

Data Path : Z:\HPCHEM1\BNA G\DATA\BG072716\
 Data File : BG023303.D
 Acq On : 28 Jul 2016 4:20
 Operator : UM/SJ
 Sample : PB92411BS
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB92411BS

Manual Integrations

sohil
 7/28/2016 4:58:36 PM

Quant Time: Jul 28 05:35:22 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG072616.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 27 17:30:33 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.13	152	180078	20.00	ng	0.00
21) Naphthalene-d8	10.97	136	797792	20.00	ng	0.00
38) Acenaphthene-d10	14.81	164	554584	20.00	ng	0.00
63) Phenanthrene-d10	17.57	188	1232544	20.00	ng	0.00
75) Chrysene-d12	21.90	240	1194033	20.00	ng	0.00
86) Perylene-d12	25.30	264	1233131	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.67	112	1403528	133.06	ng	0.00
7) Phenol-d6	7.29	99	2071606	127.92	ng	0.00
23) Nitrobenzene-d5	9.31	82	1333201	81.72	ng	0.00
41) 2,4,6-Tribromophenol	16.31	330	888222	154.74	ng	0.00
44) 2-Fluorobiphenyl	13.42	172	2628971	88.72	ng	0.00
78) Terphenyl-d14	20.18	244	3184756	86.70	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.53	88	141633	31.31	ng	# 92
3) Pyridine	3.94	79	451275	28.50	ng	# 92
4) n-Nitrosodimethylamine	3.85	42	218253	30.51	ng	# 75
6) Aniline	7.46	93	804458	37.19	ng	95
8) 2-Chlorophenol	7.70	128	536323	45.57	ng	95
9) Benzaldehyde	7.26	77	106231	9.13	ng	96
10) Phenol	7.32	94	774696	45.33	ng	81
11) bis(2-Chloroethyl)ether	7.55	93	548948	37.80	ng	90
12) 1,3-Dichlorobenzene	8.02	146	525068	39.59	ng	95
13) 1,4-Dichlorobenzene	8.17	146	534958	39.54	ng	95
14) 1,2-Dichlorobenzene	8.49	146	523446	40.10	ng	95
15) Benzyl Alcohol	8.37	79	551267	50.58	ng	93
16) 2,2'-oxybis(1-Chloropropan	8.67	45	784439	33.94	ng	90
17) 2-Methylphenol	8.59	107	529013	45.29	ng	# 85
18) Hexachloroethane	9.23	117	204121	36.21	ng	94
19) n-Nitroso-di-n-propylamine	8.94	70	532259	36.54	ng	98
20) 3+4-Methylphenols	8.92	107	731009	46.03	ng	91
22) Acetophenone	8.97	105	860665	41.32	ng	# 91
24) Nitrobenzene	9.36	77	732115	38.65	ng	# 89
25) Isophorone	9.88	82	1375709	40.83	ng	# 92
26) 2-Nitrophenol	10.07	139	314014	43.88	ng	# 75
27) 2,4-Dimethylphenol	10.13	122	584124	48.43	ng	84
28) bis(2-Chloroethoxy)methane	10.37	93	817357	43.23	ng	99
29) 2,4-Dichlorophenol	10.62	162	591634	50.67	ng	100
30) 1,2,4-Trichlorobenzene	10.83	180	551478	42.74	ng	98
31) Naphthalene	11.02	128	1612487	43.50	ng	99
32) Benzoic acid	10.27	122	366401	39.12	ng	91
33) 4-Chloroaniline	11.14	127	360014	21.02	ng	95
34) Hexachlorobutadiene	11.31	225	347938	41.14	ng	97
35) Caprolactam	11.92	113	239927m	48.32	ng	
36) 4-Chloro-3-methylphenol	12.26	107	700136	45.64	ng	86
37) 2-Methylnaphthalene	12.63	142	1258588	45.79	ng	# 92
39) 1,2,4,5-Tetrachlorobenzene	13.00	216	639310	35.61	ng	97
40) Hexachlorocyclopentadiene	12.97	237	819173	96.28	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.24	196	498265	47.25	nq	99
43) 2,4,5-Trichlorophenol	13.32	196	542417	48.48	nq	94
45) 1,1'-Biphenyl	13.64	154	1668109	43.13	nq	92
46) 2-Chloronaphthalene	13.69	162	1319971	42.59	nq	95
47) 2-Nitroaniline	13.89	65	562509	42.90	nq	91
48) Acenaphthylene	14.53	152	2071612	43.66	nq	94
49) Dimethylphthalate	14.26	163	1726862	42.19	nq	96
50) 2,6-Dinitrotoluene	14.38	165	418125	43.63	nq	# 79
51) Acenaphthene	14.87	154	1341376	43.58	nq	99
52) 3-Nitroaniline	14.72	138	377866	39.15	nq	# 83
53) 2,4-Dinitrophenol	14.92	184	525327	93.62	nq	# 85
54) Dibenzofuran	15.21	168	2049226	47.05	nq	97
55) 4-Nitrophenol	15.04	139	730435	87.24	nq	# 74
56) 2,4-Dinitrotoluene	15.17	165	599037	44.39	nq	95
57) Fluorene	15.86	166	1692471	44.01	nq	98
58) 2,3,4,6-Tetrachlorophenol	15.44	232	492493	52.00	nq	92
59) Diethylphthalate	15.62	149	1777033	42.39	nq	95
60) 4-Chlorophenyl-phenylether	15.85	204	879339	42.24	nq	98
61) 4-Nitroaniline	15.89	138	458880	45.34	nq	# 65
62) Azobenzene	16.14	77	1968890	45.30	nq	91
64) 4,6-Dinitro-2-methylphenol	15.94	198	374557	47.31	nq	94
65) n-Nitrosodiphenylamine	16.06	169	1560071	47.41	nq	# 92
66) 4-Bromophenyl-phenylether	16.75	248	600418	48.18	nq	95
67) Hexachlorobenzene	16.87	284	614262	49.02	nq	96
68) Atrazine	17.02	200	602457	47.73	nq	96
69) Pentachlorophenol	17.22	266	748692	96.81	nq	97
70) Phenanthrene	17.62	178	2551537m	47.33	nq	
71) Anthracene	17.71	178	2625271	48.19	nq	93
72) Carbazole	17.98	167	2421637	48.54	nq	97
73) Di-n-butylphthalate	18.52	149	2650024	47.61	nq	97
74) Fluoranthene	19.62	202	2798021	45.48	nq	88
76) Benzidine	19.81	184	1783183	53.10	nq	99
77) Pyrene	19.99	202	2806478	41.42	nq	93
79) Butylbenzylphthalate	20.86	149	1372394	46.04	nq	96
80) Benzo(a)anthracene	21.87	228	2798002	44.76	nq	97
81) 3,3'-Dichlorobenzidine	21.79	252	836242	32.99	nq	# 97
82) Chrysene	21.94	228	2615805	43.31	nq	93
83) Bis(2-ethylhexyl)phthalate	21.75	149	1862463	44.65	nq	# 96
84) Di-n-octyl phthalate	23.03	149	3049955	44.70	nq	99
85) Indeno(1,2,3-cd)pyrene	29.21	276	3615316	51.32	nq	# 100
87) Benzo(b)fluoranthene	24.22	252	3012990	45.99	nq	# 95
88) Benzo(k)fluoranthene	24.28	252	2884357	44.04	nq	# 92
89) Benzo(a)pyrene	25.13	252	2952366	45.99	nq	# 95
90) Dibenzo(a,h)anthracene	29.27	278	2972819	48.96	nq	# 90
91) Benzo(g,h,i)perylene	30.42	276	2992285	46.81	nq	# 87

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

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