

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022720\
 Data File : BF119116.D
 Acq On : 27 Feb 2020 18:12
 Operator : CG/JU
 Sample : L1688-02MS
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 GP-02-022420MS

Manual Integrations
 APPROVED

mohammad
 2/28/2020 2:41:40 PM

Quant Time: Feb 28 04:11:23 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 26 16:03:05 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.75	152	146028	20.00	ng	0.00
21) Naphthalene-d8	8.03	136	542837	20.00	ng	0.00
39) Acenaphthene-d10	9.78	164	289530	20.00	ng	0.00
64) Phenanthrene-d10	11.26	188	501604	20.00	ng	0.00
76) Chrysene-d12	13.90	240	322127	20.00	ng	0.00
87) Perylene-d12	15.33	264	352977	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.37	112	878938	100.59	ng	0.00
7) Phenol-d6	6.40	99	1055662	99.34	ng	0.00
23) Nitrobenzene-d5	7.31	82	702749	68.35	ng	-0.01
42) 2,4,6-Tribromophenol	10.57	330	337352	89.07	ng	0.00
45) 2-Fluorobiphenyl	9.10	172	1319663	67.15	ng	0.00
79) Terphenyl-d14	12.85	244	1374790	73.48	ng	0.00

Target Compounds					Qvalue
2) 1,4-Dioxane	2.49	88	157604	40.510 ng	99
3) Pyridine	3.22	79	398649	37.520 ng	97
4) n-Nitrosodimethylamine	3.16	42	162057	50.350 ng	96
6) Aniline	6.41	93	421515	32.145 ng	89
8) 2-Chlorophenol	6.54	128	485877	51.525 ng	97
9) Benzaldehyde	6.29	77	196477	27.672 ng	99
10) Phenol	6.41	94	554194	48.234 ng	96
11) bis(2-Chloroethyl)ether	6.48	93	432993	49.252 ng	97
12) 1,3-Dichlorobenzene	6.69	146	526239	46.560 ng	97
13) 1,4-Dichlorobenzene	6.76	146	533338	47.156 ng	99
14) 1,2-Dichlorobenzene	6.92	146	490775	47.579 ng	99
15) Benzyl Alcohol	6.90	79	389292	50.685 ng	99
16) 2,2'-oxybis(1-Chloropropan	7.03	45	346798	45.539 ng	97
17) 2-Methylphenol	7.02	107	375785	49.152 ng	100
18) Hexachloroethane	7.26	117	185561	45.858 ng	97
19) n-Nitroso-di-n-propylamine	7.17	70	306349	49.472 ng	96
20) 3+4-Methylphenols	7.17	107	466156	49.767 ng	95
22) Acetophenone	7.16	105	623716	49.660 ng	# 96
24) Nitrobenzene	7.33	77	487216	47.922 ng	99
25) Isophorone	7.57	82	863636	48.369 ng	100
26) 2-Nitrophenol	7.65	139	259220	48.121 ng	94
27) 2,4-Dimethylphenol	7.69	122	395057	54.036 ng	98
28) bis(2-Chloroethoxy)methane	7.78	93	525609	45.226 ng	99
29) 2,4-Dichlorophenol	7.90	162	401776	46.681 ng	98
30) 1,2,4-Trichlorobenzene	7.97	180	456624	44.747 ng	99
31) Naphthalene	8.05	128	1333163	47.355 ng	98
32) Benzoic acid	7.85	122	241517	45.653 ng	99
33) 4-Chloroaniline	8.10	127	171690	14.179 ng	96
34) Hexachlorobutadiene	8.17	225	279609	43.989 ng	98
35) Caprolactam	8.48	113	124579m	47.008 ng	
36) 4-Chloro-3-methylphenol	8.60	107	417418	47.922 ng	95
37) 2-Methylnaphthalene	8.74	142	903057	47.665 ng	100
38) 1-Methylnaphthalene	8.84	142	849749	47.692 ng	97
40) 1,2,4,5-Tetrachlorobenzene	8.91	216	453489	46.456 ng	98

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41) Hexachlorocyclopentadiene	8.90	237	238173	68.553	nq	100
43) 2,4,6-Trichlorophenol	9.03	196	287205	46.590	nq	99
44) 2,4,5-Trichlorophenol	9.07	196	301438	46.599	nq	97
46) 1,1'-Biphenyl	9.20	154	1081958	46.238	nq	98
47) 2-Chloronaphthalene	9.23	162	862995	45.766	nq	98
48) 2-Nitroaniline	9.33	65	235891	44.850	nq	98
49) Acenaphthylene	9.65	152	1299210	47.704	nq	100
50) Dimethylphthalate	9.51	163	1111978	50.225	nq	100
51) 2,6-Dinitrotoluene	9.57	165	233595	47.490	nq	98
52) Acenaphthene	9.82	154	766184	46.655	nq	99
53) 3-Nitroaniline	9.74	138	111540	20.454	nq	98
54) 2,4-Dinitrophenol	9.86	184	129066	56.418	nq	# 91
55) Dibenzofuran	9.99	168	1151197	46.965	nq	99
56) 4-Nitrophenol	9.92	139	267071	81.515	nq	92
57) 2,4-Dinitrotoluene	9.98	165	289370	45.386	nq	96
58) Fluorene	10.33	166	899894	47.246	nq	98
59) 2,3,4,6-Tetrachlorophenol	10.12	232	253059	46.362	nq	98
60) Diethylphthalate	10.20	149	990228	47.461	nq	99
61) 4-Chlorophenyl-phenylether	10.32	204	472486	46.865	nq	94
62) 4-Nitroaniline	10.36	138	198427	35.566	nq	98
63) Azobenzene	10.48	77	866304	47.769	nq	97
65) 4,6-Dinitro-2-methylphenol	10.39	198	119488	39.366	nq	96
66) n-Nitrosodiphenylamine	10.44	169	815447	49.649	nq	99
67) 4-Bromophenyl-phenylether	10.81	248	308776	47.094	nq	94
68) Hexachlorobenzene	10.88	284	345862	45.752	nq	96
69) Atrazine	10.97	200	287642	50.125	nq	98
70) Pentachlorophenol	11.08	266	259842	85.331	nq	100
71) Phenanthrene	11.29	178	1296946	47.412	nq	99
72) Anthracene	11.35	178	1342885	49.132	nq	98
73) Carbazole	11.50	167	1146850	47.099	nq	98
74) Di-n-butylphthalate	11.83	149	1466675	49.739	nq	98
75) Fluoranthene	12.47	202	1306384	44.991	nq	99
77) Benzidine	12.60	184	593430	45.298	nq	98
78) Pyrene	12.70	202	1281179	49.531	nq	99
80) Butylbenzylphthalate	13.32	149	561062	51.538	nq	95
81) Benzo(a)anthracene	13.89	228	1071409	48.814	nq	98
82) 3,3'-Dichlorobenzidine	13.85	252	373945	43.876	nq	99
83) Chrysene	13.93	228	1053855	48.187	nq	98
84) Bis(2-ethylhexyl)phthalate	13.88	149	769113	55.921	nq	99
85) Di-n-octyl phthalate	14.49	149	1253636	57.209	nq	99
86) Indeno(1,2,3-cd)pyrene	16.74	276	1264022	59.691	nq	100
88) Benzo(b)fluoranthene	14.92	252	1049801	45.121	nq	99
89) Benzo(k)fluoranthene	14.95	252	1069090	49.584	nq	99
90) Benzo(a)pyrene	15.27	252	982293	46.779	nq	98
91) Dibenzo(a,h)anthracene	16.74	278	1068766	57.846	nq	100
92) Benzo(g,h,i)perylene	17.16	276	1070326	58.631	nq	100

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

