

Data Path : Z:\HPCHEM1\BNA G\DATA\BG050515\
 Data File : BG016860.D
 Acq On : 5 May 2015 12:07
 Operator : TP/IZ
 Sample : SSTDICC2.5
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICC2.5

Manual Integrations
 APPROVED

apatel
 5/6/2015 5:17:37 PM

Quant Time: May 05 15:56:58 2015
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG050515.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 05 15:36:00 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.81	152	73008	20.00	ng	0.00
21) Naphthalene-d8	10.60	136	302578	20.00	ng	0.00
38) Acenaphthene-d10	14.45	164	196685	20.00	ng	0.00
63) Phenanthrene-d10	17.19	188	435700	20.00	ng	0.00
75) Chrysene-d12	21.37	240	470526	20.00	ng	0.00
86) Perylene-d12	23.68	264	458607	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.39	112	20779	5.20	ng	0.00
7) Phenol-d6	6.96	99	29090	5.20	ng	0.00
23) Nitrobenzene-d5	8.96	82	33025	6.86	ng	0.00
41) 2,4,6-Tribromophenol	15.93	330	10124	6.63	ng	0.00
44) 2-Fluorobiphenyl	13.07	172	73049	5.72	ng	0.00
78) Terphenyl-d14	19.82	244	114786	5.87	ng	0.00

Target Compounds						Qvalue
6) Aniline	7.13	93	19642	2.58	ng	91
8) 2-Chlorophenol	7.37	128	12241	2.46	ng	93
9) Benzaldehyde	6.95	77	11633	3.52	ng	91
10) Phenol	6.99	94	14999	2.59	ng	95
11) bis(2-Chloroethyl)ether	7.23	93	12978	2.81	ng	96
12) 1,3-Dichlorobenzene	7.69	146	14486	2.71	ng	91
13) 1,4-Dichlorobenzene	7.84	146	13485m	2.46	ng	
14) 1,2-Dichlorobenzene	8.16	146	13788	2.62	ng	95
15) Benzyl Alcohol	8.04	79	12580	3.38	ng	96
16) 2,2'-oxybis(1-Chloropropan	8.34	45	23907	5.09	ng	96
17) 2-Methylphenol	8.24	107	11128	2.68	ng	93
18) Hexachloroethane	8.88	117	6147	3.11	ng	# 85
19) n-Nitroso-di-n-propylamine	8.61	70	13160	3.50	ng	# 91
20) 3+4-Methylphenols	8.56	107	14175	2.49	ng	# 72
22) Acetophenone	8.62	105	18564	3.00	ng	# 86
24) Nitrobenzene	9.00	77	15822	3.43	ng	97
25) Isophorone	9.53	82	33005	3.43	ng	98
26) 2-Nitrophenol	9.71	139	6521	2.39	ng	# 80
27) 2,4-Dimethylphenol	9.77	122	11207	2.44	ng	85
28) bis(2-Chloroethoxy)methane	10.02	93	16832	3.08	ng	97
29) 2,4-Dichlorophenol	10.24	162	10699	2.66	ng	88
30) 1,2,4-Trichlorobenzene	10.46	180	11806	2.76	ng	97
31) Naphthalene	10.65	128	42026	2.80	ng	97
33) 4-Chloroaniline	10.76	127	16631	2.53	ng	94
34) Hexachlorobutadiene	10.94	225	8204	4.02	ng	93
36) 4-Chloro-3-methylphenol	11.88	107	13454	2.91	ng	# 92
37) 2-Methylnaphthalene	12.26	142	31131	2.81	ng	94
39) 1,2,4,5-Tetrachlorobenzene	12.63	216	13797	3.01	ng	# 100
40) Hexachlorocyclopentadiene	12.62	237	8448	7.03	ng	93
42) 2,4,6-Trichlorophenol	12.87	196	9381	2.72	ng	98
43) 2,4,5-Trichlorophenol	12.94	196	10302	2.95	ng	88
45) 1,1'-Biphenyl	13.28	154	38052	2.64	ng	98
46) 2-Chloronaphthalene	13.32	162	29957	2.69	ng	97
47) 2-Nitroaniline	13.52	65	9059	3.08	ng	94

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
48) Acenaphthylene	14.16	152	53001	2.68	nq	97
49) Dimethylphthalate	13.90	163	39336	2.78	nq	100
50) 2,6-Dinitrotoluene	14.02	165	7122	2.09	nq	# 84
51) Acenaphthene	14.51	154	29787	2.53	nq	97
54) Dibenzofuran	14.84	168	42235	2.51	nq	98
57) Fluorene	15.49	166	36692	2.49	nq	98
58) 2,3,4,6-Tetrachlorophenol	15.07	232	8550	3.14	nq	# 79
59) Diethylphthalate	15.27	149	41542	2.82	nq	99
60) 4-Chlorophenyl-phenylether	15.49	204	20321	3.19	nq	96
61) 4-Nitroaniline	15.50	138	9640	2.47	nq	91
62) Azobenzene	15.78	77	42630	3.38	nq	97
65) n-Nitrosodiphenylamine	15.70	169	35336	2.54	nq	94
66) 4-Bromophenyl-phenylether	16.38	248	11212	3.12	nq	89
67) Hexachlorobenzene	16.50	284	12077	3.21	nq	# 83
68) Atrazine	16.65	200	12268	2.93	nq	90
70) Phenanthrene	17.23	178	61866	2.65	nq	98
71) Anthracene	17.32	178	61263	2.63	nq	96
72) Carbazole	17.59	167	58098	2.55	nq	97
73) Di-n-butylphthalate	18.16	149	79641	2.80	nq	# 97
74) Fluoranthene	19.25	202	71207	2.79	nq	94
76) Benzidine	19.43	184	45352	2.57	nq	# 95
77) Pyrene	19.60	202	75861	2.51	nq	95
79) Butylbenzylphthalate	20.51	149	34934	2.33	nq	92
80) Benzo(a)anthracene	21.35	228	68332	2.54	nq	98
81) 3,3'-Dichlorobenzidine	21.28	252	25239	2.72	nq	# 87
82) Chrysene	21.40	228	62943	2.47	nq	95
83) Bis(2-ethylhexyl)phthalate	21.30	149	47945	2.28	nq	98
84) Di-n-octyl phthalate	22.19	149	84153	2.36	nq	# 100
85) Indeno(1,2,3-cd)pyrene	26.04	276	81617	2.81	nq	# 100
87) Benzo(b)fluoranthene	22.97	252	68648	2.64	nq	97
88) Benzo(k)fluoranthene	23.02	252	67758	2.68	nq	97
89) Benzo(a)pyrene	23.57	252	64521	2.63	nq	97
90) Dibenzo(a,h)anthracene	26.06	278	69695	2.81	nq	# 92
91) Benzo(g,h,i)perylene	26.75	276	67394	2.72	nq	95

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

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