

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG071616\
 Data File : BG023147.D
 Acq On : 16 Jul 2016 15:01
 Operator : UM/SJ
 Sample : H4018-08
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C8-B4(0-5)C

Quant Time: Jul 18 01:56:39 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	125122	20.00	ng	0.00
21) Naphthalene-d8	10.98	136	573976	20.00	ng	0.00
38) Acenaphthene-d10	14.80	164	412571	20.00	ng	0.00
63) Phenanthrene-d10	17.55	188	928631	20.00	ng	-0.01
75) Chrysene-d12	21.85	240	982609	20.00	ng	0.00
86) Perylene-d12	25.22	264	1008186	20.00	ng	0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.70	112	598808	78.27	ng	0.00
7) Phenol-d6	7.31	99	961082	80.21	ng	0.00
23) Nitrobenzene-d5	9.32	82	720014	53.71	ng	0.00
41) 2,4,6-Tribromophenol	16.29	330	436901	58.92	ng	0.00
44) 2-Fluorobiphenyl	13.42	172	1434559	54.77	ng	0.00
78) Terphenyl-d14	20.14	244	1958113	54.97	ng	-0.01
Target Compounds						Qvalue
49) Dimethylphthalate	14.24	163	378886	11.24	ng	98
70) Phenanthrene	17.59	178	220401	4.84	ng	95
74) Fluoranthene	19.60	202	866082	15.51	ng	95
77) Pyrene	19.96	202	762943	13.82	ng	92
80) Benzo(a)anthracene	21.83	228	585173	10.64	ng	93
82) Chrysene	21.90	228	570169	11.06	ng	97
85) Indeno(1,2,3-cd)pyrene	29.06	276	281058	4.27	ng	# 100
87) Benzo(b)fluoranthene	24.14	252	728674	12.53	ng	# 98
88) Benzo(k)fluoranthene	24.20	252	229356m	4.16	ng	
89) Benzo(a)pyrene	25.06	252	485605	8.71	ng	# 95
91) Benzo(g,h,i)perylene	30.27	276	282856	5.10	ng	96

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

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