

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF012821\
 Data File : BF123057.D
 Acq On : 28 Jan 2021 13:17
 Operator : JU/CG
 Sample : PB134359BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB134359BS

Manual Integrations
 APPROVED

mohammad
 1/29/2021 10:33:31 AM

Quant Time: Jan 28 14:55:54 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF012721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 28 00:48:07 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.892	152	269377	20.00	ng	0.00
21) Naphthalene-d8	8.181	136	1028673	20.00	ng	0.00
39) Acenaphthene-d10	9.939	164	548088	20.00	ng	0.00
64) Phenanthrene-d10	11.427	188	1004108	20.00	ng	# 0.00
76) Chrysene-d12	14.086	240	844673	20.00	ng	0.00
86) Perylene-d12	15.568	264	797354	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.522	112	1971909	112.92	ng	0.02
7) Phenol-d6	6.522	99	2571492	108.64	ng	0.00
23) Nitrobenzene-d5	7.457	82	1457701	71.32	ng	0.00
42) 2,4,6-Tribromophenol	10.733	330	716266	120.39	ng	0.00
45) 2-Fluorobiphenyl	9.257	172	2659128	72.12	ng	0.00
79) Terphenyl-d14	13.022	244	3086553	71.79	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.769	88	350737	40.37	ng	95
3) Pyridine	3.516	79	965570	41.56	ng	96
4) n-Nitrosodimethylamine	3.469	42	422729	43.23	ng	94
6) Aniline	6.557	93	919046	31.55	ng	99
8) 2-Chlorophenol	6.681	128	838438	44.29	ng	97
9) Benzaldehyde	6.440	77	253245	23.40	ng	98
10) Phenol	6.534	94	1022731m	43.57	ng	
11) bis(2-Chloroethyl)ether	6.628	93	832521	42.62	ng	95
12) 1,3-Dichlorobenzene	6.834	146	915526	41.47	ng	97
13) 1,4-Dichlorobenzene	6.910	146	925788	40.27	ng	97
14) 1,2-Dichlorobenzene	7.063	146	861721	40.06	ng	97
15) Benzyl Alcohol	7.034	79	732672	44.15	ng	97
16) 2,2'-oxybis(1-Chloropr...	7.169	45	1227205	40.74	ng	97
17) 2-Methylphenol	7.145	107	708490	43.99	ng	98
18) Hexachloroethane	7.410	117	344629	42.80	ng	96
19) n-Nitroso-di-n-propyla...	7.310	70	591102	41.53	ng	97
20) 3+4-Methylphenols	7.292	107	835250	44.03	ng	# 88
22) Acetophenone	7.304	105	1111846	41.48	ng	97
24) Nitrobenzene	7.475	77	911156	43.27	ng	100
25) Isophorone	7.716	82	1663887	44.44	ng	100
26) 2-Nitrophenol	7.792	139	449287	51.59	ng	95
27) 2,4-Dimethylphenol	7.828	122	809785	51.53	ng	97
28) bis(2-Chloroethoxy)met...	7.922	93	1057443	41.40	ng	99
29) 2,4-Dichlorophenol	8.034	162	725745	44.17	ng	98
30) 1,2,4-Trichlorobenzene	8.116	180	759556	41.03	ng	99
31) Naphthalene	8.198	128	2401775	42.57	ng	99
32) Benzoic acid	7.951	122	565178	44.99	ng	98
33) 4-Chloroaniline	8.245	127	657512	26.26	ng	99
34) Hexachlorobutadiene	8.322	225	433915	40.58	ng	99
35) Caprolactam	8.616	113	232824m	41.92	ng	
36) 4-Chloro-3-methylphenol	8.728	107	771620	45.16	ng	98
37) 2-Methylnaphthalene	8.892	142	1610705	43.00	ng	100
38) 1-Methylnaphthalene	8.992	142	1501703	42.34	ng	99
40) 1,2,4,5-Tetrachloroben...	9.063	216	701303	41.13	ng	97
41) Hexachlorocyclopentadiene	9.051	237	830088	102.56	ng	98
43) 2,4,6-Trichlorophenol	9.169	196	532943	45.59	ng	98

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44) 2,4,5-Trichlorophenol	9.216	196	542771	44.40	ng	99
46) 1,1'-Biphenyl	9.357	154	1906334	42.40	ng	99
47) 2-Chloronaphthalene	9.386	162	1526089	40.55	ng	98
48) 2-Nitroaniline	9.475	65	493353	43.25	ng	86
49) Acenaphthylene	9.804	152	2328772	42.28	ng	100
50) Dimethylphthalate	9.663	163	1779698	41.39	ng	99
51) 2,6-Dinitrotoluene	9.716	165	419269	45.77	ng	# 85
52) Acenaphthene	9.975	154	1612904	45.59	ng	100
53) 3-Nitroaniline	9.886	138	309453	28.57	ng	92
54) 2,4-Dinitrophenol	9.992	184	387750	91.49	ng	99
55) Dibenzofuran	10.145	168	2107488	40.94	ng	99
56) 4-Nitrophenol	10.051	139	736367	94.05	ng	95
57) 2,4-Dinitrotoluene	10.122	165	558665	46.90	ng	96
58) Fluorene	10.492	166	1591266	38.69	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.263	232	466078	44.96	ng	95
60) Diethylphthalate	10.363	149	1780502	42.11	ng	99
61) 4-Chlorophenyl-phenyle...	10.480	204	758488	39.25	ng	98
62) 4-Nitroaniline	10.504	138	421709	38.75	ng	96
63) Azobenzene	10.639	77	1760826	42.56	ng	97
65) 4,6-Dinitro-2-methylph...	10.533	198	273906	49.27	ng	98
66) n-Nitrosodiphenylamine	10.598	169	1492955	41.25	ng	99
67) 4-Bromophenyl-phenylether	10.975	248	518218	40.52	ng	94
68) Hexachlorobenzene	11.039	284	550294	40.12	ng	97
69) Atrazine	11.128	200	401053	40.13	ng	99
70) Pentachlorophenol	11.233	266	666361	89.64	ng	99
71) Phenanthrene	11.457	178	2408435	38.62	ng	99
72) Anthracene	11.510	178	2470474	41.30	ng	99
73) Carbazole	11.663	167	2269127	40.87	ng	99
74) Di-n-butylphthalate	11.992	149	2717826	41.28	ng	99
75) Fluoranthene	12.651	202	2536021	40.69	ng	99
77) Benzidine	12.769	184	1047804	54.82	ng	100
78) Pyrene	12.880	202	2576782	40.41	ng	98
80) Butylbenzylphthalate	13.498	149	1217109	44.36	ng	96
81) Benzo(a)anthracene	14.074	228	2339070	40.92	ng	100
82) 3,3'-Dichlorobenzidine	14.033	252	748485	41.60	ng	# 97
83) Chrysene	14.116	228	2356438	41.19	ng	98
84) Bis(2-ethylhexyl)phtha...	14.063	149	1588971	43.58	ng	99
85) Di-n-octyl phthalate	14.680	149	2867552	46.83	ng	98
87) Indeno(1,2,3-cd)pyrene	17.074	276	2422734	45.63	ng	96
88) Benzo(b)fluoranthene	15.133	252	2604100	45.25	ng	99
89) Benzo(k)fluoranthene	15.168	252	2224030	41.59	ng	99
90) Benzo(a)pyrene	15.510	252	2190732	43.34	ng	98
91) Dibenzo(a,h)anthracene	17.098	278	2010309	46.00	ng	98
92) Benzo(g,h,i)perylene	17.533	276	1976189	46.66	ng	98

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

