

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051917\
 Data File : BF095266.D
 Acq On : 19 May 2017 16:45
 Operator : SJ/MA
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad
 5/22/2017 1:29:40 PM

Quant Time: May 20 01:21:35 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 18 17:07:53 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.77	152	106175	20.00	ng	0.00
21) Naphthalene-d8	9.78	136	438292	20.00	ng	0.00
38) Acenaphthene-d10	12.61	164	189195	20.00	ng	0.00
63) Phenanthrene-d10	15.02	188	336598	20.00	ng	0.00
75) Chrysene-d12	18.69	240	258375	20.00	ng	0.00
86) Perylene-d12	20.36	264	187630	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.74	112	526937	82.74	ng	0.00
7) Phenol-d6	7.31	99	621288	81.31	ng	0.00
23) Nitrobenzene-d5	8.67	82	505982	72.80	ng	0.00
41) 2,4,6-Tribromophenol	13.91	330	122358	78.97	ng	0.00
44) 2-Fluorobiphenyl	11.55	172	1020331	77.62	ng	0.00
78) Terphenyl-d14	17.34	244	969503	82.65	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.23	88	113236	36.26	ng	94
3) Pyridine	2.93	79	259993	35.92	ng	99
4) n-Nitrosodimethylamine	2.86	42	99097m	32.84	ng	
6) Aniline	7.27	93	409309	42.43	ng	96
8) 2-Chlorophenol	7.46	128	293447	40.48	ng	93
9) Benzaldehyde	7.09	77	166625	35.71	ng	100
10) Phenol	7.32	94	293259	36.42	ng	92
11) bis(2-Chloroethyl)ether	7.40	93	260324	39.16	ng	96
12) 1,3-Dichlorobenzene	7.67	146	330278	40.17	ng	100
13) 1,4-Dichlorobenzene	7.79	146	329204	39.48	ng	95
14) 1,2-Dichlorobenzene	8.02	146	314492	40.41	ng	95
15) Benzyl Alcohol	8.05	79	213269	43.39	ng	94
16) 2,2'-oxybis(1-Chloropropan	8.25	45	361884	38.46	ng	79
17) 2-Methylphenol	8.26	107	244126	41.15	ng	96
18) Hexachloroethane	8.54	117	113465	40.17	ng	# 83
19) n-Nitroso-di-n-propylamine	8.47	70	202933	39.27	ng	93
20) 3+4-Methylphenols	8.52	107	325207	40.34	ng	# 59
22) Acetophenone	8.44	105	405812	38.59	ng	# 84
24) Nitrobenzene	8.70	77	270517	37.13	ng	92
25) Isophorone	9.10	82	492004	39.64	ng	# 94
26) 2-Nitrophenol	9.21	139	166318	42.31	ng	99
27) 2,4-Dimethylphenol	9.34	122	253472	39.88	ng	98
28) bis(2-Chloroethoxy)methane	9.46	93	343642	40.02	ng	99
29) 2,4-Dichlorophenol	9.62	162	233230	38.86	ng	99
30) 1,2,4-Trichlorobenzene	9.71	180	252621	39.29	ng	98
31) Naphthalene	9.81	128	884611	40.16	ng	100
32) Benzoic acid	9.67	122	134686	33.94	ng	95
33) 4-Chloroaniline	9.96	127	321112	39.12	ng	# 92
34) Hexachlorobutadiene	10.03	225	132443	39.04	ng	98
35) Caprolactam	10.60	113	64027m	35.53	ng	
36) 4-Chloro-3-methylphenol	10.81	107	254544	39.67	ng	90
37) 2-Methylnaphthalene	10.94	142	563288	38.96	ng	99
39) 1,2,4,5-Tetrachlorobenzene	11.22	216	239154	41.10	ng	98
40) Hexachlorocyclopentadiene	11.19	237	79233	36.00	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	11.44	196	155867	40.12	ng	99
43) 2,4,5-Trichlorophenol	11.54	196	157055	37.62	ng #	91
45) 1,1'-Biphenyl	11.71	154	656475	41.21	ng	98
46) 2-Chloronaphthalene	11.72	162	512917	39.88	ng	95
47) 2-Nitroaniline	11.95	65	144595	41.07	ng #	77
48) Acenaphthylene	12.38	152	823167	40.86	ng	98
49) Dimethylphthalate	12.26	163	609300	39.23	ng	99
50) 2,6-Dinitrotoluene	12.35	165	129147	40.76	ng	98
51) Acenaphthene	12.67	154	476936	39.64	ng	98
52) 3-Nitroaniline	12.63	138	120415	35.41	ng	87
53) 2,4-Dinitrophenol	12.85	184	57563	36.65	ng #	65
54) Dibenzofuran	12.95	168	685658	41.18	ng	98
55) 4-Nitrophenol	13.05	139	67001m	32.80	ng	
56) 2,4-Dinitrotoluene	13.02	165	175905	43.20	ng #	92
57) Fluorene	13.51	166	575235	39.73	ng	98
58) 2,3,4,6-Tetrachlorophenol	13.20	232	103867	39.53	ng #	95
59) Diethylphthalate	13.40	149	592552	39.80	ng	99
60) 4-Chlorophenyl-phenylether	13.53	204	241781	38.00	ng	90
61) 4-Nitroaniline	13.63	138	112271	35.96	ng	93
62) Azobenzene	13.79	77	469432	39.24	ng	94
64) 4,6-Dinitro-2-methylphenol	13.69	198	86545	38.35	ng #	68
65) n-Nitrosodiphenylamine	13.74	169	475408	38.92	ng	98
66) 4-Bromophenyl-phenylether	14.32	248	131181	36.89	ng	97
67) Hexachlorobenzene	14.41	284	143363	39.34	ng #	90
68) Atrazine	14.66	200	136073	39.45	ng	95
69) Pentachlorophenol	14.77	266	54995	36.79	ng	95
70) Phenanthrene	15.06	178	776802	40.17	ng	98
71) Anthracene	15.14	178	799362	40.46	ng	97
72) Carbazole	15.42	167	716781	39.88	ng	98
73) Di-n-butylphthalate	15.98	149	932839	41.62	ng	98
74) Fluoranthene	16.80	202	745042	37.55	ng	98
76) Benzidine	17.02	184	316386m	45.65	ng	
77) Pyrene	17.10	202	812409	42.04	ng	98
79) Butylbenzylphthalate	17.99	149	370514	41.22	ng #	88
80) Benzo(a)anthracene	18.68	228	591538	39.34	ng	98
81) 3,3'-Dichlorobenzidine	18.66	252	203970	39.27	ng #	96
82) Chrysene	18.72	228	600197	41.28	ng	100
83) Bis(2-ethylhexyl)phthalate	18.74	149	510720	39.88	ng #	99
84) Di-n-octyl phthalate	19.51	149	788788	37.57	ng	96
85) Indeno(1,2,3-cd)pyrene	21.69	276	379545	32.14	ng	100
87) Benzo(b)fluoranthene	19.95	252	442336	38.15	ng	98
88) Benzo(k)fluoranthene	19.98	252	530645m	47.45	ng	
89) Benzo(a)pyrene	20.31	252	427405	39.95	ng	99
90) Dibenzo(a,h)anthracene	21.69	278	330715	37.43	ng	98
91) Benzo(g,h,i)perylene	22.08	276	327509	37.77	ng	97

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

