

Data Path : Z:\HPCHEM1\BNA G\DATA\BG042115\
 Data File : BG016596.D
 Acq On : 21 Apr 2015 22:42
 Operator : TP/IZ
 Sample : PB82885BS
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB82885BS

Manual Integrations
 APPROVED

apatel
 4/22/2015 5:21:09 PM

Quant Time: Apr 22 04:54:56 2015
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG042115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 22 02:10:07 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	55803	20.00	ng	0.00
21) Naphthalene-d8	10.42	136	234419	20.00	ng	0.00
38) Acenaphthene-d10	14.28	164	126778	20.00	ng	0.00
63) Phenanthrene-d10	17.02	188	279783	20.00	ng	0.00
75) Chrysene-d12	21.21	240	305838	20.00	ng	0.00
86) Perylene-d12	23.42	264	291519	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.23	112	458354	131.64	ng	0.00
7) Phenol-d6	6.83	99	619434	127.34	ng	0.00
23) Nitrobenzene-d5	8.79	82	314748	82.67	ng	0.00
41) 2,4,6-Tribromophenol	15.78	330	118130	137.19	ng	0.00
44) 2-Fluorobiphenyl	12.90	172	670561	87.60	ng	0.00
78) Terphenyl-d14	19.66	244	998411	78.65	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.10	88	55579	35.50	ng	94
3) Pyridine	3.49	79	148676	34.33	ng	98
4) n-Nitrosodimethylamine	3.42	42	49560	38.22	ng	99
6) Aniline	6.97	93	165628	24.65	ng	98
8) 2-Chlorophenol	7.21	128	173734	41.80	ng	97
9) Benzaldehyde	6.78	77	14464	6.53	ng	94
10) Phenol	6.85	94	217784	42.33	ng	98
11) bis(2-Chloroethyl)ether	7.07	93	155976	38.72	ng	98
12) 1,3-Dichlorobenzene	7.52	146	169764	39.12	ng	98
13) 1,4-Dichlorobenzene	7.67	146	170358	38.56	ng	99
14) 1,2-Dichlorobenzene	7.98	146	162321	38.62	ng	99
15) Benzyl Alcohol	7.88	79	120857	39.04	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.18	45	154292	39.55	ng	97
17) 2-Methylphenol	8.09	107	141573	40.83	ng	99
18) Hexachloroethane	8.70	117	62469	39.38	ng	97
19) n-Nitroso-di-n-propylamine	8.44	70	107949	34.32	ng	97
20) 3+4-Methylphenols	8.42	107	189703	39.72	ng	99
22) Acetophenone	8.46	105	214425	41.07	ng	98
24) Nitrobenzene	8.83	77	159273	40.20	ng	98
25) Isophorone	9.35	82	290084	37.22	ng	99
26) 2-Nitrophenol	9.54	139	89350	40.84	ng	98
27) 2,4-Dimethylphenol	9.62	122	161748	43.48	ng	97
28) bis(2-Chloroethoxy)methane	9.84	93	187211	40.06	ng	100
29) 2,4-Dichlorophenol	10.08	162	137982	44.86	ng	99
30) 1,2,4-Trichlorobenzene	10.28	180	134070	40.99	ng	97
31) Naphthalene	10.47	128	486532	39.76	ng	99
32) Benzoic acid	9.77	122	29520	28.60	ng	98
33) 4-Chloroaniline	10.59	127	104556	18.68	ng	97
34) Hexachlorobutadiene	10.76	225	58255	40.08	ng	99
35) Caprolactam	11.35	113	67739m	38.03	ng	
36) 4-Chloro-3-methylphenol	11.73	107	155418	41.84	ng	99
37) 2-Methylnaphthalene	12.09	142	349589	39.90	ng	98
39) 1,2,4,5-Tetrachlorobenzene	12.46	216	121623	41.70	ng	99
40) Hexachlorocyclopentadiene	12.44	237	98218	78.41	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	12.71	196	89938	43.55	nq	97
43) 2,4,5-Trichlorophenol	12.79	196	98165	44.58	nq	97
45) 1,1'-Biphenyl	13.11	154	419259	42.94	nq	99
46) 2-Chloronaphthalene	13.15	162	316434	42.32	nq	99
47) 2-Nitroaniline	13.37	65	92283	42.31	nq	98
48) Acenaphthylene	14.00	152	547545	41.25	nq	100
49) Dimethylphthalate	13.75	163	386190	41.19	nq	99
50) 2,6-Dinitrotoluene	13.87	165	96766	42.03	nq	99
51) Acenaphthene	14.34	154	326486	41.22	nq	99
52) 3-Nitroaniline	14.20	138	75381	26.51	nq	100
53) 2,4-Dinitrophenol	14.41	184	90273	84.06	nq	97
54) Dibenzofuran	14.68	168	475726	42.66	nq	100
55) 4-Nitrophenol	14.53	139	211762	98.40	nq	98
56) 2,4-Dinitrotoluene	14.66	165	140364	44.31	nq	98
57) Fluorene	15.33	166	411256	41.85	nq	99
58) 2,3,4,6-Tetrachlorophenol	14.91	232	73063	43.97	nq	99
59) Diethylphthalate	15.11	149	408424	40.89	nq	99
60) 4-Chlorophenyl-phenylether	15.33	204	165199	40.95	nq	98
61) 4-Nitroaniline	15.36	138	136090	41.77	nq	96
62) Azobenzene	15.62	77	403703	43.55	nq	98
64) 4,6-Dinitro-2-methylphenol	15.43	198	75836	40.12	nq	93
65) n-Nitrosodiphenylamine	15.54	169	371771	39.63	nq	98
66) 4-Bromophenyl-phenylether	16.22	248	88501	40.18	nq	98
67) Hexachlorobenzene	16.34	284	92203	40.08	nq	90
68) Atrazine	16.50	200	120159	45.82	nq	96
69) Pentachlorophenol	16.68	266	97009	81.32	nq	99
70) Phenanthrene	17.07	178	663782	41.68	nq	99
71) Anthracene	17.16	178	672955	42.52	nq	99
72) Carbazole	17.43	167	752366	46.59	nq	99
73) Di-n-butylphthalate	17.99	149	896390	45.61	nq	100
74) Fluoranthene	19.09	202	797587	46.49	nq	99
76) Benzidine	19.27	184	404051	35.73	nq	99
77) Pyrene	19.45	202	845735	41.17	nq	99
79) Butylbenzylphthalate	20.35	149	465867	43.32	nq	100
80) Benzo(a)anthracene	21.19	228	761136	41.27	nq	99
81) 3,3'-Dichlorobenzidine	21.13	252	200108	32.77	nq	96
82) Chrysene	21.24	228	719975	40.66	nq	99
83) Bis(2-ethylhexyl)phthalate	21.12	149	633378	42.91	nq	99
84) Di-n-octyl phthalate	21.99	149	1056828	43.11	nq	100
85) Indeno(1,2,3-cd)pyrene	25.68	276	797811	41.06	nq	99
87) Benzo(b)fluoranthene	22.76	252	750267	43.13	nq	99
88) Benzo(k)fluoranthene	22.80	252	689110	40.39	nq	98
89) Benzo(a)pyrene	23.33	252	707759	42.85	nq	99
90) Dibenzo(a,h)anthracene	25.69	278	664481	41.40	nq	97
91) Benzo(g,h,i)perylene	26.38	276	676964	41.81	nq	99

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

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