

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041015\
 Data File : BF078420.D
 Acq On : 11 Apr 2015 1:11
 Operator : TP/IZ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040EC

Manual Integrations
 APPROVED

apatel
 4/13/2015 4:16:51 PM

Quant Time: Apr 13 11:01:25 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Apr 11 02:29:51 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	32041	20.00	ng	-0.01
21) Naphthalene-d8	8.43	136	136928	20.00	ng	-0.01
38) Acenaphthene-d10	10.59	164	71851	20.00	ng	0.00
63) Phenanthrene-d10	12.42	188	143661	20.00	ng	0.00
75) Chrysene-d12	15.68	240	164786	20.00	ng	-0.01
86) Perylene-d12	17.32	264	159568	20.00	ng	-0.13

System Monitoring Compounds

5) 2-Fluorophenol	5.10	112	158284	81.45	ng	-0.01
7) Phenol-d6	6.45	99	200696	82.41	ng	-0.01
23) Nitrobenzene-d5	7.56	82	191029	84.56	ng	0.00
41) 2,4,6-Tribromophenol	11.57	330	56502	94.82	ng	0.00
44) 2-Fluorobiphenyl	9.78	172	393536	78.29	ng	0.00
78) Terphenyl-d14	14.41	244	567850	77.87	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	1.64	88	49669m	56.52	ng	
3) Pyridine	2.22	79	92571	40.21	ng	99
4) n-Nitrosodimethylamine	2.17	42	41871	47.26	ng	90
6) Aniline	6.44	93	145159	42.03	ng	91
8) 2-Chlorophenol	6.59	128	96000	41.78	ng	98
9) Benzaldehyde	6.29	77	59708	36.99	ng	90
10) Phenol	6.48	94	114439	39.22	ng	98
11) bis(2-Chloroethyl)ether	6.56	93	85543	40.76	ng	94
12) 1,3-Dichlorobenzene	6.77	146	102783	40.72	ng	97
13) 1,4-Dichlorobenzene	6.88	146	105448	41.50	ng	99
14) 1,2-Dichlorobenzene	7.06	146	99052	41.58	ng	98
15) Benzyl Alcohol	7.06	79	76625	44.77	ng	94
16) 2,2'-oxybis(1-Chloropropan	7.24	45	120182	42.93	ng	97
17) 2-Methylphenol	7.22	107	75922	42.56	ng	97
18) Hexachloroethane	7.48	117	36332	42.40	ng	96
19) n-Nitroso-di-n-propylamine	7.39	70	68342	42.53	ng	# 85
20) 3+4-Methylphenols	7.42	107	103447	43.47	ng	92
22) Acetophenone	7.38	105	133272	41.77	ng	# 92
24) Nitrobenzene	7.58	77	104203	44.12	ng	89
25) Isophorone	7.88	82	184806	43.10	ng	# 92
26) 2-Nitrophenol	7.97	139	46883	41.66	ng	# 77
27) 2,4-Dimethylphenol	8.06	122	86867	41.51	ng	98
28) bis(2-Chloroethoxy)methane	8.18	93	117700	42.81	ng	100
29) 2,4-Dichlorophenol	8.28	162	81725	42.93	ng	91
30) 1,2,4-Trichlorobenzene	8.37	180	89077	40.81	ng	98
31) Naphthalene	8.46	128	294727	41.35	ng	99
32) Benzoic acid	8.20	122	49125	49.69	ng	99
33) 4-Chloroaniline	8.56	127	125111	42.30	ng	95
34) Hexachlorobutadiene	8.62	225	51148	42.67	ng	98
35) Caprolactam	8.98	113	26272	46.23	ng	96
36) 4-Chloro-3-methylphenol	9.17	107	85627	44.57	ng	87
37) 2-Methylnaphthalene	9.32	142	201943	41.74	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.53	216	91074	40.91	ng	99
40) Hexachlorocyclopentadiene	9.50	237	30269	35.71	ng	98

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041015\
 Data File : BF078420.D
 Acq On : 11 Apr 2015 1:11
 Operator : TP/IZ
 Sample : SSTDC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDC040EC

Manual Integrations
 APPROVED

apatel
 4/13/2015 4:16:51 PM

Quant Time: Apr 13 11:01:25 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Apr 11 02:29:51 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.68	196	62357	44.41	ng	99
43) 2,4,5-Trichlorophenol	9.72	196	63383	43.21	ng	95
45) 1,1'-Biphenyl	9.90	154	236563	40.25	ng	99
46) 2-Chloronaphthalene	9.92	162	188516	40.77	ng	95
47) 2-Nitroaniline	10.05	65	52547	45.93	ng	87
48) Acenaphthylene	10.42	152	321114	43.00	ng	99
49) Dimethylphthalate	10.29	163	227989	42.24	ng	99
50) 2,6-Dinitrotoluene	10.37	165	48112	43.32	ng	99
51) Acenaphthene	10.64	154	177330	42.18	ng	98
52) 3-Nitroaniline	10.57	138	59035	46.75	ng	93
53) 2,4-Dinitrophenol	10.70	184	14837	45.14	ng	93
54) Dibenzofuran	10.84	168	264342	41.17	ng	97
55) 4-Nitrophenol	10.81	139	36598m	40.76	ng	
56) 2,4-Dinitrotoluene	10.86	165	66619	46.31	ng	# 86
57) Fluorene	11.26	166	218481	41.98	ng	100
58) 2,3,4,6-Tetrachlorophenol	11.01	232	49142	45.19	ng	98
59) Diethylphthalate	11.17	149	236186	45.38	ng	99
60) 4-Chlorophenyl-phenylether	11.29	204	104452	43.62	ng	# 89
61) 4-Nitroaniline	11.31	138	56092	43.94	ng	94
62) Azobenzene	11.48	77	228751	48.16	ng	85
64) 4,6-Dinitro-2-methylphenol	11.36	198	29277	42.29	ng	# 70
65) n-Nitrosodiphenylamine	11.44	169	210189	42.30	ng	99
66) 4-Bromophenyl-phenylether	11.88	248	63350	41.64	ng	99
67) Hexachlorobenzene	11.94	284	68640	41.17	ng	# 92
68) Atrazine	12.11	200	61723	42.88	ng	95
69) Pentachlorophenol	12.19	266	32347	47.27	ng	98
70) Phenanthrene	12.44	178	333020	40.07	ng	99
71) Anthracene	12.51	178	353888	43.22	ng	99
72) Carbazole	12.73	167	333865	43.50	ng	98
73) Di-n-butylphthalate	13.19	149	410364	47.20	ng	99
74) Fluoranthene	13.91	202	379188	43.27	ng	97
76) Benzidine	14.10	184	216218	34.08	ng	99
77) Pyrene	14.18	202	386694	37.38	ng	99
79) Butylbenzylphthalate	15.04	149	186673	47.35	ng	# 83
80) Benzo(a)anthracene	15.67	228	388432	39.62	ng	98
81) 3,3'-Dichlorobenzidine	15.65	252	138293	42.12	ng	# 97
82) Chrysene	15.71	228	366443	39.78	ng	99
83) Bis(2-ethylhexyl)phthalate	15.76	149	287189	46.44	ng	97
84) Di-n-octyl phthalate	16.52	149	505737	49.81	ng	# 100
85) Indeno(1,2,3-cd)pyrene	18.53	276	435873m	41.70	ng	
87) Benzo(b)fluoranthene	16.91	252	404309	41.00	ng	99
88) Benzo(k)fluoranthene	16.93	252	341148m	38.93	ng	
89) Benzo(a)pyrene	17.27	252	352988	41.01	ng	98
90) Dibenzo(a,h)anthracene	18.56	278	356366m	41.36	ng	
91) Benzo(g,h,i)perylene	18.89	276	355533m	41.15	ng	

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

