

Data Path : Z:\HPCHEM1\BNA G\DATA\BG042016\
 Data File : BG021765.D
 Acq On : 19 Apr 2016 21:37
 Operator : UM/SJ
 Sample : H2620-03MS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EUT01-TU-B2(9-13)-CMS

Manual Integrations
 APPROVED

Sohil
 4/20/2016 1:38:39 PM

Quant Time: Apr 20 03:46:31 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG041316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 13 15:48:13 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.21	152	27068	20.00	ng	-0.01
21) Naphthalene-d8	11.03	136	129090	20.00	ng	0.00
38) Acenaphthene-d10	14.82	164	84427	20.00	ng	0.00
63) Phenanthrene-d10	17.56	188	186724	20.00	ng	0.00
75) Chrysene-d12	21.84	240	204917	20.00	ng	0.00
86) Perylene-d12	25.17	264	202232	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.76	112	196220	120.72	ng	0.00
7) Phenol-d6	7.37	99	300828	123.90	ng	0.01
23) Nitrobenzene-d5	9.38	82	182420	72.63	ng	0.00
41) 2,4,6-Tribromophenol	16.30	330	118570	130.31	ng	0.00
44) 2-Fluorobiphenyl	13.45	172	425537	73.85	ng	0.00
78) Terphenyl-d14	20.14	244	579062	70.51	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.62	88	17549	24.25	ng	# 83
3) Pyridine	4.03	79	57517	26.49	ng	92
4) n-Nitrosodimethylamine	3.94	42	32600	30.30	ng	# 76
6) Aniline	7.53	93	86853	27.05	ng	91
8) 2-Chlorophenol	7.78	128	75184	38.58	ng	94
9) Benzaldehyde	7.35	77	19035	13.56	ng	93
10) Phenol	7.39	94	104986	40.20	ng	96
11) bis(2-Chloroethyl)ether	7.62	93	72013	34.45	ng	91
12) 1,3-Dichlorobenzene	8.10	146	73397	33.73	ng	98
13) 1,4-Dichlorobenzene	8.25	146	74860	33.79	ng	94
14) 1,2-Dichlorobenzene	8.57	146	74002	34.82	ng	98
15) Benzyl Alcohol	8.45	79	75454	40.66	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.74	45	134310	30.00	ng	99
17) 2-Methylphenol	8.65	107	71477	39.59	ng	94
18) Hexachloroethane	9.30	117	27580	34.20	ng	# 73
19) n-Nitroso-di-n-propylamine	9.02	70	64277	34.07	ng	# 93
20) 3+4-Methylphenols	8.98	107	99484	40.27	ng	97
22) Acetophenone	9.04	105	119589	33.26	ng	# 96
24) Nitrobenzene	9.42	77	92325	34.33	ng	97
25) Isophorone	9.94	82	179977	32.74	ng	99
26) 2-Nitrophenol	10.14	139	43806	42.67	ng	# 91
27) 2,4-Dimethylphenol	10.19	122	79224	33.97	ng	98
28) bis(2-Chloroethoxy)methane	10.42	93	104868	35.90	ng	91
29) 2,4-Dichlorophenol	10.67	162	82933	41.71	ng	99
30) 1,2,4-Trichlorobenzene	10.89	180	78474	36.70	ng	94
31) Naphthalene	11.08	128	246872	35.73	ng	# 97
32) Benzoic acid	10.25	122	2920	6.78	ng	96
33) 4-Chloroaniline	11.18	127	70647	23.62	ng	98
34) Hexachlorobutadiene	11.35	225	48674	35.30	ng	99
35) Caprolactam	11.95	113	31035m	34.11	ng	
36) 4-Chloro-3-methylphenol	12.29	107	97489	38.99	ng	96
37) 2-Methylnaphthalene	12.67	142	186528	39.33	ng	99
39) 1,2,4,5-Tetrachlorobenzene	13.03	216	97118	36.81	ng	98
40) Hexachlorocyclopentadiene	13.00	237	104924	71.59	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.26	196	67358	40.03	nq	99
43) 2,4,5-Trichlorophenol	13.34	196	72395	39.88	nq	96
45) 1,1'-Biphenyl	13.66	154	245499	34.51	nq	97
46) 2-Chloronaphthalene	13.71	162	189253	35.99	nq	99
47) 2-Nitroaniline	13.90	65	69850	38.74	nq	95
48) Acenaphthylene	14.55	152	312086	35.89	nq	98
49) Dimethylphthalate	14.27	163	295428	40.47	nq	99
50) 2,6-Dinitrotoluene	14.39	165	54495	39.15	nq	98
51) Acenaphthene	14.88	154	202105	36.36	nq	99
52) 3-Nitroaniline	14.73	138	51538	33.39	nq	95
53) 2,4-Dinitrophenol	14.93	184	16770	35.86	nq	88
54) Dibenzofuran	15.22	168	306037	40.32	nq	100
55) 4-Nitrophenol	15.03	139	95060	71.92	nq	97
56) 2,4-Dinitrotoluene	15.18	165	76107	40.46	nq	93
57) Fluorene	15.87	166	249045	36.74	nq	100
58) 2,3,4,6-Tetrachlorophenol	15.44	232	65169	43.26	nq	98
59) Diethylphthalate	15.62	149	261159	34.28	nq	97
60) 4-Chlorophenyl-phenylether	15.85	204	122438	36.67	nq	98
61) 4-Nitroaniline	15.88	138	59805	35.21	nq	96
62) Azobenzene	16.14	77	244871	37.50	nq	97
64) 4,6-Dinitro-2-methylphenol	15.94	198	24127	32.94	nq	93
65) n-Nitrosodiphenylamine	16.07	169	221363	39.53	nq	95
66) 4-Bromophenyl-phenylether	16.74	248	79983	39.98	nq	94
67) Hexachlorobenzene	16.86	284	82847	39.23	nq	96
68) Atrazine	17.00	200	84113	38.03	nq	99
69) Pentachlorophenol	17.21	266	86518	71.76	nq	95
70) Phenanthrene	17.60	178	380220	36.95	nq	99
71) Anthracene	17.69	178	385070	36.86	nq	96
72) Carbazole	17.96	167	354204	36.33	nq	99
73) Di-n-butylphthalate	18.50	149	426657	34.40	nq	100
74) Fluoranthene	19.59	202	427570	34.79	nq	99
76) Benzidine	19.77	184	199152	31.21	nq	100
77) Pyrene	19.96	202	441066	37.33	nq	99
79) Butylbenzylphthalate	20.82	149	196754	35.87	nq	99
80) Benzo(a)anthracene	21.81	228	438233	37.16	nq	99
81) 3,3'-Dichlorobenzidine	21.72	252	97676	10.46	nq	96
82) Chrysene	21.88	228	399626	35.27	nq	99
83) Bis(2-ethylhexyl)phthalate	21.69	149	283923	35.39	nq	# 98
84) Di-n-octyl phthalate	22.94	149	492944	36.68	nq	97
85) Indeno(1,2,3-cd)pyrene	28.99	276	499334	37.07	nq	# 95
87) Benzo(b)fluoranthene	24.10	252	433209	36.92	nq	99
88) Benzo(k)fluoranthene	24.17	252	408366	35.59	nq	99
89) Benzo(a)pyrene	25.01	252	419756	36.92	nq	98
90) Dibenzo(a,h)anthracene	29.05	278	417455	36.62	nq	99
91) Benzo(g,h,i)perylene	30.19	276	412410	36.87	nq	98

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

