

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110122\
 Data File : BG055395.D
 Acq On : 1 Nov 2022 15:39
 Operator : CG/JU
 Sample : SSTDICC010
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

SSTDICC010

Manual Integrations

APPROVED

Reviewed By :Christian Giraldo 11/02/2022
 Supervised By :Jagrut Upadhyay 11/02/2022

Quant Time: Nov 02 04:10:59 2022

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG110122.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Wed Nov 02 04:08:03 2022

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.244	152	24973	20.000	ng	0.00
21) Naphthalene-d8	11.076	136	110806	20.000	ng	0.00
39) Acenaphthene-d10	14.860	164	83188	20.000	ng	0.00
64) Phenanthrene-d10	17.592	188	202052	20.000	ng	0.00
76) Chrysene-d12	21.870	240	205531	20.000	ng	0.00
86) Perylene-d12	25.242	264	224121	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.771	112	28612	20.061	ng	0.00
7) Phenol-d6	7.387	99	39860	19.291	ng	0.00
23) Nitrobenzene-d5	9.420	82	42457	19.566	ng	0.00
42) 2,4,6-Tribromophenol	16.341	330	23632	26.132	ng	0.00
45) 2-Fluorobiphenyl	13.485	172	119487	23.510	ng	0.00
79) Terphenyl-d14	20.166	244	236754	23.933	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.573	88	6574	9.546	ng	98
3) Pyridine	4.002	79	18491	8.916	ng	95
4) n-Nitrosodimethylamine	3.908	42	7426	8.060	ng	# 88
6) Aniline	7.557	93	26066	9.636	ng	98
8) 2-Chlorophenol	7.804	128	18020	10.868	ng	98
9) Benzaldehyde	7.375	77	12119	8.590	ng	96
10) Phenol	7.416	94	22541	9.739	ng	98
11) bis(2-Chloroethyl)ether	7.651	93	17638	10.095	ng	94
12) 1,3-Dichlorobenzene	8.133	146	20311	11.227	ng	99
13) 1,4-Dichlorobenzene	8.280	146	20500	11.094	ng	98
14) 1,2-Dichlorobenzene	8.603	146	19952	11.214	ng	95
15) Benzyl Alcohol	8.479	79	15223	8.804	ng	98
16) 2,2'-oxybis(1-Chloropr...	8.762	45	22996	8.043	ng	97
17) 2-Methylphenol	8.685	107	16374	10.004	ng	97
18) Hexachloroethane	9.337	117	7149	10.864	ng	92
19) n-Nitroso-di-n-propyla...	9.044	70	16069	9.365	ng	97
20) 3+4-Methylphenols	9.014	107	23309	9.943	ng	97
22) Acetophenone	9.073	105	29878	10.547	ng	# 99
24) Nitrobenzene	9.461	77	22265	9.854	ng	98
25) Isophorone	9.984	82	42558	9.710	ng	98
26) 2-Nitrophenol	10.178	139	10900	11.016	ng	99
27) 2,4-Dimethylphenol	10.225	122	15951m	10.326	ng	
28) bis(2-Chloroethoxy)met...	10.460	93	22744	10.278	ng	96
29) 2,4-Dichlorophenol	10.724	162	18663	11.262	ng	98
30) 1,2,4-Trichlorobenzene	10.930	180	21183	12.525	ng	93
31) Naphthalene	11.123	128	64723	10.949	ng	99
32) Benzoic acid	10.413	122	5678m	5.292	ng	
33) 4-Chloroaniline	11.229	127	26154	10.349	ng	97
34) Hexachlorobutadiene	11.400	225	13209	13.669	ng	96
35) Caprolactam	11.981	113	7658	9.957	ng	95
36) 4-Chloro-3-methylphenol	12.334	107	21066	10.103	ng	99
37) 2-Methylnaphthalene	12.710	142	47655	11.182	ng	98
38) 1-Methylnaphthalene	12.927	142	46076	10.959	ng	98
40) 1,2,4,5-Tetrachloroben...	13.068	216	24803	12.362	ng	100
41) Hexachlorocyclopentadiene	13.045	237	6564	7.975	ng	94
43) 2,4,6-Trichlorophenol	13.309	196	16330	11.020	ng	94

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.391	196	18957	11.566	ng	95
46) 1,1'-Biphenyl	13.697	154	63499	11.129	ng	99
47) 2-Chloronaphthalene	13.750	162	50252	11.169	ng	100
48) 2-Nitroaniline	13.944	65	14938	8.490	ng	99
49) Acenaphthylene	14.584	152	83544	10.757	ng	99
50) Dimethylphthalate	14.296	163	74664	11.431	ng	98
51) 2,6-Dinitrotoluene	14.425	165	14565	10.959	ng	93
52) Acenaphthene	14.925	154	49906	10.081	ng	99
53) 3-Nitroaniline	14.760	138	16653	9.969	ng	92
54) 2,4-Dinitrophenol	14.984	184	5469	8.568	ng	88
55) Dibenzofuran	15.254	168	83401	11.229	ng	97
56) 4-Nitrophenol	15.072	139	10873	9.100	ng	99
57) 2,4-Dinitrotoluene	15.213	165	20900	10.836	ng	96
58) Fluorene	15.900	166	69074	11.177	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.483	232	17019	11.647	ng	# 98
60) Diethylphthalate	15.648	149	78751	11.335	ng	99
61) 4-Chlorophenyl-phenyle...	15.883	204	35204	12.396	ng	96
62) 4-Nitroaniline	15.918	138	18453	10.512	ng	98
63) Azobenzene	16.170	77	63997	9.455	ng	96
65) 4,6-Dinitro-2-methylph...	15.977	198	12184	10.969	ng	95
66) n-Nitrosodiphenylamine	16.094	169	61337	10.907	ng	99
67) 4-Bromophenyl-phenylether	16.770	248	23975	12.213	ng	94
68) Hexachlorobenzene	16.899	284	27295	12.298	ng	98
69) Atrazine	17.028	200	21889	11.446	ng	99
70) Pentachlorophenol	17.252	266	9458m	8.636	ng	
71) Phenanthrene	17.633	178	119487	11.089	ng	100
72) Anthracene	17.727	178	118054	11.207	ng	98
73) Carbazole	17.992	167	111741	11.028	ng	99
74) Di-n-butylphthalate	18.521	149	142030	11.214	ng	99
75) Fluoranthene	19.625	202	151792	12.089	ng	99
77) Benzidine	19.796	184	40661	8.171	ng	99
78) Pyrene	19.984	202	156397	11.102	ng	99
80) Butylbenzylphthalate	20.841	149	61176	9.189	ng	99
81) Benzo(a)anthracene	21.846	228	148332	10.747	ng	98
82) 3,3'-Dichlorobenzidine	21.746	252	48452	10.779	ng	96
83) Chrysene	21.917	228	141610	10.802	ng	99
84) Bis(2-ethylhexyl)phtha...	21.711	149	89930	9.456	ng	98
85) Di-n-octyl phthalate	22.963	149	153896	9.714	ng	98
87) Indeno(1,2,3-cd)pyrene	29.126	276	184346	11.137	ng	100
88) Benzo(b)fluoranthene	24.161	252	148622	11.024	ng	99
89) Benzo(k)fluoranthene	24.232	252	149696	10.997	ng	98
90) Benzo(a)pyrene	25.084	252	128016	11.073	ng	97
91) Dibenzo(a,h)anthracene	29.179	278	153786	11.321	ng	99
92) Benzo(g,h,i)perylene	30.342	276	150730	11.187	ng	100

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

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