

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF071823\
 Data File : BF134337.D
 Acq On : 18 Jul 2023 14:32
 Operator : CG\JU
 Sample : PB154222BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB154222BS

Manual Integrations
 APPROVED

Reviewed By :Yogesh
 Patel
 07/19/2023

Supervised By :mohammad
 ahmed
 07/19/2023

Quant Time: Jul 18 15:11:47 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF062623.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 14 22:51:27 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.839	152	77780	20.000	ng	0.00
21) Naphthalene-d8	8.122	136	307856	20.000	ng	0.00
39) Acenaphthene-d10	9.880	164	162022	20.000	ng	0.00
64) Phenanthrene-d10	11.374	188	321931	20.000	ng	0.00
76) Chrysene-d12	14.039	240	277603	20.000	ng	0.00
86) Perylene-d12	15.545	264	230199	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.487	112	663925	142.786	ng	0.02
7) Phenol-d6	6.492	99	658974	115.239	ng	0.00
23) Nitrobenzene-d5	7.410	82	461961	80.889	ng	0.00
42) 2,4,6-Tribromophenol	10.680	330	261009	128.485	ng	0.00
45) 2-Fluorobiphenyl	9.198	172	908980	81.567	ng	0.00
79) Terphenyl-d14	12.974	244	1326586	76.910	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.693	88	82327	42.940	ng	97
3) Pyridine	3.481	79	166097	33.260	ng	98
4) n-Nitrosodimethylamine	3.434	42	135200	47.402	ng	98
6) Aniline	6.504	93	125460	18.030	ng	# 1
8) 2-Chlorophenol	6.628	128	213620	45.258	ng	98
9) Benzaldehyde	6.392	77	28651	6.720	ng	94
10) Phenol	6.504	94	251056	41.756	ng	85
11) bis(2-Chloroethyl)ether	6.581	93	192967	43.594	ng	97
12) 1,3-Dichlorobenzene	6.781	146	229145	45.536	ng	100
13) 1,4-Dichlorobenzene	6.857	146	227753	44.809	ng	97
14) 1,2-Dichlorobenzene	7.010	146	222039	46.645	ng	98
15) Benzyl Alcohol	6.987	79	201031	45.603	ng	96
16) 2,2'-oxybis(1-Chloropr...	7.116	45	292082	37.299	ng	82
17) 2-Methylphenol	7.104	107	175093	41.425	ng	# 92
18) Hexachloroethane	7.345	117	89265	47.365	ng	96
19) n-Nitroso-di-n-propyla...	7.257	70	159541	43.995	ng	96
20) 3+4-Methylphenols	7.257	107	221377	43.173	ng	# 88
22) Acetophenone	7.251	105	319675	45.659	ng	95
24) Nitrobenzene	7.428	77	240681	41.704	ng	97
25) Isophorone	7.663	82	422639	40.719	ng	98
26) 2-Nitrophenol	7.739	139	119069	43.008	ng	92
27) 2,4-Dimethylphenol	7.781	122	187473	42.779	ng	95
28) bis(2-Chloroethoxy)met...	7.869	93	242686	40.380	ng	98
29) 2,4-Dichlorophenol	7.986	162	185180	42.613	ng	98
30) 1,2,4-Trichlorobenzene	8.063	180	200364	44.685	ng	99
31) Naphthalene	8.145	128	619790	41.679	ng	99
32) Benzoic acid	7.928	122	137263	41.800	ng	93
33) 4-Chloroaniline	8.198	127	71335	11.214	ng	98
34) Hexachlorobutadiene	8.251	225	130984	46.308	ng	100
35) Caprolactam	8.569	113	60272m	42.123	ng	
36) 4-Chloro-3-methylphenol	8.686	107	215477	44.138	ng	96
37) 2-Methylnaphthalene	8.833	142	423257	40.250	ng	99
38) 1-Methylnaphthalene	8.933	142	399994	40.916	ng	99
40) 1,2,4,5-Tetrachloroben...	9.004	216	213512	44.480	ng	99
41) Hexachlorocyclopentadiene	8.980	237	163490	71.792	ng	97
43) 2,4,6-Trichlorophenol	9.116	196	135913	41.593	ng	99

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44) 2,4,5-Trichlorophenol	9.169	196	144958	37.713	ng	96
46) 1,1'-Biphenyl	9.298	154	549203	42.257	ng	98
47) 2-Chloronaphthalene	9.328	162	400059	42.070	ng	98
48) 2-Nitroaniline	9.428	65	136961	39.520	ng	98
49) Acenaphthylene	9.745	152	648527	39.923	ng	99
50) Dimethylphthalate	9.604	163	489790	41.645	ng	99
51) 2,6-Dinitrotoluene	9.675	165	106586	42.076	ng	91
52) Acenaphthene	9.916	154	388031	41.167	ng	98
53) 3-Nitroaniline	9.839	138	71580	23.554	ng	# 96
54) 2,4-Dinitrophenol	9.963	184	118175	80.742	ng	92
55) Dibenzofuran	10.086	168	595489	40.424	ng	96
56) 4-Nitrophenol	10.022	139	146831	69.074	ng	92
57) 2,4-Dinitrotoluene	10.080	165	144472	43.185	ng	# 86
58) Fluorene	10.433	166	484583	42.282	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.210	232	123223	40.642	ng	98
60) Diethylphthalate	10.304	149	518860	42.344	ng	98
61) 4-Chlorophenyl-phenyle...	10.422	204	235786	42.691	ng	97
62) 4-Nitroaniline	10.463	138	114433	37.364	ng	94
63) Azobenzene	10.586	77	461213	39.792	ng	98
65) 4,6-Dinitro-2-methylph...	10.498	198	82333	46.909	ng	96
66) n-Nitrosodiphenylamine	10.545	169	422456	41.072	ng	98
67) 4-Bromophenyl-phenylether	10.916	248	147323	43.055	ng	96
68) Hexachlorobenzene	10.986	284	175311	48.254	ng	# 90
69) Atrazine	11.074	200	87033	30.590	ng	99
70) Pentachlorophenol	11.180	266	149164	68.667	ng	100
71) Phenanthrene	11.404	178	698851	43.122	ng	99
72) Anthracene	11.457	178	704735	41.808	ng	99
73) Carbazole	11.616	167	675488	43.780	ng	100
74) Di-n-butylphthalate	11.939	149	845031	46.056	ng	100
75) Fluoranthene	12.598	202	797606	45.523	ng	98
77) Benzidine	12.721	184	80880	12.245	ng	98
78) Pyrene	12.827	202	835734	32.349	ng	99
80) Butylbenzylphthalate	13.445	149	404675	41.765	ng	98
81) Benzo(a)anthracene	14.027	228	815131	42.339	ng	99
82) 3,3'-Dichlorobenzidine	13.992	252	199467	32.702	ng	97
83) Chrysene	14.068	228	755054	40.492	ng	100
84) Bis(2-ethylhexyl)phtha...	14.010	149	568563	51.068	ng	99
85) Di-n-octyl phthalate	14.627	149	896471	54.799	ng	98
87) Indeno(1,2,3-cd)pyrene	17.104	276	711601	44.549	ng	96
88) Benzo(b)fluoranthene	15.104	252	709797	52.438	ng	99
89) Benzo(k)fluoranthene	15.133	252	617393	44.798	ng	97
90) Benzo(a)pyrene	15.480	252	596594	45.580	ng	98
91) Dibenzo(a,h)anthracene	17.115	278	595777	45.664	ng	98
92) Benzo(g,h,i)perylene	17.574	276	605601	45.011	ng	98

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

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[illegible]