

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112319\
 Data File : BF117963.D
 Acq On : 23 Nov 2019 18:15
 Operator : JU
 Sample : K5951-01MSD
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 IRONBOUND-4FTMSD

Manual Integrations
 APPROVED

mohammad
 11/25/2019 2:32:22 PM

Quant Time: Nov 25 08:30:23 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	136085	20.00	ng	-0.02
21) Naphthalene-d8	8.19	136	440183	20.00	ng	-0.02
39) Acenaphthene-d10	9.96	164	236218	20.00	ng	-0.02
64) Phenanthrene-d10	11.45	188	474533	20.00	ng	-0.02
76) Chrysene-d12	14.10	240	369314	20.00	ng	-0.02
87) Perylene-d12	15.59	264	373163	20.00	ng	-0.03

System Monitoring Compounds

5) 2-Fluorophenol	5.52	112	958716	124.70	ng	-0.01
7) Phenol-d6	6.53	99	1227456	132.37	ng	-0.01
23) Nitrobenzene-d5	7.47	82	756304	102.14	ng	-0.02
42) 2,4,6-Tribromophenol	10.75	330	359918	143.92	ng	-0.02
45) 2-Fluorobiphenyl	9.27	172	1237229	93.97	ng	-0.02
79) Terphenyl-d14	13.04	244	1522746	75.36	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.72	88	120909	33.645	ng	96
3) Pyridine	3.48	79	282620	29.981	ng	98
4) n-Nitrosodimethylamine	3.43	42	153988	37.421	ng	98
6) Aniline	6.56	93	338723	27.642	ng	99
8) 2-Chlorophenol	6.69	128	382601	45.864	ng	97
9) Benzaldehyde	6.45	77	166452	24.563	ng	98
10) Phenol	6.55	94	483172	50.143	ng	98
11) bis(2-Chloroethyl)ether	6.64	93	330377	42.695	ng	97
12) 1,3-Dichlorobenzene	6.85	146	420619	42.828	ng	99
13) 1,4-Dichlorobenzene	6.92	146	424228	42.734	ng	99
14) 1,2-Dichlorobenzene	7.07	146	356872	37.588	ng	100
15) Benzyl Alcohol	7.05	79	297242	41.519	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.17	45	409421	34.366	ng	98
17) 2-Methylphenol	7.15	107	291835	41.310	ng	99
18) Hexachloroethane	7.42	117	153818	42.179	ng	97
19) n-Nitroso-di-n-propylamine	7.32	70	244951	42.819	ng	97
20) 3+4-Methylphenols	7.30	107	390069	44.481	ng	# 88
22) Acetophenone	7.31	105	511289	51.700	ng	96
24) Nitrobenzene	7.49	77	403106	52.770	ng	99
25) Isophorone	7.73	82	602652	44.405	ng	98
26) 2-Nitrophenol	7.80	139	182887	49.188	ng	96
27) 2,4-Dimethylphenol	7.84	122	291021	50.371	ng	99
28) bis(2-Chloroethoxy)methane	7.93	93	354821	41.199	ng	99
29) 2,4-Dichlorophenol	8.05	162	281403	44.903	ng	98
30) 1,2,4-Trichlorobenzene	8.13	180	304132	41.838	ng	99
31) Naphthalene	8.22	128	953019	46.441	ng	98
32) Benzoic acid	7.96	122	216459	44.745	ng	97
33) 4-Chloroaniline	8.26	127	130322	14.186	ng	97
34) Hexachlorobutadiene	8.33	225	173085	40.725	ng	100
35) Caprolactam	8.63	113	93510m	46.949	ng	
36) 4-Chloro-3-methylphenol	8.75	107	296940	46.688	ng	100
37) 2-Methylnaphthalene	8.91	142	646950	46.660	ng	98
38) 1-Methylnaphthalene	9.01	142	612253	46.708	ng	96
40) 1,2,4,5-Tetrachlorobenzene	9.07	216	284866	41.526	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	9.06	237	354084	92.105	ng	99
43) 2,4,6-Trichlorophenol	9.19	196	209267	42.590	ng	100
44) 2,4,5-Trichlorophenol	9.23	196	219086	42.944	ng	96
46) 1,1'-Biphenyl	9.37	154	783419	44.702	ng	98
47) 2-Chloronaphthalene	9.40	162	610845	42.369	ng	99
48) 2-Nitroaniline	9.49	65	182323	41.888	ng	94
49) Acenaphthylene	9.82	152	992523	47.219	ng	99
50) Dimethylphthalate	9.67	163	937608	56.603	ng	98
51) 2,6-Dinitrotoluene	9.73	165	164011	45.304	ng	97
52) Acenaphthene	9.99	154	692246	51.411	ng	97
53) 3-Nitroaniline	9.90	138	96201	22.208	ng	93
54) 2,4-Dinitrophenol	10.01	184	178816	94.673	ng	# 83
55) Dibenzofuran	10.16	168	872203	46.778	ng	99
56) 4-Nitrophenol	10.07	139	308038	95.265	ng	92
57) 2,4-Dinitrotoluene	10.14	165	227782	49.462	ng	99
58) Fluorene	10.50	166	674627	46.690	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.28	232	186632	44.524	ng	97
60) Diethylphthalate	10.37	149	733089	45.394	ng	100
61) 4-Chlorophenyl-phenylether	10.50	204	332340	45.323	ng	99
62) 4-Nitroaniline	10.52	138	181142	41.763	ng	97
63) Azobenzene	10.66	77	667056	45.402	ng	97
65) 4,6-Dinitro-2-methylphenol	10.55	198	124162	56.048	ng	94
66) n-Nitrosodiphenylamine	10.62	169	602543	43.630	ng	99
67) 4-Bromophenyl-phenylether	10.99	248	210810	41.173	ng	91
68) Hexachlorobenzene	11.06	284	245235	41.076	ng	97
69) Atrazine	11.15	200	208953	45.995	ng	99
70) Pentachlorophenol	11.25	266	285662	81.897	ng	98
71) Phenanthrene	11.47	178	1062038	44.515	ng	98
72) Anthracene	11.53	178	1102066	46.590	ng	98
73) Carbazole	11.68	167	976858	47.258	ng	98
74) Di-n-butylphthalate	12.00	149	1178226	45.780	ng	99
75) Fluoranthene	12.67	202	1138306	53.879	ng	98
77) Benzidine	12.79	184	512504	42.366	ng	98
78) Pyrene	12.90	202	1171374	39.247	ng	99
80) Butylbenzylphthalate	13.51	149	530069	41.350	ng	96
81) Benzo(a)anthracene	14.09	228	1078855	44.207	ng	99
82) 3,3'-Dichlorobenzidine	14.04	252	254183	29.776	ng	99
83) Chrysene	14.13	228	1018935	43.087	ng	99
84) Bis(2-ethylhexyl)phthalate	14.07	149	723342	46.544	ng	99
85) Di-n-octyl phthalate	14.69	149	1244031	54.488	ng	98
86) Indeno(1,2,3-cd)pyrene	17.12	276	950151	47.208	ng	99
88) Benzo(b)fluoranthene	15.15	252	1042432	42.997	ng	99
89) Benzo(k)fluoranthene	15.19	252	944499	41.346	ng	99
90) Benzo(a)pyrene	15.53	252	885089	41.516	ng	100
91) Dibenzo(a,h)anthracene	17.14	278	793786	43.258	ng	98
92) Benzo(g,h,i)perylene	17.59	276	797312	43.624	ng	99

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

