

Data Path : Z:\HPCHEM1\BNA F\DATA\BF041718\
 Data File : BF104561.D
 Acq On : 17 Apr 2018 15:16
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Apr 17 17:08:38 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 17 17:08:38 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	137435	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	573384	20.00	ng	0.00
38) Acenaphthene-d10	9.85	164	263006	20.00	ng	0.00
63) Phenanthrene-d10	11.33	188	487073	20.00	ng	0.00
75) Chrysene-d12	13.98	240	368702	20.00	ng	0.00
86) Perylene-d12	15.44	264	317435	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.42	112	678572	75.50	ng	0.00
7) Phenol-d6	6.45	99	851200	77.08	ng	0.00
23) Nitrobenzene-d5	7.37	82	771576	87.87	ng	0.00
41) 2,4,6-Tribromophenol	10.64	330	224485	82.28	ng	0.00
44) 2-Fluorobiphenyl	9.17	172	1253089	75.27	ng	0.00
78) Terphenyl-d14	12.92	244	1238407	76.58	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.52	88	168674	40.274	ng	# 90
3) Pyridine	3.25	79	458467	38.935	ng	86
4) n-Nitrosodimethylamine	3.22	42	152375	37.019	ng	79
6) Aniline	6.47	93	589256	38.405	ng	98
8) 2-Chlorophenol	6.59	128	381085	40.170	ng	94
9) Benzaldehyde	6.36	77	205557	35.434	ng	92
10) Phenol	6.46	94	473777	40.433	ng	88
11) bis(2-Chloroethyl)ether	6.55	93	367381	39.289	ng	94
12) 1,3-Dichlorobenzene	6.75	146	424761	39.522	ng	99
13) 1,4-Dichlorobenzene	6.82	146	424521	39.625	ng	99
14) 1,2-Dichlorobenzene	6.97	146	395652	39.852	ng	99
15) Benzyl Alcohol	6.95	79	320793	38.245	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.08	45	485314	38.647	ng	92
17) 2-Methylphenol	7.06	107	331526	40.202	ng	97
18) Hexachloroethane	7.32	117	155262	40.647	ng	97
19) n-Nitroso-di-n-propylamine	7.22	70	249141	34.524	ng	96
20) 3+4-Methylphenols	7.22	107	378051	36.964	ng	# 70
22) Acetophenone	7.22	105	509155	36.068	ng	# 98
24) Nitrobenzene	7.39	77	409990	42.290	ng	98
25) Isophorone	7.63	82	707698	38.112	ng	98
26) 2-Nitrophenol	7.70	139	211824	53.670	ng	99
27) 2,4-Dimethylphenol	7.75	122	302259	36.883	ng	99
28) bis(2-Chloroethoxy)methane	7.84	93	434065	38.233	ng	99
29) 2,4-Dichlorophenol	7.95	162	306714	39.226	ng	99
30) 1,2,4-Trichlorobenzene	8.03	180	336059	39.162	ng	98
31) Naphthalene	8.12	128	1100088	38.530	ng	100
32) Benzoic acid	7.88	122	291982	44.655	ng	98
33) 4-Chloroaniline	8.16	127	486182	39.747	ng	98
34) Hexachlorobutadiene	8.23	225	182211	38.766	ng	99
35) Caprolactam	8.55	113	106728	39.127	ng	# 81
36) 4-Chloro-3-methylphenol	8.65	107	330731	39.447	ng	99
37) 2-Methylnaphthalene	8.80	142	711572	38.529	ng	98
39) 1,2,4,5-Tetrachlorobenzene	8.97	216	307349	40.823	ng	98
40) Hexachlorocyclopentadiene	8.95	237	167389	39.714	ng	99

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42) 2,4,6-Trichlorophenol	9.08	196	215373	41.209	nq	99
43) 2,4,5-Trichlorophenol	9.13	196	244234	41.760	nq	94
45) 1,1'-Biphenyl	9.27	154	857477	38.810	nq	100
46) 2-Chloronaphthalene	9.29	162	692373	39.674	nq	99
47) 2-Nitroaniline	9.39	65	216954	50.270	nq	98
48) Acenaphthylene	9.71	152	1063759	38.850	nq	100
49) Dimethylphthalate	9.57	163	819864	39.530	nq	99
50) 2,6-Dinitrotoluene	9.63	165	183716	47.871	nq	98
51) Acenaphthene	9.88	154	614422	36.778	nq	100
52) 3-Nitroaniline	9.80	138	218986	48.741	nq	98
53) 2,4-Dinitrophenol	9.91	184	81338	58.940	nq	# 20
54) Dibenzofuran	10.05	168	898173	38.578	nq	97
55) 4-Nitrophenol	9.97	139	155289	44.000	nq	88
56) 2,4-Dinitrotoluene	10.04	165	237393	47.158	nq	94
57) Fluorene	10.40	166	683227	40.317	nq	100
58) 2,3,4,6-Tetrachlorophenol	10.17	232	175926	40.320	nq	94
59) Diethylphthalate	10.27	149	782674	37.891	nq	98
60) 4-Chlorophenyl-phenylether	10.39	204	313671	41.562	nq	97
61) 4-Nitroaniline	10.42	138	217960	49.616	nq	96
62) Azobenzene	10.55	77	747699	37.889	nq	93
64) 4,6-Dinitro-2-methylphenol	10.45	198	130152	53.816	nq	# 75
65) n-Nitrosodiphenylamine	10.51	169	637996	37.778	nq	99
66) 4-Bromophenyl-phenylether	10.88	248	199260	38.461	nq	# 87
67) Hexachlorobenzene	10.95	284	231389	39.692	nq	# 87
68) Atrazine	11.04	200	198624	38.964	nq	97
69) Pentachlorophenol	11.14	266	133122	36.912	nq	98
70) Phenanthrene	11.36	178	1011496	37.290	nq	99
71) Anthracene	11.41	178	1041544	37.774	nq	100
72) Carbazole	11.57	167	1009347	37.462	nq	99
73) Di-n-butylphthalate	11.89	149	1197570	36.386	nq	100
74) Fluoranthene	12.55	202	1051979	37.120	nq	98
76) Benzidine	12.67	184	520799	42.365	nq	99
77) Pyrene	12.78	202	1079232	39.490	nq	100
79) Butylbenzylphthalate	13.40	149	512242	39.785	nq	96
80) Benzo(a)anthracene	13.97	228	900402	40.878	nq	99
81) 3,3'-Dichlorobenzidine	13.93	252	324731	39.540	nq	98
82) Chrysene	14.01	228	880321	38.910	nq	99
83) Bis(2-ethylhexyl)phthalate	13.95	149	664431	40.121	nq	98
84) Di-n-octyl phthalate	14.57	149	1011520	36.554	nq	96
85) Indeno(1,2,3-cd)pyrene	16.89	276	707065	36.789	nq	98
87) Benzo(b)fluoranthene	15.02	252	838526	41.825	nq	98
88) Benzo(k)fluoranthene	15.05	252	736803	38.927	nq	98
89) Benzo(a)pyrene	15.38	252	724830	40.406	nq	99
90) Dibenzo(a,h)anthracene	16.91	278	581671	39.481	nq	99
91) Benzo(g,h,i)perylene	17.33	276	580489	40.668	nq	97

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

