

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF051523\
 Data File : BF133384.D
 Acq On : 15 May 2023 13:58
 Operator : CG\JU
 Sample : PB152781BSD
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB152781BSD

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 05/16/2023
 Supervised By : Yogesh Patel 05/16/2023

Quant Time: May 15 14:46:35 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 11 16:40:17 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.722	152	236418	20.000	ng	0.00
21) Naphthalene-d8	8.010	136	980297	20.000	ng	0.00
39) Acenaphthene-d10	9.763	164	505029	20.000	ng	0.00
64) Phenanthrene-d10	11.245	188	886266	20.000	ng	0.00
76) Chrysene-d12	13.886	240	532605	20.000	ng	0.00
86) Perylene-d12	15.304	264	466293	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.351	112	1915692	122.405	ng	0.02
7) Phenol-d6	6.375	99	2287201	117.742	ng	0.00
23) Nitrobenzene-d5	7.292	82	1489898	82.036	ng	0.00
42) 2,4,6-Tribromophenol	10.557	330	604117	118.943	ng	0.00
45) 2-Fluorobiphenyl	9.086	172	2463043	81.593	ng	0.00
79) Terphenyl-d14	12.833	244	2698110	87.369	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.481	88	276674	38.113	ng	96
3) Pyridine	3.169	79	723831	37.542	ng	99
4) n-Nitrosodimethylamine	3.140	42	403479	40.402	ng	92
6) Aniline	6.387	93	784421	31.352	ng	# 19
8) 2-Chlorophenol	6.510	128	781222	49.163	ng	95
9) Benzaldehyde	6.269	77	424130	52.674	ng	98
10) Phenol	6.387	94	891723	41.611	ng	# 72
11) bis(2-Chloroethyl)ether	6.457	93	740513	46.048	ng	99
12) 1,3-Dichlorobenzene	6.663	146	806129	47.716	ng	97
13) 1,4-Dichlorobenzene	6.740	146	808134	47.729	ng	99
14) 1,2-Dichlorobenzene	6.892	146	779433	49.931	ng	97
15) Benzyl Alcohol	6.875	79	629260	46.895	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.004	45	1131456	40.509	ng	83
17) 2-Methylphenol	6.992	107	631340	43.390	ng	95
18) Hexachloroethane	7.234	117	285287	48.463	ng	99
19) n-Nitroso-di-n-propyla...	7.151	70	492473	40.722	ng	93
20) 3+4-Methylphenols	7.145	107	760027	44.354	ng	# 81
22) Acetophenone	7.140	105	1055518	46.482	ng	95
24) Nitrobenzene	7.310	77	790831	42.972	ng	98
25) Isophorone	7.551	82	1479584	42.908	ng	99
26) 2-Nitrophenol	7.628	139	442150	49.811	ng	94
27) 2,4-Dimethylphenol	7.675	122	693986	45.595	ng	98
28) bis(2-Chloroethoxy)met...	7.763	93	888547	43.558	ng	99
29) 2,4-Dichlorophenol	7.875	162	643269	46.425	ng	98
30) 1,2,4-Trichlorobenzene	7.951	180	666833	46.453	ng	99
31) Naphthalene	8.034	128	2132124	43.502	ng	99
32) Benzoic acid	7.822	122	433869	46.148	ng	99
33) 4-Chloroaniline	8.081	127	337629m	17.050	ng	
34) Hexachlorobutadiene	8.145	225	357609	44.948	ng	98
35) Caprolactam	8.469	113	232308m	49.907	ng	
36) 4-Chloro-3-methylphenol	8.581	107	696794	46.633	ng	97
37) 2-Methylnaphthalene	8.722	142	1411616	42.128	ng	98
38) 1-Methylnaphthalene	8.822	142	1310129	41.795	ng	100
40) 1,2,4,5-Tetrachloroben...	8.892	216	592268	43.617	ng	99
41) Hexachlorocyclopentadiene	8.875	237	676951	85.482	ng	99
43) 2,4,6-Trichlorophenol	9.004	196	420466	44.775	ng	96

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.051	196	462293	42.566	ng	95
46) 1,1'-Biphenyl	9.186	154	1724200	43.055	ng	99
47) 2-Chloronaphthalene	9.210	162	1322079	45.105	ng	99
48) 2-Nitroaniline	9.310	65	419486	43.303	ng	98
49) Acenaphthylene	9.628	152	2001623	42.738	ng	99
50) Dimethylphthalate	9.492	163	1590076	45.154	ng	99
51) 2,6-Dinitrotoluene	9.557	165	364577	45.998	ng	# 83
52) Acenaphthene	9.798	154	1463918m	46.669	ng	
53) 3-Nitroaniline	9.722	138	256183	27.184	ng	97
54) 2,4-Dinitrophenol	9.833	184	401884	100.876	ng	97
55) Dibenzofuran	9.969	168	1748076	40.854	ng	97
56) 4-Nitrophenol	9.904	139	574291	78.679	ng	98
57) 2,4-Dinitrotoluene	9.963	165	455091	45.024	ng	93
58) Fluorene	10.316	166	1333445	42.536	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.092	232	374823	44.517	ng	98
60) Diethylphthalate	10.192	149	1439309	43.271	ng	100
61) 4-Chlorophenyl-phenyle...	10.304	204	640870	41.950	ng	97
62) 4-Nitroaniline	10.345	138	408819	47.199	ng	96
63) Azobenzene	10.463	77	1434357	40.707	ng	99
65) 4,6-Dinitro-2-methylph...	10.375	198	281708	52.441	ng	94
66) n-Nitrosodiphenylamine	10.428	169	1216444	42.314	ng	99
67) 4-Bromophenyl-phenylether	10.792	248	414035	43.074	ng	97
68) Hexachlorobenzene	10.863	284	446832	45.367	ng	99
69) Atrazine	10.957	200	430304	51.946	ng	98
70) Pentachlorophenol	11.057	266	496825	70.827	ng	98
71) Phenanthrene	11.275	178	1998765	41.961	ng	99
72) Anthracene	11.327	178	2089787	42.869	ng	100
73) Carbazole	11.480	167	1914648	44.109	ng	99
74) Di-n-butylphthalate	11.810	149	2295539	46.732	ng	100
75) Fluoranthene	12.457	202	2046871	42.816	ng	99
77) Benzidine	12.580	184	1016098	112.817	ng	# 93
78) Pyrene	12.686	202	2115766	46.341	ng	99
80) Butylbenzylphthalate	13.304	149	915577	56.637	ng	94
81) Benzo(a)anthracene	13.874	228	1546772	42.232	ng	99
82) 3,3'-Dichlorobenzidine	13.839	252	357432	35.327	ng	99
83) Chrysene	13.916	228	1558162	42.553	ng	99
84) Bis(2-ethylhexyl)phtha...	13.868	149	1050971	51.795	ng	99
85) Di-n-octyl phthalate	14.480	149	1707138	51.598	ng	96
87) Indeno(1,2,3-cd)pyrene	16.692	276	1429483	53.895	ng	99
88) Benzo(b)fluoranthene	14.904	252	1377592	46.438	ng	100
89) Benzo(k)fluoranthene	14.933	252	1237069	42.068	ng	99
90) Benzo(a)pyrene	15.251	252	1146061	42.445	ng	98
91) Dibenzo(a,h)anthracene	16.709	278	1178548	53.275	ng	98
92) Benzo(g,h,i)perylene	17.115	276	1215917	55.674	ng	97

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

