

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG060119\
 Data File : BG041010.D
 Acq On : 1 Jun 2019 13:37
 Operator : HP/JU
 Sample : K3040-02
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 0445-FIELD-BLANK-2

Quant Time: Jun 03 01:12:01 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG052919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 29 18:39:57 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	0.00	152	0	0.00	ng	-8.13
21) Naphthalene-d8	10.98	136	54	20.00	ng	0.02
39) Acenaphthene-d10	14.81	164	56	20.00	ng	0.05
64) Phenanthrene-d10	0.00	188	0	0.00	ng	-17.50
76) Chrysene-d12	21.58	240	54	20.00	ng	-0.21
87) Perylene-d12	24.71	264	59	20.00	ng	-0.42

System Monitoring Compounds						
5) 2-Fluorophenol	5.22	112	59	0.00	ng	-0.44
7) Phenol-d6	7.33	99	58	0.00	ng	0.07
23) Nitrobenzene-d5	9.52	82	63	51.79	ng	0.22
42) 2,4,6-Tribromophenol	0.00	330	0	0.00	ng	
45) 2-Fluorobiphenyl	13.74	172	59	15.17	ng	0.36
79) Terphenyl-d14	20.11	244	57	22.40	ng	0.01

Target Compounds				Qvalue		
22) Acetophenone	9.05	105	54	35.648	ng	# 47
24) Nitrobenzene	9.42	77	62	50.016	ng	# 44
25) Isophorone	10.18	82	61	26.351	ng	# 67
26) 2-Nitrophenol	10.49	139	59	115.453	ng	# 27
28) bis(2-Chloroethoxy)methane	10.59	93	60	48.023	ng	# 49
29) 2,4-Dichlorophenol	10.69	162	77	71.844	ng	# 28
30) 1,2,4-Trichlorobenzene	10.65	180	86	70.536	ng	# 10
31) Naphthalene	10.73	128	63	21.729	ng	# 68
33) 4-Chloroaniline	10.88	127	73	54.265	ng	# 41
35) Caprolactam	12.02	113	62	175.178	ng	# 1
36) 4-Chloro-3-methylphenol	11.88	107	69	56.371	ng	# 23
38) 1-Methylnaphthalene	13.29	142	75	35.642	ng	# 6
48) 2-Nitroaniline	14.11	65	55	43.082	ng	# 11
49) Acenaphthylene	14.30	152	61	11.389	ng	# 59
50) Dimethylphthalate	14.00	163	98	20.673	ng	# 76
54) 2,4-Dinitrophenol	14.46	184	74	116.613	ng	# 18
60) Diethylphthalate	15.18	149	94	19.431	ng	# 57
61) 4-Chlorophenyl-phenylether	16.24	204	61	21.585	ng	# 30
63) Azobenzene	15.98	77	59	13.418	ng	# 47
77) Benzidine	19.07	184	53	41.143	ng	# 64
78) Pyrene	19.35	202	57	16.238	ng	# 55
80) Butylbenzylphthalate	20.59	149	55	40.261	ng	# 24
81) Benzo(a)anthracene	21.58	228	63	17.526	ng	# 49
82) 3,3'-Dichlorobenzidine	21.46	252	65	51.412	ng	# 24
83) Chrysene	21.60	228	81	23.547	ng	# 49
84) Bis(2-ethylhexyl)phthalate	21.43	149	84	43.681	ng	# 50
85) Di-n-octyl phthalate	22.64	149	58	18.109	ng	# 1
86) Indeno(1,2,3-cd)pyrene	28.40	276	57	13.527	ng	# 56
88) Benzo(b)fluoranthene	23.60	252	144	40.240	ng	# 59
89) Benzo(k)fluoranthene	23.71	252	61	17.226	ng	# 60
90) Benzo(a)pyrene	24.57	252	60	17.574	ng	# 57
91) Dibenzo(a,h)anthracene	28.89	278	63	18.467	ng	# 58

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG060119\
Data File : BG041010.D
Acq On : 1 Jun 2019 13:37
Operator : HP/JU
Sample : K3040-02
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
0445-FIELD-BLANK-2

Quant Time: Jun 03 01:12:01 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG052919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 29 18:39:57 2019
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
--------------------	------	------	----------	------	-------	----------

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

