

Data Path : Z:\HPCHEM1\BNA F\DATA\BF041817\
 Data File : BF094481.D
 Acq On : 18 Apr 2017 12:25
 Operator : SJ/MA
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad
 4/19/2017 1:27:31 PM

Quant Time: Apr 19 01:25:25 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	125726	20.00	ng	0.01
21) Naphthalene-d8	7.27	136	540765	20.00	ng	0.00
38) Acenaphthene-d10	9.40	164	245854	20.00	ng	0.00
63) Phenanthrene-d10	11.22	188	452179	20.00	ng	0.00
75) Chrysene-d12	14.47	240	392229	20.00	ng	0.00
86) Perylene-d12	16.09	264	344536	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	4.32	112	626397	82.66	ng	0.02
7) Phenol-d6	5.40	99	692741	77.54	ng	0.01
23) Nitrobenzene-d5	6.43	82	693159	79.15	ng	0.01
41) 2,4,6-Tribromophenol	10.38	330	163494	85.02	ng	0.00
44) 2-Fluorobiphenyl	8.59	172	1268224	75.90	ng	0.00
78) Terphenyl-d14	13.20	244	1154805	78.70	ng	0.00

Target Compounds						Qvalue
2) 1,4-Dioxane	2.10	88	180323	40.92	ng	96
3) Pyridine	2.62	79	399918	41.52	ng	98
4) n-Nitrosodimethylamine	2.58	42	162128	40.68	ng	88
6) Aniline	5.40	93	452892	38.15	ng	89
8) 2-Chlorophenol	5.55	128	341773	39.63	ng	96
9) Benzaldehyde	5.27	77	236448	34.79	ng	97
10) Phenol	5.42	94	425143	38.74	ng	95
11) bis(2-Chloroethyl)ether	5.49	93	318859	39.77	ng	96
12) 1,3-Dichlorobenzene	5.70	146	378713	39.64	ng	100
13) 1,4-Dichlorobenzene	5.79	146	379133	39.44	ng	96
14) 1,2-Dichlorobenzene	5.97	146	354997	40.10	ng	95
15) Benzyl Alcohol	5.95	79	269653	40.69	ng	100
16) 2,2'-oxybis(1-Chloropropan	6.10	45	453105	43.10	ng	81
17) 2-Methylphenol	6.09	107	284259	41.85	ng	96
18) Hexachloroethane	6.36	117	137096	41.61	ng	# 86
19) n-Nitroso-di-n-propylamine	6.26	70	260312	43.98	ng	95
20) 3+4-Methylphenols	6.28	107	396015	42.91	ng	# 62
22) Acetophenone	6.25	105	518125	39.41	ng	# 85
24) Nitrobenzene	6.45	77	372537	40.39	ng	95
25) Isophorone	6.74	82	660888	40.71	ng	95
26) 2-Nitrophenol	6.82	139	201480	40.67	ng	96
27) 2,4-Dimethylphenol	6.90	122	321821	39.99	ng	98
28) bis(2-Chloroethoxy)methane	7.00	93	427085	40.67	ng	98
29) 2,4-Dichlorophenol	7.12	162	304025	40.61	ng	98
30) 1,2,4-Trichlorobenzene	7.20	180	306609	39.61	ng	98
31) Naphthalene	7.30	128	1023394	39.37	ng	99
32) Benzoic acid	7.07	122	185098	43.49	ng	94
33) 4-Chloroaniline	7.37	127	432250	39.97	ng	94
34) Hexachlorobutadiene	7.45	225	154423	40.02	ng	99
35) Caprolactam	7.83	113	94939m	40.37	ng	
36) 4-Chloro-3-methylphenol	7.99	107	306157	39.02	ng	90
37) 2-Methylnaphthalene	8.13	142	710142	39.93	ng	99
39) 1,2,4,5-Tetrachlorobenzene	8.34	216	278636	39.80	ng	100
40) Hexachlorocyclopentadiene	8.33	237	128124	45.24	ng	96

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	8.49	196	195217	39.71	nq	97
43) 2,4,5-Trichlorophenol	8.55	196	202169	40.04	nq	94
45) 1,1'-Biphenyl	8.71	154	789732	39.32	nq	99
46) 2-Chloronaphthalene	8.73	162	634066	39.70	nq	96
47) 2-Nitroaniline	8.86	65	184389	40.13	nq	# 83
48) Acenaphthylene	9.23	152	968231	38.89	nq	99
49) Dimethylphthalate	9.10	163	751633	41.02	nq	99
50) 2,6-Dinitrotoluene	9.17	165	167151	40.64	nq	92
51) Acenaphthene	9.45	154	589611	39.54	nq	99
52) 3-Nitroaniline	9.37	138	185356	38.96	nq	85
53) 2,4-Dinitrophenol	9.50	184	83790	40.60	nq	# 76
54) Dibenzofuran	9.66	168	854303	39.41	nq	100
55) 4-Nitrophenol	9.63	139	132123	39.31	nq	# 78
56) 2,4-Dinitrotoluene	9.66	165	208354	41.29	nq	# 74
57) Fluorene	10.07	166	670163	39.70	nq	100
58) 2,3,4,6-Tetrachlorophenol	9.82	232	156135	43.15	nq	95
59) Diethylphthalate	9.97	149	721961	41.76	nq	99
60) 4-Chlorophenyl-phenylether	10.09	204	293810	41.83	nq	90
61) 4-Nitroaniline	10.13	138	174128	36.00	nq	98
62) Azobenzene	10.28	77	693876	41.34	nq	96
64) 4,6-Dinitro-2-methylphenol	10.17	198	119150	40.22	nq	# 66
65) n-Nitrosodiphenylamine	10.24	169	614240	39.06	nq	99
66) 4-Bromophenyl-phenylether	10.68	248	177387	39.51	nq	96
67) Hexachlorobenzene	10.75	284	179885	39.75	nq	# 85
68) Atrazine	10.92	200	194553	41.35	nq	97
69) Pentachlorophenol	11.00	266	85304	47.48	nq	96
70) Phenanthrene	11.25	178	994808	39.07	nq	99
71) Anthracene	11.32	178	1034220	39.27	nq	98
72) Carbazole	11.52	167	967566	39.14	nq	98
73) Di-n-butylphthalate	11.97	149	1163121	42.73	nq	99
74) Fluoranthene	12.71	202	1048088	39.71	nq	98
76) Benzidine	12.89	184	597330	37.61	nq	96
77) Pyrene	12.99	202	1064358	38.84	nq	99
79) Butylbenzylphthalate	13.81	149	508165	42.31	nq	# 86
80) Benzo(a)anthracene	14.46	228	905820	39.73	nq	99
81) 3,3'-Dichlorobenzidine	14.44	252	319193	39.55	nq	# 97
82) Chrysene	14.50	228	830043	39.84	nq	99
83) Bis(2-ethylhexyl)phthalate	14.53	149	713122	44.23	nq	# 98
84) Di-n-octyl phthalate	15.28	149	1189804	43.99	nq	98
85) Indeno(1,2,3-cd)pyrene	17.37	276	743876	37.52	nq	99
87) Benzo(b)fluoranthene	15.69	252	877499	41.63	nq	99
88) Benzo(k)fluoranthene	15.73	252	737796	36.99	nq	# 94
89) Benzo(a)pyrene	16.04	252	769500	39.24	nq	99
90) Dibenzo(a,h)anthracene	17.39	278	625488	38.42	nq	97
91) Benzo(g,h,i)perylene	17.74	276	594452	35.86	nq	98

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

