

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG020923\
 Data File : BG056595.D
 Acq On : 9 Feb 2023 12:22
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

SSTDCCC040

Manual Integrations

APPROVED

Reviewed By : Christian Giraldo 02/10/2023
 Supervised By : Jagrut Upadhyay 02/10/2023

Quant Time: Feb 10 00:22:25 2023

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG012723.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Wed Feb 08 03:02:11 2023

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.265	152	31132	20.000	ng	0.00
21) Naphthalene-d8	11.097	136	116984	20.000	ng	0.00
39) Acenaphthene-d10	14.887	164	95110	20.000	ng	0.00
64) Phenanthrene-d10	17.624	188	234159	20.000	ng	0.00
76) Chrysene-d12	21.914	240	213890	20.000	ng	-0.01
86) Perylene-d12	25.333	264	233697	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.785	112	130072	76.859	ng	0.00
7) Phenol-d6	7.413	99	187877	74.926	ng	0.00
23) Nitrobenzene-d5	9.440	82	164017	79.615	ng	0.00
42) 2,4,6-Tribromophenol	16.373	330	94512	74.747	ng	0.00
45) 2-Fluorobiphenyl	13.512	172	554247	80.793	ng	0.00
79) Terphenyl-d14	20.204	244	960992	78.912	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.576	88	26590	37.733	ng	97
3) Pyridine	3.999	79	77471	36.880	ng	98
4) n-Nitrosodimethylamine	3.911	42	17003	38.062	ng	# 96
6) Aniline	7.577	93	123573	37.640	ng	96
8) 2-Chlorophenol	7.824	128	70479	37.910	ng	98
9) Benzaldehyde	7.389	77	41192	34.452	ng	90
10) Phenol	7.442	94	95543	37.489	ng	98
11) bis(2-Chloroethyl)ether	7.666	93	66798	38.132	ng	95
12) 1,3-Dichlorobenzene	8.147	146	82974	38.763	ng	96
13) 1,4-Dichlorobenzene	8.294	146	83319	38.436	ng	98
14) 1,2-Dichlorobenzene	8.623	146	80466	38.259	ng	98
15) Benzyl Alcohol	8.500	79	61770	35.904	ng	97
16) 2,2'-oxybis(1-Chloropr...	8.782	45	14179	38.201	ng	92
17) 2-Methylphenol	8.706	107	73955	38.144	ng	97
18) Hexachloroethane	9.358	117	25689	39.181	ng	97
19) n-Nitroso-di-n-propyla...	9.070	70	54372	37.627	ng	98
20) 3+4-Methylphenols	9.040	107	102348	37.668	ng	99
22) Acetophenone	9.093	105	124287	39.986	ng	# 100
24) Nitrobenzene	9.481	77	78454	39.481	ng	96
25) Isophorone	10.004	82	165261	38.884	ng	100
26) 2-Nitrophenol	10.204	139	48458	40.192	ng	96
27) 2,4-Dimethylphenol	10.251	122	80963m	40.125	ng	
28) bis(2-Chloroethoxy)met...	10.474	93	92660	39.465	ng	98
29) 2,4-Dichlorophenol	10.744	162	91097	40.109	ng	96
30) 1,2,4-Trichlorobenzene	10.950	180	101446	40.194	ng	100
31) Naphthalene	11.144	128	245711	39.795	ng	98
32) Benzoic acid	10.415	122	32870m	29.117	ng	
33) 4-Chloroaniline	11.250	127	114883	39.860	ng	98
34) Hexachlorobutadiene	11.420	225	74134	41.845	ng	95
35) Caprolactam	12.031	113	27853	36.362	ng	96
36) 4-Chloro-3-methylphenol	12.360	107	87166	39.740	ng	96
37) 2-Methylnaphthalene	12.730	142	203070	39.893	ng	99
38) 1-Methylnaphthalene	12.948	142	187689	39.480	ng	98
40) 1,2,4,5-Tetrachloroben...	13.095	216	140286	40.374	ng	99
41) Hexachlorocyclopentadiene	13.065	237	60482	36.592	ng	99
43) 2,4,6-Trichlorophenol	13.335	196	87265	38.972	ng	94

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.418	196	100840	38.464	ng	96
46) 1,1'-Biphenyl	13.723	154	277552	40.234	ng	98
47) 2-Chloronaphthalene	13.770	162	216673	40.183	ng	98
48) 2-Nitroaniline	13.976	65	44532	39.950	ng	96
49) Acenaphthylene	14.616	152	347507	39.611	ng	99
50) Dimethylphthalate	14.328	163	299773	38.959	ng	99
51) 2,6-Dinitrotoluene	14.458	165	65704	39.155	ng	98
52) Acenaphthene	14.951	154	204870	39.380	ng	96
53) 3-Nitroaniline	14.793	138	62589	38.474	ng	100
54) 2,4-Dinitrophenol	15.010	184	41875	39.806	ng	96
55) Dibenzofuran	15.280	168	369999	39.657	ng	98
56) 4-Nitrophenol	15.110	139	40789	36.694	ng	95
57) 2,4-Dinitrotoluene	15.245	165	92785	39.255	ng	99
58) Fluorene	15.926	166	292868	39.449	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.509	232	84277	36.155	ng	# 96
60) Diethylphthalate	15.674	149	276131	38.217	ng	98
61) 4-Chlorophenyl-phenylether	15.909	204	178391	39.379	ng	97
62) 4-Nitroaniline	15.950	138	63332	38.906	ng	98
63) Azobenzene	16.203	77	194759	39.174	ng	99
65) 4,6-Dinitro-2-methylph...	16.015	198	58659	35.632	ng	97
66) n-Nitrosodiphenylamine	16.120	169	264478	39.893	ng	98
67) 4-Bromophenyl-phenylether	16.802	248	121139	40.411	ng	98
68) Hexachlorobenzene	16.931	284	119205	40.602	ng	98
69) Atrazine	17.055	200	101058	39.512	ng	100
70) Pentachlorophenol	17.284	266	62396	35.123	ng	98
71) Phenanthrene	17.672	178	486417	39.690	ng	98
72) Anthracene	17.760	178	497588	39.552	ng	99
73) Carbazole	18.030	167	419960	39.194	ng	99
74) Di-n-butylphthalate	18.547	149	437301	38.870	ng	100
75) Fluoranthene	19.663	202	601248	38.871	ng	99
77) Benzidine	19.834	184	261856	45.240	ng	97
78) Pyrene	20.022	202	598728	39.358	ng	99
80) Butylbenzylphthalate	20.874	149	182517	39.286	ng	98
81) Benzo(a)anthracene	21.896	228	590402	39.483	ng	99
82) 3,3'-Dichlorobenzidine	21.796	252	216149	40.140	ng	100
83) Chrysene	21.966	228	554168	39.447	ng	98
84) Bis(2-ethylhexyl)phtha...	21.749	149	259803	39.004	ng	99
85) Di-n-octyl phthalate	23.012	149	437001	39.395	ng	100
87) Indeno(1,2,3-cd)pyrene	29.270	276	684250	39.998	ng	# 94
88) Benzo(b)fluoranthene	24.246	252	574766	39.624	ng	98
89) Benzo(k)fluoranthene	24.317	252	588528	40.186	ng	99
90) Benzo(a)pyrene	25.174	252	570576	39.932	ng	99
91) Dibenzo(a,h)anthracene	29.334	278	576754	40.165	ng	99
92) Benzo(g,h,i)perylene	30.521	276	556226	40.145	ng	99

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

