

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_G\DATA\BG010318\
 Data File : BG031907.D
 Acq On : 4 Jan 2018 2:00
 Operator : SJ/JU
 Sample : J1005-04
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 OWS-4

Quant Time: Jan 04 08:24:01 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	58989	20.00	ng	0.00
21) Naphthalene-d8	11.67	136	236694	20.00	ng	-0.01
38) Acenaphthene-d10	15.42	164	161520	20.00	ng	-0.01
63) Phenanthrene-d10	18.15	188	393729	20.00	ng	-0.01
75) Chrysene-d12	22.49	240	438377	20.00	ng	-0.02
86) Perylene-d12	26.23	264	427471	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	6.21	112	424900	131.20	ng	0.00
7) Phenol-d6	7.90	99	498221	118.49	ng	0.00
23) Nitrobenzene-d5	9.99	82	349743	111.58	ng	0.00
41) 2,4,6-Tribromophenol	16.89	330	372386	151.10	ng	0.00
44) 2-Fluorobiphenyl	14.05	172	1228361	95.45	ng	0.00
78) Terphenyl-d14	20.68	244	1967993	102.56	ng	-0.01
Target Compounds						
10) Phenol	7.92	94	16992	4.003	ng	92
20) 3+4-Methylphenols	9.55	107	27235	6.309	ng	85

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

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