

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG020417\
 Data File : BG025801.D
 Acq On : 3 Feb 2017 19:16
 Operator : SJ/MA
 Sample : PB96126BSD
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB96126BSD

Manual Integrations
 APPROVED

mohammad
 2/6/2017 3:19:33 PM

Quant Time: Feb 03 22:19:26 2017
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG013017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 30 18:29:38 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.16	152	78917	20.00	ng	0.00
21) Naphthalene-d8	10.99	136	351143	20.00	ng	0.00
38) Acenaphthene-d10	14.80	164	285744	20.00	ng	0.00
63) Phenanthrene-d10	17.54	188	651196	20.00	ng	0.00
75) Chrysene-d12	21.83	240	790829	20.00	ng	0.00
86) Perylene-d12	25.21	264	780894	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.71	112	659459	149.76	ng	0.00
7) Phenol-d6	7.33	99	966707	150.32	ng	0.00
23) Nitrobenzene-d5	9.35	82	686027	82.55	ng	0.00
41) 2,4,6-Tribromophenol	16.29	330	496734	121.90	ng	0.00
44) 2-Fluorobiphenyl	13.41	172	1399357	79.54	ng	0.00
78) Terphenyl-d14	20.12	244	2419523	73.85	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.53	88	68431	34.28	ng	99
3) Pyridine	3.94	79	199378	36.81	ng	95
4) n-Nitrosodimethylamine	3.87	42	139765	44.85	ng	93
6) Aniline	7.49	93	218434	24.11	ng	93
8) 2-Chlorophenol	7.73	128	239083	47.83	ng	99
9) Benzaldehyde	7.31	77	125738	23.05	ng	92
10) Phenol	7.36	94	355996	49.25	ng	92
11) bis(2-Chloroethyl)ether	7.57	93	246122	43.06	ng	90
12) 1,3-Dichlorobenzene	8.05	146	250071	40.91	ng	95
13) 1,4-Dichlorobenzene	8.20	146	250617	40.33	ng	96
14) 1,2-Dichlorobenzene	8.52	146	243454	40.89	ng	97
15) Benzyl Alcohol	8.41	79	266792	47.97	ng	89
16) 2,2'-oxybis(1-Chloropropan	8.69	45	484591	41.99	ng	95
17) 2-Methylphenol	8.61	107	231079	46.80	ng	94
18) Hexachloroethane	9.24	117	100428	39.68	ng	93
19) n-Nitroso-di-n-propylamine	8.97	70	239328	42.17	ng	98
20) 3+4-Methylphenols	8.95	107	314473	46.67	ng	97
22) Acetophenone	9.00	105	409775	42.62	ng	# 99
24) Nitrobenzene	9.40	77	363480	40.06	ng	99
25) Isophorone	9.90	82	642346	39.67	ng	96
26) 2-Nitrophenol	10.11	139	144855	45.05	ng	91
27) 2,4-Dimethylphenol	10.15	122	264265	47.19	ng	92
28) bis(2-Chloroethoxy)methane	10.38	93	364481	42.52	ng	97
29) 2,4-Dichlorophenol	10.66	162	265468	46.19	ng	94
30) 1,2,4-Trichlorobenzene	10.85	180	275532	41.22	ng	98
31) Naphthalene	11.04	128	739247	40.55	ng	99
32) Benzoic acid	10.33	122	150248	42.54	ng	93
33) 4-Chloroaniline	11.18	127	101625	13.41	ng	92
34) Hexachlorobutadiene	11.29	225	194940	35.85	ng	98
35) Caprolactam	11.95	113	100715m	41.13	ng	
36) 4-Chloro-3-methylphenol	12.28	107	315417	45.05	ng	97
37) 2-Methylnaphthalene	12.63	142	584420	41.52	ng	97
39) 1,2,4,5-Tetrachlorobenzene	13.00	216	371364	38.81	ng	99
40) Hexachlorocyclopentadiene	12.96	237	224386	63.27	ng	93

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.24	196	235530	42.04	ng	97
43) 2,4,5-Trichlorophenol	13.33	196	245608	40.61	ng	99
45) 1,1'-Biphenyl	13.63	154	829762	41.74	ng	98
46) 2-Chloronaphthalene	13.68	162	628202	40.24	ng	96
47) 2-Nitroaniline	13.90	65	277034	41.24	ng	98
48) Acenaphthylene	14.52	152	992362	38.21	ng	97
49) Dimethylphthalate	14.24	163	835044	38.81	ng	98
50) 2,6-Dinitrotoluene	14.38	165	192909	39.59	ng	96
51) Acenaphthene	14.86	154	649520	40.48	ng	97
52) 3-Nitroaniline	14.73	138	105264	22.91	ng #	98
53) 2,4-Dinitrophenol	14.95	184	214842	74.44	ng #	89
54) Dibenzofuran	15.19	168	1035007	42.69	ng	96
55) 4-Nitrophenol	15.05	139	263802	74.66	ng	87
56) 2,4-Dinitrotoluene	15.18	165	284330	39.88	ng #	77
57) Fluorene	15.84	166	842289	39.10	ng	99
58) 2,3,4,6-Tetrachlorophenol	15.42	232	244119	43.20	ng	98
59) Diethylphthalate	15.59	149	881150	38.63	ng	99
60) 4-Chlorophenyl-phenylether	15.82	204	450054	38.54	ng	98
61) 4-Nitroaniline	15.89	138	185038	40.10	ng	95
62) Azobenzene	16.12	77	1032349	44.47	ng	96
64) 4,6-Dinitro-2-methylphenol	15.95	198	181747	39.27	ng	80
65) n-Nitrosodiphenylamine	16.04	169	781394	41.80	ng	99
66) 4-Bromophenyl-phenylether	16.72	248	314237	39.68	ng	98
67) Hexachlorobenzene	16.85	284	349531	39.25	ng	94
68) Atrazine	16.98	200	327053	35.44	ng	97
69) Pentachlorophenol	17.20	266	303684	77.60	ng	99
70) Phenanthrene	17.59	178	1425557	40.29	ng	98
71) Anthracene	17.68	178	1493564	40.93	ng	99
72) Carbazole	17.96	167	1322204	37.41	ng	96
73) Di-n-butylphthalate	18.46	149	1615786	39.51	ng	98
74) Fluoranthene	19.58	202	1785790	36.82	ng	97
76) Benzidine	19.77	184	412976	15.21	ng	95
77) Pyrene	19.95	202	1818914	41.29	ng	98
79) Butylbenzylphthalate	20.79	149	798545	43.66	ng	95
80) Benzo(a)anthracene	21.81	228	1915443	40.90	ng	99
81) 3,3'-Dichlorobenzidine	21.72	252	398554	21.68	ng	99
82) Chrysene	21.88	228	1728155	39.04	ng	99
83) Bis(2-ethylhexyl)phthalate	21.65	149	1122571	44.07	ng	96
84) Di-n-octyl phthalate	22.89	149	1879547	45.42	ng	98
85) Indeno(1,2,3-cd)pyrene	29.12	276	2260636	41.89	ng #	98
87) Benzo(b)fluoranthene	24.13	252	1931822	42.80	ng	98
88) Benzo(k)fluoranthene	24.20	252	1850876	41.90	ng	96
89) Benzo(a)pyrene	25.05	252	1858073	43.28	ng #	97
90) Dibenzo(a,h)anthracene	29.15	278	1895585	42.70	ng #	96
91) Benzo(g,h,i)perylene	30.34	276	1867268	42.33	ng #	93

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

