

Data Path : Z:\HPCHEM1\BNA G\DATA\BG022618\
 Data File : BG032823.D
 Acq On : 26 Feb 2018 15:53
 Operator : SJ/JU
 Sample : J1688-01MSD
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 T-3-ROADMSD

Manual Integrations
 APPROVED

Sohil
 2/27/2018 10:23:12 AM

Quant Time: Feb 26 18:10:04 2018
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 26 17:50:45 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.93	152	74800	20.00	ng	0.00
21) Naphthalene-d8	11.81	136	328606	20.00	ng	0.00
38) Acenaphthene-d10	15.54	164	184117	20.00	ng	0.00
63) Phenanthrene-d10	18.27	188	425349	20.00	ng	0.00
75) Chrysene-d12	22.63	240	408603	20.00	ng	0.00
86) Perylene-d12	26.44	264	452669	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	6.34	112	492797	105.40	ng	0.00
7) Phenol-d6	8.02	99	657547	100.90	ng	0.00
23) Nitrobenzene-d5	10.13	82	368620	78.68	ng	0.00
41) 2,4,6-Tribromophenol	17.01	330	224277	115.85	ng	0.00
44) 2-Fluorobiphenyl	14.17	172	778307	60.97	ng	0.00
78) Terphenyl-d14	20.79	244	1144183	62.91	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.96	88	63454	31.272	ng	87
3) Pyridine	4.43	79	178785	31.968	ng	94
4) n-Nitrosodimethylamine	4.33	42	87433	38.994	ng	# 91
6) Aniline	8.22	93	204926	23.231	ng	96
8) 2-Chlorophenol	8.47	128	196967	38.489	ng	93
9) Benzaldehyde	8.02	77	88204	23.484	ng	90
10) Phenol	8.05	94	268637	40.892	ng	88
11) bis(2-Chloroethyl)ether	8.30	93	179896	33.489	ng	94
12) 1,3-Dichlorobenzene	8.81	146	198719	33.153	ng	95
13) 1,4-Dichlorobenzene	8.96	146	202089	33.495	ng	93
14) 1,2-Dichlorobenzene	9.30	146	189826	32.832	ng	97
15) Benzyl Alcohol	9.16	79	167692	35.002	ng	88
16) 2,2'-oxybis(1-Chloropropan	9.45	45	289053	31.301	ng	96
17) 2-Methylphenol	9.35	107	169948	35.082	ng	93
18) Hexachloroethane	10.06	117	71836	34.969	ng	# 83
19) n-Nitroso-di-n-propylamine	9.74	70	146917	31.933	ng	# 97
20) 3+4-Methylphenols	9.69	107	235158	34.912	ng	93
22) Acetophenone	9.77	105	280239	33.768	ng	# 95
24) Nitrobenzene	10.17	77	201233	39.676	ng	# 89
25) Isophorone	10.70	82	405390	35.148	ng	# 94
26) 2-Nitrophenol	10.90	139	95369	50.602	ng	90
27) 2,4-Dimethylphenol	10.92	122	219505	43.809	ng	87
28) bis(2-Chloroethoxy)methane	11.17	93	246458	36.680	ng	# 96
29) 2,4-Dichlorophenol	11.44	162	179130	39.391	ng	98
30) 1,2,4-Trichlorobenzene	11.66	180	174054	32.652	ng	98
31) Naphthalene	11.86	128	590721	34.402	ng	99
32) Benzoic acid	11.02	122	11746	10.246	ng	# 81
33) 4-Chloroaniline	11.95	127	184383	24.016	ng	96
34) Hexachlorobutadiene	12.13	225	92168	30.825	ng	99
35) Caprolactam	12.69	113	72824	39.994	ng	# 84
36) 4-Chloro-3-methylphenol	13.01	107	212723	40.198	ng	91
37) 2-Methylnaphthalene	13.42	142	425031	34.214	ng	92
39) 1,2,4,5-Tetrachlorobenzene	13.76	216	175832	32.171	ng	# 96
40) Hexachlorocyclopentadiene	13.74	237	202740	72.785	ng	93

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.99	196	135582	39.336	nq	98
43) 2,4,5-Trichlorophenol	14.06	196	154302	38.679	nq	95
45) 1,1'-Biphenyl	14.38	154	525109	32.638	nq	96
46) 2-Chloronaphthalene	14.44	162	401848	33.436	nq	98
47) 2-Nitroaniline	14.62	65	134607	45.146	nq #	87
48) Acenaphthylene	15.27	152	655593	33.318	nq	98
49) Dimethylphthalate	14.97	163	588964	37.674	nq	99
50) 2,6-Dinitrotoluene	15.09	165	114482	44.940	nq	88
51) Acenaphthene	15.61	154	422975	34.516	nq	98
52) 3-Nitroaniline	15.42	138	109316	36.326	nq #	83
53) 2,4-Dinitrophenol	15.62	184	54506	70.126	nq #	1
54) Dibenzofuran	15.93	168	607243	34.801	nq	95
55) 4-Nitrophenol	15.69	139	229502	88.794	nq #	78
56) 2,4-Dinitrotoluene	15.87	165	161513	52.296	nq	86
57) Fluorene	16.58	166	493692	34.280	nq	99
58) 2,3,4,6-Tetrachlorophenol	16.14	232	132641m	42.826	nq	
59) Diethylphthalate	16.31	149	561418	34.048	nq	98
60) 4-Chlorophenyl-phenylether	16.55	204	237471	33.707	nq	94
61) 4-Nitroaniline	16.58	138	137257	41.378	nq #	80
62) Azobenzene	16.85	77	506843	34.359	nq	97
64) 4,6-Dinitro-2-methylphenol	16.62	198	67397	52.446	nq	96
65) n-Nitrosodiphenylamine	16.76	169	475158	34.339	nq	99
66) 4-Bromophenyl-phenylether	17.45	248	146433	32.558	nq #	90
67) Hexachlorobenzene	17.57	284	155664	32.027	nq	94
68) Atrazine	17.68	200	169515	37.218	nq	95
69) Pentachlorophenol	17.90	266	217540	71.058	nq	99
70) Phenanthrene	18.31	178	806282	33.977	nq	98
71) Anthracene	18.40	178	820946	34.459	nq	99
72) Carbazole	18.66	167	806824	34.359	nq	97
73) Di-n-butylphthalate	19.16	149	983793	34.150	nq	99
74) Fluoranthene	20.26	202	914574	34.890	nq	94
76) Benzidine	20.42	184	443713	31.858	nq	98
77) Pyrene	20.62	202	911668	31.639	nq	97
79) Butylbenzylphthalate	21.49	149	466296	36.499	nq	93
80) Benzo(a)anthracene	22.61	228	894295	34.131	nq	99
81) 3,3'-Dichlorobenzidine	22.49	252	264190	27.224	nq #	97
82) Chrysene	22.69	228	835989	33.401	nq	98
83) Bis(2-ethylhexyl)phthalate	22.45	149	680690	35.995	nq #	97
84) Di-n-octyl phthalate	23.88	149	1173322	40.705	nq	98
85) Indeno(1,2,3-cd)pyrene	30.88	276	1046936	35.866	nq	98
87) Benzo(b)fluoranthene	25.22	252	891719	33.317	nq #	97
88) Benzo(k)fluoranthene	25.30	252	901037	33.874	nq #	96
89) Benzo(a)pyrene	26.26	252	882635	34.671	nq #	97
90) Dibenzo(a,h)anthracene	30.96	278	872933	34.067	nq #	94
91) Benzo(g,h,i)perylene	32.28	276	872261	34.532	nq #	91

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

