

Data Path : Z:\HPCHEM1\BNA G\DATA\BG061616\
 Data File : BG022387.D
 Acq On : 17 Jun 2016 5:00
 Operator : UM/SJ
 Sample : H3569-11
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-03-COMP

Quant Time: Jun 17 07:27:13 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	36607	20.00	ng	-0.04
21) Naphthalene-d8	10.88	136	185846	20.00	ng	-0.03
38) Acenaphthene-d10	14.69	164	107786	20.00	ng	-0.03
63) Phenanthrene-d10	17.44	188	209703	20.00	ng	-0.03
75) Chrysene-d12	21.71	240	176964	20.00	ng	-0.03
86) Perylene-d12	24.99	264	165692	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.62	112	212940	112.47	ng	-0.02
7) Phenol-d6	7.24	99	314314	105.59	ng	0.00
23) Nitrobenzene-d5	9.23	82	175507	65.84	ng	-0.03
41) 2,4,6-Tribromophenol	16.19	330	90119	73.99	ng	-0.03
44) 2-Fluorobiphenyl	13.31	172	452807	64.02	ng	-0.03
78) Terphenyl-d14	20.04	244	457176	61.33	ng	-0.02
Target Compounds						Qvalue
49) Dimethylphthalate	14.14	163	98506	11.02	ng	99
70) Phenanthrene	17.48	178	113600	9.97	ng	99
74) Fluoranthene	19.48	202	134684	10.29	ng	96
77) Pyrene	19.85	202	129711	11.79	ng	99
80) Benzo(a)anthracene	21.69	228	58199	5.86	ng	99
82) Chrysene	21.75	228	55755	5.77	ng	98
85) Indeno(1,2,3-cd)pyrene	28.75	276	24828m	2.29	ng	
87) Benzo(b)fluoranthene	23.94	252	55187m	5.71	ng	
88) Benzo(k)fluoranthene	23.99	252	19928m	2.09	ng	
89) Benzo(a)pyrene	24.84	252	35969	3.89	ng	98
91) Benzo(g,h,i)perylene	29.94	276	23742m	2.62	ng	

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

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