

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110419\
 Data File : BG043485.D
 Acq On : 4 Nov 2019 17:50
 Operator : JU
 Sample : K5623-02
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PAT-TOTE

Quant Time: Nov 05 01:19:55 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	51035	20.00	ng	0.00
21) Naphthalene-d8	10.82	136	194114	20.00	ng	0.00
39) Acenaphthene-d10	14.64	164	134677	20.00	ng	0.00
64) Phenanthrene-d10	17.39	188	290620	20.00	ng	0.00
76) Chrysene-d12	21.65	240	303275	20.00	ng	0.00
87) Perylene-d12	24.85	264	368617	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	246719	88.59	ng	0.00
7) Phenol-d6	7.20	99	237813	59.35	ng	0.00
23) Nitrobenzene-d5	9.17	82	267190	70.96	ng	0.00
42) 2,4,6-Tribromophenol	16.13	330	256658	100.14	ng	0.00
45) 2-Fluorobiphenyl	13.26	172	701595	71.98	ng	0.00
79) Terphenyl-d14	20.00	244	1214531	75.13	ng	0.00
Target Compounds						
10) Phenol	7.22	94	43764	11.952	ng	94
50) Dimethylphthalate	14.09	163	239195	22.930	ng	97

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

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