

Data Path : Z:\HPCHEM1\BNA G\DATA\BG071316\
 Data File : BG023091.D
 Acq On : 14 Jul 2016 6:33
 Operator : UM/SJ
 Sample : H4014-01
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SYSDIS-071216

Quant Time: Jul 14 07:36:27 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	102348	20.00	ng	0.00
21) Naphthalene-d8	10.98	136	468412	20.00	ng	0.00
38) Acenaphthene-d10	14.80	164	352709	20.00	ng	0.00
63) Phenanthrene-d10	17.56	188	875585	20.00	ng	0.00
75) Chrysene-d12	21.86	240	983969	20.00	ng	0.01
86) Perylene-d12	25.24	264	987861	20.00	ng	0.06
System Monitoring Compounds						
5) 2-Fluorophenol	5.70	112	30942m	4.94	ng	0.00
7) Phenol-d6	7.35	99	51580	5.26	ng	0.04
23) Nitrobenzene-d5	9.33	82	671942	61.42	ng	0.00
41) 2,4,6-Tribromophenol	16.30	330	34011	5.37	ng	0.00
44) 2-Fluorobiphenyl	13.42	172	1395934	62.34	ng	0.00
78) Terphenyl-d14	20.16	244	2367688	71.97	ng	0.00
Target Compounds						
3) Pyridine	3.99	79	23609	2.84	ng	# 81
6) Aniline	7.53	93	10433408	768.11	ng	# 25
8) 2-Chlorophenol	7.73	128	21884	3.29	ng	95
12) 1,3-Dichlorobenzene	8.05	146	85424	10.48	ng	95
13) 1,4-Dichlorobenzene	8.20	146	1270311	155.66	ng	97
14) 1,2-Dichlorobenzene	8.52	146	1482620	182.74	ng	98
20) 3+4-Methylphenols	8.96	107	88131	9.63	ng	# 50
33) 4-Chloroaniline	11.14	127	125857	10.79	ng	97
49) Dimethylphthalate	14.24	163	148652	5.16	ng	99
76) Benzidine	19.78	184	114248	4.05	ng	98

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

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