

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082121\
 Data File : BG049929.D
 Acq On : 21 Aug 2021 5:02
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040EC

Manual Integrations
 APPROVED

mohammad
 8/23/2021 4:29:09 PM

Quant Time: Aug 23 06:02:30 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG080321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Aug 21 05:13:22 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.992	152	36288	20.000	ng	0.01
21) Naphthalene-d8	10.812	136	147295	20.000	ng	# 0.01
39) Acenaphthene-d10	14.653	164	95214	20.000	ng	0.01
64) Phenanthrene-d10	17.408	188	176705	20.000	ng	# 0.02
76) Chrysene-d12	21.685	240	123786	20.000	ng	0.03
86) Perylene-d12	24.939	264	58941	20.000	ng	0.06
System Monitoring Compounds						
5) 2-Fluorophenol	5.548	112	143275	70.777	ng	0.02
7) Phenol-d6	7.169	99	222272	72.536	ng	0.03
23) Nitrobenzene-d5	9.167	82	223773	67.227	ng	0.01
42) 2,4,6-Tribromophenol	16.151	330	96866	69.214	ng	0.02
45) 2-Fluorobiphenyl	13.267	172	474036	72.443	ng	0.01
79) Terphenyl-d14	20.017	244	553287	81.179	ng	0.02
Target Compounds						Qvalue
2) 1,4-Dioxane	3.386	88	22800	25.774	ng	99
3) Pyridine	3.803	79	70169	28.938	ng	# 84
4) n-Nitrosodimethylamine	3.715	42	39083	29.975	ng	# 92
6) Aniline	7.316	93	114546	35.271	ng	95
8) 2-Chlorophenol	7.557	128	83549	37.856	ng	93
9) Benzaldehyde	7.122	77	64959	31.770	ng	92
10) Phenol	7.199	94	107766	36.296	ng	91
11) bis(2-Chloroethyl)ether	7.404	93	68268	34.053	ng	91
12) 1,3-Dichlorobenzene	7.880	146	93116	36.838	ng	98
13) 1,4-Dichlorobenzene	8.027	146	92264	36.060	ng	96
14) 1,2-Dichlorobenzene	8.350	146	87670	36.169	ng	95
15) Benzyl Alcohol	8.239	79	91565	35.581	ng	95
16) 2,2'-oxybis(1-Chloropr...	8.520	45	78497	34.460	ng	97
17) 2-Methylphenol	8.450	107	73398	37.540	ng	97
18) Hexachloroethane	9.073	117	27047	27.499	ng	98
19) n-Nitroso-di-n-propyla...	8.802	70	69213	33.753	ng	# 88
20) 3+4-Methylphenols	8.785	107	103613	37.733	ng	95
22) Acetophenone	8.820	105	131931	32.969	ng	# 97
24) Nitrobenzene	9.208	77	116454	33.155	ng	98
25) Isophorone	9.731	82	193627	32.599	ng	97
26) 2-Nitrophenol	9.924	139	18952	14.139	ng	96
27) 2,4-Dimethylphenol	9.989	122	72800	36.746	ng	95
28) bis(2-Chloroethoxy)met...	10.206	93	90200	34.715	ng	# 86
29) 2,4-Dichlorophenol	10.471	162	90811	37.353	ng	96
30) 1,2,4-Trichlorobenzene	10.676	180	98901	36.946	ng	95
31) Naphthalene	10.864	128	267593	34.689	ng	97
32) Benzoic acid	10.171	122	27908m	22.839	ng	
33) 4-Chloroaniline	10.976	127	108989	35.787	ng	96
34) Hexachlorobutadiene	11.140	225	71180	36.591	ng	97
35) Caprolactam	11.787	113	25362	33.763	ng	# 82
36) 4-Chloro-3-methylphenol	12.116	107	98373	35.030	ng	# 89
37) 2-Methylnaphthalene	12.474	142	203631	35.041	ng	94
38) 1-Methylnaphthalene	12.697	142	188996	34.605	ng	92
40) 1,2,4,5-Tetrachloroben...	12.844	216	108166	38.556	ng	96
43) 2,4,6-Trichlorophenol	13.091	196	74298	39.283	ng	94
44) 2,4,5-Trichlorophenol	13.179	196	77822	37.927	ng	90

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
46) 1,1'-Biphenyl	13.478	154	245876	36.257	ng	99
47) 2-Chloronaphthalene	13.531	162	199459	36.825	ng	98
48) 2-Nitroaniline	13.737	65	84918	42.697	ng	94
49) Acenaphthylene	14.377	152	301854	35.162	ng	98
50) Dimethylphthalate	14.101	163	241503	33.618	ng	99
51) 2,6-Dinitrotoluene	14.230	165	35060	22.166	ng	# 86
52) Acenaphthene	14.718	154	181570	35.110	ng	94
53) 3-Nitroaniline	14.559	138	68920	44.682	ng	93
55) Dibenzofuran	15.053	168	288717	34.414	ng	97
56) 4-Nitrophenol	14.900	139	29540	26.214	ng	89
57) 2,4-Dinitrotoluene	15.023	165	43467	19.566	ng	95
58) Fluorene	15.705	166	230851	33.855	ng	97
59) 2,3,4,6-Tetrachlorophenol	15.288	232	63462	33.301	ng	98
60) Diethylphthalate	15.458	149	239592	33.161	ng	97
61) 4-Chlorophenyl-phenyle...	15.687	204	121223	33.308	ng	95
62) 4-Nitroaniline	15.728	138	71085	44.802	ng	89
63) Azobenzene	15.981	77	237148	31.011	ng	87
66) n-Nitrosodiphenylamine	15.905	169	197679	39.592	ng	98
67) 4-Bromophenyl-phenylether	16.586	248	79872	38.380	ng	96
68) Hexachlorobenzene	16.715	284	86443	37.387	ng	95
69) Atrazine	16.856	200	71759	35.929	ng	98
70) Pentachlorophenol	17.074	266	23655	19.080	ng	94
71) Phenanthrene	17.450	178	325933	35.059	ng	96
72) Anthracene	17.544	178	324279	35.517	ng	99
73) Carbazole	17.814	167	301956	34.919	ng	99
74) Di-n-butylphthalate	18.354	149	330405	32.777	ng	99
75) Fluoranthene	19.464	202	336529	29.417	ng	99
77) Benzidine	19.641	184	114110	35.978	ng	98
78) Pyrene	19.829	202	334481	39.635	ng	97
80) Butylbenzylphthalate	20.698	149	123596	38.456	ng	97
81) Benzo(a)anthracene	21.667	228	270520	33.557	ng	98
82) 3,3'-Dichlorobenzidine	21.573	252	88001	37.387	ng	99
83) Chrysene	21.732	228	286363	37.161	ng	98
84) Bis(2-ethylhexyl)phtha...	21.550	149	174896	38.443	ng	99
85) Di-n-octyl phthalate	22.766	149	297944	38.766	ng	100
87) Indeno(1,2,3-cd)pyrene	28.670	276	86793m	20.418	ng	
88) Benzo(b)fluoranthene	23.906	252	149066	38.562	ng	# 96
89) Benzo(k)fluoranthene	23.970	252	190875	52.936	ng	# 96
90) Benzo(a)pyrene	24.787	252	145297	41.533	ng	# 96
91) Dibenzo(a,h)anthracene	28.723	278	73326	20.471	ng	# 97
92) Benzo(g,h,i)perylene	29.815	276	71323m	20.324	ng	

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

