

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG112118\
 Data File : BG038126.D
 Acq On : 21 Nov 2018 20:54
 Operator : JU/SJ
 Sample : J5765-04MS
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 NB-08-112018MS

Manual Integrations
 APPROVED

mohammad
 11/23/2018 11:50:17 AM

Quant Time: Nov 22 01:38:55 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG111318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 13 16:39:28 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.09	152	39335	20.00	ng	0.00
21) Naphthalene-d8	10.91	136	173245	20.00	ng	0.00
39) Acenaphthene-d10	14.72	164	109948	20.00	ng	0.00
64) Phenanthrene-d10	17.47	188	249589	20.00	ng	0.00
76) Chrysene-d12	21.75	240	230428	20.00	ng	0.00
87) Perylene-d12	25.07	264	245304	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.63	112	256605	112.63	ng	0.00
7) Phenol-d6	7.24	99	329783	101.41	ng	0.00
23) Nitrobenzene-d5	9.26	82	251404	77.68	ng	0.00
42) 2,4,6-Tribromophenol	16.21	330	139664	111.38	ng	0.00
45) 2-Fluorobiphenyl	13.34	172	555078	73.55	ng	0.00
79) Terphenyl-d14	20.07	244	840966	85.87	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.53	88	34468	32.031	ng	# 1
3) Pyridine	3.93	79	99886	32.540	ng	94
4) n-Nitrosodimethylamine	3.85	42	54120	48.597	ng	93
6) Aniline	7.41	93	65090	15.508	ng	95
8) 2-Chlorophenol	7.65	128	120976	45.763	ng	97
9) Benzaldehyde	7.22	77	60526	25.736	ng	99
10) Phenol	7.26	94	134301	38.989	ng	94
11) bis(2-Chloroethyl)ether	7.51	93	120982	43.021	ng	99
12) 1,3-Dichlorobenzene	7.97	146	124346	40.831	ng	96
13) 1,4-Dichlorobenzene	8.12	146	125735	40.872	ng	97
14) 1,2-Dichlorobenzene	8.44	146	123231	41.957	ng	99
15) Benzyl Alcohol	8.32	79	108762	46.220	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.60	45	246232	47.154	ng	98
17) 2-Methylphenol	8.52	107	105429	45.271	ng	98
18) Hexachloroethane	9.17	117	47024	42.001	ng	98
19) n-Nitroso-di-n-propylamine	8.89	70	101017	42.929	ng	96
20) 3+4-Methylphenols	8.85	107	149734	45.355	ng	99
22) Acetophenone	8.92	105	189356	43.930	ng	# 98
24) Nitrobenzene	9.30	77	151714	45.825	ng	95
25) Isophorone	9.82	82	288675	49.556	ng	99
26) 2-Nitrophenol	10.02	139	69988	42.345	ng	99
27) 2,4-Dimethylphenol	10.06	122	128388	51.557	ng	99
28) bis(2-Chloroethoxy)methane	10.30	93	173209	46.500	ng	96
29) 2,4-Dichlorophenol	10.54	162	130251	46.501	ng	100
30) 1,2,4-Trichlorobenzene	10.76	180	132511	42.217	ng	98
31) Naphthalene	10.95	128	410909	48.223	ng	99
32) Benzoic acid	10.18	122	43324m	32.410	ng	
33) 4-Chloroaniline	11.07	127	14704	3.710	ng	96
34) Hexachlorobutadiene	11.22	225	78556	40.665	ng	97
35) Caprolactam	11.85	113	33804	30.151	ng	# 87
36) 4-Chloro-3-methylphenol	12.17	107	143056	46.748	ng	93
37) 2-Methylnaphthalene	12.55	142	295244	48.237	ng	97
38) 1-Methylnaphthalene	12.78	142	283690	48.151	ng	97
40) 1,2,4,5-Tetrachlorobenzene	12.92	216	150776	41.039	ng	99

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41) Hexachlorocyclopentadiene	12.88	237	168669	85.735	nq	100
43) 2,4,6-Trichlorophenol	13.15	196	108031	43.151	nq	97
44) 2,4,5-Trichlorophenol	13.23	196	114348	43.409	nq	98
46) 1,1'-Biphenyl	13.55	154	380163	41.157	nq	99
47) 2-Chloronaphthalene	13.60	162	300394	42.618	nq	99
48) 2-Nitroaniline	13.81	65	104950	47.310	nq	100
49) Acenaphthylene	14.44	152	500064	47.178	nq	99
50) Dimethylphthalate	14.17	163	395293	45.351	nq	100
51) 2,6-Dinitrotoluene	14.30	165	90308	45.778	nq	97
52) Acenaphthene	14.79	154	291473	45.678	nq	100
53) 3-Nitroaniline	14.63	138	28648	13.160	nq	# 99
54) 2,4-Dinitrophenol	14.84	184	36445	39.542	nq	# 86
55) Dibenzofuran	15.12	168	468823	44.432	nq	98
56) 4-Nitrophenol	14.93	139	111006	66.118	nq	93
57) 2,4-Dinitrotoluene	15.09	165	125561	46.779	nq	96
58) Fluorene	15.77	166	385518	48.970	nq	97
59) 2,3,4,6-Tetrachlorophenol	15.34	232	100342	45.522	nq	97
60) Diethylphthalate	15.52	149	406913	43.952	nq	97
61) 4-Chlorophenyl-phenylether	15.76	204	193789	42.068	nq	99
62) 4-Nitroaniline	15.80	138	93812	43.381	nq	97
63) Azobenzene	16.05	77	379132	45.801	nq	98
65) 4,6-Dinitro-2-methylphenol	15.85	198	44007	27.876	nq	98
66) n-Nitrosodiphenylamine	15.97	169	340701	45.121	nq	98
67) 4-Bromophenyl-phenylether	16.65	248	119715	43.700	nq	93
68) Hexachlorobenzene	16.76	284	124441	44.008	nq	97
69) Atrazine	16.91	200	124983	44.902	nq	99
70) Pentachlorophenol	17.11	266	113129	58.908	nq	98
71) Phenanthrene	17.51	178	637803	49.912	nq	100
72) Anthracene	17.61	178	629692	49.990	nq	99
73) Carbazole	17.88	167	575308	45.523	nq	100
74) Di-n-butylphthalate	18.40	149	681608	48.046	nq	99
75) Fluoranthene	19.52	202	720600	49.426	nq	99
77) Benzidine	19.70	184	55231	6.831	nq	97
78) Pyrene	19.88	202	733910	48.714	nq	99
80) Butylbenzylphthalate	20.75	149	316311	48.010	nq	99
81) Benzo(a)anthracene	21.73	228	680964	48.902	nq	99
82) 3,3'-Dichlorobenzidine	21.65	252	90882	16.755	nq	99
83) Chrysene	21.80	228	634899	47.654	nq	99
84) Bis(2-ethylhexyl)phthalate	21.60	149	443961	47.251	nq	99
85) Di-n-octyl phthalate	22.84	149	757517	49.835	nq	99
86) Indeno(1,2,3-cd)pyrene	28.87	276	665849	44.998	nq	97
88) Benzo(b)fluoranthene	24.01	252	669456	49.592	nq	99
89) Benzo(k)fluoranthene	24.08	252	664172	49.861	nq	98
90) Benzo(a)pyrene	24.91	252	658914	51.075	nq	99
91) Dibenzo(a,h)anthracene	28.95	278	566238	45.617	nq	97
92) Benzo(g,h,i)perylene	30.09	276	559061	45.296	nq	98

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

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