

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112222\
 Data File : BG055741.D
 Acq On : 22 Nov 2022 15:44
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 11/23/2022
 Supervised By : Jagrut Upadhyay 11/23/2022

Quant Time: Nov 23 04:13:31 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 23 01:35:20 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.244	152	42605	20.000	ng	0.00
21) Naphthalene-d8	11.076	136	183421	20.000	ng	0.00
39) Acenaphthene-d10	14.872	164	134085	20.000	ng	0.00
64) Phenanthrene-d10	17.610	188	306950	20.000	ng	0.00
76) Chrysene-d12	21.911	240	268431	20.000	ng	0.00
86) Perylene-d12	25.318	264	285385	20.000	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.777	112	180924	76.757	ng	0.00
7) Phenol-d6	7.392	99	249084	79.385	ng	0.00
23) Nitrobenzene-d5	9.419	82	270369	77.910	ng	0.00
42) 2,4,6-Tribromophenol	16.358	330	142540	75.494	ng	0.01
45) 2-Fluorobiphenyl	13.497	172	675511	75.475	ng	0.00
79) Terphenyl-d14	20.195	244	1030719	76.801	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.603	88	40925	40.628	ng	98
3) Pyridine	4.020	79	118360	38.302	ng	98
4) n-Nitrosodimethylamine	3.932	42	62919	38.458	ng	98
6) Aniline	7.563	93	157189	39.885	ng	98
8) 2-Chlorophenol	7.810	128	106732	39.610	ng	98
9) Benzaldehyde	7.375	77	78598	36.885	ng	98
10) Phenol	7.416	94	139867	39.812	ng	100
11) bis(2-Chloroethyl)ether	7.651	93	106672	39.620	ng	100
12) 1,3-Dichlorobenzene	8.133	146	122553	38.684	ng	98
13) 1,4-Dichlorobenzene	8.285	146	123457	38.564	ng	98
14) 1,2-Dichlorobenzene	8.609	146	118423	38.603	ng	97
15) Benzyl Alcohol	8.479	79	107299	42.078	ng	99
16) 2,2'-oxybis(1-Chloropr...	8.767	45	195168	40.043	ng	98
17) 2-Methylphenol	8.685	107	102227	40.985	ng	97
18) Hexachloroethane	9.343	117	44324	40.227	ng	99
19) n-Nitroso-di-n-propyla...	9.049	70	98267	40.761	ng	99
20) 3+4-Methylphenols	9.014	107	145326	41.709	ng	98
22) Acetophenone	9.073	105	181044	37.985	ng	# 98
24) Nitrobenzene	9.466	77	139680	38.535	ng	99
25) Isophorone	9.983	82	261635	39.791	ng	99
26) 2-Nitrophenol	10.177	139	69300	41.414	ng	96
27) 2,4-Dimethylphenol	10.230	122	101454m	39.837	ng	
28) bis(2-Chloroethoxy)met...	10.459	93	139019	39.383	ng	99
29) 2,4-Dichlorophenol	10.718	162	120930	39.672	ng	99
30) 1,2,4-Trichlorobenzene	10.935	180	132950	37.542	ng	99
31) Naphthalene	11.129	128	378136	38.136	ng	99
32) Benzoic acid	10.371	122	76808	36.658	ng	97
33) 4-Chloroaniline	11.229	127	161827	39.564	ng	97
34) Hexachlorobutadiene	11.399	225	90365	37.369	ng	97
35) Caprolactam	11.999	113	40491	38.386	ng	98
36) 4-Chloro-3-methylphenol	12.334	107	135736	40.515	ng	96
37) 2-Methylnaphthalene	12.716	142	289911	39.050	ng	99
38) 1-Methylnaphthalene	12.933	142	281778	39.096	ng	99
40) 1,2,4,5-Tetrachloroben...	13.080	216	160788	38.159	ng	100
41) Hexachlorocyclopentadiene	13.050	237	93392	43.976	ng	95
43) 2,4,6-Trichlorophenol	13.309	196	113672	39.547	ng	100

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.385	196	124116	39.294	ng	99
46) 1,1'-Biphenyl	13.709	154	373334	38.208	ng	98
47) 2-Chloronaphthalene	13.756	162	293141	38.598	ng	98
48) 2-Nitroaniline	13.955	65	101018	40.457	ng	95
49) Acenaphthylene	14.596	152	482035	38.544	ng	100
50) Dimethylphthalate	14.314	163	407017	38.079	ng	98
51) 2,6-Dinitrotoluene	14.437	165	86676	39.659	ng	98
52) Acenaphthene	14.936	154	322947m	39.511	ng	
53) 3-Nitroaniline	14.772	138	94229	39.276	ng	96
54) 2,4-Dinitrophenol	14.978	184	57824	46.017	ng	98
55) Dibenzofuran	15.266	168	480687	38.121	ng	99
56) 4-Nitrophenol	15.072	139	76737	40.753	ng	100
57) 2,4-Dinitrotoluene	15.224	165	124076	40.267	ng	97
58) Fluorene	15.912	166	383399	38.070	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.489	232	114695	37.474	ng	97
60) Diethylphthalate	15.665	149	411132	38.044	ng	99
61) 4-Chlorophenyl-phenylether	15.894	204	209397	37.825	ng	96
62) 4-Nitroaniline	15.929	138	95486	38.074	ng	97
63) Azobenzene	16.188	77	366000	38.793	ng	99
65) 4,6-Dinitro-2-methylph...	15.988	198	75696	44.164	ng	96
66) n-Nitrosodiphenylamine	16.112	169	339209	40.099	ng	98
67) 4-Bromophenyl-phenylether	16.793	248	139412	39.862	ng	97
68) Hexachlorobenzene	16.917	284	150210	38.734	ng	99
69) Atrazine	17.046	200	127987	38.255	ng	99
70) Pentachlorophenol	17.257	266	90084	37.974	ng	98
71) Phenanthrene	17.657	178	628303	37.931	ng	100
72) Anthracene	17.745	178	616788	38.324	ng	100
73) Carbazole	18.009	167	546030	37.739	ng	100
74) Di-n-butylphthalate	18.544	149	672343	37.424	ng	100
75) Fluoranthene	19.649	202	707548	35.111	ng	99
77) Benzidine	19.819	184	214371	34.520	ng	99
78) Pyrene	20.013	202	713050	39.244	ng	99
80) Butylbenzylphthalate	20.877	149	279153	39.246	ng	99
81) Benzo(a)anthracene	21.887	228	703096	38.029	ng	100
82) 3,3'-Dichlorobenzidine	21.793	252	244174	37.585	ng	99
83) Chrysene	21.964	228	667995	37.952	ng	98
84) Bis(2-ethylhexyl)phtha...	21.764	149	394597	38.601	ng	99
85) Di-n-octyl phthalate	23.033	149	682842	39.033	ng	99
87) Indeno(1,2,3-cd)pyrene	29.249	276	842442	36.033	ng	# 94
88) Benzo(b)fluoranthene	24.231	252	699071	37.613	ng	99
89) Benzo(k)fluoranthene	24.302	252	687759	37.123	ng	99
90) Benzo(a)pyrene	25.160	252	594832	37.296	ng	99
91) Dibenzo(a,h)anthracene	29.314	278	695447	36.401	ng	99
92) Benzo(g,h,i)perylene	30.489	276	683835	35.513	ng	99

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

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