

Data Path : Z:\HPCHEM1\BNA G\DATA\BG120715\
 Data File : BG019985.D
 Acq On : 7 Dec 2015 17:39
 Operator : UM/NP
 Sample : PB87126BS
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB87126BS

Manual Integrations
 APPROVED

Mmdadoda
 12/8/2015 6:30:10 PM

Quant Time: Dec 07 18:22:53 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	17063	20.00	ng	-0.02
21) Naphthalene-d8	10.93	136	70991	20.00	ng	-0.02
38) Acenaphthene-d10	14.74	164	49064	20.00	ng	-0.02
63) Phenanthrene-d10	17.48	188	127187	20.00	ng	-0.02
75) Chrysene-d12	21.76	240	159012	20.00	ng	-0.03
86) Perylene-d12	25.03	264	148120	20.00	ng	-0.06

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.66	112	85615	90.69	ng	-0.01
7) Phenol-d6	7.26	99	126140	88.49	ng	-0.01
23) Nitrobenzene-d5	9.28	82	122556	88.11	ng	-0.02
41) 2,4,6-Tribromophenol	16.23	330	63952	85.19	ng	-0.02
44) 2-Fluorobiphenyl	13.37	172	294290	81.90	ng	-0.02
78) Terphenyl-d14	20.08	244	548520	84.14	ng	-0.03

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.52	88	12457	33.87	ng	# 100
3) Pyridine	3.92	79	40028	35.84	ng	# 87
4) n-Nitrosodimethylamine	3.84	42	21802	37.11	ng	# 81
6) Aniline	7.43	93	43939	23.80	ng	# 86
8) 2-Chlorophenol	7.67	128	48466	42.10	ng	# 96
9) Benzaldehyde	7.24	77	13997	14.73	ng	# 91
10) Phenol	7.29	94	66234	44.15	ng	# 79
11) bis(2-Chloroethyl)ether	7.53	93	44025	39.62	ng	# 87
12) 1,3-Dichlorobenzene	8.00	146	49570	37.99	ng	# 94
13) 1,4-Dichlorobenzene	8.15	146	51693	38.36	ng	# 94
14) 1,2-Dichlorobenzene	8.47	146	49762	38.71	ng	# 94
15) Benzyl Alcohol	8.35	79	50696	40.53	ng	# 90
16) 2,2'-oxybis(1-Chloropropan	8.65	45	67497	37.25	ng	# 96
17) 2-Methylphenol	8.55	107	45250	42.73	ng	# 93
18) Hexachloroethane	9.21	117	19587	38.82	ng	# 99
19) n-Nitroso-di-n-propylamine	8.92	70	42554	37.85	ng	# 88
20) 3+4-Methylphenols	8.88	107	61811	41.45	ng	# 94
22) Acetophenone	8.94	105	73763	37.91	ng	# 94
24) Nitrobenzene	9.32	77	62810	41.63	ng	# 94
25) Isophorone	9.85	82	112997	37.89	ng	# 97
26) 2-Nitrophenol	10.03	139	27008	46.35	ng	# 76
27) 2,4-Dimethylphenol	10.09	122	51421	42.37	ng	# 90
28) bis(2-Chloroethoxy)methane	10.33	93	62493	42.88	ng	# 98
29) 2,4-Dichlorophenol	10.57	162	56156	45.30	ng	# 97
30) 1,2,4-Trichlorobenzene	10.79	180	54801	39.11	ng	# 97
31) Naphthalene	10.98	128	155390	40.86	ng	# 99
32) Benzoic acid	10.21	122	36919	45.50	ng	# 85
33) 4-Chloroaniline	11.08	127	29338	17.51	ng	# 91
34) Hexachlorobutadiene	11.27	225	39301	38.03	ng	# 98
35) Caprolactam	11.86	113	19616m	42.51	ng	# 96
36) 4-Chloro-3-methylphenol	12.20	107	65814	42.63	ng	# 97
37) 2-Methylnaphthalene	12.58	142	118391	42.48	ng	# 99
39) 1,2,4,5-Tetrachlorobenzene	12.94	216	74331	39.92	ng	# 99
40) Hexachlorocyclopentadiene	12.93	237	78323	73.77	ng	# 99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.18	196	51175	41.51	nq	95
43) 2,4,5-Trichlorophenol	13.25	196	58216	44.68	nq	95
45) 1,1'-Biphenyl	13.58	154	157394	39.09	nq	99
46) 2-Chloronaphthalene	13.62	162	126971	40.91	nq	95
47) 2-Nitroaniline	13.82	65	46543	47.48	nq	87
48) Acenaphthylene	14.47	152	214685	40.61	nq	99
49) Dimethylphthalate	14.19	163	176225	39.93	nq	98
50) 2,6-Dinitrotoluene	14.31	165	38124	45.94	nq #	75
51) Acenaphthene	14.81	154	129856	40.48	nq	99
52) 3-Nitroaniline	14.64	138	27387	31.77	nq	95
53) 2,4-Dinitrophenol	14.84	184	37053	83.46	nq #	85
54) Dibenzofuran	15.14	168	213333	44.70	nq	93
55) 4-Nitrophenol	14.94	139	64607	86.05	nq #	66
56) 2,4-Dinitrotoluene	15.09	165	55092	47.59	nq #	85
57) Fluorene	15.79	166	177458	41.55	nq	98
58) 2,3,4,6-Tetrachlorophenol	15.36	232	56700	47.02	nq	98
59) Diethylphthalate	15.55	149	186600	39.07	nq	97
60) 4-Chlorophenyl-phenylether	15.78	204	95695	39.02	nq	97
61) 4-Nitroaniline	15.80	138	42252	43.58	nq #	72
62) Azobenzene	16.07	77	177260	44.60	nq	94
64) 4,6-Dinitro-2-methylphenol	15.86	198	31300	44.34	nq	91
65) n-Nitrosodiphenylamine	15.99	169	162068	44.03	nq	94
66) 4-Bromophenyl-phenylether	16.67	248	63589	40.24	nq	97
67) Hexachlorobenzene	16.79	284	69852	42.43	nq	98
68) Atrazine	16.93	200	67017	41.91	nq	96
69) Pentachlorophenol	17.13	266	89006	92.63	nq	95
70) Phenanthrene	17.53	178	306834	42.64	nq	97
71) Anthracene	17.62	178	312288	42.61	nq	95
72) Carbazole	17.88	167	275856	42.72	nq	96
73) Di-n-butylphthalate	18.44	149	329477	41.63	nq	97
74) Fluoranthene	19.53	202	391204	42.50	nq	94
76) Benzidine	19.71	184	151270	28.96	nq	95
77) Pyrene	19.89	202	400607	42.82	nq	95
79) Butylbenzylphthalate	20.77	149	155680	42.46	nq	99
80) Benzo(a)anthracene	21.74	228	406795	41.79	nq	98
81) 3,3'-Dichlorobenzidine	21.65	252	104709	28.25	nq	98
82) Chrysene	21.80	228	369573	39.99	nq	96
83) Bis(2-ethylhexyl)phthalate	21.64	149	225108	41.83	nq	98
84) Di-n-octyl phthalate	22.87	149	385632	44.08	nq #	94
85) Indeno(1,2,3-cd)pyrene	28.80	276	449260	44.13	nq #	100
87) Benzo(b)fluoranthene	23.99	252	396195	42.20	nq #	95
88) Benzo(k)fluoranthene	24.07	252	381347	41.02	nq #	95
89) Benzo(a)pyrene	24.89	252	379048	42.53	nq #	96
90) Dibenzo(a,h)anthracene	28.86	278	370315	42.06	nq #	94
91) Benzo(g,h,i)perylene	29.96	276	365187	42.24	nq #	92

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

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