

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG070820\
 Data File : BG045982.D
 Acq On : 8 Jul 2020 15:56
 Operator : CG/JU
 Sample : L3216-08
 Misc : LOO-WTR-10PPM
 ALS Vial : 10 Sample Multiplier: 1

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

Quant Time: Jul 08 16:33:04 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG070620.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 08 15:53:46 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.01	152	145437	20.00	ng	0.00
21) Naphthalene-d8	10.81	136	590423	20.00	ng	0.00
39) Acenaphthene-d10	14.63	164	354144	20.00	ng	0.00
64) Phenanthrene-d10	17.37	188	730172	20.00	ng	0.00
76) Chrysene-d12	21.63	240	673365	20.00	ng	0.00
86) Perylene-d12	24.78	264	677406	20.00	ng	-0.01

System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	879734	98.09	ng	0.00
7) Phenol-d6	7.17	99	1276175	98.84	ng	0.00
23) Nitrobenzene-d5	9.16	82	856739	83.52	ng	0.00
42) 2,4,6-Tribromophenol	16.11	330	334523	101.78	ng	0.00
45) 2-Fluorobiphenyl	13.26	172	1660363	73.13	ng	0.00
79) Terphenyl-d14	19.98	244	1792112	59.23	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.51	88	45174	10.918	ng	94
3) Pyridine	3.91	79	97005	7.754	ng	95
4) n-Nitrosodimethylamine	3.82	42	44233	10.456	ng	93
6) Aniline	7.33	93	157562	9.539	ng	98
8) 2-Chlorophenol	7.57	128	87676	9.082	ng	97
9) Benzaldehyde	7.14	77	140328	17.154	ng	96
10) Phenol	7.19	94	118613	9.158	ng	96
11) bis(2-Chloroethyl)ether	7.42	93	96773	9.035	ng	100
12) 1,3-Dichlorobenzene	7.90	146	101575	9.128	ng	98
13) 1,4-Dichlorobenzene	8.04	146	103105	9.419	ng	98
14) 1,2-Dichlorobenzene	8.36	146	93113	8.838	ng	99
15) Benzyl Alcohol	8.23	79	90518	8.941	ng	96
16) 2,2'-oxybis(1-Chloropropan	8.54	45	126809	9.075	ng	98
17) 2-Methylphenol	8.44	107	81983	8.436	ng	98
18) Hexachloroethane	9.09	117	38631	9.172	ng	93
19) n-Nitroso-di-n-propylamine	8.80	70	84927	9.363	ng	98
20) 3+4-Methylphenols	8.76	107	111594	8.425	ng	98
22) Acetophenone	8.82	105	225429	14.292	ng	# 96
24) Nitrobenzene	9.20	77	114489	10.112	ng	97
25) Isophorone	9.72	82	218741	8.735	ng	99
26) 2-Nitrophenol	9.91	139	35291	9.563	ng	95
27) 2,4-Dimethylphenol	9.97	122	86382	9.691	ng	96
28) bis(2-Chloroethoxy)methane	10.21	93	129366	9.200	ng	94
29) 2,4-Dichlorophenol	10.44	162	80719	8.748	ng	98
30) 1,2,4-Trichlorobenzene	10.67	180	91452	8.945	ng	96
31) Naphthalene	10.85	128	292171	9.231	ng	97
32) Benzoic acid	10.05	122	50128	10.386	ng	92
33) 4-Chloroaniline	10.95	127	123098	8.703	ng	97
34) Hexachlorobutadiene	11.15	225	50266	8.420	ng	96
35) Caprolactam	11.69	113	30176	8.614	ng	93
36) 4-Chloro-3-methylphenol	12.07	107	93696	8.915	ng	100
37) 2-Methylnaphthalene	12.46	142	211168	9.385	ng	98
38) 1-Methylnaphthalene	12.68	142	200085	9.423	ng	98
40) 1,2,4,5-Tetrachlorobenzene	12.83	216	149821	15.008	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.82	237	52036	8.785	nq	92
43) 2,4,6-Trichlorophenol	13.06	196	58807	8.710	nq	99
44) 2,4,5-Trichlorophenol	13.13	196	64959	8.510	nq	97
46) 1,1'-Biphenyl	13.46	154	421376	15.817	nq	98
47) 2-Chloronaphthalene	13.50	162	191902	9.130	nq	99
48) 2-Nitroaniline	13.69	65	53533	10.882	nq	98
49) Acenaphthylene	14.35	152	322861	9.520	nq	96
50) Dimethylphthalate	14.08	163	240435	9.047	nq	99
51) 2,6-Dinitrotoluene	14.19	165	43774	10.615	nq	94
52) Acenaphthene	14.69	154	196361	9.179	nq	99
53) 3-Nitroaniline	14.51	138	50980	9.976	nq	# 91
54) 2,4-Dinitrophenol	14.72	184	16847	11.581	nq	# 87
55) Dibenzofuran	15.03	168	298130	9.508	nq	96
56) 4-Nitrophenol	14.81	139	39541	9.132	nq	92
57) 2,4-Dinitrotoluene	14.97	165	56793	11.145	nq	88
58) Fluorene	15.67	166	240884	9.364	nq	99
59) 2,3,4,6-Tetrachlorophenol	15.25	232	55605	9.434	nq	92
60) Diethylphthalate	15.45	149	255888	9.181	nq	98
61) 4-Chlorophenyl-phenylether	15.67	204	111170	8.687	nq	98
62) 4-Nitroaniline	15.67	138	56462	10.279	nq	89
63) Azobenzene	15.97	77	286008	9.560	nq	99
65) 4,6-Dinitro-2-methylphenol	15.74	198	17649	8.253	nq	89
66) n-Nitrosodiphenylamine	15.88	169	215112	9.507	nq	96
67) 4-Bromophenyl-phenylether	16.56	248	70612	8.659	nq	96
68) Hexachlorobenzene	16.68	284	74597	9.116	nq	98
69) Atrazine	16.83	200	73699	11.017	nq	96
70) Pentachlorophenol	17.02	266	36468	8.015	nq	93
71) Phenanthrene	17.41	178	375838	9.736	nq	97
72) Anthracene	17.50	178	384343	9.811	nq	94
73) Carbazole	17.76	167	366646	9.337	nq	97
74) Di-n-butylphthalate	18.34	149	452191	9.530	nq	95
75) Fluoranthene	19.42	202	435361	9.756	nq	96
77) Benzidine	19.60	184	225724	11.724	nq	100
78) Pyrene	19.78	202	446341	9.711	nq	95
80) Butylbenzylphthalate	20.68	149	186582	8.590	nq	98
81) Benzo(a)anthracene	21.60	228	412626	9.319	nq	93
82) 3,3'-Dichlorobenzidine	21.52	252	155653	9.514	nq	98
83) Chrysene	21.67	228	389341	9.230	nq	95
84) Bis(2-ethylhexyl)phthalate	21.55	149	281665	8.933	nq	93
85) Di-n-octyl phthalate	22.75	149	455028	8.316	nq	98
87) Indeno(1,2,3-cd)pyrene	28.35	276	453462	9.150	nq	# 92
88) Benzo(b)fluoranthene	23.78	252	410625	9.549	nq	98
89) Benzo(k)fluoranthene	23.85	252	394085	9.692	nq	97
90) Benzo(a)pyrene	24.64	252	367394	9.086	nq	95
91) Dibenzo(a,h)anthracene	28.42	278	375631	9.398	nq	98
92) Benzo(g,h,i)perylene	29.46	276	378407	9.407	nq	97

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

