

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG112519\
 Data File : BG043771.D
 Acq On : 25 Nov 2019 16:17
 Operator : JU
 Sample : K5668-11
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 HD-01-112119

Quant Time: Nov 26 00:44:44 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG112219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Nov 23 00:54:00 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.02	152	73345	20.00	ng	-0.01
21) Naphthalene-d8	10.83	136	357885	20.00	ng	0.00
39) Acenaphthene-d10	14.65	164	278805	20.00	ng	0.00
64) Phenanthrene-d10	17.40	188	807474	20.00	ng	0.00
76) Chrysene-d12	21.66	240	817972	20.00	ng	-0.01
87) Perylene-d12	24.83	264	752630	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	651117	159.60	ng	0.00
7) Phenol-d6	7.18	99	973154	164.09	ng	0.00
23) Nitrobenzene-d5	9.18	82	551873	84.79	ng	0.00
42) 2,4,6-Tribromophenol	16.13	330	487510	144.97	ng	-0.01
45) 2-Fluorobiphenyl	13.28	172	1354510	79.86	ng	-0.01
79) Terphenyl-d14	20.01	244	2948261	81.45	ng	0.00
Target Compounds						
10) Phenol	7.21	94	22922	3.765	ng	96
50) Dimethylphthalate	14.10	163	345046	15.623	ng	99
77) Benzidine	19.61	184	120810	5.056	ng	98

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

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