

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG031320\
 Data File : BG044787.D
 Acq On : 14 Mar 2020 6:50
 Operator : CG/JU
 Sample : L1761-10MS
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SU-01-031220MS

Manual Integrations
 APPROVED

mohammad
 3/18/2020 8:43:46 AM

Quant Time: Mar 16 02:11:00 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG031020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 10 16:27:39 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.06	152	106954	20.00	ng	0.00
21) Naphthalene-d8	10.86	136	437950	20.00	ng	0.00
39) Acenaphthene-d10	14.69	164	292095	20.00	ng	0.00
64) Phenanthrene-d10	17.43	188	627015	20.00	ng	0.00
76) Chrysene-d12	21.70	240	499191	20.00	ng	0.00
87) Perylene-d12	24.92	264	580096	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.62	112	860429	125.23	ng	0.00
7) Phenol-d6	7.21	99	1108340	111.03	ng	0.00
23) Nitrobenzene-d5	9.21	82	881163	88.29	ng	0.00
42) 2,4,6-Tribromophenol	16.18	330	526100	137.56	ng	0.00
45) 2-Fluorobiphenyl	13.31	172	1754344	86.01	ng	0.00
79) Terphenyl-d14	20.04	244	2407252	90.20	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.54	88	110043	33.260	ng	# 23
3) Pyridine	3.94	79	216185	22.072	ng	100
4) n-Nitrosodimethylamine	3.83	42	154371	41.540	ng	99
6) Aniline	7.37	93	240995	19.402	ng	96
8) 2-Chlorophenol	7.62	128	352116	51.339	ng	96
9) Benzaldehyde	7.18	77	277393	48.376	ng	98
10) Phenol	7.24	94	427558	43.262	ng	99
11) bis(2-Chloroethyl)ether	7.47	93	351332	44.543	ng	99
12) 1,3-Dichlorobenzene	7.94	146	322219	38.138	ng	94
13) 1,4-Dichlorobenzene	8.09	146	319686	38.273	ng	98
14) 1,2-Dichlorobenzene	8.41	146	307004	38.754	ng	95
15) Benzyl Alcohol	8.28	79	376344	46.988	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.58	45	546543	39.266	ng	98
17) 2-Methylphenol	8.49	107	319142	47.672	ng	95
18) Hexachloroethane	9.15	117	120835	36.745	ng	95
19) n-Nitroso-di-n-propylamine	8.86	70	311148	44.143	ng	95
20) 3+4-Methylphenols	8.82	107	431350	46.742	ng	99
22) Acetophenone	8.87	105	549262	44.270	ng	98
24) Nitrobenzene	9.25	77	472018	45.580	ng	97
25) Isophorone	9.78	82	861497	44.228	ng	97
26) 2-Nitrophenol	9.97	139	193556	53.167	ng	97
27) 2,4-Dimethylphenol	10.03	122	337263	53.169	ng	95
28) bis(2-Chloroethoxy)methane	10.26	93	488297	42.006	ng	99
29) 2,4-Dichlorophenol	10.51	162	360392	48.642	ng	96
30) 1,2,4-Trichlorobenzene	10.72	180	360383	40.680	ng	99
31) Naphthalene	10.91	128	1030530	42.478	ng	99
32) Benzoic acid	10.16	122	180712	38.872	ng	97
33) 4-Chloroaniline	11.01	127	71156	6.510	ng	99
34) Hexachlorobutadiene	11.21	225	228680	39.252	ng	98
35) Caprolactam	11.78	113	95620m	34.087	ng	
36) 4-Chloro-3-methylphenol	12.13	107	417302	47.225	ng	98
37) 2-Methylnaphthalene	12.52	142	788769	43.464	ng	98
38) 1-Methylnaphthalene	12.73	142	749678	43.941	ng	96
40) 1,2,4,5-Tetrachlorobenzene	12.89	216	434597	45.178	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.87	237	207672	39.503	nq	96
43) 2,4,6-Trichlorophenol	13.12	196	319902	50.110	nq	99
44) 2,4,5-Trichlorophenol	13.19	196	334900	50.584	nq	96
46) 1,1'-Biphenyl	13.52	154	1004167	43.020	nq	99
47) 2-Chloronaphthalene	13.56	162	770567	43.214	nq	99
48) 2-Nitroaniline	13.75	65	297220	47.611	nq	94
49) Acenaphthylene	14.41	152	1272249	45.543	nq	100
50) Dimethylphthalate	14.14	163	1083519	46.575	nq	99
51) 2,6-Dinitrotoluene	14.25	165	238218	50.800	nq	96
52) Acenaphthene	14.75	154	804633	43.981	nq	98
53) 3-Nitroaniline	14.58	138	129793	24.075	nq	99
54) 2,4-Dinitrophenol	14.78	184	64717	31.986	nq	92
55) Dibenzofuran	15.08	168	1214124	45.764	nq	99
56) 4-Nitrophenol	14.88	139	352634	85.704	nq	96
57) 2,4-Dinitrotoluene	15.04	165	319452	49.672	nq	89
58) Fluorene	15.73	166	980749	44.939	nq	97
59) 2,3,4,6-Tetrachlorophenol	15.31	232	303868m	49.852	nq	
60) Diethylphthalate	15.51	149	1043816	43.019	nq	99
61) 4-Chlorophenyl-phenylether	15.73	204	537703	45.757	nq	98
62) 4-Nitroaniline	15.74	138	206149	35.859	nq	98
63) Azobenzene	16.02	77	1102800	43.767	nq	98
65) 4,6-Dinitro-2-methylphenol	15.81	198	56757	23.867	nq	98
66) n-Nitrosodiphenylamine	15.94	169	888269	47.054	nq	99
67) 4-Bromophenyl-phenylether	16.62	248	370450	47.260	nq	96
68) Hexachlorobenzene	16.75	284	399568	46.986	nq	97
69) Atrazine	16.89	200	342565	49.531	nq	98
70) Pentachlorophenol	17.09	266	476063	101.744	nq	99
71) Phenanthrene	17.47	178	1532337	45.732	nq	99
72) Anthracene	17.57	178	1570401	47.531	nq	100
73) Carbazole	17.83	167	1441444	45.419	nq	99
74) Di-n-butylphthalate	18.40	149	1674549	43.405	nq	100
75) Fluoranthene	19.48	202	1756116	44.018	nq	99
77) Benzidine	19.65	184	599048	36.965	nq	98
78) Pyrene	19.84	202	1753405	47.876	nq	98
80) Butylbenzylphthalate	20.73	149	758590	46.559	nq	99
81) Benzo(a)anthracene	21.69	228	1674397	46.582	nq	99
82) 3,3'-Dichlorobenzidine	21.60	252	474241	34.103	nq	99
83) Chrysene	21.75	228	1581777	46.068	nq	97
84) Bis(2-ethylhexyl)phthalate	21.61	149	1034728	46.015	nq	100
85) Di-n-octyl phthalate	22.83	149	1763334	46.861	nq	100
86) Indeno(1,2,3-cd)pyrene	28.59	276	1950632	43.333	nq	98
88) Benzo(b)fluoranthene	23.91	252	1733867	45.275	nq	98
89) Benzo(k)fluoranthene	23.98	252	1706307	47.081	nq	98
90) Benzo(a)pyrene	24.78	252	1607687	44.864	nq	99
91) Dibenzo(a,h)anthracene	28.66	278	1605525	45.415	nq	99
92) Benzo(g,h,i)perylene	29.72	276	1513493	42.374	nq	98

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

