

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071522\
 Data File : BG054297.D
 Acq On : 15 Jul 2022 11:09
 Operator : CG/JU
 Sample : SSTDICC010
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICC010

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 07/18/2022
 Supervised By : Jagrut Upadhyay 07/18/2022

Quant Time: Jul 15 14:20:41 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 15 14:17:51 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.999	152	14580	20.000	ng	0.00
21) Naphthalene-d8	10.807	136	56053	20.000	ng	# 0.00
39) Acenaphthene-d10	14.632	164	48649	20.000	ng	0.00
64) Phenanthrene-d10	17.376	188	135914	20.000	ng	# 0.00
76) Chrysene-d12	21.642	240	172890	20.000	ng	# 0.00
86) Perylene-d12	24.838	264	182097	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.566	112	13861	19.100	ng	0.00
7) Phenol-d6	7.165	99	18392	18.375	ng	0.00
23) Nitrobenzene-d5	9.156	82	19320	18.704	ng	0.00
42) 2,4,6-Tribromophenol	16.119	330	19707	20.912	ng	0.00
45) 2-Fluorobiphenyl	13.252	172	66712	20.149	ng	0.00
79) Terphenyl-d14	19.985	244	179068	21.299	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.428	88	2882	9.081	ng	94
3) Pyridine	3.833	79	6559	8.275	ng	# 93
4) n-Nitrosodimethylamine	3.751	42	3559	8.888	ng	# 71
6) Aniline	7.317	93	10945	9.302	ng	98
8) 2-Chlorophenol	7.564	128	8086	9.941	ng	89
9) Benzaldehyde	7.129	77	6025	10.024	ng	# 80
10) Phenol	7.188	94	9727	9.632	ng	93
11) bis(2-Chloroethyl)ether	7.417	93	6630	8.903	ng	87
12) 1,3-Dichlorobenzene	7.887	146	9846m	9.969	ng	
13) 1,4-Dichlorobenzene	8.034	146	10078	9.958	ng	96
14) 1,2-Dichlorobenzene	8.351	146	9422m	9.695	ng	
15) Benzyl Alcohol	8.240	79	7369	9.133	ng	96
16) 2,2'-oxybis(1-Chloropr...	8.528	45	8634	8.349	ng	91
17) 2-Methylphenol	8.451	107	6867	9.517	ng	# 86
18) Hexachloroethane	9.086	117	3221	9.354	ng	# 64
19) n-Nitroso-di-n-propyla...	8.798	70	6424	9.401	ng	92
20) 3+4-Methylphenols	8.774	107	9494	9.378	ng	96
22) Acetophenone	8.822	105	13349	9.854	ng	# 97
24) Nitrobenzene	9.203	77	9874	9.724	ng	93
25) Isophorone	9.726	82	17802	9.527	ng	# 95
26) 2-Nitrophenol	9.920	139	4645	9.285	ng	# 82
27) 2,4-Dimethylphenol	9.979	122	6827	9.756	ng	84
28) bis(2-Chloroethoxy)met...	10.208	93	8834	9.288	ng	92
29) 2,4-Dichlorophenol	10.461	162	10236	10.093	ng	84
30) 1,2,4-Trichlorobenzene	10.672	180	12532	10.015	ng	92
31) Naphthalene	10.854	128	27070	9.604	ng	98
32) Benzoic acid	10.190	122	683m	2.399	ng	
33) 4-Chloroaniline	10.966	127	11457	9.774	ng	# 92
34) Hexachlorobutadiene	11.148	225	11220	10.806	ng	92
35) Caprolactam	11.706	113	2655	9.192	ng	# 66
36) 4-Chloro-3-methylphenol	12.094	107	9615	9.854	ng	96
37) 2-Methylnaphthalene	12.464	142	21374	9.895	ng	98
38) 1-Methylnaphthalene	12.682	142	21520m	10.261	ng	
40) 1,2,4,5-Tetrachloroben...	12.834	216	20141	10.199	ng	# 94
41) Hexachlorocyclopentadiene	12.811	237	1413	2.861	ng	83
43) 2,4,6-Trichlorophenol	13.069	196	10522	9.278	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.152	196	12381	9.980	ng	# 91
46) 1,1'-Biphenyl	13.469	154	31256	9.554	ng	95
47) 2-Chloronaphthalene	13.510	162	26477	9.681	ng	96
48) 2-Nitroaniline	13.710	65	6649	8.590	ng	# 88
49) Acenaphthylene	14.356	152	40464	9.753	ng	99
50) Dimethylphthalate	14.080	163	37241	9.750	ng	# 98
51) 2,6-Dinitrotoluene	14.203	165	7799	9.394	ng	# 74
52) Acenaphthene	14.697	154	24050	9.584	ng	98
53) 3-Nitroaniline	14.532	138	6582	8.944	ng	90
54) 2,4-Dinitrophenol	14.756	184	2344	6.022	ng	# 80
55) Dibenzofuran	15.032	168	45173	10.248	ng	93
56) 4-Nitrophenol	14.861	139	3714	7.856	ng	88
57) 2,4-Dinitrotoluene	14.997	165	11488	9.556	ng	# 81
58) Fluorene	15.678	166	36314	9.917	ng	89
59) 2,3,4,6-Tetrachlorophenol	15.261	232	10964	8.824	ng	# 87
60) Diethylphthalate	15.443	149	36786	9.874	ng	95
61) 4-Chlorophenyl-phenyle...	15.666	204	23813	10.238	ng	# 88
62) 4-Nitroaniline	15.690	138	7207	9.376	ng	96
63) Azobenzene	15.960	77	26052	9.279	ng	82
65) 4,6-Dinitro-2-methylph...	15.766	198	6613	8.271	ng	75
66) n-Nitrosodiphenylamine	15.884	169	32232	9.402	ng	96
67) 4-Bromophenyl-phenylether	16.559	248	17747	9.790	ng	# 90
68) Hexachlorobenzene	16.689	284	20764	10.035	ng	# 89
69) Atrazine	16.824	200	15883	10.490	ng	92
70) Pentachlorophenol	17.035	266	6058	7.288	ng	# 88
71) Phenanthrene	17.417	178	66843	9.832	ng	98
72) Anthracene	17.511	178	65882	9.895	ng	98
73) Carbazole	17.776	167	54722	9.705	ng	98
74) Di-n-butylphthalate	18.334	149	63916	9.573	ng	97
75) Fluoranthene	19.427	202	95331	10.311	ng	95
77) Benzidine	19.603	184	37321	12.297	ng	97
78) Pyrene	19.791	202	97358	9.785	ng	97
80) Butylbenzylphthalate	20.672	149	29398	9.118	ng	# 86
81) Benzo(a)anthracene	21.624	228	109088	9.975	ng	99
82) 3,3'-Dichlorobenzidine	21.530	252	33742	10.094	ng	# 98
83) Chrysene	21.683	228	103042	9.962	ng	100
84) Bis(2-ethylhexyl)phtha...	21.524	149	41400	9.051	ng	# 98
85) Di-n-octyl phthalate	22.723	149	70434	8.944	ng	# 95
87) Indeno(1,2,3-cd)pyrene	28.493	276	133923	9.776	ng	# 99
88) Benzo(b)fluoranthene	23.822	252	102528	9.495	ng	# 98
89) Benzo(k)fluoranthene	23.886	252	106128	9.907	ng	# 97
90) Benzo(a)pyrene	24.685	252	89834	9.723	ng	98
91) Dibenzo(a,h)anthracene	28.551	278	110822	9.875	ng	# 96
92) Benzo(g,h,i)perylene	29.638	276	109160	9.752	ng	# 92

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

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