

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG071616\
 Data File : BG023153.D
 Acq On : 16 Jul 2016 19:04
 Operator : UM/SJ
 Sample : H3951-09
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 LSS-SB-SB03(5-7ft)

Quant Time: Jul 18 14:55:22 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	133351	20.00	ng	0.00
21) Naphthalene-d8	10.98	136	603172	20.00	ng	0.00
38) Acenaphthene-d10	14.80	164	403763	20.00	ng	0.00
63) Phenanthrene-d10	17.55	188	854377	20.00	ng	0.00
75) Chrysene-d12	21.85	240	879399	20.00	ng	0.00
86) Perylene-d12	25.23	264	892453	20.00	ng	0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.70	112	1024904	125.70	ng	0.00
7) Phenol-d6	7.31	99	1587923	124.34	ng	0.00
23) Nitrobenzene-d5	9.33	82	1166941	82.84	ng	0.00
41) 2,4,6-Tribromophenol	16.29	330	695933	95.90	ng	0.00
44) 2-Fluorobiphenyl	13.42	172	2092554	81.63	ng	0.00
78) Terphenyl-d14	20.15	244	2434840	89.67	ng	0.00
Target Compounds						
10) Phenol	7.33	94	27923	2.13	ng	Qvalue # 25
49) Dimethylphthalate	14.24	163	659617	20.00	ng	98
70) Phenanthrene	17.59	178	218484	5.22	ng	97
74) Fluoranthene	19.60	202	727685	14.16	ng	93
77) Pyrene	19.96	202	707650	14.32	ng	93
80) Benzo(a)anthracene	21.84	228	484476	9.85	ng	95
82) Chrysene	21.91	228	433360	9.39	ng	98
85) Indeno(1,2,3-cd)pyrene	29.07	276	250433	4.26	ng	# 100
87) Benzo(b)fluoranthene	24.16	252	628991	12.22	ng	# 97
88) Benzo(k)fluoranthene	24.21	252	199315m	4.08	ng	
89) Benzo(a)pyrene	25.07	252	480576	9.74	ng	# 99
91) Benzo(g,h,i)perylene	30.29	276	222680m	4.54	ng	

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

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