

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG021022\
 Data File : BG052357.D
 Acq On : 10 Feb 2022 17:00
 Operator : CG/JU
 Sample : N1432-19
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SP-7

Quant Time: Feb 10 17:50:51 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG020722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 08 01:21:20 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.224	152	27831	20.000	ng	0.00
21) Naphthalene-d8	11.067	136	111034	20.000	ng	# 0.00
39) Acenaphthene-d10	14.873	164	77688	20.000	ng	0.00
64) Phenanthrene-d10	17.623	188	194323	20.000	ng	0.00
76) Chrysene-d12	21.940	240	200919	20.000	ng	0.00
86) Perylene-d12	25.412	264	209229	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.756	112	197596	123.741	ng	0.00
7) Phenol-d6	7.372	99	274227	111.300	ng	0.00
23) Nitrobenzene-d5	9.398	82	192914	75.511	ng	0.00
42) 2,4,6-Tribromophenol	16.360	330	141903	127.140	ng	0.00
45) 2-Fluorobiphenyl	13.493	172	422831	75.628	ng	0.00
79) Terphenyl-d14	20.213	244	812647	71.423	ng	0.00
Target Compounds						
						Qvalue
3) Pyridine	4.059	79	61	0.029	ng	# 23
4) n-Nitrosodimethylamine	3.853	42	61	0.055	ng	# 1
15) Benzyl Alcohol	8.218	79	973	0.476	ng	# 48
16) 2,2'-oxybis(1-Chloropr...	8.271	45	53	0.020	ng	# 68
24) Nitrobenzene	9.387	77	611	0.252	ng	# 48
25) Isophorone	9.751	82	62	0.014	ng	# 68
26) 2-Nitrophenol	10.180	139	58	0.056	ng	# 23
31) Naphthalene	11.108	128	65	0.011	ng	# 66
35) Caprolactam	11.971	113	55	0.083	ng	# 1
46) 1,1'-Biphenyl	13.499	154	751	0.132	ng	# 1
48) 2-Nitroaniline	13.745	65	210	0.134	ng	# 10
49) Acenaphthylene	14.603	152	628	0.088	ng	# 60
50) Dimethylphthalate	14.309	163	298	0.050	ng	# 76
52) Acenaphthene	14.885	154	55	0.012	ng	# 4
58) Fluorene	16.360	166	4529	0.771	ng	# 84
60) Diethylphthalate	15.661	149	2037	0.339	ng	# 89
63) Azobenzene	16.213	77	129	0.021	ng	# 48
65) 4,6-Dinitro-2-methylph...	16.095	198	56	0.048	ng	# 29
66) n-Nitrosodiphenylamine	16.360	169	4184	0.781	ng	# 48
71) Phenanthrene	17.664	178	3844	0.384	ng	# 98
72) Anthracene	17.664	178	3844	0.392	ng	# 95
74) Di-n-butylphthalate	18.557	149	1642	0.157	ng	# 77
75) Fluoranthene	19.667	202	5085	0.399	ng	# 86
78) Pyrene	20.025	202	9205	0.685	ng	# 95
80) Butylbenzylphthalate	20.895	149	139	0.030	ng	# 74
81) Benzo(a)anthracene	21.917	228	4381	0.330	ng	# 93
82) 3,3'-Dichlorobenzidine	21.717	252	57	0.012	ng	# 24
83) Chrysene	21.917	228	4381	0.348	ng	# 96
84) Bis(2-ethylhexyl)phtha...	21.793	149	1271	0.196	ng	# 86
85) Di-n-octyl phthalate	23.098	149	393	0.036	ng	# 71
87) Indeno(1,2,3-cd)pyrene	29.401	276	472	0.031	ng	# 17
88) Benzo(b)fluoranthene	24.290	252	1850	0.142	ng	# 86
89) Benzo(k)fluoranthene	24.314	252	1840	0.143	ng	# 79
90) Benzo(a)pyrene	25.248	252	4266	0.387	ng	# 80
91) Dibenzo(a,h)anthracene	29.483	278	218	0.017	ng	# 57
92) Benzo(g,h,i)perylene	30.664	276	2477	0.200	ng	# 66

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

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