

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080620\
 Data File : BF121199.D
 Acq On : 6 Aug 2020 12:28
 Operator : JU/CG
 Sample : L3521-07MSD
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 COMP-CDMSD

Manual Integrations
 APPROVED

mohammad
 8/7/2020 1:50:42 PM

Quant Time: Aug 06 13:52:20 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071620.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 29 14:55:43 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	141595	20.00	ng	-0.01
21) Naphthalene-d8	8.08	136	559008	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	299956	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	593094	20.00	ng	0.00
76) Chrysene-d12	13.97	240	486450	20.00	ng	0.00
86) Perylene-d12	15.41	264	508256	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.42	112	808266	93.58	ng	0.00
7) Phenol-d6	6.44	99	1021633	91.57	ng	0.00
23) Nitrobenzene-d5	7.36	82	680489	70.44	ng	-0.01
42) 2,4,6-Tribromophenol	10.63	330	330479	100.65	ng	0.00
45) 2-Fluorobiphenyl	9.16	172	1252027	62.32	ng	0.00
79) Terphenyl-d14	12.91	244	1371497	61.29	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.52	88	149517	37.613	ng	99
3) Pyridine	3.26	79	366807	34.688	ng	98
4) n-Nitrosodimethylamine	3.21	42	197640	43.708	ng	93
6) Aniline	6.46	93	520224	35.720	ng	# 81
8) 2-Chlorophenol	6.58	128	464954	48.865	ng	99
9) Benzaldehyde	6.35	77	285334	60.321	ng	98
10) Phenol	6.45	94	539395	43.949	ng	95
11) bis(2-Chloroethyl)ether	6.53	93	432639	45.996	ng	100
12) 1,3-Dichlorobenzene	6.73	146	446998	43.324	ng	98
13) 1,4-Dichlorobenzene	6.81	146	452009	44.058	ng	99
14) 1,2-Dichlorobenzene	6.96	146	439519	45.284	ng	99
15) Benzyl Alcohol	6.95	79	380054	48.168	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.07	45	599856	44.246	ng	99
17) 2-Methylphenol	7.06	107	368103	43.648	ng	100
18) Hexachloroethane	7.30	117	166402	44.812	ng	100
19) n-Nitroso-di-n-propylamine	7.22	70	282101	44.465	ng	98
20) 3+4-Methylphenols	7.21	107	419388	39.092	ng	94
22) Acetophenone	7.20	105	568033	45.277	ng	97
24) Nitrobenzene	7.38	77	465831	44.738	ng	99
25) Isophorone	7.62	82	878347	45.555	ng	98
26) 2-Nitrophenol	7.70	139	255739	51.714	ng	98
27) 2,4-Dimethylphenol	7.74	122	393388	48.995	ng	99
28) bis(2-Chloroethoxy)methane	7.83	93	556661	47.297	ng	100
29) 2,4-Dichlorophenol	7.94	162	391531	47.721	ng	97
30) 1,2,4-Trichlorobenzene	8.02	180	405719	44.571	ng	99
31) Naphthalene	8.10	128	1255448	43.783	ng	100
32) Benzoic acid	7.89	122	304983	50.601	ng	99
33) 4-Chloroaniline	8.15	127	381702	29.840	ng	99
34) Hexachlorobutadiene	8.22	225	228493	42.581	ng	98
35) Caprolactam	8.55	113	137248m	52.164	ng	
36) 4-Chloro-3-methylphenol	8.65	107	415067	49.327	ng	97
37) 2-Methylnaphthalene	8.79	142	873035	46.891	ng	100
38) 1-Methylnaphthalene	8.89	142	818377	46.410	ng	100
40) 1,2,4,5-Tetrachlorobenzene	8.96	216	401660	45.959	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.95	237	383447	79.387	nq	100
43) 2,4,6-Trichlorophenol	9.07	196	291244	47.331	nq	98
44) 2,4,5-Trichlorophenol	9.12	196	316038	45.278	nq	99
46) 1,1'-Biphenyl	9.26	154	1062344	46.430	nq	99
47) 2-Chloronaphthalene	9.28	162	822159	44.073	nq	98
48) 2-Nitroaniline	9.38	65	275191	48.814	nq	98
49) Acenaphthylene	9.70	152	1323861	45.681	nq	100
50) Dimethylphthalate	9.56	163	1151559	53.214	nq	100
51) 2,6-Dinitrotoluene	9.63	165	229687	49.404	nq	98
52) Acenaphthene	9.87	154	919263m	49.892	nq	
53) 3-Nitroaniline	9.80	138	182324	31.269	nq	96
54) 2,4-Dinitrophenol	9.91	184	265912	97.713	nq	97
55) Dibenzofuran	10.05	168	1151088	43.202	nq	99
56) 4-Nitrophenol	9.97	139	391698	88.434	nq	95
57) 2,4-Dinitrotoluene	10.03	165	296624	48.585	nq	# 87
58) Fluorene	10.39	166	878909	44.229	nq	98
59) 2,3,4,6-Tetrachlorophenol	10.17	232	269945	50.795	nq	100
60) Diethylphthalate	10.26	149	1021163	46.653	nq	100
61) 4-Chlorophenyl-phenylether	10.37	204	428943	40.469	nq	99
62) 4-Nitroaniline	10.42	138	267565	47.108	nq	97
63) Azobenzene	10.54	77	948500	45.518	nq	98
65) 4,6-Dinitro-2-methylphenol	10.45	198	183246	52.953	nq	98
66) n-Nitrosodiphenylamine	10.50	169	854599	45.804	nq	99
67) 4-Bromophenyl-phenylether	10.87	248	292864	44.289	nq	96
68) Hexachlorobenzene	10.93	284	334534	46.768	nq	# 92
69) Atrazine	11.03	200	250501	54.213	nq	99
70) Pentachlorophenol	11.13	266	370212	90.343	nq	98
71) Phenanthrene	11.35	178	1344877	44.090	nq	100
72) Anthracene	11.40	178	1397859	45.638	nq	100
73) Carbazole	11.56	167	1368192	45.011	nq	100
74) Di-n-butylphthalate	11.89	149	1598586	45.572	nq	100
75) Fluoranthene	12.54	202	1572261	46.503	nq	99
77) Benzidine	12.66	184	737385	57.655	nq	99
78) Pyrene	12.77	202	1604181	46.644	nq	99
80) Butylbenzylphthalate	13.38	149	738768	48.186	nq	94
81) Benzo(a)anthracene	13.96	228	1390891	47.218	nq	99
82) 3,3'-Dichlorobenzidine	13.92	252	499388	44.936	nq	99
83) Chrysene	14.00	228	1395639	45.084	nq	100
84) Bis(2-ethylhexyl)phthalate	13.94	149	945621	50.018	nq	100
85) Di-n-octyl phthalate	14.56	149	1743500	46.354	nq	100
87) Indeno(1,2,3-cd)pyrene	16.85	276	1536194	43.460	nq	100
88) Benzo(b)fluoranthene	15.00	252	1508519	49.102	nq	99
89) Benzo(k)fluoranthene	15.03	252	1378669	49.206	nq	99
90) Benzo(a)pyrene	15.36	252	1331018	47.486	nq	100
91) Dibenzo(a,h)anthracene	16.87	278	1250356	43.305	nq	99
92) Benzo(g,h,i)perylene	17.29	276	1312426	46.100	nq	100

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

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