

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122023\
 Data File : BG060104.D
 Acq On : 20 Dec 2023 17:44
 Operator : MA/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

SSTDCCC040

Manual Integrations

APPROVED

Reviewed By :Yogesh Patel 12/21/2023

Supervised By :mohammad ahmed 12/21/2023

Quant Time: Dec 21 02:38:59 2023

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG121323.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Tue Dec 19 04:37:23 2023

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.946	152	32710	20.000	ng	#-0.01
21) Naphthalene-d8	10.755	136	124401	20.000	ng	# 0.00
39) Acenaphthene-d10	14.585	164	129252	20.000	ng	-0.01
64) Phenanthrene-d10	17.335	188	431083	20.000	ng	#-0.01
76) Chrysene-d12	21.607	240	567996	20.000	ng	#-0.01
86) Perylene-d12	24.797	264	635863	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.514	112	106546	55.760	ng	-0.01
7) Phenol-d6	7.112	99	186385	69.344	ng	0.00
23) Nitrobenzene-d5	9.110	82	252386	92.125	ng	-0.01
42) 2,4,6-Tribromophenol	16.078	330	309820	102.856	ng	0.00
45) 2-Fluorobiphenyl	13.211	172	920554	87.168	ng	0.00
79) Terphenyl-d14	19.956	244	2771589	89.087	ng	-0.01
Target Compounds						Qvalue
2) 1,4-Dioxane	3.422	88	29476m	32.600	ng	
3) Pyridine	3.822	79	68096	26.892	ng	# 75
4) n-Nitrosodimethylamine	3.733	42	48344	38.814	ng	# 51
6) Aniline	7.276	93	121686	44.448	ng	# 88
8) 2-Chlorophenol	7.511	128	60724	33.536	ng	95
9) Benzaldehyde	7.082	77	60760	37.178	ng	94
10) Phenol	7.135	94	100192	37.107	ng	# 73
11) bis(2-Chloroethyl)ether	7.370	93	59093	30.571	ng	85
12) 1,3-Dichlorobenzene	7.835	146	81678	33.732	ng	# 90
13) 1,4-Dichlorobenzene	7.981	146	84583	33.569	ng	92
14) 1,2-Dichlorobenzene	8.299	146	82685	34.109	ng	95
15) Benzyl Alcohol	8.181	79	108549	50.248	ng	96
16) 2,2'-oxybis(1-Chloropr...	8.475	45	126584	36.596	ng	97
17) 2-Methylphenol	8.387	107	75965	42.199	ng	# 82
18) Hexachloroethane	9.027	117	26977	35.344	ng	83
19) n-Nitroso-di-n-propyla...	8.751	70	81070	46.750	ng	# 81
20) 3+4-Methylphenols	8.716	107	103766	41.268	ng	91
22) Acetophenone	8.769	105	153068	42.224	ng	# 91
24) Nitrobenzene	9.151	77	125116	44.984	ng	# 86
25) Isophorone	9.674	82	221613	42.265	ng	# 94
26) 2-Nitrophenol	9.867	139	44792	43.789	ng	# 52
27) 2,4-Dimethylphenol	9.920	122	75028	53.510	ng	# 82
28) bis(2-Chloroethoxy)met...	10.155	93	94372	34.860	ng	99
29) 2,4-Dichlorophenol	10.396	162	115886	46.018	ng	87
30) 1,2,4-Trichlorobenzene	10.614	180	157484	46.200	ng	# 91
31) Naphthalene	10.802	128	252128	38.932	ng	98
32) Benzoic acid	10.061	122	47453m	37.206	ng	
33) 4-Chloroaniline	10.907	127	113476	46.051	ng	# 90
34) Hexachlorobutadiene	11.090	225	162871	52.205	ng	88
35) Caprolactam	11.689	113	26917m	42.076	ng	
36) 4-Chloro-3-methylphenol	12.036	107	108564	47.402	ng	98
37) 2-Methylnaphthalene	12.412	142	232303	45.735	ng	# 91
38) 1-Methylnaphthalene	12.629	142	211284	42.111	ng	# 99
40) 1,2,4,5-Tetrachloroben...	12.776	216	314269	46.639	ng	97
41) Hexachlorocyclopentadiene	12.758	237	175340	62.980	ng	95
43) 2,4,6-Trichlorophenol	13.023	196	163061	43.159	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.087	196	190962	46.469	ng	# 98
46) 1,1'-Biphenyl	13.422	154	387846	40.462	ng	97
47) 2-Chloronaphthalene	13.463	162	312448	37.638	ng	96
48) 2-Nitroaniline	13.669	65	98841	55.157	ng	# 80
49) Acenaphthylene	14.309	152	469485	41.345	ng	99
50) Dimethylphthalate	14.045	163	451872	45.040	ng	# 97
51) 2,6-Dinitrotoluene	14.162	165	96989	46.668	ng	# 81
52) Acenaphthene	14.656	154	287787	36.818	ng	95
53) 3-Nitroaniline	14.497	138	64462	38.816	ng	86
54) 2,4-Dinitrophenol	14.697	184	68413	44.679	ng	# 74
55) Dibenzofuran	14.985	168	554808	43.609	ng	95
56) 4-Nitrophenol	14.803	139	52663	37.731	ng	# 54
57) 2,4-Dinitrotoluene	14.950	165	140068	47.944	ng	# 89
58) Fluorene	15.637	166	459025	42.787	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.208	232	206897	48.366	ng	95
60) Diethylphthalate	15.408	149	401705	43.516	ng	99
61) 4-Chlorophenyl-phenyle...	15.631	204	378953	49.046	ng	97
62) 4-Nitroaniline	15.661	138	73517	40.654	ng	# 42
63) Azobenzene	15.919	77	351024	38.569	ng	87
65) 4,6-Dinitro-2-methylph...	15.708	198	116388	43.061	ng	85
66) n-Nitrosodiphenylamine	15.843	169	441298	41.904	ng	94
67) 4-Bromophenyl-phenylether	16.524	248	283796	46.660	ng	95
68) Hexachlorobenzene	16.642	284	292068	41.077	ng	98
69) Atrazine	16.789	200	132127	34.344	ng	92
70) Pentachlorophenol	16.983	266	173700	37.710	ng	96
71) Phenanthrene	17.382	178	856324	41.011	ng	98
72) Anthracene	17.470	178	894280	42.596	ng	98
73) Carbazole	17.741	167	668865	37.492	ng	99
74) Di-n-butylphthalate	18.299	149	628010	38.588	ng	98
75) Fluoranthene	19.392	202	1312963	43.090	ng	97
77) Benzidine	19.574	184	245366	33.631	ng	97
78) Pyrene	19.756	202	1328936	38.597	ng	99
80) Butylbenzylphthalate	20.649	149	261359	42.265	ng	89
81) Benzo(a)anthracene	21.589	228	1577146	41.336	ng	97
82) 3,3'-Dichlorobenzidine	21.507	252	581117	46.287	ng	# 96
83) Chrysene	21.654	228	1480447	40.956	ng	99
84) Bis(2-ethylhexyl)phtha...	21.501	149	381498	40.703	ng	99
85) Di-n-octyl phthalate	22.699	149	610497	37.655	ng	# 94
87) Indeno(1,2,3-cd)pyrene	28.434	276	1870468	41.218	ng	# 97
88) Benzo(b)fluoranthene	23.781	252	1518744	38.674	ng	# 99
89) Benzo(k)fluoranthene	23.851	252	1551581	40.733	ng	# 98
90) Benzo(a)pyrene	24.650	252	1467949	42.191	ng	# 99
91) Dibenzo(a,h)anthracene	28.510	278	1602066	42.218	ng	# 97
92) Benzo(g,h,i)perylene	29.568	276	1514839	40.133	ng	# 93

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

