

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110419\
 Data File : BG043494.D
 Acq On : 4 Nov 2019 23:36
 Operator : JU
 Sample : K5602-08
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 20190942-AMENDED-COMPOSITE-

Quant Time: Nov 05 17:32:49 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	38599	20.00	ng	0.00
21) Naphthalene-d8	10.81	136	143878	20.00	ng	0.00
39) Acenaphthene-d10	14.64	164	113334	20.00	ng	0.00
64) Phenanthrene-d10	17.39	188	274854	20.00	ng	0.00
76) Chrysene-d12	21.66	240	299891	20.00	ng	0.00
87) Perylene-d12	24.85	264	365630	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	151605	71.97	ng	0.00
7) Phenol-d6	7.20	99	137385	45.33	ng	0.00
23) Nitrobenzene-d5	9.17	82	211798	75.89	ng	0.00
42) 2,4,6-Tribromophenol	16.13	330	258328	119.78	ng	0.00
45) 2-Fluorobiphenyl	13.27	172	604392	73.69	ng	0.00
79) Terphenyl-d14	20.00	244	1299419	81.29	ng	0.00
Target Compounds						
10) Phenol	7.23	94	12862	4.644	ng	96
19) n-Nitroso-di-n-propylamine	9.17	70	28936	14.776	ng	# 73
32) Benzoic acid	10.17	122	452	9.310	ng	91
50) Dimethylphthalate	14.09	163	89625	10.210	ng	99
57) 2,4-Dinitrotoluene	14.64	165	15227	5.587	ng	# 27
67) 4-Bromophenyl-phenylether	16.13	248	17060	4.612	ng	# 1
77) Benzidine	20.00	184	19370	3.209	ng	# 88

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

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