

Data Path : Z:\HPCHEM1\BNA G\DATA\BG042715\
 Data File : BG016748.D
 Acq On : 28 Apr 2015 2:16
 Operator : TP/IZ
 Sample : G1942-02
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-2

Quant Time: Apr 28 03:30:32 2015
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG042315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 28 01:49:03 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	70717	20.00	ng	0.00
21) Naphthalene-d8	10.36	136	304336	20.00	ng	0.00
38) Acenaphthene-d10	14.23	164	180746	20.00	ng	0.00
63) Phenanthrene-d10	16.97	188	374840	20.00	ng	0.00
75) Chrysene-d12	21.16	240	339937	20.00	ng	0.00
86) Perylene-d12	23.36	264	322014	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.17	112	469305	107.69	ng	0.00
7) Phenol-d6	6.78	99	644599	105.44	ng	0.00
23) Nitrobenzene-d5	8.73	82	317852	65.99	ng	0.00
41) 2,4,6-Tribromophenol	15.72	330	147224	109.80	ng	0.00
44) 2-Fluorobiphenyl	12.84	172	703033	62.49	ng	0.00
78) Terphenyl-d14	19.61	244	820303	55.59	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	13.69	163	35593	2.63	ng	98
70) Phenanthrene	17.01	178	52832	2.45	ng	98
74) Fluoranthene	19.03	202	147012	6.44	ng	96
77) Pyrene	19.40	202	129750	5.43	ng	97
80) Benzo(a)anthracene	21.14	228	63476m	3.08	ng	
82) Chrysene	21.20	228	62922m	3.18	ng	
87) Benzo(b)fluoranthene	22.69	252	85309m	4.36	ng	
89) Benzo(a)pyrene	23.26	252	48269	2.62	ng	94

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

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