

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112122\
 Data File : BG055715.D
 Acq On : 21 Nov 2022 14:49
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040EC

Manual Integrations
 APPROVED

Reviewed By : Christian
 Giraldo

Quant Time: Nov 22 02:03:25 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 17 02:02:46 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.250	152	35982	20.000	ng	0.00
21) Naphthalene-d8	11.082	136	159830	20.000	ng	0.00
39) Acenaphthene-d10	14.871	164	119840	20.000	ng	0.00
64) Phenanthrene-d10	17.615	188	276466	20.000	ng	0.00
76) Chrysene-d12	21.910	240	249860	20.000	ng	0.00
86) Perylene-d12	25.324	264	267262	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.776	112	147372	74.031	ng	0.00
7) Phenol-d6	7.398	99	208406	78.646	ng	0.00
23) Nitrobenzene-d5	9.425	82	227150	75.117	ng	0.00
42) 2,4,6-Tribromophenol	16.358	330	125256	74.226	ng	0.00
45) 2-Fluorobiphenyl	13.496	172	583344	72.924	ng	0.00
79) Terphenyl-d14	20.200	244	940979	75.325	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.614	88	29969	35.227	ng	98
3) Pyridine	4.025	79	100898	38.661	ng	99
4) n-Nitrosodimethylamine	3.937	42	53809	38.943	ng	98
6) Aniline	7.562	93	132901	39.929	ng	97
8) 2-Chlorophenol	7.809	128	87553	38.473	ng	96
9) Benzaldehyde	7.374	77	65881	36.608	ng	99
10) Phenol	7.421	94	115956	39.081	ng	99
11) bis(2-Chloroethyl)ether	7.656	93	88639	38.982	ng	97
12) 1,3-Dichlorobenzene	8.138	146	98616	36.858	ng	99
13) 1,4-Dichlorobenzene	8.285	146	101100	37.393	ng	97
14) 1,2-Dichlorobenzene	8.608	146	97235	37.530	ng	98
15) Benzyl Alcohol	8.485	79	89770	41.683	ng	99
16) 2,2'-oxybis(1-Chloropr...	8.773	45	161826	39.313	ng	99
17) 2-Methylphenol	8.684	107	86072	40.860	ng	99
18) Hexachloroethane	9.343	117	35318	37.953	ng	98
19) n-Nitroso-di-n-propyla...	9.055	70	85478	41.983	ng	97
20) 3+4-Methylphenols	9.014	107	122000	41.459	ng	98
22) Acetophenone	9.078	105	154195	37.127	ng	99
24) Nitrobenzene	9.466	77	117137	37.085	ng	99
25) Isophorone	9.989	82	222293	38.797	ng	99
26) 2-Nitrophenol	10.183	139	58484	40.109	ng	93
27) 2,4-Dimethylphenol	10.230	122	85442m	38.502	ng	
28) bis(2-Chloroethoxy)met...	10.465	93	115780	37.641	ng	99
29) 2,4-Dichlorophenol	10.717	162	102412	38.556	ng	98
30) 1,2,4-Trichlorobenzene	10.935	180	114021	36.949	ng	97
31) Naphthalene	11.129	128	316337	36.612	ng	99
32) Benzoic acid	10.371	122	71468	39.143	ng	95
33) 4-Chloroaniline	11.229	127	135498	38.017	ng	99
34) Hexachlorobutadiene	11.405	225	78403	37.208	ng	97
35) Caprolactam	12.004	113	34562	37.601	ng	99
36) 4-Chloro-3-methylphenol	12.333	107	114375	39.178	ng	99
37) 2-Methylnaphthalene	12.715	142	244232	37.753	ng	99
38) 1-Methylnaphthalene	12.932	142	237838	37.871	ng	99
40) 1,2,4,5-Tetrachloroben...	13.079	216	140130	37.209	ng	100
41) Hexachlorocyclopentadiene	13.056	237	70811	37.306	ng	98
43) 2,4,6-Trichlorophenol	13.314	196	98093	38.184	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
44) 2,4,5-Trichlorophenol	13.385	196	109271	38.706	ng	98	11/22/2022
46) 1,1'-Biphenyl	13.708	154	319625	36.600	ng	99	Supervised By :Jagrut Upadhyay
47) 2-Chloronaphthalene	13.755	162	249708	36.788	ng	99	
48) 2-Nitroaniline	13.955	65	85348	38.244	ng	97	
49) Acenaphthylene	14.595	152	410602	36.734	ng	99	
50) Dimethylphthalate	14.313	163	355031	37.164	ng	99	
51) 2,6-Dinitrotoluene	14.437	165	72839	37.290	ng	98	11/22/2022
52) Acenaphthene	14.936	154	271052m	37.104	ng		
53) 3-Nitroaniline	14.771	138	79207	36.939	ng	99	
54) 2,4-Dinitrophenol	14.977	184	43884	39.075	ng	98	
55) Dibenzofuran	15.265	168	408842	36.278	ng	99	
56) 4-Nitrophenol	15.071	139	60242	35.796	ng	97	
57) 2,4-Dinitrotoluene	15.224	165	105322	38.243	ng	93	
58) Fluorene	15.917	166	326532	36.277	ng	98	
59) 2,3,4,6-Tetrachlorophenol	15.488	232	103154	37.710	ng	100	
60) Diethylphthalate	15.665	149	350200	36.258	ng	99	
61) 4-Chlorophenyl-phenyle...	15.900	204	181042	36.590	ng	99	
62) 4-Nitroaniline	15.929	138	81751	36.472	ng	97	
63) Azobenzene	16.193	77	303900	36.040	ng	98	
65) 4,6-Dinitro-2-methylph...	15.988	198	64154	41.557	ng	97	
66) n-Nitrosodiphenylamine	16.111	169	288102	37.813	ng	98	
67) 4-Bromophenyl-phenylether	16.793	248	120430	38.232	ng	98	
68) Hexachlorobenzene	16.916	284	133226	38.142	ng	100	
69) Atrazine	17.051	200	111655	37.054	ng	97	
70) Pentachlorophenol	17.263	266	78363	36.675	ng	99	
71) Phenanthrene	17.656	178	544755	36.513	ng	99	
72) Anthracene	17.750	178	531356	36.656	ng	99	
73) Carbazole	18.009	167	467651	35.886	ng	100	
74) Di-n-butylphthalate	18.549	149	578869	35.774	ng	99	
75) Fluoranthene	19.654	202	620963	34.212	ng	100	
77) Benzidine	19.824	184	202322	35.001	ng	98	
78) Pyrene	20.012	202	628416	37.157	ng	99	
80) Butylbenzylphthalate	20.882	149	247434	37.372	ng	97	
81) Benzo(a)anthracene	21.893	228	630514	36.638	ng	100	
82) 3,3'-Dichlorobenzidine	21.793	252	224303	37.093	ng	99	
83) Chrysene	21.963	228	598863	36.554	ng	98	
84) Bis(2-ethylhexyl)phtha...	21.763	149	352977	37.096	ng	99	
85) Di-n-octyl phthalate	23.038	149	601906	36.964	ng	100	
87) Indeno(1,2,3-cd)pyrene	29.260	276	753289	34.405	ng	# 95	
88) Benzo(b)fluoranthene	24.231	252	614619	35.312	ng	99	
89) Benzo(k)fluoranthene	24.307	252	634253	36.556	ng	100	
90) Benzo(a)pyrene	25.165	252	536134	35.895	ng	100	
91) Dibenzo(a,h)anthracene	29.325	278	627218	35.056	ng	99	
92) Benzo(g,h,i)perylene	30.494	276	612712	33.978	ng	99	

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

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