

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG030518\
 Data File : BG033009.D
 Acq On : 6 Mar 2018 4:38
 Operator : SJ/JU
 Sample : J1783-04
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SLUDGE

Quant Time: Mar 06 07:08:53 2018
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 01 17:38:52 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.91	152	60194	20.00	ng	0.00
21) Naphthalene-d8	11.80	136	272960	20.00	ng	0.00
38) Acenaphthene-d10	15.53	164	157766	20.00	ng	0.00
63) Phenanthrene-d10	18.25	188	336401	20.00	ng	0.00
75) Chrysene-d12	22.61	240	301613	20.00	ng	0.00
86) Perylene-d12	26.41	264	313802	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	6.33	112	361746	96.15	ng	0.00
7) Phenol-d6	8.01	99	587772	112.08	ng	0.00
23) Nitrobenzene-d5	10.11	82	464473	114.23	ng	0.00
41) 2,4,6-Tribromophenol	17.00	330	234087	141.11	ng	0.00
44) 2-Fluorobiphenyl	14.16	172	1042005	95.26	ng	0.00
78) Terphenyl-d14	20.77	244	1386505	103.28	ng	0.00
Target Compounds						
20) 3+4-Methylphenols	9.67	107	13009	2.400	ng	95
76) Benzidine	20.40	184	53114	5.166	ng	97

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

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