

Data Path : Z:\HPCHEM1\BNA G\DATA\BG061616\
 Data File : BG022390.D
 Acq On : 17 Jun 2016 7:00
 Operator : UM/SJ
 Sample : H3569-13
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-05-COMP

Quant Time: Jun 17 07:54:44 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG060616.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 13 17:40:23 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.05	152	33877	20.00	ng	-0.04
21) Naphthalene-d8	10.88	136	164648	20.00	ng	-0.03
38) Acenaphthene-d10	14.69	164	96206	20.00	ng	-0.03
63) Phenanthrene-d10	17.44	188	185062	20.00	ng	-0.03
75) Chrysene-d12	21.71	240	154772	20.00	ng	-0.03
86) Perylene-d12	25.00	264	142081	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.62	112	207994	118.71	ng	-0.02
7) Phenol-d6	7.24	99	298137	108.22	ng	0.00
23) Nitrobenzene-d5	9.23	82	168972	71.55	ng	-0.03
41) 2,4,6-Tribromophenol	16.19	330	84544	77.77	ng	-0.03
44) 2-Fluorobiphenyl	13.31	172	422466	66.92	ng	-0.03
78) Terphenyl-d14	20.03	244	432557	66.35	ng	-0.02
Target Compounds						Qvalue
49) Dimethylphthalate	14.14	163	100345	12.58	ng	99
74) Fluoranthene	19.49	202	66206	5.73	ng	96
77) Pyrene	19.85	202	51519	5.35	ng	99
80) Benzo(a)anthracene	21.69	228	25712	2.96	ng	97
82) Chrysene	21.75	228	25394	3.01	ng	97
87) Benzo(b)fluoranthene	23.95	252	32161m	3.88	ng	
88) Benzo(k)fluoranthene	23.99	252	16800m	2.06	ng	
89) Benzo(a)pyrene	24.84	252	20327	2.57	ng	# 93

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

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