

Data Path : Z:\HPCHEM1\BNA G\DATA\BG021318\
 Data File : BG032668.D
 Acq On : 13 Feb 2018 19:50
 Operator : SJ/JU
 Sample : J1397-02MS
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 CL-01-021218-AMS

Manual Integrations
 APPROVED

Sohil
 2/14/2018 10:31:10 AM

Quant Time: Feb 14 03:28:33 2018
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG021218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 12 18:08:01 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.67	152	36109	20.00	ng	0.00
21) Naphthalene-d8	11.55	136	153856	20.00	ng	0.00
38) Acenaphthene-d10	15.31	164	109877	20.00	ng	0.00
63) Phenanthrene-d10	18.04	188	281495	20.00	ng	0.00
75) Chrysene-d12	22.37	240	346078	20.00	ng	0.00
86) Perylene-d12	26.01	264	366639	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	6.12	112	285583	158.68	ng	0.00
7) Phenol-d6	7.80	99	324800	131.98	ng	0.00
23) Nitrobenzene-d5	9.87	82	245604	101.45	ng	0.00
41) 2,4,6-Tribromophenol	16.79	330	250023	147.78	ng	0.00
44) 2-Fluorobiphenyl	13.93	172	822198	93.58	ng	0.00
78) Terphenyl-d14	20.59	244	1524303	100.39	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.80	88	29659	39.394	ng	# 70
3) Pyridine	4.25	79	67772	35.095	ng	# 87
4) n-Nitrosodimethylamine	4.15	42	49591	54.652	ng	# 73
6) Aniline	7.97	93	31956	10.987	ng	# 88
8) 2-Chlorophenol	8.22	128	114810	55.476	ng	# 85
9) Benzaldehyde	7.78	77	17349	11.929	ng	# 85
10) Phenol	7.82	94	107630	44.969	ng	# 86
11) bis(2-Chloroethyl)ether	8.06	93	89988	46.898	ng	# 79
12) 1,3-Dichlorobenzene	8.56	146	133472	45.728	ng	# 98
13) 1,4-Dichlorobenzene	8.71	146	130849	45.376	ng	# 98
14) 1,2-Dichlorobenzene	9.04	146	130894	45.004	ng	# 96
15) Benzyl Alcohol	8.91	79	93169	50.511	ng	# 92
16) 2,2'-oxybis(1-Chloropropan	9.20	45	92254	44.673	ng	# 72
17) 2-Methylphenol	9.11	107	80553	47.523	ng	# 92
18) Hexachloroethane	9.78	117	48204	45.781	ng	# 79
19) n-Nitroso-di-n-propylamine	9.48	70	76570	46.282	ng	# 94
20) 3+4-Methylphenols	9.45	107	110038	42.952	ng	# 95
22) Acetophenone	9.51	105	159293	48.681	ng	# 96
24) Nitrobenzene	9.91	77	116804	46.082	ng	# 95
25) Isophorone	10.44	82	223182	48.355	ng	# 85
26) 2-Nitrophenol	10.64	139	65303	48.171	ng	# 87
27) 2,4-Dimethylphenol	10.68	122	108433	60.881	ng	# 98
28) bis(2-Chloroethoxy)methane	10.92	93	120324	50.997	ng	# 96
29) 2,4-Dichlorophenol	11.18	162	131981	50.629	ng	# 93
30) 1,2,4-Trichlorobenzene	11.40	180	163015	44.606	ng	# 97
31) Naphthalene	11.60	128	361936	48.447	ng	# 99
32) Benzoic acid	10.77	122	7611m	9.229	ng	# 92
33) 4-Chloroaniline	11.72	127	14055	4.741	ng	# 92
34) Hexachlorobutadiene	11.86	225	129265	43.158	ng	# 96
35) Caprolactam	12.44	113	24655	33.305	ng	# 57
36) 4-Chloro-3-methylphenol	12.79	107	117400	50.068	ng	# 93
37) 2-Methylnaphthalene	13.17	142	287725	47.481	ng	# 93
39) 1,2,4,5-Tetrachlorobenzene	13.52	216	234502	45.187	ng	# 96
40) Hexachlorocyclopentadiene	13.50	237	290319	88.399	ng	# 92

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.75	196	130239	50.282	nq	98
43) 2,4,5-Trichlorophenol	13.83	196	145147	44.966	nq #	91
45) 1,1'-Biphenyl	14.15	154	411999	45.571	nq	95
46) 2-Chloronaphthalene	14.20	162	327549	45.785	nq	95
47) 2-Nitroaniline	14.39	65	75697	48.643	nq	89
48) Acenaphthylene	15.04	152	474051	45.040	nq	99
49) Dimethylphthalate	14.74	163	423122	44.550	nq #	98
50) 2,6-Dinitrotoluene	14.87	165	93206	47.650	nq #	76
51) Acenaphthene	15.37	154	309291	43.029	nq	99
52) 3-Nitroaniline	15.21	138	22635	13.104	nq #	79
53) 2,4-Dinitrophenol	15.41	184	13035	22.044	nq #	90
54) Dibenzofuran	15.70	168	508275	47.528	nq	98
55) 4-Nitrophenol	15.50	139	96849	75.900	nq #	77
56) 2,4-Dinitrotoluene	15.65	165	132519	49.169	nq #	75
57) Fluorene	16.35	166	407068	46.286	nq	95
58) 2,3,4,6-Tetrachlorophenol	15.92	232	145792	47.973	nq #	83
59) Diethylphthalate	16.09	149	414933	44.899	nq	98
60) 4-Chlorophenyl-phenylether	16.32	204	269175	46.356	nq	94
61) 4-Nitroaniline	16.37	138	65647	36.862	nq	86
62) Azobenzene	16.62	77	278824	46.960	nq	86
64) 4,6-Dinitro-2-methylphenol	16.41	198	22676	14.939	nq	77
65) n-Nitrosodiphenylamine	16.54	169	368507	44.636	nq	98
66) 4-Bromophenyl-phenylether	17.22	248	188705	46.841	nq	96
67) Hexachlorobenzene	17.34	284	192439	44.475	nq	93
68) Atrazine	17.47	200	166975	47.155	nq	93
69) Pentachlorophenol	17.69	266	114851	43.357	nq	98
70) Phenanthrene	18.09	178	681808	45.089	nq	99
71) Anthracene	18.18	178	718353	46.429	nq	98
72) Carbazole	18.44	167	603565	45.753	nq	99
73) Di-n-butylphthalate	18.95	149	676864	45.570	nq	99
74) Fluoranthene	20.05	202	920493	45.598	nq	95
76) Benzidine	20.22	184	63639	7.809	nq	97
77) Pyrene	20.41	202	929195	45.131	nq	98
79) Butylbenzylphthalate	21.27	149	309521	45.636	nq #	77
80) Benzo(a)anthracene	22.35	228	1003158	46.296	nq	99
81) 3,3'-Dichlorobenzidine	22.24	252	85111	10.518	nq #	97
82) Chrysene	22.42	228	940940	44.039	nq	98
83) Bis(2-ethylhexyl)phthalate	22.19	149	436296	44.608	nq #	99
84) Di-n-octyl phthalate	23.54	149	762576	49.695	nq #	90
85) Indeno(1,2,3-cd)pyrene	30.24	276	1208889	47.762	nq #	90
87) Benzo(b)fluoranthene	24.85	252	1026439	46.977	nq #	95
88) Benzo(k)fluoranthene	24.92	252	1051162	46.485	nq #	96
89) Benzo(a)pyrene	25.85	252	1027166	48.255	nq #	95
90) Dibenzo(a,h)anthracene	30.31	278	1014810	47.888	nq #	93
91) Benzo(g,h,i)perylene	31.56	276	989029	48.280	nq #	93

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

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