

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG071816\
 Data File : BG023170.D
 Acq On : 18 Jul 2016 22:37
 Operator : UM/SJ
 Sample : H4044-07
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 F9-B2(6-11)C

Quant Time: Jul 19 13:56:25 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG070816.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 19 01:16:03 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.11	152	151454	20.00	ng	0.00
21) Naphthalene-d8	10.94	136	671522	20.00	ng	0.00
38) Acenaphthene-d10	14.75	164	434471	20.00	ng	0.00
63) Phenanthrene-d10	17.49	188	911535	20.00	ng	0.00
75) Chrysene-d12	21.75	240	894327	20.00	ng	0.00
86) Perylene-d12	25.04	264	873825	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.67	112	1088295	117.52	ng	0.00
7) Phenol-d6	7.28	99	1642491	113.24	ng	0.00
23) Nitrobenzene-d5	9.29	82	1237318	78.90	ng	0.00
41) 2,4,6-Tribromophenol	16.23	330	626639	80.25	ng	0.00
44) 2-Fluorobiphenyl	13.37	172	1983515	71.91	ng	0.00
78) Terphenyl-d14	20.08	244	2185345	73.64	ng	0.00
Target Compounds						
49) Dimethylphthalate	14.19	163	514271	14.49	ng	Qvalue 98
74) Fluoranthene	19.53	202	166440	3.04	ng	97
77) Pyrene	19.89	202	169107m	3.36	ng	
80) Benzo(a)anthracene	21.73	228	104226m	2.08	ng	
82) Chrysene	21.80	228	94769m	2.02	ng	
87) Benzo(b)fluoranthene	24.00	252	114275	2.27	ng	# 93

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

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