

Data Path : Z:\HPCHEM1\BNA G\DATA\BG120515\
 Data File : BG019944.D
 Acq On : 5 Dec 2015 16:21
 Operator : UM/NP
 Sample : PB87048BS
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB87048BS

Manual Integrations
 APPROVED

Mmdadoda
 12/7/2015 6:31:37 PM

Quant Time: Dec 07 02:21:46 2015
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG120215.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 02 18:50:01 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.11	152	32350	20.00	ng	-0.02
21) Naphthalene-d8	10.93	136	148461	20.00	ng	-0.02
38) Acenaphthene-d10	14.75	164	104189	20.00	ng	-0.02
63) Phenanthrene-d10	17.49	188	265343	20.00	ng	-0.02
75) Chrysene-d12	21.76	240	316722	20.00	ng	-0.03
86) Perylene-d12	25.04	264	294340	20.00	ng	-0.05

System Monitoring Compounds						
5) 2-Fluorophenol	5.66	112	233957	130.71	ng	-0.01
7) Phenol-d6	7.26	99	335902	124.29	ng	-0.01
23) Nitrobenzene-d5	9.28	82	218917	75.26	ng	-0.02
41) 2,4,6-Tribromophenol	16.23	330	170198	106.76	ng	-0.02
44) 2-Fluorobiphenyl	13.37	172	532004	69.72	ng	-0.02
78) Terphenyl-d14	20.09	244	925384	71.27	ng	-0.02

Target Compounds						Qvalue
2) 1,4-Dioxane	3.52	88	25357	36.36	ng	# 100
3) Pyridine	3.93	79	83272	39.32	ng	91
4) n-Nitrosodimethylamine	3.84	42	41433	37.20	ng	91
6) Aniline	7.43	93	103967	29.71	ng	# 87
8) 2-Chlorophenol	7.67	128	99438	45.56	ng	93
9) Benzaldehyde	7.24	77	34930	19.38	ng	91
10) Phenol	7.29	94	136286	47.92	ng	83
11) bis(2-Chloroethyl)ether	7.53	93	88945	42.22	ng	88
12) 1,3-Dichlorobenzene	8.00	146	97696	39.49	ng	# 92
13) 1,4-Dichlorobenzene	8.15	146	102905m	40.28	ng	
14) 1,2-Dichlorobenzene	8.47	146	98289	40.33	ng	# 94
15) Benzyl Alcohol	8.35	79	105299	44.40	ng	90
16) 2,2'-oxybis(1-Chloropropan	8.65	45	140504	40.90	ng	92
17) 2-Methylphenol	8.55	107	94461	47.04	ng	# 94
18) Hexachloroethane	9.21	117	38264	40.00	ng	92
19) n-Nitroso-di-n-propylamine	8.93	70	84394	39.60	ng	92
20) 3+4-Methylphenols	8.88	107	129574	45.83	ng	95
22) Acetophenone	8.94	105	149855	36.82	ng	# 91
24) Nitrobenzene	9.32	77	128250	40.64	ng	91
25) Isophorone	9.85	82	227029	36.40	ng	# 96
26) 2-Nitrophenol	10.04	139	55076	45.19	ng	# 70
27) 2,4-Dimethylphenol	10.09	122	105242	41.47	ng	90
28) bis(2-Chloroethoxy)methane	10.33	93	128234	42.08	ng	100
29) 2,4-Dichlorophenol	10.58	162	113393	43.74	ng	96
30) 1,2,4-Trichlorobenzene	10.79	180	109614	37.41	ng	95
31) Naphthalene	10.99	128	313761	39.45	ng	99
32) Benzoic acid	10.22	122	76244	45.01	ng	83
33) 4-Chloroaniline	11.09	127	59724	17.04	ng	# 92
34) Hexachlorobutadiene	11.27	225	77665	35.93	ng	95
35) Caprolactam	11.87	113	41155m	42.64	ng	
36) 4-Chloro-3-methylphenol	12.20	107	133840	41.45	ng	95
37) 2-Methylnaphthalene	12.59	142	240674	41.29	ng	95
39) 1,2,4,5-Tetrachlorobenzene	12.94	216	146360	37.02	ng	97
40) Hexachlorocyclopentadiene	12.93	237	169193	75.05	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.18	196	105770	40.41	nq	95
43) 2,4,5-Trichlorophenol	13.25	196	118664	42.89	nq	96
45) 1,1'-Biphenyl	13.58	154	317806	37.17	nq	99
46) 2-Chloronaphthalene	13.63	162	257397	39.05	nq	94
47) 2-Nitroaniline	13.82	65	96485	46.35	nq	88
48) Acenaphthylene	14.47	152	431291	38.42	nq	99
49) Dimethylphthalate	14.20	163	356324	38.02	nq	98
50) 2,6-Dinitrotoluene	14.31	165	78175	44.37	nq #	72
51) Acenaphthene	14.81	154	301136m	44.21	nq	
52) 3-Nitroaniline	14.64	138	58349	31.87	nq	92
53) 2,4-Dinitrophenol	14.85	184	82744	87.29	nq	94
54) Dibenzofuran	15.14	168	433734	42.80	nq	93
55) 4-Nitrophenol	14.94	139	133159	83.52	nq #	63
56) 2,4-Dinitrotoluene	15.09	165	112583	45.79	nq #	87
57) Fluorene	15.79	166	349047	38.48	nq	99
58) 2,3,4,6-Tetrachlorophenol	15.36	232	115615	45.15	nq	97
59) Diethylphthalate	15.56	149	371627	36.64	nq	99
60) 4-Chlorophenyl-phenylether	15.78	204	190888	36.65	nq	96
61) 4-Nitroaniline	15.81	138	86018	41.78	nq #	77
62) Azobenzene	16.08	77	353540	41.89	nq	95
64) 4,6-Dinitro-2-methylphenol	15.86	198	65779	44.62	nq	96
65) n-Nitrosodiphenylamine	15.99	169	318076	41.42	nq	97
66) 4-Bromophenyl-phenylether	16.67	248	128096	38.85	nq	97
67) Hexachlorobenzene	16.79	284	137790	40.12	nq	95
68) Atrazine	16.94	200	133012	39.87	nq	95
69) Pentachlorophenol	17.13	266	177677	88.63	nq	96
70) Phenanthrene	17.53	178	605550	40.34	nq	97
71) Anthracene	17.62	178	612159	40.03	nq	96
72) Carbazole	17.89	167	545726	40.51	nq	97
73) Di-n-butylphthalate	18.44	149	652793	39.54	nq	98
74) Fluoranthene	19.53	202	745273	38.81	nq	95
76) Benzidine	19.71	184	244088	23.46	nq	97
77) Pyrene	19.89	202	774379	41.56	nq	97
79) Butylbenzylphthalate	20.78	149	297923	40.79	nq	97
80) Benzo(a)anthracene	21.74	228	775066	39.98	nq	99
81) 3,3'-Dichlorobenzidine	21.65	252	183162	24.81	nq	98
82) Chrysene	21.81	228	709356	38.54	nq	98
83) Bis(2-ethylhexyl)phthalate	21.65	149	434378	40.53	nq	97
84) Di-n-octyl phthalate	22.89	149	737132	42.30	nq #	94
85) Indeno(1,2,3-cd)pyrene	28.79	276	869861	42.90	nq #	100
87) Benzo(b)fluoranthene	24.00	252	757736	40.62	nq #	95
88) Benzo(k)fluoranthene	24.07	252	737737	39.93	nq #	95
89) Benzo(a)pyrene	24.89	252	737224	41.63	nq #	95
90) Dibenzo(a,h)anthracene	28.87	278	715356	40.88	nq #	94
91) Benzo(g,h,i)perylene	29.97	276	718262	41.80	nq #	94

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

