

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG031920\
 Data File : BG044905.D
 Acq On : 19 Mar 2020 20:15
 Operator : CG/JU
 Sample : L1964-03MS
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 2000448388MS

Manual Integrations
 APPROVED

mohammad
 3/20/2020 11:48:04 AM

Quant Time: Mar 20 02:15:34 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG031020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 16 17:37:15 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.04	152	81306	20.00	ng	0.00
21) Naphthalene-d8	10.85	136	308044	20.00	ng	0.00
39) Acenaphthene-d10	14.68	164	211922	20.00	ng	0.00
64) Phenanthrene-d10	17.42	188	487279	20.00	ng	0.00
76) Chrysene-d12	21.69	240	452614	20.00	ng	-0.01
87) Perylene-d12	24.90	264	539770	20.00	ng	-0.01

System Monitoring Compounds						
5) 2-Fluorophenol	5.62	112	619689	118.65	ng	0.00
7) Phenol-d6	7.20	99	781254	102.95	ng	0.00
23) Nitrobenzene-d5	9.20	82	624947	89.02	ng	0.00
42) 2,4,6-Tribromophenol	16.16	330	418700	150.89	ng	0.00
45) 2-Fluorobiphenyl	13.30	172	1276227	86.24	ng	0.00
79) Terphenyl-d14	20.03	244	2261791	93.47	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.53	88	86156	34.255	ng	# 24
3) Pyridine	3.93	79	167977	22.560	ng	99
4) n-Nitrosodimethylamine	3.83	42	120706	42.727	ng	99
6) Aniline	7.36	93	132802	14.064	ng	94
8) 2-Chlorophenol	7.61	128	254200	48.754	ng	98
9) Benzaldehyde	7.17	77	203954	46.036	ng	95
10) Phenol	7.23	94	285603	38.015	ng	96
11) bis(2-Chloroethyl)ether	7.46	93	247376	41.257	ng	93
12) 1,3-Dichlorobenzene	7.94	146	267614	41.667	ng	96
13) 1,4-Dichlorobenzene	8.08	146	269105	42.381	ng	97
14) 1,2-Dichlorobenzene	8.40	146	248137	41.204	ng	99
15) Benzyl Alcohol	8.28	79	259397	42.603	ng	93
16) 2,2'-oxybis(1-Chloropropan	8.57	45	368185	34.796	ng	98
17) 2-Methylphenol	8.49	107	220657	43.358	ng	98
18) Hexachloroethane	9.14	117	102112	40.847	ng	93
19) n-Nitroso-di-n-propylamine	8.85	70	208959	38.996	ng	98
20) 3+4-Methylphenols	8.81	107	303842	43.311	ng	97
22) Acetophenone	8.86	105	380980	43.656	ng	# 98
24) Nitrobenzene	9.24	77	322579	44.286	ng	100
25) Isophorone	9.77	82	596330	43.525	ng	99
26) 2-Nitrophenol	9.96	139	137502	53.643	ng	98
27) 2,4-Dimethylphenol	10.02	122	239336	53.642	ng	94
28) bis(2-Chloroethoxy)methane	10.25	93	328926	40.229	ng	99
29) 2,4-Dichlorophenol	10.49	162	253703	48.683	ng	98
30) 1,2,4-Trichlorobenzene	10.71	180	272823	43.783	ng	96
31) Naphthalene	10.90	128	753683	44.167	ng	98
32) Benzoic acid	10.15	122	109475	34.821	ng	97
33) 4-Chloroaniline	11.01	127	25061	3.260	ng	# 91
34) Hexachlorobutadiene	11.20	225	182726	44.591	ng	97
35) Caprolactam	11.76	113	65664m	33.280	ng	
36) 4-Chloro-3-methylphenol	12.12	107	292907	47.127	ng	96
37) 2-Methylnaphthalene	12.50	142	567366	44.448	ng	95
38) 1-Methylnaphthalene	12.73	142	541553	45.128	ng	95
40) 1,2,4,5-Tetrachlorobenzene	12.87	216	300096	42.998	ng	99

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG031920\
 Data File : BG044905.D
 Acq On : 19 Mar 2020 20:15
 Operator : CG/JU
 Sample : L1964-03MS
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 2000448388MS

Manual Integrations
 APPROVED

mohammad
 3/20/2020 11:48:04 AM

Quant Time: Mar 20 02:15:34 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG031020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 16 17:37:15 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.86	237	368666	96.656	nq	99
43) 2,4,6-Trichlorophenol	13.11	196	225965	48.786	nq	98
44) 2,4,5-Trichlorophenol	13.18	196	242301	50.443	nq	94
46) 1,1'-Biphenyl	13.51	154	709904	41.919	nq	99
47) 2-Chloronaphthalene	13.56	162	548022	42.361	nq	99
48) 2-Nitroaniline	13.74	65	202789	44.774	nq	98
49) Acenaphthylene	14.40	152	917867	45.287	nq	98
50) Dimethylphthalate	14.13	163	794112	47.048	nq	99
51) 2,6-Dinitrotoluene	14.24	165	173550	51.011	nq	97
52) Acenaphthene	14.74	154	660255	49.742	nq	99
53) 3-Nitroaniline	14.57	138	53695	13.727	nq	99
54) 2,4-Dinitrophenol	14.77	184	179113	98.589	nq	92
55) Dibenzofuran	15.08	168	878626	45.647	nq	99
56) 4-Nitrophenol	14.88	139	261943	87.747	nq	92
57) 2,4-Dinitrotoluene	15.03	165	246226	52.770	nq	97
58) Fluorene	15.72	166	727551	45.949	nq	98
59) 2,3,4,6-Tetrachlorophenol	15.30	232	229105	51.806	nq	# 93
60) Diethylphthalate	15.49	149	786151	44.657	nq	98
61) 4-Chlorophenyl-phenylether	15.72	204	396539	46.510	nq	98
62) 4-Nitroaniline	15.74	138	173052	41.490	nq	95
63) Azobenzene	16.01	77	780180	42.677	nq	99
65) 4,6-Dinitro-2-methylphenol	15.79	198	138587	58.435	nq	96
66) n-Nitrosodiphenylamine	15.93	169	665266	45.347	nq	98
67) 4-Bromophenyl-phenylether	16.61	248	275230	45.181	nq	98
68) Hexachlorobenzene	16.73	284	297515	45.018	nq	99
69) Atrazine	16.88	200	268820	50.014	nq	98
70) Pentachlorophenol	17.08	266	368548	101.353	nq	98
71) Phenanthrene	17.46	178	1204226	46.246	nq	100
72) Anthracene	17.56	178	1237590	48.200	nq	99
73) Carbazole	17.82	167	1170047	47.440	nq	99
74) Di-n-butylphthalate	18.39	149	1388598	46.315	nq	99
75) Fluoranthene	19.47	202	1534152	49.482	nq	99
77) Benzidine	19.64	184	185375	12.616	nq	98
78) Pyrene	19.83	202	1529736	46.067	nq	99
80) Butylbenzylphthalate	20.72	149	655480	44.370	nq	96
81) Benzo(a)anthracene	21.67	228	1498273	45.971	nq	98
82) 3,3'-Dichlorobenzidine	21.58	252	186113	14.761	nq	97
83) Chrysene	21.74	228	1408144	45.231	nq	99
84) Bis(2-ethylhexyl)phthalate	21.59	149	907743	44.522	nq	99
85) Di-n-octyl phthalate	22.81	149	1560395	45.735	nq	100
86) Indeno(1,2,3-cd)pyrene	28.54	276	1968233	48.223	nq	# 98
88) Benzo(b)fluoranthene	23.88	252	1628403	45.698	nq	99
89) Benzo(k)fluoranthene	23.95	252	1570453	46.570	nq	99
90) Benzo(a)pyrene	24.75	252	1504788	45.130	nq	96
91) Dibenzo(a,h)anthracene	28.60	278	1580274	48.040	nq	98
92) Benzo(g,h,i)perylene	29.68	276	1623158	48.839	nq	98

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

