

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040216\
 Data File : BF085967.D
 Acq On : 1 Apr 2016 17:50
 Operator : UM/SJ
 Sample : SSTDICC050
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC050

Manual Integrations
 APPROVED

Sohil
 4/4/2016 7:31:50 PM

Quant Time: Apr 01 18:45:55 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040216.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 01 18:28:07 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.77	152	98960	20.00	ng	0.00
21) Naphthalene-d8	8.06	136	397477	20.00	ng	0.00
38) Acenaphthene-d10	9.81	164	184459	20.00	ng	0.00
63) Phenanthrene-d10	11.30	188	341668	20.00	ng	0.00
75) Chrysene-d12	13.93	240	263398	20.00	ng	0.00
86) Perylene-d12	15.33	264	225023	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.37	112	612465	103.07	ng	0.00
7) Phenol-d6	6.42	99	714308	94.55	ng	0.00
23) Nitrobenzene-d5	7.34	82	639710	90.63	ng	0.00
41) 2,4,6-Tribromophenol	10.60	330	168631	96.22	ng	0.00
44) 2-Fluorobiphenyl	9.14	172	1092035	98.91	ng	0.00
78) Terphenyl-d14	12.88	244	1096426	85.61	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.29	88	137098	48.32	ng	98
3) Pyridine	3.00	79	348992	51.30	ng	100
4) n-Nitrosodimethylamine	2.96	42	142546	52.35	ng	97
6) Aniline	6.44	93	504640	49.62	ng	99
8) 2-Chlorophenol	6.56	128	335366	48.58	ng	91
9) Benzaldehyde	6.33	77	192758	42.35	ng	91
10) Phenol	6.43	94	432315	50.51	ng	85
11) bis(2-Chloroethyl)ether	6.51	93	309287	46.96	ng	92
12) 1,3-Dichlorobenzene	6.72	146	368064m	49.51	ng	
13) 1,4-Dichlorobenzene	6.80	146	368730	48.90	ng	93
14) 1,2-Dichlorobenzene	6.94	146	341264	48.57	ng	97
15) Benzyl Alcohol	6.92	79	228810	45.39	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.06	45	395403	45.24	ng	94
17) 2-Methylphenol	7.04	107	272408	49.29	ng	98
18) Hexachloroethane	7.29	117	137194	49.69	ng	96
19) n-Nitroso-di-n-propylamine	7.20	70	194937	43.09	ng	94
20) 3+4-Methylphenols	7.20	107	308636	46.44	ng	91
22) Acetophenone	7.20	105	457183	49.37	ng	# 94
24) Nitrobenzene	7.37	77	393628	54.10	ng	92
25) Isophorone	7.61	82	657382	48.35	ng	97
26) 2-Nitrophenol	7.69	139	178562	49.06	ng	# 87
27) 2,4-Dimethylphenol	7.72	122	283384	44.54	ng	99
28) bis(2-Chloroethoxy)methane	7.81	93	354096	45.71	ng	98
29) 2,4-Dichlorophenol	7.93	162	269179	50.33	ng	96
30) 1,2,4-Trichlorobenzene	8.01	180	298778	50.32	ng	95
31) Naphthalene	8.09	128	981919	49.03	ng	99
32) Benzoic acid	7.87	122	241351	57.12	ng	99
33) 4-Chloroaniline	8.14	127	380687	46.55	ng	95
34) Hexachlorobutadiene	8.20	225	162010	50.20	ng	99
35) Caprolactam	8.52	113	85271	45.01	ng	98
36) 4-Chloro-3-methylphenol	8.62	107	292544	46.82	ng	95
37) 2-Methylnaphthalene	8.77	142	589492	46.74	ng	98
39) 1,2,4,5-Tetrachlorobenzene	8.94	216	285369	52.17	ng	99
40) Hexachlorocyclopentadiene	8.92	237	87193	48.24	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.06	196	184090	49.83	nq	98
43) 2,4,5-Trichlorophenol	9.10	196	178532	46.08	nq	98
45) 1,1'-Biphenyl	9.24	154	803534	51.62	nq	95
46) 2-Chloronaphthalene	9.26	162	602787	52.67	nq	96
47) 2-Nitroaniline	9.37	65	187945	53.70	nq	88
48) Acenaphthylene	9.68	152	933629	50.06	nq	100
49) Dimethylphthalate	9.54	163	683697	48.85	nq	99
50) 2,6-Dinitrotoluene	9.61	165	142946	47.45	nq	95
51) Acenaphthene	9.85	154	536054	47.06	nq	99
52) 3-Nitroaniline	9.78	138	187894	54.49	nq	93
53) 2,4-Dinitrophenol	9.89	184	79254	61.40	nq	# 74
54) Dibenzofuran	10.02	168	693228	47.02	nq	98
55) 4-Nitrophenol	9.95	139	110862	49.35	nq	93
56) 2,4-Dinitrotoluene	10.02	165	196245	51.25	nq	# 81
57) Fluorene	10.36	166	616042	50.26	nq	100
58) 2,3,4,6-Tetrachlorophenol	10.14	232	141040	52.38	nq	96
59) Diethylphthalate	10.24	149	641989	46.43	nq	97
60) 4-Chlorophenyl-phenylether	10.35	204	263980	44.05	nq	95
61) 4-Nitroaniline	10.40	138	190287	57.70	nq	96
62) Azobenzene	10.51	77	640530	48.21	nq	97
64) 4,6-Dinitro-2-methylphenol	10.43	198	111036	55.49	nq	# 52
65) n-Nitrosodiphenylamine	10.48	169	540606	49.96	nq	99
66) 4-Bromophenyl-phenylether	10.84	248	182261	52.14	nq	99
67) Hexachlorobenzene	10.91	284	191953	52.51	nq	94
68) Atrazine	11.00	200	154185	42.20	nq	97
69) Pentachlorophenol	11.10	266	92184	51.37	nq	98
70) Phenanthrene	11.32	178	940725	53.62	nq	100
71) Anthracene	11.37	178	919616	49.93	nq	99
72) Carbazole	11.53	167	768990	49.73	nq	99
73) Di-n-butylphthalate	11.86	149	1061704	50.04	nq	100
74) Fluoranthene	12.51	202	966559	51.83	nq	88
76) Benzidine	12.63	184	504625	54.40	nq	98
77) Pyrene	12.73	202	953376	44.03	nq	99
79) Butylbenzylphthalate	13.36	149	467154	51.52	nq	99
80) Benzo(a)anthracene	13.92	228	788449	52.50	nq	99
81) 3,3'-Dichlorobenzidine	13.89	252	649473	62.31	nq	# 95
82) Chrysene	13.96	228	825173	54.55	nq	99
83) Bis(2-ethylhexyl)phthalate	13.91	149	582676	53.67	nq	100
84) Di-n-octyl phthalate	14.52	149	998973	61.88	nq	99
85) Indeno(1,2,3-cd)pyrene	16.69	276	611457	52.34	nq	99
87) Benzo(b)fluoranthene	14.95	252	820382m	57.87	nq	
88) Benzo(k)fluoranthene	14.97	252	572340m	44.79	nq	
89) Benzo(a)pyrene	15.28	252	633036	50.29	nq	98
90) Dibenzo(a,h)anthracene	16.71	278	515720	44.10	nq	99
91) Benzo(g,h,i)perylene	17.11	276	490752	41.44	nq	# 94

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

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