

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF050522\
 Data File : BF128093.D
 Acq On : 05 May 2022 14:33
 Operator : CG\JU
 Sample : PB144601BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB144601BS

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 05/06/2022
 Supervised By : mohammad ahmed 05/09/2022

Quant Time: May 05 18:00:08 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 05 17:56:23 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.904	152	104912	20.000	ng	0.00
21) Naphthalene-d8	8.192	136	416268	20.000	ng	0.00
39) Acenaphthene-d10	9.951	164	236591	20.000	ng	0.00
64) Phenanthrene-d10	11.445	188	442412	20.000	ng	# 0.00
76) Chrysene-d12	14.092	240	310461	20.000	ng	0.00
86) Perylene-d12	15.592	264	244094	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.546	112	715702	108.659	ng	0.01
7) Phenol-d6	6.545	99	929216	108.722	ng	0.00
23) Nitrobenzene-d5	7.475	82	678323	76.627	ng	0.00
42) 2,4,6-Tribromophenol	10.745	330	288799	121.258	ng	0.00
45) 2-Fluorobiphenyl	9.269	172	1104487	78.976	ng	0.00
79) Terphenyl-d14	13.027	244	1360429	80.127	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.834	88	90653	28.846	ng	98
3) Pyridine	3.587	79	246473	30.609	ng	97
4) n-Nitrosodimethylamine	3.552	42	154499	39.639	ng	99
6) Aniline	6.575	93	375461	36.975	ng	100
8) 2-Chlorophenol	6.693	128	278373	40.720	ng	99
9) Benzaldehyde	6.463	77	183266	34.662	ng	96
10) Phenol	6.557	94	344619m	39.979	ng	
11) bis(2-Chloroethyl)ether	6.645	93	280596	39.241	ng	98
12) 1,3-Dichlorobenzene	6.845	146	299760	38.093	ng	95
13) 1,4-Dichlorobenzene	6.928	146	304813	38.806	ng	99
14) 1,2-Dichlorobenzene	7.075	146	293604	40.378	ng	99
15) Benzyl Alcohol	7.051	79	272805	46.938	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.175	45	410015	37.517	ng	99
17) 2-Methylphenol	7.157	107	229054	38.833	ng	97
18) Hexachloroethane	7.416	117	116462	40.331	ng	95
19) n-Nitroso-di-n-propyla...	7.322	70	211336	40.394	ng	98
20) 3+4-Methylphenols	7.310	107	274317	38.498	ng	90
22) Acetophenone	7.322	105	389210	38.362	ng	97
24) Nitrobenzene	7.492	77	333730	39.124	ng	99
25) Isophorone	7.734	82	600036	40.274	ng	99
26) 2-Nitrophenol	7.810	139	147837	40.672	ng	92
27) 2,4-Dimethylphenol	7.840	122	248512	41.161	ng	96
28) bis(2-Chloroethoxy)met...	7.939	93	343322	35.956	ng	97
29) 2,4-Dichlorophenol	8.051	162	231872	36.659	ng	98
30) 1,2,4-Trichlorobenzene	8.128	180	262310	36.509	ng	99
31) Naphthalene	8.216	128	785096	37.031	ng	100
32) Benzoic acid	7.975	122	133675	30.890	ng	95
33) 4-Chloroaniline	8.263	127	251140	29.938	ng	99
34) Hexachlorobutadiene	8.322	225	174076	38.465	ng	97
35) Caprolactam	8.639	113	76350m	37.395	ng	
36) 4-Chloro-3-methylphenol	8.745	107	255447	37.815	ng	98
37) 2-Methylnaphthalene	8.904	142	531305	37.919	ng	97
38) 1-Methylnaphthalene	9.004	142	504842	37.068	ng	100
40) 1,2,4,5-Tetrachloroben...	9.069	216	260777	37.395	ng	98
41) Hexachlorocyclopentadiene	9.051	237	311993	82.577	ng	99
43) 2,4,6-Trichlorophenol	9.181	196	169213	36.877	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.228	196	181023	36.291	ng	95
46) 1,1'-Biphenyl	9.369	154	654225	37.330	ng	99
47) 2-Chloronaphthalene	9.398	162	484128	36.287	ng	99
48) 2-Nitroaniline	9.498	65	178696	39.596	ng	97
49) Acenaphthylene	9.816	152	782968	38.366	ng	100
50) Dimethylphthalate	9.669	163	581069	36.621	ng	99
51) 2,6-Dinitrotoluene	9.739	165	136514	39.737	ng	94
52) Acenaphthene	9.986	154	566719m	41.305	ng	
53) 3-Nitroaniline	9.910	138	112406	29.465	ng	97
54) 2,4-Dinitrophenol	10.022	184	149324	82.535	ng	91
55) Dibenzofuran	10.157	168	712265	35.982	ng	98
56) 4-Nitrophenol	10.075	139	197665	67.855	ng	91
57) 2,4-Dinitrotoluene	10.145	165	182441	40.092	ng	# 89
58) Fluorene	10.504	166	558288	37.796	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.275	232	152825	36.434	ng	98
60) Diethylphthalate	10.369	149	581052	37.345	ng	100
61) 4-Chlorophenyl-phenyle...	10.492	204	272752	37.159	ng	99
62) 4-Nitroaniline	10.528	138	141051	37.819	ng	95
63) Azobenzene	10.651	77	624670	36.486	ng	99
65) 4,6-Dinitro-2-methylph...	10.551	198	102650	39.049	ng	89
66) n-Nitrosodiphenylamine	10.610	169	491391	35.696	ng	100
67) 4-Bromophenyl-phenylether	10.980	248	178639	36.424	ng	99
68) Hexachlorobenzene	11.051	284	199750	38.657	ng	# 88
69) Atrazine	11.139	200	178646	41.968	ng	99
70) Pentachlorophenol	11.245	266	200090	65.822	ng	98
71) Phenanthrene	11.469	178	850447	36.723	ng	100
72) Anthracene	11.522	178	860442	37.299	ng	99
73) Carbazole	11.675	167	784848	35.300	ng	99
74) Di-n-butylphthalate	11.998	149	938523	37.540	ng	99
75) Fluoranthene	12.663	202	913289	37.471	ng	96
77) Benzidine	12.780	184	452550	77.038	ng	99
78) Pyrene	12.892	202	934318	37.280	ng	99
80) Butylbenzylphthalate	13.498	149	386415	39.337	ng	99
81) Benzo(a)anthracene	14.080	228	789112	38.134	ng	99
82) 3,3'-Dichlorobenzidine	14.039	252	199077	30.299	ng	99
83) Chrysene	14.121	228	731071	36.992	ng	100
84) Bis(2-ethylhexyl)phtha...	14.051	149	517833	39.752	ng	98
85) Di-n-octyl phthalate	14.668	149	794903	39.376	ng	100
87) Indeno(1,2,3-cd)pyrene	17.121	276	576495	35.384	ng	97
88) Benzo(b)fluoranthene	15.145	252	660476	43.080	ng	99
89) Benzo(k)fluoranthene	15.180	252	596597	38.625	ng	98
90) Benzo(a)pyrene	15.527	252	581414	45.719	ng	98
91) Dibenzo(a,h)anthracene	17.139	278	497660	38.086	ng	97
92) Benzo(g,h,i)perylene	17.592	276	507092	36.825	ng	96

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

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