

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081319\
 Data File : BF116224.D
 Acq On : 13 Aug 2019 11:50
 Operator : HP/JU
 Sample : PB122181BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB122181BS

Quant Time: Aug 13 23:43:30 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 07 11:13:18 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	127110	20.00	ng	-0.01
21) Naphthalene-d8	8.12	136	523692	20.00	ng	-0.01
39) Acenaphthene-d10	9.87	164	277676	20.00	ng	-0.01
64) Phenanthrene-d10	11.36	188	480915	20.00	ng	0.00
76) Chrysene-d12	14.00	240	335416	20.00	ng	0.00
87) Perylene-d12	15.47	264	295416	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.47	112	885634	116.64	ng	0.00
7) Phenol-d6	6.48	99	1081798	112.92	ng	0.00
23) Nitrobenzene-d5	7.40	82	715057	78.82	ng	-0.01
42) 2,4,6-Tribromophenol	10.67	330	302639	112.42	ng	0.00
45) 2-Fluorobiphenyl	9.19	172	1221254	82.38	ng	-0.01
79) Terphenyl-d14	12.94	244	1365225	85.77	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.70	88	155806	40.618	ng	99
3) Pyridine	3.45	79	370156	34.545	ng	97
4) n-Nitrosodimethylamine	3.40	42	159641	37.753	ng	100
6) Aniline	6.50	93	429466	31.304	ng	# 65
8) 2-Chlorophenol	6.63	128	375005	44.664	ng	98
9) Benzaldehyde	6.39	77	153879	23.280	ng	98
10) Phenol	6.49	94	465002	41.219	ng	94
11) bis(2-Chloroethyl)ether	6.57	93	362801	43.742	ng	97
12) 1,3-Dichlorobenzene	6.77	146	403473	41.580	ng	100
13) 1,4-Dichlorobenzene	6.85	146	413132	42.522	ng	98
14) 1,2-Dichlorobenzene	7.00	146	377286	42.173	ng	98
15) Benzyl Alcohol	6.98	79	304075	41.826	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.11	45	474753	41.965	ng	93
17) 2-Methylphenol	7.10	107	310564	43.683	ng	97
18) Hexachloroethane	7.35	117	148630	41.550	ng	94
19) n-Nitroso-di-n-propylamine	7.25	70	257528	43.381	ng	98
20) 3+4-Methylphenols	7.25	107	383357	45.566	ng	96
22) Acetophenone	7.25	105	507670	44.202	ng	98
24) Nitrobenzene	7.42	77	416193	42.485	ng	97
25) Isophorone	7.66	82	759852	43.801	ng	99
26) 2-Nitrophenol	7.73	139	207877	45.017	ng	99
27) 2,4-Dimethylphenol	7.77	122	333754	48.041	ng	99
28) bis(2-Chloroethoxy)methane	7.86	93	467563	43.484	ng	100
29) 2,4-Dichlorophenol	7.98	162	329232	44.883	ng	98
30) 1,2,4-Trichlorobenzene	8.06	180	348546	41.684	ng	99
31) Naphthalene	8.14	128	1084729	44.016	ng	99
32) Benzoic acid	7.93	122	174817	35.691	ng	97
33) 4-Chloroaniline	8.19	127	291702	26.492	ng	97
34) Hexachlorobutadiene	8.25	225	197747	41.058	ng	99
35) Caprolactam	8.57	113	100557m	42.385	ng	
36) 4-Chloro-3-methylphenol	8.69	107	343058	44.955	ng	96
37) 2-Methylnaphthalene	8.83	142	715309	44.073	ng	99
38) 1-Methylnaphthalene	8.93	142	666536	43.410	ng	100
40) 1,2,4,5-Tetrachlorobenzene	9.00	216	331507	44.657	ng	99

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41) Hexachlorocyclopentadiene	8.98	237	133675	43.064	ng	99
43) 2,4,6-Trichlorophenol	9.12	196	212758	41.639	ng	98
44) 2,4,5-Trichlorophenol	9.17	196	219975	41.648	ng	97
46) 1,1'-Biphenyl	9.29	154	876393	44.705	ng	100
47) 2-Chloronaphthalene	9.32	162	688810	42.216	ng	99
48) 2-Nitroaniline	9.42	65	216518	40.147	ng	95
49) Acenaphthylene	9.73	152	1031857	43.348	ng	99
50) Dimethylphthalate	9.59	163	813916	43.049	ng	100
51) 2,6-Dinitrotoluene	9.66	165	178914	43.545	ng	90
52) Acenaphthene	9.91	154	618968	42.386	ng	98
53) 3-Nitroaniline	9.83	138	141264	27.825	ng	98
54) 2,4-Dinitrophenol	9.96	184	135840	66.467	ng	98
55) Dibenzofuran	10.08	168	887011	41.767	ng	99
56) 4-Nitrophenol	10.02	139	206043	63.354	ng	95
57) 2,4-Dinitrotoluene	10.07	165	216595	39.594	ng	91
58) Fluorene	10.42	166	716368	43.844	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.20	232	199572	44.911	ng	98
60) Diethylphthalate	10.29	149	788889	44.692	ng	99
61) 4-Chlorophenyl-phenylether	10.41	204	368695	44.555	ng	100
62) 4-Nitroaniline	10.46	138	174973	33.349	ng	98
63) Azobenzene	10.57	77	809024	44.560	ng	96
65) 4,6-Dinitro-2-methylphenol	10.49	198	109545	42.984	ng	97
66) n-Nitrosodiphenylamine	10.53	169	650560	44.944	ng	99
67) 4-Bromophenyl-phenylether	10.90	248	224928	43.691	ng	97
68) Hexachlorobenzene	10.98	284	248048	41.667	ng	95
69) Atrazine	11.06	200	202299	42.482	ng	98
70) Pentachlorophenol	11.17	266	225670	78.081	ng	98
71) Phenanthrene	11.39	178	1047984	42.626	ng	100
72) Anthracene	11.44	178	1091554	43.870	ng	99
73) Carbazole	11.59	167	998789	44.246	ng	99
74) Di-n-butylphthalate	11.91	149	1236584	46.959	ng	99
75) Fluoranthene	12.57	202	1115631	43.345	ng	98
77) Benzidine	12.69	184	382666	33.769	ng	99
78) Pyrene	12.80	202	1123889	44.450	ng	100
80) Butylbenzylphthalate	13.41	149	519402	48.564	ng	98
81) Benzo(a)anthracene	14.00	228	925912	43.189	ng	99
82) 3,3'-Dichlorobenzidine	13.95	252	282815	37.971	ng	98
83) Chrysene	14.03	228	898239	43.156	ng	99
84) Bis(2-ethylhexyl)phthalate	13.97	149	718963	51.730	ng	98
85) Di-n-octyl phthalate	14.58	149	1142311	51.745	ng	99
86) Indeno(1,2,3-cd)pyrene	16.96	276	797349	45.612	ng	99
88) Benzo(b)fluoranthene	15.05	252	793411	46.449	ng	100
89) Benzo(k)fluoranthene	15.08	252	800681	48.125	ng	98
90) Benzo(a)pyrene	15.42	252	765838	50.155	ng	98
91) Dibenzo(a,h)anthracene	16.96	278	671756	50.824	ng	99
92) Benzo(g,h,i)perylene	17.40	276	667832	51.448	ng	99

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

