

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012723\
 Data File : BG056455.D
 Acq On : 27 Jan 2023 12:17
 Operator : CG/JU
 Sample : SSTDICC010
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICC010

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 01/30/2023
 Supervised By : Jagrut Upadhyay 01/30/2023

Quant Time: Jan 27 23:45:35 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG012723.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 27 23:42:35 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.281	152	35254	20.000	ng	0.00
21) Naphthalene-d8	11.113	136	137497	20.000	ng	0.00
39) Acenaphthene-d10	14.903	164	111623	20.000	ng	0.00
64) Phenanthrene-d10	17.641	188	285777	20.000	ng	0.00
76) Chrysene-d12	21.930	240	271461	20.000	ng	0.00
86) Perylene-d12	25.361	264	300281	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.802	112	37582	19.611	ng	0.00
7) Phenol-d6	7.423	99	54399	19.158	ng	0.00
23) Nitrobenzene-d5	9.456	82	47695	19.698	ng	0.00
42) 2,4,6-Tribromophenol	16.383	330	28649	19.306	ng	0.00
45) 2-Fluorobiphenyl	13.528	172	163278	20.280	ng	0.00
79) Terphenyl-d14	20.214	244	317295	20.529	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.604	88	8417	10.548	ng	92
3) Pyridine	4.027	79	22435	9.488	ng	96
4) n-Nitrosodimethylamine	3.939	42	5066	10.015	ng	# 84
6) Aniline	7.594	93	35857	9.645	ng	98
8) 2-Chlorophenol	7.840	128	21356	10.144	ng	95
9) Benzaldehyde	7.406	77	16041	11.516	ng	90
10) Phenol	7.453	94	28326	9.815	ng	95
11) bis(2-Chloroethyl)ether	7.682	93	20061	10.113	ng	97
12) 1,3-Dichlorobenzene	8.169	146	23957	9.883	ng	97
13) 1,4-Dichlorobenzene	8.316	146	25227	10.277	ng	100
14) 1,2-Dichlorobenzene	8.639	146	24380	10.236	ng	97
15) Benzyl Alcohol	8.516	79	17971	9.224	ng	99
16) 2,2'-oxybis(1-Chloropr...	8.804	45	4122	9.807	ng	84
17) 2-Methylphenol	8.716	107	20742	9.447	ng	97
18) Hexachloroethane	9.380	117	7453	10.038	ng	97
19) n-Nitroso-di-n-propyla...	9.080	70	15477	9.458	ng	95
20) 3+4-Methylphenols	9.045	107	29896	9.716	ng	96
22) Acetophenone	9.110	105	35801	9.800	ng	# 94
24) Nitrobenzene	9.497	77	22819	9.770	ng	# 96
25) Isophorone	10.020	82	48408	9.691	ng	97
26) 2-Nitrophenol	10.214	139	13456	9.496	ng	95
27) 2,4-Dimethylphenol	10.261	122	23007	9.766	ng	96
28) bis(2-Chloroethoxy)met...	10.496	93	28010	10.150	ng	99
29) 2,4-Dichlorophenol	10.761	162	24916	9.334	ng	92
30) 1,2,4-Trichlorobenzene	10.972	180	29514	9.949	ng	# 93
31) Naphthalene	11.166	128	70781	9.753	ng	97
32) Benzoic acid	10.367	122	9869m	7.368	ng	
33) 4-Chloroaniline	11.266	127	32906	9.714	ng	98
34) Hexachlorobutadiene	11.442	225	21111	10.138	ng	94
35) Caprolactam	12.018	113	8346	9.270	ng	97
36) 4-Chloro-3-methylphenol	12.370	107	25544	9.908	ng	97
37) 2-Methylnaphthalene	12.752	142	57696	9.643	ng	97
38) 1-Methylnaphthalene	12.970	142	53797	9.628	ng	89
40) 1,2,4,5-Tetrachloroben...	13.111	216	40619	9.961	ng	99
41) Hexachlorocyclopentadiene	13.087	237	11355	6.265	ng	88
43) 2,4,6-Trichlorophenol	13.346	196	25106	9.553	ng	96

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.422	196	29921m	9.783	ng	
46) 1,1'-Biphenyl	13.739	154	79881	9.866	ng	100
47) 2-Chloronaphthalene	13.786	162	62624	9.896	ng	99
48) 2-Nitroaniline	13.980	65	12500	9.555	ng	91
49) Acenaphthylene	14.627	152	101803	9.888	ng	99
50) Dimethylphthalate	14.339	163	90121	9.980	ng	99
51) 2,6-Dinitrotoluene	14.468	165	19087	9.692	ng	96
52) Acenaphthene	14.967	154	60312	9.878	ng	96
53) 3-Nitroaniline	14.803	138	18781	9.837	ng	94
54) 2,4-Dinitrophenol	15.014	184	9101	7.371	ng	# 91
55) Dibenzofuran	15.296	168	108926	9.948	ng	99
56) 4-Nitrophenol	15.102	139	11166	8.604	ng	94
57) 2,4-Dinitrotoluene	15.255	165	27324	9.850	ng	98
58) Fluorene	15.943	166	86788	9.961	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.526	232	25203	9.213	ng	# 99
60) Diethylphthalate	15.684	149	85917	10.132	ng	98
61) 4-Chlorophenyl-phenyle...	15.925	204	52458	9.867	ng	96
62) 4-Nitroaniline	15.954	138	18654	9.764	ng	98
63) Azobenzene	16.219	77	58044	9.948	ng	98
65) 4,6-Dinitro-2-methylph...	16.019	198	17973	8.946	ng	98
66) n-Nitrosodiphenylamine	16.137	169	79646	9.844	ng	99
67) 4-Bromophenyl-phenylether	16.818	248	35117	9.599	ng	97
68) Hexachlorobenzene	16.947	284	34333	9.582	ng	94
69) Atrazine	17.071	200	31591	10.120	ng	98
70) Pentachlorophenol	17.294	266	16952	7.843	ng	95
71) Phenanthrene	17.682	178	147144	9.838	ng	99
72) Anthracene	17.776	178	151353	9.858	ng	99
73) Carbazole	18.040	167	130062	9.946	ng	99
74) Di-n-butylphthalate	18.563	149	138746	10.105	ng	99
75) Fluoranthene	19.674	202	187710	9.944	ng	99
77) Benzidine	19.844	184	90731	12.351	ng	98
78) Pyrene	20.038	202	195759	10.139	ng	98
80) Butylbenzylphthalate	20.890	149	57662	9.779	ng	99
81) Benzo(a)anthracene	21.906	228	187330	9.871	ng	99
82) 3,3'-Dichlorobenzidine	21.812	252	69799	10.213	ng	98
83) Chrysene	21.977	228	179116	10.046	ng	97
84) Bis(2-ethylhexyl)phtha...	21.771	149	83013	9.820	ng	99
85) Di-n-octyl phthalate	23.040	149	139710	9.924	ng	99
87) Indeno(1,2,3-cd)pyrene	29.292	276	214129	9.742	ng	# 95
88) Benzo(b)fluoranthene	24.256	252	181066	9.715	ng	96
89) Benzo(k)fluoranthene	24.327	252	188311	10.010	ng	100
90) Benzo(a)pyrene	25.191	252	180629	9.838	ng	99
91) Dibenzo(a,h)anthracene	29.345	278	181616	9.843	ng	99
92) Benzo(g,h,i)perylene	30.537	276	173701	9.757	ng	99

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

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