

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_G\DATA\BG010318\
 Data File : BG031906.D
 Acq On : 4 Jan 2018 1:22
 Operator : SJ/JU
 Sample : J1005-03
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 OWS-3

Quant Time: Jan 04 08:21:47 2018
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG122117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 29 15:51:46 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.79	152	56148	20.00	ng	0.00
21) Naphthalene-d8	11.67	136	231318	20.00	ng	0.00
38) Acenaphthene-d10	15.41	164	156154	20.00	ng	-0.01
63) Phenanthrene-d10	18.15	188	388298	20.00	ng	-0.01
75) Chrysene-d12	22.49	240	445601	20.00	ng	-0.02
86) Perylene-d12	26.23	264	464383	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	6.21	112	369475	119.85	ng	0.00
7) Phenol-d6	7.89	99	420974	105.18	ng	0.00
23) Nitrobenzene-d5	9.99	82	303176	98.97	ng	0.00
41) 2,4,6-Tribromophenol	16.89	330	339340	142.42	ng	0.00
44) 2-Fluorobiphenyl	14.05	172	1064551	85.57	ng	0.00
78) Terphenyl-d14	20.68	244	1824975	93.57	ng	0.00
Target Compounds						
10) Phenol	7.92	94	24060	5.954	ng	91
20) 3+4-Methylphenols	9.54	107	74703	18.179	ng	92

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

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