

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112219\
 Data File : BF117927.D
 Acq On : 22 Nov 2019 19:31
 Operator : JU
 Sample : PB124996BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB124996BS

Manual Integrations
 APPROVED

mohammad
 11/25/2019 8:19:31 AM

Quant Time: Nov 25 03:44:24 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 13 18:36:07 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	241182	20.00	ng	-0.02
21) Naphthalene-d8	8.19	136	1019278	20.00	ng	-0.02
39) Acenaphthene-d10	9.96	164	532645	20.00	ng	-0.02
64) Phenanthrene-d10	11.45	188	1015406	20.00	ng	-0.02
76) Chrysene-d12	14.10	240	637818	20.00	ng	-0.01
87) Perylene-d12	15.60	264	464585	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.53	112	1676582	123.05	ng	0.00
7) Phenol-d6	6.54	99	2054548	125.02	ng	0.00
23) Nitrobenzene-d5	7.47	82	1136232	66.27	ng	-0.02
42) 2,4,6-Tribromophenol	10.76	330	712959	126.43	ng	-0.01
45) 2-Fluorobiphenyl	9.27	172	2223557	74.89	ng	-0.02
79) Terphenyl-d14	13.04	244	2359800	67.62	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.78	88	196999	30.931	ng	98
3) Pyridine	3.54	79	512047	30.649	ng	98
4) n-Nitrosodimethylamine	3.49	42	289150	39.648	ng	100
6) Aniline	6.57	93	737594	33.964	ng	99
8) 2-Chlorophenol	6.70	128	685339	46.355	ng	97
9) Benzaldehyde	6.45	77	297773	24.862	ng	98
10) Phenol	6.55	94	840834	49.236	ng	96
11) bis(2-Chloroethyl)ether	6.65	93	614369	44.799	ng	95
12) 1,3-Dichlorobenzene	6.85	146	738631	42.435	ng	96
13) 1,4-Dichlorobenzene	6.92	146	759895	43.191	ng	98
14) 1,2-Dichlorobenzene	7.08	146	717024	42.612	ng	98
15) Benzyl Alcohol	7.05	79	636077	50.132	ng	100
16) 2,2'-oxybis(1-Chloropropan	7.18	45	865664	40.999	ng	97
17) 2-Methylphenol	7.16	107	619303	49.464	ng	99
18) Hexachloroethane	7.42	117	275220	42.582	ng	94
19) n-Nitroso-di-n-propylamine	7.32	70	467294	46.090	ng	98
20) 3+4-Methylphenols	7.31	107	708392	45.579	ng	# 85
22) Acetophenone	7.32	105	937129	40.922	ng	# 97
24) Nitrobenzene	7.49	77	743791	42.049	ng	100
25) Isophorone	7.73	82	1385833	44.097	ng	98
26) 2-Nitrophenol	7.81	139	418730	48.635	ng	97
27) 2,4-Dimethylphenol	7.85	122	686121	51.286	ng	100
28) bis(2-Chloroethoxy)methane	7.94	93	833218	41.781	ng	99
29) 2,4-Dichlorophenol	8.05	162	649517	44.758	ng	98
30) 1,2,4-Trichlorobenzene	8.13	180	678680	40.320	ng	99
31) Naphthalene	8.22	128	2046308	43.064	ng	100
32) Benzoic acid	7.98	122	558698	49.876	ng	98
33) 4-Chloroaniline	8.26	127	573149	26.943	ng	99
34) Hexachlorobutadiene	8.33	225	382270	38.843	ng	99
35) Caprolactam	8.65	113	213521m	46.297	ng	
36) 4-Chloro-3-methylphenol	8.75	107	669310	45.447	ng	95
37) 2-Methylnaphthalene	8.91	142	1404733	43.753	ng	99
38) 1-Methylnaphthalene	9.01	142	1347123	44.382	ng	100
40) 1,2,4,5-Tetrachlorobenzene	9.08	216	609186	39.382	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	9.06	237	665158	76.732	ng	99
43) 2,4,6-Trichlorophenol	9.19	196	483274	43.619	ng	98
44) 2,4,5-Trichlorophenol	9.23	196	495511	43.075	ng	98
46) 1,1'-Biphenyl	9.38	154	1704882	43.143	ng	99
47) 2-Chloronaphthalene	9.40	162	1342226	41.288	ng	99
48) 2-Nitroaniline	9.50	65	410934	41.870	ng	93
49) Acenaphthylene	9.82	152	2095498	44.212	ng	100
50) Dimethylphthalate	9.68	163	1569862	42.030	ng	100
51) 2,6-Dinitrotoluene	9.74	165	358920	43.968	ng	99
52) Acenaphthene	9.99	154	1264950	41.662	ng	98
53) 3-Nitroaniline	9.91	138	283837	29.058	ng	95
54) 2,4-Dinitrophenol	10.02	184	405563	95.184	ng	96
55) Dibenzofuran	10.17	168	1804545	42.920	ng	99
56) 4-Nitrophenol	10.07	139	639428	87.699	ng	98
57) 2,4-Dinitrotoluene	10.15	165	486048	46.807	ng	93
58) Fluorene	10.51	166	1402616	43.050	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.28	232	412540	43.646	ng	99
60) Diethylphthalate	10.38	149	1571066	43.143	ng	100
61) 4-Chlorophenyl-phenylether	10.50	204	701606	42.433	ng	95
62) 4-Nitroaniline	10.53	138	365075	37.328	ng	98
63) Azobenzene	10.66	77	1393858	42.073	ng	96
65) 4,6-Dinitro-2-methylphenol	10.56	198	277577	58.557	ng	84
66) n-Nitrosodiphenylamine	10.62	169	1272969	43.077	ng	99
67) 4-Bromophenyl-phenylether	10.99	248	455087	41.537	ng	94
68) Hexachlorobenzene	11.06	284	521069	40.787	ng	95
69) Atrazine	11.15	200	432746	44.517	ng	98
70) Pentachlorophenol	11.26	266	606493	81.258	ng	99
71) Phenanthrene	11.48	178	2100773	41.150	ng	99
72) Anthracene	11.53	178	2143912	42.356	ng	100
73) Carbazole	11.68	167	1952437	44.142	ng	99
74) Di-n-butylphthalate	12.01	149	2294118	41.657	ng	99
75) Fluoranthene	12.67	202	2093596	45.041	ng	98
77) Benzidine	12.79	184	446419	21.368	ng	98
78) Pyrene	12.90	202	2144481	41.604	ng	99
80) Butylbenzylphthalate	13.52	149	936367	42.295	ng	98
81) Benzo(a)anthracene	14.09	228	1763061	41.830	ng	99
82) 3,3'-Dichlorobenzidine	14.05	252	538234	36.508	ng	98
83) Chrysene	14.13	228	1661603	40.684	ng	98
84) Bis(2-ethylhexyl)phthalate	14.07	149	1203356	44.834	ng	100
85) Di-n-octyl phthalate	14.69	149	1861334	47.205	ng	99
86) Indeno(1,2,3-cd)pyrene	17.13	276	1373828	39.523	ng	99
88) Benzo(b)fluoranthene	15.16	252	1563045	51.783	ng	100
89) Benzo(k)fluoranthene	15.19	252	1391985	48.944	ng	99
90) Benzo(a)pyrene	15.54	252	1317028	49.620	ng	100
91) Dibenzo(a,h)anthracene	17.16	278	1092936	47.840	ng	97
92) Benzo(g,h,i)perylene	17.60	276	1205009	52.957	ng	98

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

