

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG062119\
 Data File : BG041387.D
 Acq On : 21 Jun 2019 21:51
 Operator : HP/JU
 Sample : K3433-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 OILY-WATER

Quant Time: Jun 22 01:40:38 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG061719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 17 14:58:02 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.06	152	39105	20.00	ng	-0.03
21) Naphthalene-d8	10.88	136	146629	20.00	ng	-0.03
39) Acenaphthene-d10	14.70	164	97932	20.00	ng	-0.03
64) Phenanthrene-d10	17.45	188	242822	20.00	ng	-0.01
76) Chrysene-d12	21.73	240	246192	20.00	ng	0.00
87) Perylene-d12	25.02	264	244742	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.68	112	97258	45.70	ng	0.06
7) Phenol-d6	7.22	99	84676	27.45	ng	-0.02
23) Nitrobenzene-d5	9.24	82	257876	73.96	ng	-0.03
42) 2,4,6-Tribromophenol	16.19	330	162578	108.13	ng	-0.01
45) 2-Fluorobiphenyl	13.31	172	610662	84.75	ng	-0.03
79) Terphenyl-d14	20.06	244	1033355	83.18	ng	0.00
Target Compounds						
10) Phenol	7.25	94	7277	2.194	ng	93
50) Dimethylphthalate	14.14	163	48678	5.586	ng	99
84) Bis(2-ethylhexyl)phthalate	21.57	149	18345	2.143	ng	# 77

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

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