

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082521\
 Data File : BG049963.D
 Acq On : 25 Aug 2021 11:02
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Quant Time: Aug 25 12:48:18 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG080321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 23 16:13:24 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.979	152	23338	20.000	ng	# 0.00
21) Naphthalene-d8	10.804	136	91398	20.000	ng	# 0.00
39) Acenaphthene-d10	14.640	164	64686	20.000	ng	0.00
64) Phenanthrene-d10	17.395	188	144370	20.000	ng	# 0.00
76) Chrysene-d12	21.672	240	125164	20.000	ng	0.00
86) Perylene-d12	24.915	264	127581	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.529	112	104681	80.406	ng	0.00
7) Phenol-d6	7.150	99	159312	80.839	ng	0.00
23) Nitrobenzene-d5	9.153	82	161463	78.174	ng	0.00
42) 2,4,6-Tribromophenol	16.138	330	78284	82.335	ng	0.00
45) 2-Fluorobiphenyl	13.254	172	361926	81.413	ng	0.00
79) Terphenyl-d14	20.004	244	592859	86.027	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.367	88	19236	33.812	ng	94
3) Pyridine	3.784	79	55470	35.569	ng	# 83
4) n-Nitrosodimethylamine	3.702	42	30141	35.944	ng	# 75
6) Aniline	7.297	93	85946	41.149	ng	# 95
8) 2-Chlorophenol	7.544	128	58244	41.034	ng	99
9) Benzaldehyde	7.115	77	45991	34.975	ng	86
10) Phenol	7.180	94	78212	40.959	ng	90
11) bis(2-Chloroethyl)ether	7.397	93	52606	40.801	ng	91
12) 1,3-Dichlorobenzene	7.867	146	64431	39.634	ng	96
13) 1,4-Dichlorobenzene	8.020	146	66750	40.564	ng	96
14) 1,2-Dichlorobenzene	8.337	146	63768	40.906	ng	93
15) Benzyl Alcohol	8.225	79	66484	40.170	ng	95
16) 2,2'-oxybis(1-Chloropr...	8.507	45	55432	37.838	ng	96
17) 2-Methylphenol	8.437	107	55006	43.744	ng	96
18) Hexachloroethane	9.065	117	26454	41.820	ng	98
19) n-Nitroso-di-n-propyla...	8.789	70	52146	39.541	ng	89
20) 3+4-Methylphenols	8.766	107	73489	41.613	ng	96
22) Acetophenone	8.807	105	99418	40.038	ng	# 95
24) Nitrobenzene	9.195	77	82516	37.861	ng	92
25) Isophorone	9.717	82	145578	39.499	ng	97
26) 2-Nitrophenol	9.911	139	36118	43.424	ng	98
27) 2,4-Dimethylphenol	9.976	122	51369	41.786	ng	96
28) bis(2-Chloroethoxy)met...	10.199	93	67654	41.962	ng	# 91
29) 2,4-Dichlorophenol	10.458	162	64345	42.653	ng	95
30) 1,2,4-Trichlorobenzene	10.663	180	69559	41.877	ng	95
31) Naphthalene	10.851	128	193973	40.524	ng	99
32) Benzoic acid	10.129	122	24270	30.023	ng	83
33) 4-Chloroaniline	10.963	127	80262	42.472	ng	94
34) Hexachlorobutadiene	11.127	225	48208	39.938	ng	97
35) Caprolactam	11.744	113	20426	43.822	ng	# 79
36) 4-Chloro-3-methylphenol	12.102	107	76162	43.707	ng	# 92
37) 2-Methylnaphthalene	12.461	142	149444	41.445	ng	96
38) 1-Methylnaphthalene	12.684	142	141057	41.623	ng	98
40) 1,2,4,5-Tetrachloroben...	12.831	216	80361	42.163	ng	99
41) Hexachlorocyclopentadiene	12.807	237	37748	38.705	ng	98
43) 2,4,6-Trichlorophenol	13.078	196	52965	41.220	ng	91

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44) 2,4,5-Trichlorophenol	13.160	196	56958	40.860	ng	93
46) 1,1'-Biphenyl	13.471	154	186805	40.547	ng	99
47) 2-Chloronaphthalene	13.518	162	150397	40.871	ng	98
48) 2-Nitroaniline	13.724	65	53448	39.556	ng	92
49) Acenaphthylene	14.364	152	236101	40.482	ng	94
50) Dimethylphthalate	14.088	163	198892	40.753	ng	# 98
51) 2,6-Dinitrotoluene	14.217	165	45034	41.910	ng	87
52) Acenaphthene	14.705	154	144843	41.227	ng	96
53) 3-Nitroaniline	14.552	138	44745	42.699	ng	95
54) 2,4-Dinitrophenol	14.775	184	21951	35.983	ng	92
55) Dibenzofuran	15.040	168	230670	40.472	ng	98
56) 4-Nitrophenol	14.887	139	29062	37.962	ng	# 76
57) 2,4-Dinitrotoluene	15.010	165	67298	44.591	ng	# 94
58) Fluorene	15.692	166	191641	41.368	ng	94
59) 2,3,4,6-Tetrachlorophenol	15.275	232	48713	37.625	ng	96
60) Diethylphthalate	15.451	149	204497	41.662	ng	96
61) 4-Chlorophenyl-phenyle...	15.680	204	99380	40.193	ng	97
62) 4-Nitroaniline	15.721	138	48772	45.246	ng	# 82
63) Azobenzene	15.974	77	196999	37.919	ng	88
65) 4,6-Dinitro-2-methylph...	15.786	198	38740	41.746	ng	92
66) n-Nitrosodiphenylamine	15.897	169	169045	41.440	ng	95
67) 4-Bromophenyl-phenylether	16.573	248	71004	41.760	ng	96
68) Hexachlorobenzene	16.702	284	75377	39.902	ng	96
69) Atrazine	16.843	200	70729	43.345	ng	94
70) Pentachlorophenol	17.055	266	37107	36.633	ng	95
71) Phenanthrene	17.436	178	309282	40.719	ng	96
72) Anthracene	17.530	178	304616	40.836	ng	97
73) Carbazole	17.801	167	294000	41.614	ng	98
74) Di-n-butylphthalate	18.347	149	342731	41.615	ng	# 97
75) Fluoranthene	19.451	202	369219	39.503	ng	97
77) Benzidine	19.628	184	149775	46.703	ng	99
78) Pyrene	19.816	202	366856	42.992	ng	98
80) Butylbenzylphthalate	20.685	149	136943	42.139	ng	97
81) Benzo(a)anthracene	21.654	228	332931	40.845	ng	98
82) 3,3'-Dichlorobenzidine	21.566	252	97621	41.018	ng	100
83) Chrysene	21.719	228	321701	41.287	ng	98
84) Bis(2-ethylhexyl)phtha...	21.543	149	181930	39.549	ng	92
85) Di-n-octyl phthalate	22.753	149	297057	38.225	ng	99
87) Indeno(1,2,3-cd)pyrene	28.627	276	369018	40.107	ng	98
88) Benzo(b)fluoranthene	23.881	252	335233	40.064	ng	94
89) Benzo(k)fluoranthene	23.951	252	313494	40.166	ng	97
90) Benzo(a)pyrene	24.762	252	308748	40.773	ng	# 99
91) Dibenzo(a,h)anthracene	28.680	278	321466	41.462	ng	# 96
92) Benzo(g,h,i)perylene	29.796	276	307717	40.511	ng	99

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

