenol	16.15	330	1578	0.66	ng	0.02
45) 2-Fluorobiphenyl	13.27	172	87992	9.74	ng	0.00
79) Terphenyl-d14	19.99	244	166802	10.53	ng	0.00
Target Compounds						
32) Benzoic acid	9.96	122	55	9.074	ng	# 1
51) 2,6-Dinitrotoluene	14.64	165	17502	8.332	ng	# 24
54) 2,4-Dinitrophenol	15.18	184	73	7.333	ng	# 17
57) 2,4-Dinitrotoluene	14.64	165	17502	5.830	ng	# 27

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

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