

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010419\
 Data File : BG038927.D
 Acq On : 4 Jan 2019 12:39
 Operator : JU/SJ
 Sample : K1019-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TEST-PIT-D-11-12-A

Manual Integrations
 APPROVED

Sohil
 1/7/2019 12:48:03 PM

Quant Time: Jan 04 14:09:26 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 12 15:26:27 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.05	152	29022	20.00	ng	-0.01
21) Naphthalene-d8	10.87	136	118249	20.00	ng	-0.01
39) Acenaphthene-d10	14.69	164	89260	20.00	ng	-0.01
64) Phenanthrene-d10	17.43	188	224424	20.00	ng	0.00
76) Chrysene-d12	21.71	240	223948	20.00	ng	-0.01
87) Perylene-d12	25.01	264	238606	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	187957	114.87	ng	0.00
7) Phenol-d6	7.21	99	265304	118.11	ng	0.00
23) Nitrobenzene-d5	9.23	82	159365	68.85	ng	-0.01
42) 2,4,6-Tribromophenol	16.18	330	151884	123.47	ng	-0.01
45) 2-Fluorobiphenyl	13.30	172	446947	68.69	ng	-0.01
79) Terphenyl-d14	20.04	244	773349	75.15	ng	0.00
Target Compounds						
10) Phenol	7.23	94	11907	4.989	ng	94
50) Dimethylphthalate	14.13	163	64447	8.841	ng	99
71) Phenanthrene	17.48	178	59375	5.102	ng	98
75) Fluoranthene	19.49	202	103515	7.189	ng	97
78) Pyrene	19.85	202	94169	6.365	ng	99
81) Benzo(a)anthracene	21.69	228	43279	3.171	ng	99
83) Chrysene	21.76	228	36784	2.799	ng	97
88) Benzo(b)fluoranthene	23.96	252	42844	3.270	ng	97
90) Benzo(a)pyrene	24.85	252	33184	2.626	ng	# 95

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

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