

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF031723\
 Data File : BF132828.D
 Acq On : 17 Mar 2023 17:52
 Operator : CG\JU
 Sample : 01908-02MS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HILLBOR-DRUMS-0314

Quant Time: Mar 18 01:03:37 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF031623.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 17 03:35:59 2023
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By : Christian Giraldo 03/20/2023
 Supervised By : Jagrut Upadhyay 03/20/2023

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.957	152	226351	20.000	ng	0.00
21) Naphthalene-d8	8.245	136	852298	20.000	ng	0.00
39) Acenaphthene-d10	10.004	164	452056	20.000	ng	0.00
64) Phenanthrene-d10	11.498	188	855565	20.000	ng	0.00
76) Chrysene-d12	14.157	240	516188	20.000	ng	0.00
86) Perylene-d12	15.674	264	480878	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.622	112	1549450	115.508	ng	0.04
7) Phenol-d6	6.604	99	1861737	106.927	ng	0.00
23) Nitrobenzene-d5	7.528	82	1410045	84.664	ng	0.00
42) 2,4,6-Tribromophenol	10.804	330	600394	124.938	ng	0.00
45) 2-Fluorobiphenyl	9.322	172	2174226	79.007	ng	0.00
79) Terphenyl-d14	13.086	244	2572568	89.992	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.075	88	290130m	38.375	ng	Qvalue
3) Pyridine	3.775	79	672673	36.155	ng	99
4) n-Nitrosodimethylamine	3.710	42	308270	44.630	ng	98
6) Aniline	6.634	93	410849	20.543	ng	# 91
8) 2-Chlorophenol	6.757	128	705860	49.312	ng	98
9) Benzaldehyde	6.516	77	249896	34.877	ng	97
10) Phenol	6.616	94	818750	41.394	ng	99
11) bis(2-Chloroethyl)ether	6.698	93	863850	47.224	ng	100
12) 1,3-Dichlorobenzene	6.904	146	728925	46.986	ng	99
13) 1,4-Dichlorobenzene	6.981	146	728772	47.066	ng	99
14) 1,2-Dichlorobenzene	7.128	146	697815	49.391	ng	99
15) Benzyl Alcohol	7.110	79	645314	64.081	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.234	45	811439	45.590	ng	77
17) 2-Methylphenol	7.222	107	577547	44.399	ng	# 91
18) Hexachloroethane	7.469	117	286849	48.410	ng	98
19) n-Nitroso-di-n-propyla...	7.381	70	460209	46.325	ng	93
20) 3+4-Methylphenols	7.375	107	675114	43.103	ng	# 82
22) Acetophenone	7.369	105	917895	45.106	ng	# 96
24) Nitrobenzene	7.551	77	818014	45.431	ng	99
25) Isophorone	7.787	82	1480329	45.214	ng	99
26) 2-Nitrophenol	7.863	139	402593	48.002	ng	97
27) 2,4-Dimethylphenol	7.898	122	596702	45.921	ng	99
28) bis(2-Chloroethoxy)met...	7.987	93	916767	46.619	ng	99
29) 2,4-Dichlorophenol	8.110	162	656603	48.808	ng	99
30) 1,2,4-Trichlorobenzene	8.187	180	673098	46.778	ng	99
31) Naphthalene	8.269	128	1928295	45.099	ng	99
32) Benzoic acid	8.051	122	409522	50.351	ng	97
33) 4-Chloroaniline	8.345	127	171342m	9.192	ng	
34) Hexachlorobutadiene	8.375	225	424103	46.455	ng	98
35) Caprolactam	8.704	113	184948m	44.842	ng	
36) 4-Chloro-3-methylphenol	8.816	107	643212	47.011	ng	98
37) 2-Methylnaphthalene	8.957	142	1266112	44.156	ng	100
38) 1-Methylnaphthalene	9.057	142	1181233	43.754	ng	98
40) 1,2,4,5-Tetrachloroben...	9.128	216	674881	45.509	ng	99
41) Hexachlorocyclopentadiene	9.110	237	722955	99.829	ng	99
43) 2,4,6-Trichlorophenol	9.245	196	453693	46.986	ng	100

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.292	196	496674	44.180	ng	98
46) 1,1'-Biphenyl	9.422	154	1495434	44.591	ng	99
47) 2-Chloronaphthalene	9.451	162	1236802	46.575	ng	100
48) 2-Nitroaniline	9.557	65	405656	46.515	ng	95
49) Acenaphthylene	9.869	152	1790239	43.870	ng	99
50) Dimethylphthalate	9.728	163	1529082	47.178	ng	100
51) 2,6-Dinitrotoluene	9.792	165	338156	47.185	ng	90
52) Acenaphthene	10.039	154	1084243	44.139	ng	99
53) 3-Nitroaniline	9.963	138	211757	26.433	ng	94
54) 2,4-Dinitrophenol	10.081	184	372625	90.482	ng	97
55) Dibenzofuran	10.216	168	1611098	42.667	ng	100
56) 4-Nitrophenol	10.139	139	439924	92.106	ng	93
57) 2,4-Dinitrotoluene	10.204	165	418087	45.670	ng	97
58) Fluorene	10.557	166	1310067	43.630	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.334	232	409981	50.320	ng	99
60) Diethylphthalate	10.428	149	1386590	45.144	ng	99
61) 4-Chlorophenyl-phenyle...	10.545	204	676115	42.896	ng	97
62) 4-Nitroaniline	10.592	138	345042	49.256	ng	97
63) Azobenzene	10.704	77	1391131	44.301	ng	99
65) 4,6-Dinitro-2-methylph...	10.616	198	248887	47.464	ng	99
66) n-Nitrosodiphenylamine	10.669	169	1154759	43.673	ng	99
67) 4-Bromophenyl-phenylether	11.033	248	444758	44.964	ng	98
68) Hexachlorobenzene	11.110	284	496368	47.386	ng	95
69) Atrazine	11.192	200	398920	52.389	ng	98
70) Pentachlorophenol	11.310	266	475184	106.917	ng	99
71) Phenanthrene	11.528	178	1877670	43.601	ng	99
72) Anthracene	11.581	178	1890824	42.861	ng	100
73) Carbazole	11.733	167	1751385	44.217	ng	99
74) Di-n-butylphthalate	12.045	149	2019978	42.807	ng	100
75) Fluoranthene	12.722	202	1971746	42.072	ng	97
77) Benzidine	12.839	184	367669	114.372	ng	97
78) Pyrene	12.951	202	2051505	47.820	ng	99
80) Butylbenzylphthalate	13.557	149	862537	49.145	ng	92
81) Benzo(a)anthracene	14.145	228	1736177	46.468	ng	99
82) 3,3'-Dichlorobenzidine	14.098	252	323961	33.403	ng	97
83) Chrysene	14.180	228	1620282	45.891	ng	99
84) Bis(2-ethylhexyl)phtha...	14.110	149	1128982	51.156	ng	# 98
85) Di-n-octyl phthalate	14.727	149	1856946	53.710	ng	100
87) Indeno(1,2,3-cd)pyrene	17.257	276	1563870	51.880	ng	99
88) Benzo(b)fluoranthene	15.227	252	1484546	47.327	ng	99
89) Benzo(k)fluoranthene	15.257	252	1445523	47.157	ng	100
90) Benzo(a)pyrene	15.616	252	1307189	44.702	ng	99
91) Dibenzo(a,h)anthracene	17.268	278	1266166	51.471	ng	99
92) Benzo(g,h,i)perylene	17.739	276	1265858	49.248	ng	97

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

