

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061219\
 Data File : BG041222.D
 Acq On : 13 Jun 2019 1:12
 Operator : HP/JU
 Sample : K3303-04
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BAT-B02

Quant Time: Jun 13 08:26:49 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG052919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Jun 09 23:08:34 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.11	152	43274	20.00	ng	0.00
21) Naphthalene-d8	10.93	136	177437	20.00	ng	0.00
39) Acenaphthene-d10	14.73	164	133425	20.00	ng	0.00
64) Phenanthrene-d10	17.48	188	306454	20.00	ng	0.00
76) Chrysene-d12	21.75	240	296699	20.00	ng	0.00
87) Perylene-d12	25.09	264	316246	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.64	112	231607	96.51	ng	0.00
7) Phenol-d6	7.25	99	359429	96.98	ng	0.00
23) Nitrobenzene-d5	9.28	82	251247	62.86	ng	0.00
42) 2,4,6-Tribromophenol	16.23	330	184479	93.13	ng	0.00
45) 2-Fluorobiphenyl	13.36	172	557828	60.21	ng	0.00
79) Terphenyl-d14	20.07	244	850504	60.84	ng	0.00
Target Compounds						
50) Dimethylphthalate	14.18	163	119179	10.552	ng	Qvalue 100
75) Fluoranthene	19.52	202	95015	4.819	ng	95
78) Pyrene	19.89	202	115478	5.987	ng	95
81) Benzo(a)anthracene	21.74	228	68859	3.487	ng	96
83) Chrysene	21.80	228	60730	3.213	ng	97
88) Benzo(b)fluoranthene	24.02	252	95113	4.959	ng	99
90) Benzo(a)pyrene	24.92	252	67414	3.684	ng	96

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

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