

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG051119\
 Data File : BG040909.D
 Acq On : 12 May 2019 14:47
 Operator : HP/JU
 Sample : K2766-07
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 P026-SS007-0612-01

Quant Time: May 13 05:16:29 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG042319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 24 05:35:33 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.13	152	54344	20.00	ng	-0.03
21) Naphthalene-d8	10.96	136	226395	20.00	ng	-0.03
39) Acenaphthene-d10	14.76	164	170298	20.00	ng	-0.03
64) Phenanthrene-d10	17.51	188	426518	20.00	ng	-0.03
76) Chrysene-d12	21.79	240	431994	20.00	ng	-0.04
87) Perylene-d12	25.14	264	470419	20.00	ng	-0.06
System Monitoring Compounds						
5) 2-Fluorophenol	5.67	112	407544	132.20	ng	-0.02
7) Phenol-d6	7.28	99	621270	130.88	ng	-0.02
23) Nitrobenzene-d5	9.31	82	424245	86.59	ng	-0.03
42) 2,4,6-Tribromophenol	16.25	330	326193	139.35	ng	-0.02
45) 2-Fluorobiphenyl	13.39	172	852094	72.26	ng	-0.03
79) Terphenyl-d14	20.10	244	1332589	68.34	ng	-0.03
Target Compounds						
50) Dimethylphthalate	14.20	163	123031	8.635	ng	Qvalue 99
71) Phenanthrene	17.55	178	82980	3.612	ng	96
75) Fluoranthene	19.55	202	162322	5.670	ng	95
78) Pyrene	19.92	202	297163	10.975	ng	97
81) Benzo(a)anthracene	21.77	228	115063	4.214	ng	97
83) Chrysene	21.83	228	136986	5.276	ng	98
86) Indeno(1,2,3-cd)pyrene	28.96	276	74137	2.362	ng	# 82
88) Benzo(b)fluoranthene	24.06	252	139218	4.843	ng	97
90) Benzo(a)pyrene	24.97	252	99210	3.646	ng	97
92) Benzo(g,h,i)perylene	30.19	276	90013	3.284	ng	93

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

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