

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF053015\
 Data File : BF079599.D
 Acq On : 30 May 2015 13:14
 Operator : TP/IZ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

apatel
 6/1/2015 8:46:00 PM

Quant Time: May 31 06:32:59 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF052615.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 29 17:35:14 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.94	152	73369	20.00	ng	-0.02
21) Naphthalene-d8	8.51	136	300167	20.00	ng	-0.02
38) Acenaphthene-d10	10.67	164	141148	20.00	ng	-0.03
63) Phenanthrene-d10	12.51	188	258513	20.00	ng	-0.02
75) Chrysene-d12	15.78	240	261353	20.00	ng	-0.06
86) Perylene-d12	17.52	264	283938	20.00	ng	-0.15

System Monitoring Compounds						
5) 2-Fluorophenol	5.20	112	356115	84.19	ng	-0.03
7) Phenol-d6	6.54	99	450471	83.55	ng	-0.02
23) Nitrobenzene-d5	7.63	82	418325	80.56	ng	-0.03
41) 2,4,6-Tribromophenol	11.66	330	129476	83.52	ng	-0.03
44) 2-Fluorobiphenyl	9.86	172	820665	82.95	ng	-0.02
78) Terphenyl-d14	14.50	244	908405	81.57	ng	-0.03

Target Compounds				Qvalue		
2) 1,4-Dioxane	1.74	88	111315	55.32	ng	# 100
3) Pyridine	2.36	79	184992	41.75	ng	96
4) n-Nitrosodimethylamine	2.30	42	94873	43.49	ng	84
6) Aniline	6.52	93	311961	40.77	ng	95
8) 2-Chlorophenol	6.67	128	200291	40.68	ng	93
9) Benzaldehyde	6.38	77	123889	41.55	ng	93
10) Phenol	6.55	94	266396	41.96	ng	99
11) bis(2-Chloroethyl)ether	6.64	93	196153	42.72	ng	100
12) 1,3-Dichlorobenzene	6.86	146	222542	40.79	ng	95
13) 1,4-Dichlorobenzene	6.96	146	226720	40.74	ng	98
14) 1,2-Dichlorobenzene	7.14	146	206060	39.87	ng	96
15) Benzyl Alcohol	7.14	79	163244	41.20	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.31	45	267914	39.86	ng	99
17) 2-Methylphenol	7.29	107	155935	40.31	ng	98
18) Hexachloroethane	7.55	117	77231	40.32	ng	# 85
19) n-Nitroso-di-n-propylamine	7.47	70	152996	40.33	ng	# 86
20) 3+4-Methylphenols	7.50	107	215849	40.60	ng	99
22) Acetophenone	7.45	105	274826	40.87	ng	# 99
24) Nitrobenzene	7.66	77	201712	39.41	ng	99
25) Isophorone	7.96	82	386388	40.82	ng	98
26) 2-Nitrophenol	8.06	139	105085	43.37	ng	98
27) 2,4-Dimethylphenol	8.14	122	179999	41.32	ng	98
28) bis(2-Chloroethoxy)methane	8.25	93	247439	42.18	ng	97
29) 2,4-Dichlorophenol	8.36	162	161957	40.59	ng	98
30) 1,2,4-Trichlorobenzene	8.44	180	168611	40.74	ng	97
31) Naphthalene	8.54	128	573962	39.84	ng	99
32) Benzoic acid	8.28	122	108380	40.10	ng	97
33) 4-Chloroaniline	8.63	127	253903	42.19	ng	99
34) Hexachlorobutadiene	8.70	225	97599	41.46	ng	97
35) Caprolactam	9.06	113	49849	39.62	ng	# 83
36) 4-Chloro-3-methylphenol	9.26	107	162558	39.28	ng	# 73
37) 2-Methylnaphthalene	9.39	142	395084	39.50	ng	97
39) 1,2,4,5-Tetrachlorobenzene	9.60	216	163442	41.40	ng	# 100
40) Hexachlorocyclopentadiene	9.59	237	27736	35.10	ng	99

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42) 2,4,6-Trichlorophenol	9.76	196	114727	41.62	ng	99
43) 2,4,5-Trichlorophenol	9.82	196	117052	42.66	ng	93
45) 1,1'-Biphenyl	9.98	154	455182	41.29	ng	97
46) 2-Chloronaphthalene	10.00	162	357064	41.01	ng	95
47) 2-Nitroaniline	10.14	65	100987	39.01	ng	# 86
48) Acenaphthylene	10.50	152	601148	41.59	ng	99
49) Dimethylphthalate	10.38	163	406252	39.06	ng	99
50) 2,6-Dinitrotoluene	10.46	165	88230	42.46	ng	98
51) Acenaphthene	10.72	154	344540	41.69	ng	99
52) 3-Nitroaniline	10.65	138	109965	40.69	ng	91
53) 2,4-Dinitrophenol	10.80	184	28659	44.65	ng	97
54) Dibenzofuran	10.94	168	506353	41.54	ng	93
55) 4-Nitrophenol	10.90	139	49657m	33.13	ng	
56) 2,4-Dinitrotoluene	10.95	165	117829	42.28	ng	99
57) Fluorene	11.36	166	407723	40.58	ng	99
58) 2,3,4,6-Tetrachlorophenol	11.10	232	86743	43.44	ng	# 100
59) Diethylphthalate	11.26	149	414884	40.75	ng	99
60) 4-Chlorophenyl-phenylether	11.37	204	182786	42.10	ng	90
61) 4-Nitroaniline	11.40	138	98586	39.16	ng	95
62) Azobenzene	11.56	77	383201	38.70	ng	94
64) 4,6-Dinitro-2-methylphenol	11.45	198	51697	40.79	ng	95
65) n-Nitrosodiphenylamine	11.52	169	352734	41.08	ng	97
66) 4-Bromophenyl-phenylether	11.98	248	115808	43.50	ng	# 89
67) Hexachlorobenzene	12.02	284	126243	41.59	ng	# 79
68) Atrazine	12.20	200	97627	42.09	ng	98
69) Pentachlorophenol	12.28	266	47110	40.52	ng	95
70) Phenanthrene	12.54	178	578239	40.77	ng	99
71) Anthracene	12.61	178	597801	41.52	ng	99
72) Carbazole	12.82	167	567110	40.91	ng	97
73) Di-n-butylphthalate	13.28	149	695250	44.85	ng	99
74) Fluoranthene	14.01	202	642522	42.25	ng	98
76) Benzidine	14.21	184	277551	49.58	ng	98
77) Pyrene	14.29	202	647662	40.00	ng	100
79) Butylbenzylphthalate	15.13	149	316513	43.47	ng	92
80) Benzo(a)anthracene	15.77	228	606500	40.34	ng	100
81) 3,3'-Dichlorobenzidine	15.76	252	230865	41.93	ng	# 98
82) Chrysene	15.82	228	579878	40.58	ng	99
83) Bis(2-ethylhexyl)phthalate	15.85	149	471574	45.21	ng	99
84) Di-n-octyl phthalate	16.65	149	849104	46.77	ng	99
85) Indeno(1,2,3-cd)pyrene	18.81	276	786783	42.80	ng	98
87) Benzo(b)fluoranthene	17.07	252	753779	46.55	ng	99
88) Benzo(k)fluoranthene	17.11	252	512437m	35.39	ng	
89) Benzo(a)pyrene	17.45	252	604121	42.43	ng	# 97
90) Dibenzo(a,h)anthracene	18.83	278	658611	44.31	ng	98
91) Benzo(g,h,i)perylene	19.19	276	646367	43.58	ng	96

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

