

Data Path : Z:\HPCHEM1\BNA G\DATA\BG022716\
 Data File : BG021129.D
 Acq On : 28 Feb 2016 00:16
 Operator : UM/SJ
 Sample : PB88567BS
 Misc :
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB88567BS

Manual Integrations
 APPROVED

Sohil
 2/29/2016 12:39:04 PM

Quant Time: Feb 29 03:00:48 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022616.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 26 16:39:44 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.17	152	6838	20.00	ng	-0.01
21) Naphthalene-d8	10.98	136	32748	20.00	ng	-0.02
38) Acenaphthene-d10	14.77	164	21355	20.00	ng	-0.01
63) Phenanthrene-d10	17.51	188	53501	20.00	ng	-0.01
75) Chrysene-d12	21.77	240	60333	20.00	ng	-0.02
86) Perylene-d12	25.05	264	54237	20.00	ng	-0.03

System Monitoring Compounds

5) 2-Fluorophenol	5.72	112	57517	147.26	ng	0.00
7) Phenol-d6	7.32	99	89366	153.76	ng	-0.01
23) Nitrobenzene-d5	9.33	82	61926	90.10	ng	-0.02
41) 2,4,6-Tribromophenol	16.25	330	31716	113.28	ng	-0.01
44) 2-Fluorobiphenyl	13.40	172	125672	75.31	ng	-0.01
78) Terphenyl-d14	20.10	244	225367	87.30	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.61	88	6304	34.85	ng	# 86
3) Pyridine	4.01	79	19341	36.56	ng	95
4) n-Nitrosodimethylamine	3.93	42	10479	39.07	ng	96
6) Aniline	7.49	93	17225	23.96	ng	# 88
8) 2-Chlorophenol	7.73	128	22044	45.25	ng	89
9) Benzaldehyde	7.30	77	3663	11.64	ng	84
10) Phenol	7.34	94	31660	53.17	ng	79
11) bis(2-Chloroethyl)ether	7.58	93	20230	44.33	ng	89
12) 1,3-Dichlorobenzene	8.06	146	20309	39.17	ng	# 90
13) 1,4-Dichlorobenzene	8.20	146	20701	37.38	ng	99
14) 1,2-Dichlorobenzene	8.52	146	20844	39.75	ng	95
15) Benzyl Alcohol	8.40	79	25625	50.47	ng	90
16) 2,2'-oxybis(1-Chloropropan	8.70	45	30577	58.43	ng	90
17) 2-Methylphenol	8.60	107	21096	50.20	ng	# 86
18) Hexachloroethane	9.25	117	8581	40.96	ng	95
19) n-Nitroso-di-n-propylamine	8.97	70	22019	49.11	ng	96
20) 3+4-Methylphenols	8.93	107	30077	51.62	ng	91
22) Acetophenone	8.98	105	34667	38.35	ng	# 92
24) Nitrobenzene	9.37	77	31909	44.94	ng	# 86
25) Isophorone	9.89	82	58553	45.45	ng	# 94
26) 2-Nitrophenol	10.08	139	11620	38.71	ng	# 60
27) 2,4-Dimethylphenol	10.13	122	24505	46.28	ng	89
28) bis(2-Chloroethoxy)methane	10.37	93	30056	47.87	ng	95
29) 2,4-Dichlorophenol	10.61	162	22611	43.62	ng	96
30) 1,2,4-Trichlorobenzene	10.83	180	20991	35.73	ng	95
31) Naphthalene	11.03	128	67964	40.40	ng	99
32) Benzoic acid	10.25	122	18272	37.58	ng	# 75
33) 4-Chloroaniline	11.13	127	6758	9.62	ng	# 87
34) Hexachlorobutadiene	11.30	225	12891	32.79	ng	95
35) Caprolactam	11.90	113	9208m	52.57	ng	
36) 4-Chloro-3-methylphenol	12.23	107	30290	47.67	ng	86
37) 2-Methylnaphthalene	12.62	142	50054	42.37	ng	# 94
39) 1,2,4,5-Tetrachlorobenzene	12.98	216	25388	32.17	ng	# 97
40) Hexachlorocyclopentadiene	12.96	237	35925	69.23	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.21	196	21098	42.63	ng	98
43) 2,4,5-Trichlorophenol	13.28	196	22684	44.61	ng	92
45) 1,1'-Biphenyl	13.61	154	67978	37.80	ng	97
46) 2-Chloronaphthalene	13.66	162	52392	38.40	ng	95
47) 2-Nitroaniline	13.85	65	24667	55.79	ng #	74
48) Acenaphthylene	14.50	152	88590	41.04	ng	98
49) Dimethylphthalate	14.22	163	76612	41.63	ng #	99
50) 2,6-Dinitrotoluene	14.34	165	16827	44.44	ng #	67
51) Acenaphthene	14.84	154	54594	37.07	ng	99
52) 3-Nitroaniline	14.67	138	7008	19.02	ng #	61
53) 2,4-Dinitrophenol	14.87	184	22334	84.60	ng #	78
54) Dibenzofuran	15.17	168	85442	45.80	ng	93
55) 4-Nitrophenol	14.97	139	32966	100.88	ng #	60
56) 2,4-Dinitrotoluene	15.12	165	24661	45.42	ng #	77
57) Fluorene	15.82	166	76933	42.39	ng	95
58) 2,3,4,6-Tetrachlorophenol	15.39	232	19868	44.96	ng #	95
59) Diethylphthalate	15.58	149	85263	43.95	ng	99
60) 4-Chlorophenyl-phenylether	15.81	204	38014	38.08	ng	97
61) 4-Nitroaniline	15.83	138	19896	47.42	ng #	52
62) Azobenzene	16.09	77	95465	56.80	ng	93
64) 4,6-Dinitro-2-methylphenol	15.88	198	15533	39.58	ng	92
65) n-Nitrosodiphenylamine	16.02	169	69552	44.06	ng	93
66) 4-Bromophenyl-phenylether	16.70	248	23738	37.69	ng #	90
67) Hexachlorobenzene	16.81	284	23300	35.06	ng #	89
68) Atrazine	16.96	200	27615	47.22	ng	98
69) Pentachlorophenol	17.15	266	34192	76.22	ng	90
70) Phenanthrene	17.55	178	125326	43.54	ng	100
71) Anthracene	17.64	178	126665	43.69	ng	99
72) Carbazole	17.91	167	118734	47.44	ng	97
73) Di-n-butylphthalate	18.46	149	156792	47.52	ng #	97
74) Fluoranthene	19.54	202	153940	42.63	ng	98
76) Benzidine	19.72	184	31975	20.14	ng	100
77) Pyrene	19.91	202	160448	48.03	ng	99
79) Butylbenzylphthalate	20.78	149	72892	52.28	ng	90
80) Benzo(a)anthracene	21.75	228	153745	44.29	ng	98
81) 3,3'-Dichlorobenzidine	21.66	252	18392	13.54	ng	96
82) Chrysene	21.82	228	138714	41.49	ng	99
83) Bis(2-ethylhexyl)phthalate	21.65	149	103138	47.90	ng	99
84) Di-n-octyl phthalate	22.89	149	180481	50.02	ng	98
85) Indeno(1,2,3-cd)pyrene	28.81	276	159439	39.55	ng #	100
87) Benzo(b)fluoranthene	24.01	252	150869	44.50	ng	99
88) Benzo(k)fluoranthene	24.08	252	146865	43.95	ng	98
89) Benzo(a)pyrene	24.90	252	144564	44.42	ng	96
90) Dibenzo(a,h)anthracene	28.87	278	138227	42.13	ng	98
91) Benzo(g,h,i)perylene	29.98	276	131258	41.54	ng	96

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

