

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG051822\
 Data File : BG053601.D
 Acq On : 18 May 2022 12:15
 Operator : CG/JU
 Sample : PB144886BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB144886BS

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 05/19/2022
 Supervised By : Jagrut Upadhyay 05/19/2022

Quant Time: May 18 14:04:50 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG050522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 11 06:00:46 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.076	152	21802	20.000	ng	-0.01
21) Naphthalene-d8	10.890	136	101132	20.000	ng	# 0.00
39) Acenaphthene-d10	14.708	164	83116	20.000	ng	0.00
64) Phenanthrene-d10	17.452	188	209762	20.000	ng	0.00
76) Chrysene-d12	21.734	240	197619	20.000	ng	0.00
86) Perylene-d12	25.012	264	202827	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.650	112	184583	143.685	ng	0.00
7) Phenol-d6	7.248	99	271430	139.221	ng	0.00
23) Nitrobenzene-d5	9.245	82	188607	79.388	ng	0.00
42) 2,4,6-Tribromophenol	16.200	330	152295	124.235	ng	0.00
45) 2-Fluorobiphenyl	13.328	172	438022	77.010	ng	0.00
79) Terphenyl-d14	20.054	244	914381	86.176	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.518	88	18368	27.251	ng	94
3) Pyridine	3.929	79	49231	29.908	ng	97
4) n-Nitrosodimethylamine	3.835	42	33104	41.284	ng	98
6) Aniline	7.406	93	97973	45.371	ng	99
8) 2-Chlorophenol	7.647	128	66676	49.648	ng	96
9) Benzaldehyde	7.213	77	43228	34.901	ng	98
10) Phenol	7.271	94	94038	49.317	ng	92
11) bis(2-Chloroethyl)ether	7.489	93	63062	44.344	ng	96
12) 1,3-Dichlorobenzene	7.970	146	66865	42.205	ng	91
13) 1,4-Dichlorobenzene	8.117	146	68337	42.744	ng	97
14) 1,2-Dichlorobenzene	8.434	146	66602	44.363	ng	96
15) Benzyl Alcohol	8.317	79	72985	44.777	ng	99
16) 2,2'-oxybis(1-Chloropr...	8.605	45	101901	44.026	ng	95
17) 2-Methylphenol	8.528	107	65064	50.485	ng	95
18) Hexachloroethane	9.163	117	25162	41.445	ng	92
19) n-Nitroso-di-n-propyla...	8.881	70	63570	45.445	ng	99
20) 3+4-Methylphenols	8.857	107	88615	48.685	ng	99
22) Acetophenone	8.899	105	108194	39.090	ng	# 97
24) Nitrobenzene	9.286	77	95498	41.531	ng	97
25) Isophorone	9.809	82	176751	44.207	ng	97
26) 2-Nitrophenol	10.003	139	40721	43.724	ng	93
27) 2,4-Dimethylphenol	10.056	122	69617	49.671	ng	96
28) bis(2-Chloroethoxy)met...	10.285	93	92974	40.432	ng	# 96
29) 2,4-Dichlorophenol	10.543	162	76992	45.516	ng	94
30) 1,2,4-Trichlorobenzene	10.755	180	77567	40.359	ng	97
31) Naphthalene	10.943	128	209708	40.765	ng	98
32) Benzoic acid	10.220	122	27077	33.928	ng	86
33) 4-Chloroaniline	11.054	127	57384	26.636	ng	96
34) Hexachlorobutadiene	11.225	225	56743	39.816	ng	97
35) Caprolactam	11.818	113	27415m	44.598	ng	
36) 4-Chloro-3-methylphenol	12.171	107	90282	44.619	ng	99
37) 2-Methylnaphthalene	12.541	142	156600	40.791	ng	99
38) 1-Methylnaphthalene	12.758	142	151219	40.372	ng	98
40) 1,2,4,5-Tetrachloroben...	12.911	216	102816	39.244	ng	99
41) Hexachlorocyclopentadiene	12.887	237	91325	84.962	ng	94
43) 2,4,6-Trichlorophenol	13.152	196	72174	42.203	ng	96

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.234	196	77238	42.436	ng	99
46) 1,1'-Biphenyl	13.539	154	222824	39.330	ng	97
47) 2-Chloronaphthalene	13.586	162	174744	39.368	ng	99
48) 2-Nitroaniline	13.792	65	72717	43.965	ng	99
49) Acenaphthylene	14.432	152	303768	42.886	ng	99
50) Dimethylphthalate	14.156	163	255945	40.523	ng	99
51) 2,6-Dinitrotoluene	14.280	165	55911	43.450	ng	96
52) Acenaphthene	14.773	154	171336	40.590	ng	99
53) 3-Nitroaniline	14.614	138	48450	34.847	ng	98
54) 2,4-Dinitrophenol	14.826	184	57431	71.068	ng	98
55) Dibenzofuran	15.108	168	297564	39.246	ng	98
56) 4-Nitrophenol	14.932	139	82834	83.413	ng	89
57) 2,4-Dinitrotoluene	15.073	165	82476	44.218	ng	# 93
58) Fluorene	15.754	166	246321	40.483	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.337	232	74945	40.583	ng	99
60) Diethylphthalate	15.513	149	267495	41.000	ng	99
61) 4-Chlorophenyl-phenyle...	15.742	204	138186	40.700	ng	96
62) 4-Nitroaniline	15.777	138	59709	41.405	ng	92
63) Azobenzene	16.036	77	268191	40.143	ng	98
65) 4,6-Dinitro-2-methylph...	15.842	198	49964	39.974	ng	92
66) n-Nitrosodiphenylamine	15.954	169	224344	40.735	ng	99
67) 4-Bromophenyl-phenylether	16.635	248	99170	40.481	ng	99
68) Hexachlorobenzene	16.764	284	107799	40.179	ng	100
69) Atrazine	16.899	200	96325	41.546	ng	96
70) Pentachlorophenol	17.111	266	105820	73.632	ng	94
71) Phenanthrene	17.499	178	433767	40.668	ng	97
72) Anthracene	17.587	178	433130	41.314	ng	99
73) Carbazole	17.857	167	397737	38.710	ng	100
74) Di-n-butylphthalate	18.403	149	461085	39.901	ng	99
75) Fluoranthene	19.502	202	553024	40.044	ng	99
77) Benzidine	19.678	184	248360	58.308	ng	98
78) Pyrene	19.866	202	545002	42.602	ng	100
80) Butylbenzylphthalate	20.735	149	201645	42.985	ng	98
81) Benzo(a)anthracene	21.716	228	531176	42.142	ng	98
82) 3,3'-Dichlorobenzidine	21.622	252	156327	48.908	ng	99
83) Chrysene	21.781	228	487525	41.413	ng	98
84) Bis(2-ethylhexyl)phtha...	21.605	149	282907	43.694	ng	98
85) Di-n-octyl phthalate	22.827	149	477510	43.613	ng	100
87) Indeno(1,2,3-cd)pyrene	28.772	276	583854	41.331	ng	# 95
88) Benzo(b)fluoranthene	23.972	252	555224	46.247	ng	99
89) Benzo(k)fluoranthene	24.037	252	508454	43.098	ng	98
90) Benzo(a)pyrene	24.859	252	521029	50.988	ng	97
91) Dibenzo(a,h)anthracene	28.825	278	504404	43.333	ng	98
92) Benzo(g,h,i)perylene	29.947	276	506866	43.929	ng	98

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

