

Data Path : Z:\HPCHEM1\BNA G\DATA\BG092116\
 Data File : BG024058.D
 Acq On : 21 Sep 2016 17:04
 Operator : UM/SJ
 Sample : H4820-11
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 YORK-SB4-02

Quant Time: Sep 21 18:05:25 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG083116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 20 19:01:00 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.05	152	70367	20.00	ng	0.00
21) Naphthalene-d8	10.88	136	303397	20.00	ng	0.00
38) Acenaphthene-d10	14.72	164	222630	20.00	ng	0.00
63) Phenanthrene-d10	17.49	188	411526	20.00	ng	0.00
75) Chrysene-d12	21.80	240	440501	20.00	ng	0.00
86) Perylene-d12	25.16	264	434039	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	473307	118.09	ng	0.00
7) Phenol-d6	7.22	99	746457	120.99	ng	0.00
23) Nitrobenzene-d5	9.23	82	589511	86.74	ng	0.00
41) 2,4,6-Tribromophenol	16.23	330	311243	77.60	ng	0.00
44) 2-Fluorobiphenyl	13.34	172	1077126	69.37	ng	0.00
78) Terphenyl-d14	20.10	244	1345426	69.56	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	14.17	163	315918	16.55	ng	99
70) Phenanthrene	17.53	178	284524	13.29	ng	97
71) Anthracene	17.62	178	58707	2.69	ng	96
74) Fluoranthene	19.55	202	441697	17.14	ng	94
77) Pyrene	19.92	202	446530	16.48	ng	# 91
80) Benzo(a)anthracene	21.78	228	258141	10.47	ng	98
82) Chrysene	21.85	228	256816	10.94	ng	98
83) Bis(2-ethylhexyl)phthalate	21.65	149	30848	2.16	ng	99
85) Indeno(1,2,3-cd)pyrene	28.99	276	121872	4.69	ng	# 100
87) Benzo(b)fluoranthene	24.09	252	294275	11.94	ng	# 97
88) Benzo(k)fluoranthene	24.14	252	115340m	4.72	ng	
89) Benzo(a)pyrene	24.99	252	233873	9.99	ng	# 97
91) Benzo(a,h,i)perylene	30.18	276	91602	4.14	ng	# 91

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

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