

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062723\
 Data File : BG057866.D
 Acq On : 27 Jun 2023 14:23
 Operator : CG/JU
 Sample : SSTDICC005
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

SSTDICC005

Manual Integrations

APPROVED

Reviewed By :Yogesh Patel 06/29/2023

Supervised By :mohammad ahmed 06/29/2023

Quant Time: Jun 28 00:38:25 2023

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG062723.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Wed Jun 28 00:34:47 2023

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.924	152	39625	20.000	ng	0.00
21) Naphthalene-d8	10.727	136	175080	20.000	ng	0.00
39) Acenaphthene-d10	14.557	164	130865	20.000	ng	0.00
64) Phenanthrene-d10	17.301	188	348997	20.000	ng	0.00
76) Chrysene-d12	21.555	240	315828	20.000	ng	0.00
86) Perylene-d12	24.710	264	385179	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.492	112	27038	10.440	ng	0.00
7) Phenol-d6	7.084	99	39594	10.224	ng	0.00
23) Nitrobenzene-d5	9.082	82	36273	10.153	ng	0.00
42) 2,4,6-Tribromophenol	16.044	330	21945	9.984	ng	0.00
45) 2-Fluorobiphenyl	13.183	172	94582	11.029	ng	0.00
79) Terphenyl-d14	19.916	244	196300	11.331	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.400	88	7158	6.030	ng	88
3) Pyridine	3.800	79	17114	4.989	ng	95
4) n-Nitrosodimethylamine	3.711	42	7342	5.200	ng	90
6) Aniline	7.248	93	25284	5.226	ng	94
8) 2-Chlorophenol	7.483	128	13356	5.137	ng	96
10) Phenol	7.107	94	19094	5.074	ng	97
11) bis(2-Chloroethyl)ether	7.342	93	14847	5.233	ng	98
12) 1,3-Dichlorobenzene	7.812	146	14676	5.488	ng	94
13) 1,4-Dichlorobenzene	7.959	146	14555	5.419	ng	91
14) 1,2-Dichlorobenzene	8.277	146	13862	5.333	ng	96
15) Benzyl Alcohol	8.159	79	14380	5.040	ng	93
16) 2,2'-oxybis(1-Chloropr...	8.453	45	25533m	5.405	ng	
17) 2-Methylphenol	8.359	107	14187	5.085	ng	93
18) Hexachloroethane	9.005	117	4666	4.985	ng	92
19) n-Nitroso-di-n-propyla...	8.723	70	13146	5.073	ng	# 94
20) 3+4-Methylphenols	8.682	107	19392	4.989	ng	98
22) Acetophenone	8.741	105	25299	5.192	ng	# 99
24) Nitrobenzene	9.123	77	17661	4.925	ng	95
25) Isophorone	9.646	82	39808	5.161	ng	95
26) 2-Nitrophenol	9.839	139	6387	4.371	ng	98
27) 2,4-Dimethylphenol	9.892	122	14460	5.133	ng	94
28) bis(2-Chloroethoxy)met...	10.133	93	20827	5.024	ng	98
29) 2,4-Dichlorophenol	10.368	162	13191	4.774	ng	98
30) 1,2,4-Trichlorobenzene	10.592	180	14639	4.922	ng	98
31) Naphthalene	10.780	128	49100	5.254	ng	99
33) 4-Chloroaniline	10.885	127	21563	4.904	ng	98
34) Hexachlorobutadiene	11.062	225	9026	4.920	ng	95
35) Caprolactam	11.626	113	5944	4.964	ng	# 83
36) 4-Chloro-3-methylphenol	12.002	107	18395	5.015	ng	95
37) 2-Methylnaphthalene	12.384	142	35439	5.144	ng	94
38) 1-Methylnaphthalene	12.601	142	33467	5.141	ng	100
40) 1,2,4,5-Tetrachloroben...	12.754	216	18758	5.123	ng	99
41) Hexachlorocyclopentadiene	12.736	237	8657	4.398	ng	91
43) 2,4,6-Trichlorophenol	12.989	196	13115	4.860	ng	97
44) 2,4,5-Trichlorophenol	13.059	196	15435	4.805	ng	99
46) 1,1'-Biphenyl	13.394	154	47044	5.278	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
47) 2-Chloronaphthalene	13.435	162	35387	5.169	ng	95
48) 2-Nitroaniline	13.635	65	12335	4.522	ng	97
49) Acenaphthylene	14.281	152	62315	5.191	ng	98
50) Dimethylphthalate	14.011	163	53923	5.199	ng	99
51) 2,6-Dinitrotoluene	14.129	165	9573	4.445	ng	98
52) Acenaphthene	14.622	154	35790	5.103	ng	98
53) 3-Nitroaniline	14.458	138	11254	4.597	ng	# 99
55) Dibenzofuran	14.957	168	63399	5.271	ng	99
56) 4-Nitrophenol	14.757	139	8980	4.474	ng	100
57) 2,4-Dinitrotoluene	14.916	165	14196	4.543	ng	# 97
58) Fluorene	15.603	166	52541	5.412	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.180	232	14422	4.953	ng	# 96
60) Diethylphthalate	15.374	149	58372	5.339	ng	99
61) 4-Chlorophenyl-phenyle...	15.597	204	27099	5.204	ng	94
62) 4-Nitroaniline	15.615	138	12334	4.803	ng	97
63) Azobenzene	15.891	77	53446	5.112	ng	99
66) n-Nitrosodiphenylamine	15.809	169	46686	5.242	ng	97
67) 4-Bromophenyl-phenylether	16.491	248	18982	5.153	ng	97
68) Hexachlorobenzene	16.608	284	20205	5.194	ng	98
69) Atrazine	16.755	200	18421	5.569	ng	97
70) Pentachlorophenol	16.949	266	13241m	4.451	ng	
71) Phenanthrene	17.342	178	90236	5.386	ng	99
72) Anthracene	17.436	178	91997	5.324	ng	98
73) Carbazole	17.701	167	88862	5.409	ng	98
74) Di-n-butylphthalate	18.265	149	108976	5.367	ng	98
75) Fluoranthene	19.352	202	116858	5.537	ng	98
78) Pyrene	19.716	202	122592	5.436	ng	98
80) Butylbenzylphthalate	20.609	149	46635	5.100	ng	97
81) Benzo(a)anthracene	21.538	228	115948	5.305	ng	97
82) 3,3'-Dichlorobenzidine	21.455	252	41081	5.282	ng	98
83) Chrysene	21.602	228	108809	5.260	ng	99
84) Bis(2-ethylhexyl)phtha...	21.449	149	69904	5.384	ng	98
85) Di-n-octyl phthalate	22.636	149	116480	5.084	ng	99
87) Indeno(1,2,3-cd)pyrene	28.283	276	130457	5.113	ng	# 92
88) Benzo(b)fluoranthene	23.700	252	107918	5.137	ng	98
89) Benzo(k)fluoranthene	23.770	252	110274	5.194	ng	99
90) Benzo(a)pyrene	24.552	252	107687	5.175	ng	99
91) Dibenzo(a,h)anthracene	28.347	278	109668	5.181	ng	97
92) Benzo(g,h,i)perylene	29.405	276	106364	5.056	ng	97

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

