

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF070920\
 Data File : BF120816.D
 Acq On : 9 Jul 2020 13:57
 Operator : JU/CG
 Sample : L3216-08
 Misc : LOQ-WATER-10PPM
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LOQ-WATER-02-QT3-2020

Manual Integrations
 APPROVED

mohammad
 7/10/2020 11:49:44 AM

Quant Time: Jul 09 14:35:09 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF070220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 09 11:36:34 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	134145	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	506629	20.00	ng	0.00
39) Acenaphthene-d10	9.89	164	250614	20.00	ng	0.00
64) Phenanthrene-d10	11.37	188	445263	20.00	ng	0.00
76) Chrysene-d12	14.00	240	334955	20.00	ng	0.00
86) Perylene-d12	15.46	264	271616	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.46	112	846789	97.44	ng	0.00
7) Phenol-d6	6.47	99	1079081	97.24	ng	0.00
23) Nitrobenzene-d5	7.41	82	694070	79.37	ng	0.00
42) 2,4,6-Tribromophenol	10.67	330	262762	104.36	ng	0.00
45) 2-Fluorobiphenyl	9.21	172	1268925	76.60	ng	0.00
79) Terphenyl-d14	12.96	244	1050237	61.39	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.58	88	40102	10.102	ng	100
3) Pyridine	3.36	79	78010	7.175	ng	98
4) n-Nitrosodimethylamine	3.26	42	47242	8.685	ng	96
6) Aniline	6.50	93	132628	9.306	ng	100
8) 2-Chlorophenol	6.63	128	84084	9.088	ng	95
9) Benzaldehyde	6.39	77	73834	11.187	ng	98
10) Phenol	6.48	94	107037	9.376	ng	95
11) bis(2-Chloroethyl)ether	6.58	93	84344	9.079	ng	99
12) 1,3-Dichlorobenzene	6.79	146	93607	9.192	ng	96
13) 1,4-Dichlorobenzene	6.86	146	94941	9.308	ng	98
14) 1,2-Dichlorobenzene	7.02	146	89901	9.264	ng	99
15) Benzyl Alcohol	6.98	79	81081	9.724	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.12	45	125147	8.637	ng	100
17) 2-Methylphenol	7.09	107	68099	8.228	ng	96
18) Hexachloroethane	7.36	117	34764	9.068	ng	97
19) n-Nitroso-di-n-propylamine	7.25	70	59950	8.972	ng	97
20) 3+4-Methylphenols	7.24	107	86738	8.509	ng	93
22) Acetophenone	7.25	105	126896	10.077	ng	95
24) Nitrobenzene	7.42	77	96261	9.804	ng	94
25) Isophorone	7.66	82	161546	8.566	ng	97
26) 2-Nitrophenol	7.74	139	31053	8.652	ng	94
27) 2,4-Dimethylphenol	7.77	122	71455	9.492	ng	99
28) bis(2-Chloroethoxy)methane	7.87	93	101530	9.097	ng	98
29) 2,4-Dichlorophenol	7.98	162	65142	8.668	ng	98
30) 1,2,4-Trichlorobenzene	8.07	180	74195	8.790	ng	99
31) Naphthalene	8.15	128	254287	9.392	ng	98
32) Benzoic acid	7.83	122	25591m	6.281	ng	
33) 4-Chloroaniline	8.19	127	98052	8.534	ng	94
34) Hexachlorobutadiene	8.27	225	43011	8.472	ng	98
35) Caprolactam	8.53	113	17470	7.978	ng	91
36) 4-Chloro-3-methylphenol	8.67	107	66091	8.454	ng	97
37) 2-Methylnaphthalene	8.84	142	162275	9.220	ng	98
38) 1-Methylnaphthalene	8.94	142	151796	9.207	ng	98
40) 1,2,4,5-Tetrachlorobenzene	9.00	216	71176	9.397	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	9.00	237	34634	7.739	ng	98
43) 2,4,6-Trichlorophenol	9.12	196	40107	7.771	ng	99
44) 2,4,5-Trichlorophenol	9.15	196	44736	7.810	ng	97
46) 1,1'-Biphenyl	9.30	154	204504	10.116	ng	98
47) 2-Chloronaphthalene	9.33	162	144104	8.898	ng	99
48) 2-Nitroaniline	9.42	65	40085	9.251	ng	87
49) Acenaphthylene	9.75	152	230986	9.097	ng	98
50) Dimethylphthalate	9.60	163	161387	8.828	ng	99
51) 2,6-Dinitrotoluene	9.66	165	30001	9.084	ng	88
52) Acenaphthene	9.92	154	136974	8.696	ng	98
53) 3-Nitroaniline	9.83	138	34433	8.439	ng	95
54) 2,4-Dinitrophenol	9.93	184	4883	4.477	ng	# 22
55) Dibenzofuran	10.09	168	208908	9.201	ng	100
56) 4-Nitrophenol	9.97	139	23024	7.682	ng	95
57) 2,4-Dinitrotoluene	10.06	165	35851	9.041	ng	95
58) Fluorene	10.43	166	160408	9.491	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.20	232	36135	8.478	ng	89
60) Diethylphthalate	10.30	149	162714	8.544	ng	99
61) 4-Chlorophenyl-phenylether	10.42	204	76654	9.022	ng	93
62) 4-Nitroaniline	10.43	138	33873	8.720	ng	93
63) Azobenzene	10.58	77	175807	9.024	ng	93
65) 4,6-Dinitro-2-methylphenol	10.46	198	8813	6.160	ng	99
66) n-Nitrosodiphenylamine	10.54	169	133342	8.935	ng	98
67) 4-Bromophenyl-phenylether	10.92	248	41389	8.013	ng	97
68) Hexachlorobenzene	10.97	284	46326	8.547	ng	99
69) Atrazine	11.06	200	37167	9.667	ng	97
70) Pentachlorophenol	11.17	266	19264	6.282	ng	98
71) Phenanthrene	11.39	178	219841	9.100	ng	98
72) Anthracene	11.44	178	222402	9.110	ng	99
73) Carbazole	11.59	167	200266	8.536	ng	98
74) Di-n-butylphthalate	11.93	149	228443	7.865	ng	98
75) Fluoranthene	12.57	202	219310	8.521	ng	97
77) Benzidine	12.70	184	89231	9.438	ng	# 95
78) Pyrene	12.80	202	228184	8.530	ng	98
80) Butylbenzylphthalate	13.43	149	76948	6.662	ng	99
81) Benzo(a)anthracene	13.99	228	191637	8.925	ng	99
82) 3,3'-Dichlorobenzidine	13.96	252	57841	7.983	ng	# 98
83) Chrysene	14.03	228	181933	8.307	ng	98
84) Bis(2-ethylhexyl)phthalate	13.99	149	119163	7.868	ng	100
85) Di-n-octyl phthalate	14.60	149	154958	5.783	ng	98
87) Indeno(1,2,3-cd)pyrene	16.90	276	133351	7.654	ng	97
88) Benzo(b)fluoranthene	15.04	252	144813	8.110	ng	98
89) Benzo(k)fluoranthene	15.06	252	154808	8.910	ng	98
90) Benzo(a)pyrene	15.40	252	123565	7.757	ng	100
91) Dibenzo(a,h)anthracene	16.93	278	115443	7.950	ng	99
92) Benzo(g,h,i)perylene	17.34	276	110968	8.362	ng	96

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

