

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101322\
 Data File : BG055149.D
 Acq On : 12 Oct 2022 20:29
 Operator : CG/JU
 Sample : SSTDICC080
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

SSTDICC080

Manual Integrations

APPROVED

Reviewed By :Christian Giraldo 10/13/2022
 Supervised By :Jagrut Upadhyay 10/13/2022

Quant Time: Oct 13 04:28:13 2022

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG101322.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Thu Oct 13 04:19:34 2022

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.302	152	27623	20.000	ng	0.00
21) Naphthalene-d8	11.134	136	118912	20.000	ng	0.00
39) Acenaphthene-d10	14.912	164	99785	20.000	ng	0.00
64) Phenanthrene-d10	17.644	188	254443	20.000	ng	0.00
76) Chrysene-d12	21.927	240	242974	20.000	ng	0.00
86) Perylene-d12	25.341	264	263739	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.817	112	222933	179.258	ng	0.00
7) Phenol-d6	7.439	99	320667	172.090	ng	0.00
23) Nitrobenzene-d5	9.477	82	347563	174.653	ng	0.00
42) 2,4,6-Tribromophenol	16.393	330	231361	261.300	ng	0.00
45) 2-Fluorobiphenyl	13.543	172	878026	137.366	ng	0.00
79) Terphenyl-d14	20.218	244	1730234	143.930	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.643	88	47530	67.547	ng	99
3) Pyridine	4.060	79	151038m	71.380	ng	
4) n-Nitrosodimethylamine	3.966	42	77158	63.763	ng	92
6) Aniline	7.621	93	208537	76.488	ng	98
8) 2-Chlorophenol	7.862	128	132954	92.943	ng	99
9) Benzaldehyde	7.427	77	90815	60.668	ng	99
10) Phenol	7.462	94	178056	84.089	ng	98
11) bis(2-Chloroethyl)ether	7.709	93	122662	70.678	ng	100
12) 1,3-Dichlorobenzene	8.191	146	151402	76.892	ng	98
13) 1,4-Dichlorobenzene	8.337	146	153687	77.127	ng	98
14) 1,2-Dichlorobenzene	8.666	146	145761	75.292	ng	97
15) Benzyl Alcohol	8.537	79	146836	87.733	ng	99
16) 2,2'-oxybis(1-Chloropr...	8.831	45	201255	61.221	ng	98
17) 2-Methylphenol	8.731	107	128914	86.411	ng	97
18) Hexachloroethane	9.401	117	52383	82.905	ng	99
19) n-Nitroso-di-n-propyla...	9.113	70	126053	69.836	ng	96
20) 3+4-Methylphenols	9.066	107	186003	89.526	ng	96
22) Acetophenone	9.131	105	226901	75.245	ng	# 99
24) Nitrobenzene	9.518	77	178950	82.817	ng	98
25) Isophorone	10.047	82	340572	80.523	ng	98
26) 2-Nitrophenol	10.235	139	90027	142.695	ng	99
27) 2,4-Dimethylphenol	10.276	122	131773	90.884	ng	97
28) bis(2-Chloroethoxy)met...	10.517	93	164456	75.989	ng	99
29) 2,4-Dichlorophenol	10.770	162	161046	104.492	ng	99
30) 1,2,4-Trichlorobenzene	10.993	180	165961	79.914	ng	98
31) Naphthalene	11.187	128	486391	77.501	ng	99
32) Benzoic acid	10.435	122	121125m	135.773	ng	
33) 4-Chloroaniline	11.281	127	220659	85.028	ng	97
34) Hexachlorobutadiene	11.463	225	105451	78.308	ng	96
35) Caprolactam	12.068	113	66500m	117.900	ng	
36) 4-Chloro-3-methylphenol	12.380	107	184585	101.336	ng	96
37) 2-Methylnaphthalene	12.768	142	372645	80.572	ng	99
38) 1-Methylnaphthalene	12.985	142	361288	80.309	ng	98
40) 1,2,4,5-Tetrachloroben...	13.126	216	208515	71.568	ng	99
41) Hexachlorocyclopentadiene	13.102	237	110620	88.172	ng	99
43) 2,4,6-Trichlorophenol	13.355	196	158611	108.314	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.426	196	177439	103.980	ng	98
46) 1,1'-Biphenyl	13.755	154	491104	71.782	ng	100
47) 2-Chloronaphthalene	13.802	162	402945	73.090	ng	99
48) 2-Nitroaniline	13.996	65	154318	119.010	ng	97
49) Acenaphthylene	14.642	152	683133	74.830	ng	98
50) Dimethylphthalate	14.360	163	602706	88.991	ng	99
51) 2,6-Dinitrotoluene	14.483	165	132776	108.716	ng	97
52) Acenaphthene	14.977	154	457157	80.650	ng	98
53) 3-Nitroaniline	14.812	138	152990	105.670	ng	99
54) 2,4-Dinitrophenol	15.012	184	90484	143.736	ng	89
55) Dibenzofuran	15.306	168	684895	75.731	ng	99
56) 4-Nitrophenol	15.100	139	123540	112.836	ng	97
57) 2,4-Dinitrotoluene	15.259	165	198023	117.952	ng	99
58) Fluorene	15.952	166	567540	78.421	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.523	232	180737	113.012	ng	100
60) Diethylphthalate	15.705	149	644172	92.180	ng	99
61) 4-Chlorophenyl-phenylether	15.934	204	307319	78.120	ng	98
62) 4-Nitroaniline	15.970	138	159296	97.771	ng	94
63) Azobenzene	16.228	77	549683	69.840	ng	98
65) 4,6-Dinitro-2-methylphthalate	16.023	198	133779	141.058	ng	98
66) n-Nitrosodiphenylamine	16.152	169	517177	78.047	ng	97
67) 4-Bromophenyl-phenylether	16.828	248	222606	82.724	ng	98
68) Hexachlorobenzene	16.951	284	250837	80.625	ng	98
69) Atrazine	17.080	200	180505	83.609	ng	99
70) Pentachlorophenol	17.286	266	163080	109.466	ng	96
71) Phenanthrene	17.691	178	999110	74.752	ng	99
72) Anthracene	17.779	178	968727	74.521	ng	97
73) Carbazole	18.044	167	895003	78.440	ng	99
74) Di-n-butylphthalate	18.578	149	1089218	91.448	ng	99
75) Fluoranthene	19.677	202	1190060	71.372	ng	97
77) Benzidine	19.842	184	340498	81.347	ng	100
78) Pyrene	20.036	202	1201672	74.378	ng	96
80) Butylbenzylphthalate	20.899	149	489209	113.977	ng	99
81) Benzo(a)anthracene	21.910	228	1180206	76.463	ng	97
82) 3,3'-Dichlorobenzidine	21.810	252	405757	90.429	ng	98
83) Chrysene	21.980	228	1114905	74.945	ng	97
84) Bis(2-ethylhexyl)phthalate	21.786	149	679475	103.764	ng	98
85) Di-n-octyl phthalate	23.067	149	1135880	103.956	ng	100
87) Indeno(1,2,3-cd)pyrene	29.295	276	1530058	80.390	ng	# 96
88) Benzo(b)fluoranthene	24.260	252	1221738	78.621	ng	99
89) Benzo(k)fluoranthene	24.336	252	1185521	75.727	ng	99
90) Benzo(a)pyrene	25.188	252	1043419	78.342	ng	100
91) Dibenzo(a,h)anthracene	29.366	278	1254897	81.164	ng	99
92) Benzo(g,h,i)perylene	30.541	276	1240877	79.950	ng	98

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

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