

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG071616\
 Data File : BG023151.D
 Acq On : 16 Jul 2016 17:43
 Operator : UM/SJ
 Sample : H3951-03
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 LSS-SB-SB01(7-9ft)

Quant Time: Jul 18 02:06:47 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG070816.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 12 18:30:58 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.15	152	119117	20.00	ng	0.00
21) Naphthalene-d8	10.98	136	562015	20.00	ng	0.00
38) Acenaphthene-d10	14.80	164	388981	20.00	ng	0.00
63) Phenanthrene-d10	17.55	188	862558	20.00	ng	0.00
75) Chrysene-d12	21.86	240	889224	20.00	ng	0.00
86) Perylene-d12	25.23	264	909766	20.00	ng	0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.70	112	918008	126.04	ng	0.00
7) Phenol-d6	7.31	99	1468530	128.74	ng	0.00
23) Nitrobenzene-d5	9.32	82	1096301	83.53	ng	0.00
41) 2,4,6-Tribromophenol	16.29	330	690486	98.77	ng	0.00
44) 2-Fluorobiphenyl	13.42	172	2122947	85.97	ng	0.00
78) Terphenyl-d14	20.15	244	2601342	98.61	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	14.24	163	1101728	34.67	ng	99
74) Fluoranthene	19.60	202	167256	3.22	ng	93
77) Pyrene	19.96	202	160408	3.21	ng	# 91
80) Benzo(a)anthracene	21.83	228	108875m	2.19	ng	
82) Chrysene	21.90	228	108873	2.33	ng	93
87) Benzo(b)fluoranthene	24.15	252	139163	2.65	ng	# 97

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

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