

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071615\
 Data File : BF080513.D
 Acq On : 16 Jul 2015 16:35
 Operator : TP/UM
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Jul 17 12:22:45 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF071615.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 16 17:11:01 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.07	152	18408	20.00	ng	0.00
21) Naphthalene-d8	8.36	136	79770	20.00	ng	0.00
38) Acenaphthene-d10	10.12	164	36764	20.00	ng	0.00
63) Phenanthrene-d10	11.62	188	62835	20.00	ng	0.01
75) Chrysene-d12	14.58	240	59954	20.00	ng	0.15
86) Perylene-d12	16.42	264	57607	20.00	ng	0.27

System Monitoring Compounds

5) 2-Fluorophenol	5.70	112	92142	85.33	ng	0.00
7) Phenol-d6	6.72	99	123783	84.42	ng	0.00
23) Nitrobenzene-d5	7.64	82	117138	86.20	ng	0.00
41) 2,4,6-Tribromophenol	10.91	330	24990	87.36	ng	0.00
44) 2-Fluorobiphenyl	9.44	172	228256	84.49	ng	0.00
78) Terphenyl-d14	13.30	244	223801	77.40	ng	0.07

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.76	88	20915	39.22	ng	# 79
3) Pyridine	3.71	79	45294	37.64	ng	89
4) n-Nitrosodimethylamine	3.55	42	29979	44.42	ng	79
6) Aniline	6.75	93	77719	41.11	ng	# 89
8) 2-Chlorophenol	6.86	128	48660	40.18	ng	# 80
9) Benzaldehyde	6.62	77	41937	42.13	ng	# 68
10) Phenol	6.73	94	66146	41.93	ng	# 62
11) bis(2-Chloroethyl)ether	6.81	93	44472	38.72	ng	92
12) 1,3-Dichlorobenzene	7.01	146	62287	41.11	ng	# 91
13) 1,4-Dichlorobenzene	7.25	146	54520	34.80	ng	96
14) 1,2-Dichlorobenzene	7.25	146	54520	39.66	ng	99
15) Benzyl Alcohol	7.22	79	42084	41.12	ng	# 68
16) 2,2'-oxybis(1-Chloropropan	7.34	45	83289	40.44	ng	59
17) 2-Methylphenol	7.33	107	39916	41.17	ng	# 76
18) Hexachloroethane	7.60	117	18556	39.60	ng	# 63
19) n-Nitroso-di-n-propylamine	7.48	70	36245	40.72	ng	# 80
20) 3+4-Methylphenols	7.49	107	55119	42.07	ng	# 79
22) Acetophenone	7.48	105	73432	41.60	ng	# 79
24) Nitrobenzene	7.66	77	54672	39.85	ng	# 75
25) Isophorone	7.89	82	99943	41.64	ng	# 90
26) 2-Nitrophenol	7.98	139	23397	39.29	ng	# 75
27) 2,4-Dimethylphenol	8.02	122	48612	42.27	ng	# 80
28) bis(2-Chloroethoxy)methane	8.10	93	57487	40.78	ng	# 91
29) 2,4-Dichlorophenol	8.22	162	44354	42.31	ng	96
30) 1,2,4-Trichlorobenzene	8.30	180	47980	39.87	ng	98
31) Naphthalene	8.38	128	159491	40.71	ng	99
32) Benzoic acid	8.10	122	22815	38.38	ng	# 49
33) 4-Chloroaniline	8.44	127	63179	41.48	ng	# 90
34) Hexachlorobutadiene	8.50	225	27398	39.32	ng	99
35) Caprolactam	8.78	113	10118	40.85	ng	87
36) 4-Chloro-3-methylphenol	8.92	107	46420	41.93	ng	# 78
37) 2-Methylnaphthalene	9.07	142	97055	41.45	ng	# 90
39) 1,2,4,5-Tetrachlorobenzene	9.24	216	49141	41.68	ng	98
40) Hexachlorocyclopentadiene	9.23	237	22585	39.03	ng	94

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42) 2,4,6-Trichlorophenol	9.36	196	30252	42.87	ng	99
43) 2,4,5-Trichlorophenol	9.40	196	32546	43.33	ng	96
45) 1,1'-Biphenyl	9.54	154	124645	41.48	ng	94
46) 2-Chloronaphthalene	9.56	162	90268	40.11	ng	# 92
47) 2-Nitroaniline	9.66	65	24886	42.91	ng	# 50
48) Acenaphthylene	9.98	152	158643	42.08	ng	98
49) Dimethylphthalate	9.84	163	108892	43.19	ng	# 98
50) 2,6-Dinitrotoluene	9.90	165	20946	42.45	ng	93
51) Acenaphthene	10.16	154	93245	41.86	ng	88
52) 3-Nitroaniline	10.08	138	22226	40.78	ng	# 27
53) 2,4-Dinitrophenol	10.18	184	8744	37.90	ng	# 49
54) Dibenzofuran	10.33	168	134670	41.82	ng	93
55) 4-Nitrophenol	10.25	139	16620	42.34	ng	# 16
56) 2,4-Dinitrotoluene	10.30	165	30707	42.94	ng	# 88
57) Fluorene	10.67	166	107701	42.60	ng	93
58) 2,3,4,6-Tetrachlorophenol	10.45	232	23721	43.28	ng	96
59) Diethylphthalate	10.53	149	99377	41.79	ng	97
60) 4-Chlorophenyl-phenylether	10.66	204	46440	41.04	ng	96
61) 4-Nitroaniline	10.69	138	25166	46.64	ng	# 43
62) Azobenzene	10.82	77	107915	43.20	ng	95
64) 4,6-Dinitro-2-methylphenol	10.72	198	12744	41.62	ng	# 71
65) n-Nitrosodiphenylamine	10.77	169	94526	44.23	ng	92
66) 4-Bromophenyl-phenylether	11.15	248	26607	43.48	ng	# 87
67) Hexachlorobenzene	11.22	284	29400	43.45	ng	# 89
68) Atrazine	11.30	200	25182	43.52	ng	84
69) Pentachlorophenol	11.42	266	13480	40.28	ng	98
70) Phenanthrene	11.64	178	148883	42.08	ng	97
71) Anthracene	11.69	178	150644	44.90	ng	99
72) Carbazole	11.86	167	128728	43.74	ng	98
73) Di-n-butylphthalate	12.18	149	154467	44.79	ng	# 98
74) Fluoranthene	12.90	202	146874	44.32	ng	88
76) Benzidine	13.04	184	81126	41.09	ng	# 97
77) Pyrene	13.15	202	155263	39.71	ng	97
79) Butylbenzylphthalate	13.86	149	56715	40.51	ng	93
80) Benzo(a)anthracene	14.57	228	142799	42.45	ng	98
81) 3,3'-Dichlorobenzidine	14.53	252	42213	38.37	ng	# 93
82) Chrysene	14.61	228	133205	40.71	ng	96
83) Bis(2-ethylhexyl)phthalate	14.55	149	82635	40.37	ng	# 90
84) Di-n-octyl phthalate	15.37	149	120996	36.97	ng	100
85) Indeno(1,2,3-cd)pyrene	18.04	276	173605	43.99	ng	# 92
87) Benzo(b)fluoranthene	15.93	252	148430	44.52	ng	# 88
88) Benzo(k)fluoranthene	15.93	252	148430	43.80	ng	# 89
89) Benzo(a)pyrene	16.35	252	120638	38.70	ng	# 86
90) Dibenzo(a,h)anthracene	18.05	278	143210	41.73	ng	# 84
91) Benzo(g,h,i)perylene	18.52	276	147970	41.52	ng	# 81

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

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