

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG071816\
 Data File : BG023168.D
 Acq On : 18 Jul 2016 21:16
 Operator : UM/SJ
 Sample : H4044-13
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 E9-B2(0-6)C

Quant Time: Jul 19 13:43:42 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	8.11	152	145536	20.00	ng	-0.01
21) Naphthalene-d8	10.93	136	645644	20.00	ng	0.00
38) Acenaphthene-d10	14.75	164	425181	20.00	ng	0.00
63) Phenanthrene-d10	17.49	188	886289	20.00	ng	0.00
75) Chrysene-d12	21.75	240	872549	20.00	ng	0.00
86) Perylene-d12	25.04	264	866187	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.66	112	778457	87.48	ng	-0.01
7) Phenol-d6	7.27	99	1182707	84.86	ng	-0.01
23) Nitrobenzene-d5	9.28	82	982212	65.14	ng	-0.01
41) 2,4,6-Tribromophenol	16.23	330	441664	57.80	ng	0.00
44) 2-Fluorobiphenyl	13.37	172	1814212	67.21	ng	0.00
78) Terphenyl-d14	20.07	244	2101068	72.04	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	14.19	163	342542	9.86	ng	99
70) Phenanthrene	17.53	178	202063	4.65	ng	95
74) Fluoranthene	19.52	202	426332	8.00	ng	94
77) Pyrene	19.88	202	398175	8.12	ng	93
80) Benzo(a)anthracene	21.73	228	230282	4.72	ng	94
82) Chrysene	21.79	228	235672	5.15	ng	100
83) Bis(2-ethylhexyl)phthalate	21.60	149	92901m	3.44	ng	
85) Indeno(1,2,3-cd)pyrene	28.78	276	120438	2.06	ng	# 100
87) Benzo(b)fluoranthene	23.99	252	370132	7.41	ng	# 94
89) Benzo(a)pyrene	24.88	252	214385	4.48	ng	# 96
91) Benzo(g,h,i)perylene	29.95	276	106353	2.23	ng	# 74

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

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