

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG071816\
 Data File : BG023158.D
 Acq On : 18 Jul 2016 13:15
 Operator : UM/SJ
 Sample : H4018-12
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 F9-B4(0-12)C

Quant Time: Jul 19 01:20:30 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	8.10	152	134977	20.00	ng	-0.02
21) Naphthalene-d8	10.93	136	627171	20.00	ng	0.00
38) Acenaphthene-d10	14.74	164	478874	20.00	ng	0.00
63) Phenanthrene-d10	17.48	188	1078580	20.00	ng	0.00
75) Chrysene-d12	21.74	240	1073880	20.00	ng	0.00
86) Perylene-d12	25.02	264	1064778	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.65	112	803744	97.39	ng	-0.03
7) Phenol-d6	7.26	99	1316000	101.81	ng	-0.02
23) Nitrobenzene-d5	9.28	82	995841	67.99	ng	-0.01
41) 2,4,6-Tribromophenol	16.23	330	649854	75.51	ng	0.00
44) 2-Fluorobiphenyl	13.37	172	1972008	64.86	ng	0.00
78) Terphenyl-d14	20.07	244	2684314	76.22	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	14.18	163	568581	14.54	ng	98
74) Fluoranthene	19.52	202	191214	2.95	ng	94
77) Pyrene	19.88	202	178756	2.96	ng	92
80) Benzo(a)anthracene	21.72	228	156505	2.60	ng	92
82) Chrysene	21.79	228	146187m	2.59	ng	
87) Benzo(b)fluoranthene	23.98	252	252899	4.12	ng	# 90
89) Benzo(a)pyrene	24.86	252	163296	2.77	ng	# 97

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

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