

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_G\DATA\BG010918\
 Data File : BG031974.D
 Acq On : 10 Jan 2018 9:33
 Operator : SJ/JU
 Sample : J1057-04
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 OWS-2C

Manual Integrations
 APPROVED

Sohil
 1/10/2018 11:48:23 AM

Quant Time: Jan 10 10:48:44 2018
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG122117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 05 15:31:34 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.78	152	62761	20.00	ng	0.00
21) Naphthalene-d8	11.66	136	259359	20.00	ng	-0.01
38) Acenaphthene-d10	15.41	164	181862	20.00	ng	-0.02
63) Phenanthrene-d10	18.14	188	433674	20.00	ng	-0.02
75) Chrysene-d12	22.48	240	500493	20.00	ng	-0.04
86) Perylene-d12	26.20	264	507217	20.00	ng	-0.10
System Monitoring Compounds						
5) 2-Fluorophenol	6.21	112	429231	124.57	ng	0.00
7) Phenol-d6	7.89	99	511738	114.39	ng	0.00
23) Nitrobenzene-d5	9.98	82	366408	106.68	ng	0.00
41) 2,4,6-Tribromophenol	16.88	330	438443	158.00	ng	-0.02
44) 2-Fluorobiphenyl	14.04	172	1312348	90.57	ng	-0.01
78) Terphenyl-d14	20.67	244	2106005	96.13	ng	-0.03
Target Compounds						
10) Phenol	7.92	94	14579	3.228	ng	94
20) 3+4-Methylphenols	9.54	107	128701	28.020	ng	89
31) Naphthalene	11.71	128	50825	3.883	ng	99
32) Benzoic acid	10.84	122	8333m	9.387	ng	

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

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