

Data Path : Z:\HPCHEM1\BNA G\DATA\BG061716\
 Data File : BG022443.D
 Acq On : 19 Jun 2016 3:58
 Operator : UM/SJ
 Sample : H3489-05
 Misc :
 ALS Vial : 52 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 S6-B2(0-10)C

Quant Time: Jun 20 04:20:03 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	41808	20.00	ng	-0.05
21) Naphthalene-d8	10.87	136	189967	20.00	ng	-0.04
38) Acenaphthene-d10	14.68	164	93089	20.00	ng	-0.04
63) Phenanthrene-d10	17.43	188	164079	20.00	ng	-0.04
75) Chrysene-d12	21.70	240	129419	20.00	ng	-0.04
86) Perylene-d12	24.96	264	96210	20.00	ng	-0.06
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	141453	65.42	ng	-0.03
7) Phenol-d6	7.23	99	452091	132.98	ng	-0.02
23) Nitrobenzene-d5	9.23	82	266831	97.93	ng	-0.03
41) 2,4,6-Tribromophenol	16.18	330	24817	23.59	ng	-0.03
44) 2-Fluorobiphenyl	13.31	172	569869	93.29	ng	-0.03
78) Terphenyl-d14	20.03	244	443229	81.31	ng	-0.03
Target Compounds						Qvalue
37) 2-Methylnaphthalene	12.52	142	180714	25.82	ng	95
49) Dimethylphthalate	14.13	163	161043	20.86	ng	100
70) Phenanthrene	17.48	178	39457	4.43	ng	98
74) Fluoranthene	19.48	202	51240	5.00	ng	94
77) Pyrene	19.84	202	46374	5.76	ng	98
80) Benzo(a)anthracene	21.68	228	20509	2.83	ng	# 89
82) Chrysene	21.75	228	20151	2.85	ng	# 90
87) Benzo(b)fluoranthene	23.92	252	19221	3.43	ng	# 69
89) Benzo(a)pyrene	24.80	252	10951	2.04	ng	# 51

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

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