

Data Path : Z:\HPCHEM1\BNA G\DATA\BG012318\
 Data File : BG032227.D
 Acq On : 23 Jan 2018 12:13
 Operator : SJ/JU
 Sample : J1239-07
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 WC-12R12-02

Manual Integrations
 APPROVED

Sohil
 1/24/2018 2:49:07 PM

Quant Time: Jan 24 09:49:23 2018
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG011218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 18 10:13:16 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.73	152	45709	20.00	ng	-0.01
21) Naphthalene-d8	11.61	136	187501	20.00	ng	-0.01
38) Acenaphthene-d10	15.36	164	131186	20.00	ng	-0.02
63) Phenanthrene-d10	18.10	188	322364	20.00	ng	-0.02
75) Chrysene-d12	22.43	240	418967	20.00	ng	-0.02
86) Perylene-d12	26.12	264	452111	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	6.16	112	315999	126.44	ng	0.00
7) Phenol-d6	7.84	99	368867	117.19	ng	0.00
23) Nitrobenzene-d5	9.93	82	281466	101.56	ng	-0.01
41) 2,4,6-Tribromophenol	16.84	330	315252	145.22	ng	-0.01
44) 2-Fluorobiphenyl	13.99	172	976798	92.33	ng	-0.01
78) Terphenyl-d14	20.63	244	1765865	93.06	ng	-0.02
Target Compounds						
3) Pyridine	4.29	79	23819	9.294	ng	Qvalue # 83
17) 2-Methylphenol	9.17	107	6476	2.733	ng	98
20) 3+4-Methylphenols	9.51	107	11975m	3.648	ng	
22) Acetophenone	9.57	105	12646m	2.969	ng	
27) 2,4-Dimethylphenol	10.74	122	10540m	4.055	ng	
31) Naphthalene	11.67	128	2070169	222.878	ng	96
37) 2-Methylnaphthalene	13.22	142	192311	26.079	ng	93
45) 1,1'-Biphenyl	14.20	154	27556	2.450	ng	96
48) Acenaphthylene	15.09	152	98433	7.388	ng	97
51) Acenaphthene	15.43	154	45145	5.125	ng	98
57) Fluorene	16.40	166	50705	4.704	ng	95
70) Phenanthrene	18.14	178	108913	6.068	ng	98
72) Carbazole	18.49	167	92700	5.969	ng	96

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

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