

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041917\
 Data File : BF094513.D
 Acq On : 19 Apr 2017 13:40
 Operator : SJ/MA
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Apr 20 04:02:37 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF041717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 17 14:59:23 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	5.77	152	129304	20.00	ng	0.00
21) Naphthalene-d8	7.27	136	526256	20.00	ng	0.00
38) Acenaphthene-d10	9.40	164	238213	20.00	ng	0.00
63) Phenanthrene-d10	11.22	188	454217	20.00	ng	0.00
75) Chrysene-d12	14.47	240	416895	20.00	ng	0.00
86) Perylene-d12	16.09	264	369138	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	4.32	112	638769	81.96	ng	0.02
7) Phenol-d6	5.40	99	735868	80.09	ng	0.00
23) Nitrobenzene-d5	6.42	82	664231	77.94	ng	0.00
41) 2,4,6-Tribromophenol	10.38	330	159462	85.58	ng	0.00
44) 2-Fluorobiphenyl	8.59	172	1201839	74.23	ng	0.00
78) Terphenyl-d14	13.20	244	1200792	76.99	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	2.10	88	190210	41.96	ng	95
3) Pyridine	2.62	79	406029	40.99	ng	98
4) n-Nitrosodimethylamine	2.57	42	157039	38.31	ng	87
6) Aniline	5.40	93	488769	40.03	ng	92
8) 2-Chlorophenol	5.54	128	367295	41.41	ng	98
9) Benzaldehyde	5.27	77	240251	34.37	ng	99
10) Phenol	5.41	94	464290	41.13	ng	98
11) bis(2-Chloroethyl)ether	5.48	93	337671	40.95	ng	98
12) 1,3-Dichlorobenzene	5.71	146	384045	39.09	ng	98
13) 1,4-Dichlorobenzene	5.79	146	385794	39.03	ng	95
14) 1,2-Dichlorobenzene	5.97	146	353467	38.82	ng	96
15) Benzyl Alcohol	5.95	79	265894	39.01	ng	98
16) 2,2'-oxybis(1-Chloropropan	6.10	45	429285	39.71	ng	63
17) 2-Methylphenol	6.10	107	270082	38.67	ng	94
18) Hexachloroethane	6.35	117	134052	39.56	ng	# 86
19) n-Nitroso-di-n-propylamine	6.26	70	242551	39.84	ng	96
20) 3+4-Methylphenols	6.28	107	380462	40.08	ng	# 63
22) Acetophenone	6.25	105	502151	39.24	ng	# 86
24) Nitrobenzene	6.44	77	351299	39.14	ng	96
25) Isophorone	6.73	82	659099	41.72	ng	96
26) 2-Nitrophenol	6.82	139	206098	42.74	ng	99
27) 2,4-Dimethylphenol	6.90	122	317776	40.57	ng	98
28) bis(2-Chloroethoxy)methane	7.00	93	424243	41.51	ng	99
29) 2,4-Dichlorophenol	7.12	162	301584	41.39	ng	98
30) 1,2,4-Trichlorobenzene	7.20	180	289948	38.49	ng	97
31) Naphthalene	7.30	128	967838	38.26	ng	100
32) Benzoic acid	7.08	122	196314	47.40	ng	94
33) 4-Chloroaniline	7.37	127	420582	39.96	ng	94
34) Hexachlorobutadiene	7.45	225	151587	40.36	ng	99
35) Caprolactam	7.82	113	103252m	45.11	ng	
36) 4-Chloro-3-methylphenol	8.00	107	306020	40.08	ng	95
37) 2-Methylnaphthalene	8.13	142	690531	39.90	ng	99
39) 1,2,4,5-Tetrachlorobenzene	8.34	216	278721	41.09	ng	99
40) Hexachlorocyclopentadiene	8.32	237	118082	43.03	ng	99

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42) 2,4,6-Trichlorophenol	8.49	196	196313	41.21	ng	100
43) 2,4,5-Trichlorophenol	8.55	196	192063	39.26	ng	93
45) 1,1'-Biphenyl	8.71	154	744025	38.23	ng	98
46) 2-Chloronaphthalene	8.72	162	589791	38.11	ng	93
47) 2-Nitroaniline	8.86	65	179964	40.42	ng	87
48) Acenaphthylene	9.23	152	923540	38.28	ng	99
49) Dimethylphthalate	9.10	163	714939	40.27	ng	100
50) 2,6-Dinitrotoluene	9.17	165	162477	40.78	ng	89
51) Acenaphthene	9.44	154	558902	38.68	ng	99
52) 3-Nitroaniline	9.37	138	179099	38.85	ng	89
53) 2,4-Dinitrophenol	9.51	184	85131	42.30	ng	# 70
54) Dibenzofuran	9.65	168	831135	39.58	ng	100
55) 4-Nitrophenol	9.62	139	116131	35.66	ng	# 67
56) 2,4-Dinitrotoluene	9.67	165	208350	42.61	ng	92
57) Fluorene	10.08	166	648524	39.66	ng	99
58) 2,3,4,6-Tetrachlorophenol	9.82	232	148513	42.36	ng	96
59) Diethylphthalate	9.97	149	691980	41.31	ng	98
60) 4-Chlorophenyl-phenylether	10.08	204	282171	41.46	ng	93
61) 4-Nitroaniline	10.12	138	172863	36.88	ng	95
62) Azobenzene	10.28	77	669589	41.17	ng	96
64) 4,6-Dinitro-2-methylphenol	10.17	198	123541	41.51	ng	# 72
65) n-Nitrosodiphenylamine	10.24	169	595279	37.68	ng	99
66) 4-Bromophenyl-phenylether	10.68	248	173452	38.46	ng	94
67) Hexachlorobenzene	10.75	284	176185	38.76	ng	# 88
68) Atrazine	10.91	200	186629	39.49	ng	97
69) Pentachlorophenol	11.01	266	82435	45.68	ng	97
70) Phenanthrene	11.25	178	971961	38.00	ng	99
71) Anthracene	11.31	178	1026144	38.78	ng	99
72) Carbazole	11.52	167	975607	39.29	ng	97
73) Di-n-butylphthalate	11.97	149	1197906	43.81	ng	99
74) Fluoranthene	12.71	202	1090895	41.15	ng	98
76) Benzidine	12.88	184	629806	37.31	ng	# 95
77) Pyrene	12.99	202	1089845	37.42	ng	99
79) Butylbenzylphthalate	13.81	149	533997	41.83	ng	# 87
80) Benzo(a)anthracene	14.46	228	947702	39.11	ng	99
81) 3,3'-Dichlorobenzidine	14.44	252	341295	39.79	ng	# 96
82) Chrysene	14.50	228	853187	38.53	ng	100
83) Bis(2-ethylhexyl)phthalate	14.52	149	739310	43.14	ng	# 99
84) Di-n-octyl phthalate	15.28	149	1247620	43.40	ng	98
85) Indeno(1,2,3-cd)pyrene	17.37	276	811339	38.51	ng	100
87) Benzo(b)fluoranthene	15.69	252	861480	38.15	ng	99
88) Benzo(k)fluoranthene	15.72	252	761316	35.63	ng	# 94
89) Benzo(a)pyrene	16.04	252	822679	39.16	ng	99
90) Dibenzo(a,h)anthracene	17.38	278	671905	38.52	ng	98
91) Benzo(g,h,i)perylene	17.74	276	658117	37.06	ng	98

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

