

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062119\
 Data File : BF115101.D
 Acq On : 22 Jun 2019 1:50
 Operator : HP/JU
 Sample : PB120818BS
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB120818BS

Manual Integrations
 APPROVED

mohammad
 6/25/2019 12:41:35 PM

Quant Time: Jun 22 05:38:31 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF062119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 21 17:55:32 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.61	152	282611	20.00	ng	0.00
21) Naphthalene-d8	7.90	136	1123951	20.00	ng	0.00
39) Acenaphthene-d10	9.65	164	558476	20.00	ng	0.00
64) Phenanthrene-d10	11.12	188	1105002	20.00	ng	0.00
76) Chrysene-d12	13.74	240	769518	20.00	ng	0.00
87) Perylene-d12	15.11	264	806977	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.22	112	2189694	119.57	ng	0.02
7) Phenol-d6	6.26	99	2690828	121.05	ng	0.00
23) Nitrobenzene-d5	7.18	82	1672716	91.55	ng	0.00
42) 2,4,6-Tribromophenol	10.44	330	755894	128.35	ng	0.00
45) 2-Fluorobiphenyl	8.98	172	3022698	91.94	ng	0.00
79) Terphenyl-d14	12.71	244	3211871	88.74	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.29	88	294212	34.040	ng	99
3) Pyridine	2.94	79	842497	34.584	ng	98
4) n-Nitrosodimethylamine	2.89	42	368584	38.595	ng	100
6) Aniline	6.28	93	712701	23.329	ng	# 15
8) 2-Chlorophenol	6.40	128	834452	40.522	ng	99
9) Benzaldehyde	6.15	77	227466	15.685	ng	98
10) Phenol	6.27	94	973815	37.400	ng	96
11) bis(2-Chloroethyl)ether	6.36	93	766973	39.960	ng	99
12) 1,3-Dichlorobenzene	6.55	146	897232	39.259	ng	98
13) 1,4-Dichlorobenzene	6.63	146	902115	39.364	ng	98
14) 1,2-Dichlorobenzene	6.79	146	830157	39.116	ng	98
15) Benzyl Alcohol	6.77	79	639356	39.501	ng	99
16) 2,2'-oxybis(1-Chloropropan	6.90	45	1115536	40.073	ng	99
17) 2-Methylphenol	6.88	107	724943	42.649	ng	100
18) Hexachloroethane	7.13	117	329467	39.877	ng	99
19) n-Nitroso-di-n-propylamine	7.05	70	567256	39.861	ng	98
20) 3+4-Methylphenols	7.04	107	812052	41.374	ng	# 87
22) Acetophenone	7.03	105	1122628	43.629	ng	# 99
24) Nitrobenzene	7.20	77	864893	42.968	ng	97
25) Isophorone	7.44	82	1621076	43.311	ng	100
26) 2-Nitrophenol	7.51	139	466218	46.901	ng	97
27) 2,4-Dimethylphenol	7.57	122	794590	48.821	ng	100
28) bis(2-Chloroethoxy)methane	7.66	93	1007906	42.605	ng	99
29) 2,4-Dichlorophenol	7.75	162	750972	44.711	ng	96
30) 1,2,4-Trichlorobenzene	7.84	180	794031	42.799	ng	99
31) Naphthalene	7.92	128	2386446	43.255	ng	99
32) Benzoic acid	7.71	122	614901	52.928	ng	99
33) 4-Chloroaniline	7.97	127	397056	15.747	ng	99
34) Hexachlorobutadiene	8.05	225	426278	41.928	ng	99
35) Caprolactam	8.35	113	248704m	44.042	ng	
36) 4-Chloro-3-methylphenol	8.46	107	750531	44.123	ng	99
37) 2-Methylnaphthalene	8.61	142	1657948	44.569	ng	99
38) 1-Methylnaphthalene	8.71	142	1564158	44.026	ng	99
40) 1,2,4,5-Tetrachlorobenzene	8.78	216	665123	42.146	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.77	237	724239	77.393	nq	98
43) 2,4,6-Trichlorophenol	8.89	196	552990	45.387	nq	99
44) 2,4,5-Trichlorophenol	8.93	196	555365	44.262	nq	98
46) 1,1'-Biphenyl	9.08	154	1875004	41.960	nq	98
47) 2-Chloronaphthalene	9.10	162	1529223	42.189	nq	97
48) 2-Nitroaniline	9.19	65	458036	43.343	nq	95
49) Acenaphthylene	9.51	152	2382251	42.183	nq	99
50) Dimethylphthalate	9.39	163	1756641	42.753	nq	99
51) 2,6-Dinitrotoluene	9.44	165	428482	45.170	nq	99
52) Acenaphthene	9.68	154	1648590	46.651	nq	99
53) 3-Nitroaniline	9.60	138	307745	26.585	nq	92
54) 2,4-Dinitrophenol	9.71	184	447282	90.566	nq	97
55) Dibenzofuran	9.85	168	2099463	41.393	nq	98
56) 4-Nitrophenol	9.77	139	789331	85.052	nq	95
57) 2,4-Dinitrotoluene	9.84	165	571988	46.699	nq	# 87
58) Fluorene	10.20	166	1428420	41.836	nq	98
59) 2,3,4,6-Tetrachlorophenol	9.97	232	480752	45.694	nq	99
60) Diethylphthalate	10.09	149	1789128	43.318	nq	98
61) 4-Chlorophenyl-phenylether	10.19	204	688013	42.494	nq	95
62) 4-Nitroaniline	10.22	138	480638	41.388	nq	99
63) Azobenzene	10.35	77	1671122	43.318	nq	97
65) 4,6-Dinitro-2-methylphenol	10.24	198	269862	46.116	nq	86
66) n-Nitrosodiphenylamine	10.31	169	1487925	43.221	nq	99
67) 4-Bromophenyl-phenylether	10.68	248	502154	41.914	nq	96
68) Hexachlorobenzene	10.74	284	550237	42.313	nq	# 92
69) Atrazine	10.84	200	476295	43.009	nq	99
70) Pentachlorophenol	10.93	266	654011	78.944	nq	97
71) Phenanthrene	11.15	178	2344421	39.405	nq	99
72) Anthracene	11.20	178	2449045	40.779	nq	98
73) Carbazole	11.35	167	2239551	41.761	nq	98
74) Di-n-butylphthalate	11.70	149	2608776	42.097	nq	98
75) Fluoranthene	12.32	202	2470341	41.616	nq	99
77) Benzidine	12.44	184	521221	15.254	nq	99
78) Pyrene	12.55	202	2506916	42.391	nq	98
80) Butylbenzylphthalate	13.19	149	1179289	45.187	nq	94
81) Benzo(a)anthracene	13.73	228	2287771	41.434	nq	99
82) 3,3'-Dichlorobenzidine	13.70	252	575464	28.861	nq	99
83) Chrysene	13.77	228	2142752	42.365	nq	98
84) Bis(2-ethylhexyl)phthalate	13.75	149	1560718	45.639	nq	100
85) Di-n-octyl phthalate	14.37	149	2626706	45.716	nq	100
86) Indeno(1,2,3-cd)pyrene	16.39	276	2684613	44.856	nq	99
88) Benzo(b)fluoranthene	14.74	252	2413295	47.928	nq	98
89) Benzo(k)fluoranthene	14.77	252	2235391	49.504	nq	100
90) Benzo(a)pyrene	15.06	252	2301694	50.716	nq	99
91) Dibenzo(a,h)anthracene	16.42	278	2202073	51.974	nq	99
92) Benzo(g,h,i)perylene	16.78	276	2261370	52.735	nq	99

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

