

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG061215\
 Data File : BG017645.D
 Acq On : 12 Jun 2015 20:07
 Operator : TP/IZ
 Sample : SSTDICC080
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICC080

Manual Integrations
 APPROVED

apatel
 6/15/2015 4:40:04 PM

Quant Time: Jun 13 01:48:39 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG061215.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jun 13 00:03:37 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.52	152	21061	20.00	ng	0.00
21) Naphthalene-d8	10.32	136	82319	20.00	ng	0.00
38) Acenaphthene-d10	14.21	164	61856	20.00	ng	0.00
63) Phenanthrene-d10	16.96	188	172827	20.00	ng	0.00
75) Chrysene-d12	21.17	240	258387	20.00	ng	0.00
86) Perylene-d12	23.36	264	246495	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.11	112	183131	155.06	ng	0.00
7) Phenol-d6	6.74	99	265593	163.88	ng	0.00
23) Nitrobenzene-d5	8.70	82	335333	198.63	ng	0.00
41) 2,4,6-Tribromophenol	15.71	330	185442	207.88	ng	0.00
44) 2-Fluorobiphenyl	12.82	172	768691	185.07	ng	0.00
78) Terphenyl-d14	19.61	244	1917437	153.07	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.95	88	43982	81.29	ng	99
3) Pyridine	3.35	79	117927	82.49	ng	98
4) n-Nitrosodimethylamine	3.27	42	68952	75.27	ng	99
6) Aniline	6.86	93	173813	79.47	ng	94
8) 2-Chlorophenol	7.10	128	106608	76.28	ng	89
9) Benzaldehyde	6.67	77	69915	64.08	ng	96
10) Phenol	6.76	94	136586	82.65	ng	93
11) bis(2-Chloroethyl)ether	6.96	93	98296	78.07	ng	98
12) 1,3-Dichlorobenzene	7.41	146	125403	79.69	ng	96
13) 1,4-Dichlorobenzene	7.56	146	133366	80.70	ng	93
14) 1,2-Dichlorobenzene	7.88	146	120204	77.41	ng	# 83
15) Benzyl Alcohol	7.79	79	151059	97.17	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.08	45	40796	21.35	ng	88
17) 2-Methylphenol	8.00	107	99228	78.99	ng	95
18) Hexachloroethane	8.60	117	55871	84.49	ng	100
19) n-Nitroso-di-n-propylamine	8.35	70	115959	83.33	ng	# 90
20) 3+4-Methylphenols	8.34	107	141637	80.98	ng	92
22) Acetophenone	8.36	105	186154	91.45	ng	# 95
24) Nitrobenzene	8.74	77	177285	100.26	ng	# 96
25) Isophorone	9.26	82	309199	88.65	ng	99
26) 2-Nitrophenol	9.45	139	68538	86.97	ng	88
27) 2,4-Dimethylphenol	9.53	122	113065	87.26	ng	93
28) bis(2-Chloroethoxy)methane	9.75	93	137288	87.29	ng	98
29) 2,4-Dichlorophenol	10.00	162	119841	93.12	ng	95
30) 1,2,4-Trichlorobenzene	10.19	180	153105	99.93	ng	94
31) Naphthalene	10.37	128	367017	85.07	ng	# 94
32) Benzoic acid	9.75	122	68426	84.62	ng	# 82
33) 4-Chloroaniline	10.50	127	157184	82.74	ng	94
34) Hexachlorobutadiene	10.65	225	132868	128.39	ng	90
35) Caprolactam	11.31	113	53696	78.02	ng	87
36) 4-Chloro-3-methylphenol	11.67	107	149082	90.01	ng	92
37) 2-Methylnaphthalene	12.00	142	294448	88.40	ng	98
39) 1,2,4,5-Tetrachlorobenzene	12.38	216	211345	106.54	ng	99
40) Hexachlorocyclopentadiene	12.35	237	86693	97.86	ng	96

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42) 2,4,6-Trichlorophenol	12.63	196	133100m	97.76	ng	
43) 2,4,5-Trichlorophenol	12.72	196	140018	92.90	ng	91
45) 1,1'-Biphenyl	13.03	154	390545	86.39	ng	98
46) 2-Chloronaphthalene	13.08	162	322523	85.88	ng	100
47) 2-Nitroaniline	13.30	65	113116	81.33	ng	96
48) Acenaphthylene	13.93	152	549445	82.88	ng	100
49) Dimethylphthalate	13.69	163	441630	82.02	ng	# 97
50) 2,6-Dinitrotoluene	13.80	165	101254	84.14	ng	97
51) Acenaphthene	14.27	154	326023	83.09	ng	96
52) 3-Nitroaniline	14.14	138	89592	70.20	ng	98
53) 2,4-Dinitrophenol	14.36	184	57846	93.15	ng	95
54) Dibenzofuran	14.61	168	533232	90.93	ng	99
55) 4-Nitrophenol	14.49	139	63971	71.83	ng	98
56) 2,4-Dinitrotoluene	14.60	165	152344	88.46	ng	95
57) Fluorene	15.27	166	467327	88.63	ng	96
58) 2,3,4,6-Tetrachlorophenol	14.86	232	148400	105.00	ng	98
59) Diethylphthalate	15.05	149	482316	82.54	ng	98
60) 4-Chlorophenyl-phenylether	15.26	204	297465	107.73	ng	96
61) 4-Nitroaniline	15.31	138	99688	71.85	ng	91
62) Azobenzene	15.56	77	529080	93.20	ng	96
64) 4,6-Dinitro-2-methylphenol	15.37	198	118507	92.17	ng	95
65) n-Nitrosodiphenylamine	15.48	169	422675	82.51	ng	97
66) 4-Bromophenyl-phenylether	16.16	248	195215	98.49	ng	96
67) Hexachlorobenzene	16.27	284	213406	99.99	ng	96
68) Atrazine	16.45	200	196855	85.64	ng	99
69) Pentachlorophenol	16.63	266	92376	84.94	ng	99
70) Phenanthrene	17.01	178	785929	81.51	ng	99
71) Anthracene	17.10	178	782647	81.05	ng	98
72) Carbazole	17.38	167	735992	77.89	ng	97
73) Di-n-butylphthalate	17.94	149	885478	73.78	ng	99
74) Fluoranthene	19.04	202	1139624	88.90	ng	99
76) Benzidine	19.23	184	471825	55.94	ng	96
77) Pyrene	19.40	202	1165106	74.13	ng	99
79) Butylbenzylphthalate	20.31	149	438543	66.16	ng	98
80) Benzo(a)anthracene	21.15	228	1251548	78.83	ng	97
81) 3,3'-Dichlorobenzidine	21.09	252	486474	82.48	ng	98
82) Chrysene	21.21	228	1141212	79.35	ng	96
83) Bis(2-ethylhexyl)phthalate	21.08	149	654726	68.37	ng	94
84) Di-n-octyl phthalate	21.95	149	1053011	65.60	ng	99
85) Indeno(1,2,3-cd)pyrene	25.61	276	1533235	84.79	ng	99
87) Benzo(b)fluoranthene	22.71	252	1295068	83.12	ng	99
88) Benzo(k)fluoranthene	22.75	252	1255987	83.15	ng	98
89) Benzo(a)pyrene	23.27	252	1207207	84.02	ng	98
90) Dibenzo(a,h)anthracene	25.62	278	1279011	85.45	ng	99
91) Benzo(g,h,i)perylene	26.29	276	1203195	84.29	ng	# 97

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

