

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040423\
 Data File : BG056970.D
 Acq On : 4 Apr 2023 18:10
 Operator : CG/JU
 Sample : PB151693BSD
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB151693BSD

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 04/05/2023
 Supervised By : Jagrut Upadhyay 04/05/2023

Quant Time: Apr 05 04:09:25 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG033023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 31 03:23:11 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.413	152	14182	20.000	ng	0.00
21) Naphthalene-d8	11.256	136	54031	20.000	ng	# 0.00
39) Acenaphthene-d10	15.028	164	41787	20.000	ng	0.00
64) Phenanthrene-d10	17.765	188	108329	20.000	ng	0.00
76) Chrysene-d12	22.089	240	102084	20.000	ng	0.00
86) Perylene-d12	25.625	264	108432	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.922	112	123161	138.988	ng	0.00
7) Phenol-d6	7.544	99	177411	133.863	ng	0.00
23) Nitrobenzene-d5	9.594	82	114832	86.962	ng	0.00
42) 2,4,6-Tribromophenol	16.508	330	96023	143.842	ng	0.00
45) 2-Fluorobiphenyl	13.653	172	264424	85.807	ng	0.00
79) Terphenyl-d14	20.332	244	598206	99.160	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.725	88	13854	35.665	ng	# 92
3) Pyridine	4.148	79	39511	31.580	ng	93
4) n-Nitrosodimethylamine	4.054	42	25086	43.610	ng	# 91
6) Aniline	7.720	93	54864	32.382	ng	97
8) 2-Chlorophenol	7.967	128	41424	46.932	ng	93
9) Benzaldehyde	7.538	77	30755	36.998	ng	94
10) Phenol	7.573	94	57604	43.865	ng	91
11) bis(2-Chloroethyl)ether	7.814	93	37822	40.433	ng	99
12) 1,3-Dichlorobenzene	8.301	146	45228	46.375	ng	96
13) 1,4-Dichlorobenzene	8.448	146	46074	46.809	ng	97
14) 1,2-Dichlorobenzene	8.777	146	43246	45.476	ng	96
15) Benzyl Alcohol	8.642	79	47114	42.783	ng	97
16) 2,2'-oxybis(1-Chloropr...	8.936	45	48010m	42.586	ng	
17) 2-Methylphenol	8.842	107	39135	41.771	ng	93
18) Hexachloroethane	9.517	117	15666	43.545	ng	# 84
19) n-Nitroso-di-n-propyla...	9.218	70	37233	39.277	ng	95
20) 3+4-Methylphenols	9.171	107	51869	39.592	ng	96
22) Acetophenone	9.241	105	66811	41.555	ng	# 97
24) Nitrobenzene	9.635	77	58238	43.138	ng	98
25) Isophorone	10.158	82	99515	39.576	ng	99
26) 2-Nitrophenol	10.352	139	23095	44.320	ng	96
27) 2,4-Dimethylphenol	10.393	122	41948	45.593	ng	99
28) bis(2-Chloroethoxy)met...	10.628	93	49744	39.515	ng	94
29) 2,4-Dichlorophenol	10.886	162	44210	44.680	ng	98
30) 1,2,4-Trichlorobenzene	11.115	180	46742	44.459	ng	98
31) Naphthalene	11.303	128	123299	41.866	ng	99
32) Benzoic acid	10.510	122	27732m	43.880	ng	
33) 4-Chloroaniline	11.397	127	34101	25.638	ng	93
34) Hexachlorobutadiene	11.579	225	35331	44.737	ng	93
35) Caprolactam	12.161	113	14009m	41.054	ng	
36) 4-Chloro-3-methylphenol	12.490	107	48257	43.610	ng	97
37) 2-Methylnaphthalene	12.878	142	92698	39.659	ng	99
38) 1-Methylnaphthalene	13.095	142	85871	39.125	ng	91
40) 1,2,4,5-Tetrachloroben...	13.242	216	63685	43.949	ng	# 99
41) Hexachlorocyclopentadiene	13.218	237	92608	115.993	ng	95
43) 2,4,6-Trichlorophenol	13.465	196	40894	44.224	ng	94

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.541	196	44927	41.012	ng	96
46) 1,1'-Biphenyl	13.864	154	135305	44.233	ng	96
47) 2-Chloronaphthalene	13.917	162	100505	44.425	ng	97
48) 2-Nitroaniline	14.105	65	37117	45.882	ng	92
49) Acenaphthylene	14.757	152	159651	42.950	ng	98
50) Dimethylphthalate	14.464	163	139624	42.406	ng	# 98
51) 2,6-Dinitrotoluene	14.587	165	30923	43.469	ng	98
52) Acenaphthene	15.092	154	129312m	50.441	ng	
53) 3-Nitroaniline	14.922	138	22398	31.809	ng	92
54) 2,4-Dinitrophenol	15.122	184	44261	110.539	ng	# 88
55) Dibenzofuran	15.421	168	166076	43.162	ng	99
56) 4-Nitrophenol	15.210	139	54973	105.578	ng	94
57) 2,4-Dinitrotoluene	15.374	165	44332	44.185	ng	99
58) Fluorene	16.067	166	140447	44.112	ng	97
59) 2,3,4,6-Tetrachlorophenol	15.639	232	46489	46.610	ng	# 99
60) Diethylphthalate	15.809	149	148024	44.846	ng	99
61) 4-Chlorophenyl-phenyle...	16.044	204	80616	43.125	ng	97
62) 4-Nitroaniline	16.079	138	32524	44.581	ng	97
63) Azobenzene	16.338	77	143686	43.681	ng	97
65) 4,6-Dinitro-2-methylph...	16.132	198	32062	48.869	ng	97
66) n-Nitrosodiphenylamine	16.261	169	121595	41.814	ng	99
67) 4-Bromophenyl-phenylether	16.943	248	55792	41.624	ng	98
68) Hexachlorobenzene	17.072	284	59550	44.938	ng	95
69) Atrazine	17.195	200	57948	48.005	ng	99
70) Pentachlorophenol	17.413	266	81411	85.633	ng	97
71) Phenanthrene	17.812	178	241600	43.743	ng	99
72) Anthracene	17.900	178	249771	44.114	ng	99
73) Carbazole	18.165	167	227947	46.176	ng	98
74) Di-n-butylphthalate	18.687	149	266061	46.544	ng	99
75) Fluoranthene	19.798	202	317362	45.328	ng	97
77) Benzidine	19.962	184	166890	59.811	ng	98
78) Pyrene	20.156	202	318020	45.041	ng	99
80) Butylbenzylphthalate	21.019	149	111059	45.709	ng	97
81) Benzo(a)anthracene	22.065	228	312804	44.280	ng	99
82) 3,3'-Dichlorobenzidine	21.965	252	78556	32.932	ng	96
83) Chrysene	22.142	228	286530	43.320	ng	99
84) Bis(2-ethylhexyl)phtha...	21.918	149	156745	44.552	ng	99
85) Di-n-octyl phthalate	23.240	149	259025	43.688	ng	99
87) Indeno(1,2,3-cd)pyrene	29.714	276	352721	44.015	ng	# 96
88) Benzo(b)fluoranthene	24.497	252	326674	48.178	ng	99
89) Benzo(k)fluoranthene	24.568	252	310861	45.923	ng	99
90) Benzo(a)pyrene	25.466	252	280552	42.106	ng	98
91) Dibenzo(a,h)anthracene	29.790	278	292292	44.374	ng	97
92) Benzo(g,h,i)perylene	31.006	276	283353	43.931	ng	99

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

