

Data Path : Z:\HPCHEM1\BNA F\DATA\BF052617\
 Data File : BF095455.D
 Acq On : 26 May 2017 12:01
 Operator : SJ/MA
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad
 5/30/2017 1:35:26 PM

Quant Time: May 26 17:41:47 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	7.74	152	129095	20.00	ng	-0.02
21) Naphthalene-d8	9.76	136	519414	20.00	ng	-0.02
38) Acenaphthene-d10	12.59	164	233626	20.00	ng	-0.02
63) Phenanthrene-d10	15.00	188	412678	20.00	ng	-0.02
75) Chrysene-d12	18.68	240	331529	20.00	ng	-0.02
86) Perylene-d12	20.35	264	261591	20.00	ng	-0.01

System Monitoring Compounds						
5) 2-Fluorophenol	5.71	112	611043	78.91	ng	-0.03
7) Phenol-d6	7.30	99	736710	79.30	ng	-0.02
23) Nitrobenzene-d5	8.65	82	642423	77.99	ng	-0.02
41) 2,4,6-Tribromophenol	13.90	330	149526	78.15	ng	-0.02
44) 2-Fluorobiphenyl	11.54	172	1176379	72.48	ng	-0.02
78) Terphenyl-d14	17.32	244	1160282	77.09	ng	-0.02

Target Compounds						Qvalue
2) 1,4-Dioxane	2.18	88	135560	35.70	ng	94
3) Pyridine	2.88	79	325977	37.04	ng	97
4) n-Nitrosodimethylamine	2.81	42	118052	32.18	ng	82
6) Aniline	7.25	93	491137	41.87	ng	96
8) 2-Chlorophenol	7.44	128	355169	40.30	ng	96
9) Benzaldehyde	7.07	77	186212	32.82	ng	96
10) Phenol	7.31	94	343873	35.12	ng	88
11) bis(2-Chloroethyl)ether	7.38	93	311298	38.52	ng	96
12) 1,3-Dichlorobenzene	7.65	146	398337	39.84	ng	99
13) 1,4-Dichlorobenzene	7.77	146	407335	40.18	ng	97
14) 1,2-Dichlorobenzene	8.00	146	372017	39.31	ng	96
15) Benzyl Alcohol	8.04	79	263565	44.11	ng	95
16) 2,2'-oxybis(1-Chloropropan	8.23	45	443840	38.80	ng	90
17) 2-Methylphenol	8.24	107	295043	40.91	ng	96
18) Hexachloroethane	8.52	117	132125	38.47	ng	88
19) n-Nitroso-di-n-propylamine	8.46	70	244042	38.84	ng	91
20) 3+4-Methylphenols	8.51	107	381022	38.88	ng	# 59
22) Acetophenone	8.42	105	480518	38.56	ng	# 83
24) Nitrobenzene	8.68	77	337448	39.08	ng	92
25) Isophorone	9.08	82	598626	40.70	ng	96
26) 2-Nitrophenol	9.19	139	197342	42.36	ng	98
27) 2,4-Dimethylphenol	9.32	122	307742	40.86	ng	97
28) bis(2-Chloroethoxy)methane	9.44	93	406505	39.95	ng	99
29) 2,4-Dichlorophenol	9.61	162	276779	38.92	ng	99
30) 1,2,4-Trichlorobenzene	9.69	180	299474	39.30	ng	98
31) Naphthalene	9.80	128	1002665	38.41	ng	100
32) Benzoic acid	9.68	122	170214	35.86	ng	88
33) 4-Chloroaniline	9.94	127	409501	42.09	ng	# 91
34) Hexachlorobutadiene	10.01	225	155276	38.62	ng	98
35) Caprolactam	10.59	113	88866m	41.61	ng	
36) 4-Chloro-3-methylphenol	10.80	107	311855	41.01	ng	89
37) 2-Methylnaphthalene	10.92	142	675148	39.41	ng	99
39) 1,2,4,5-Tetrachlorobenzene	11.20	216	277498	38.62	ng	99
40) Hexachlorocyclopentadiene	11.17	237	95609	35.31	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	11.43	196	187921	39.18	nq	99
43) 2,4,5-Trichlorophenol	11.52	196	178919	34.70	nq	95
45) 1,1'-Biphenyl	11.69	154	762641	38.77	nq	98
46) 2-Chloronaphthalene	11.71	162	611148	38.48	nq	96
47) 2-Nitroaniline	11.93	65	176531	40.61	nq	# 80
48) Acenaphthylene	12.37	152	954383	38.36	nq	99
49) Dimethylphthalate	12.24	163	720676	37.58	nq	99
50) 2,6-Dinitrotoluene	12.34	165	153192	39.15	nq	95
51) Acenaphthene	12.65	154	570101	38.37	nq	99
52) 3-Nitroaniline	12.61	138	174414	41.53	nq	94
53) 2,4-Dinitrophenol	12.84	184	69772	36.09	nq	# 61
54) Dibenzofuran	12.94	168	809471	39.37	nq	99
55) 4-Nitrophenol	13.04	139	90661m	35.94	nq	
56) 2,4-Dinitrotoluene	13.01	165	211372	42.04	nq	# 90
57) Fluorene	13.49	166	681554	38.12	nq	99
58) 2,3,4,6-Tetrachlorophenol	13.18	232	127724	39.36	nq	# 95
59) Diethylphthalate	13.38	149	708735	38.55	nq	98
60) 4-Chlorophenyl-phenylether	13.51	204	301759	38.40	nq	90
61) 4-Nitroaniline	13.62	138	149083	38.67	nq	97
62) Azobenzene	13.77	77	593374	40.17	nq	94
64) 4,6-Dinitro-2-methylphenol	13.69	198	111710	40.38	nq	# 68
65) n-Nitrosodiphenylamine	13.72	169	592445	39.56	nq	99
66) 4-Bromophenyl-phenylether	14.30	248	170557	39.12	nq	97
67) Hexachlorobenzene	14.39	284	170960	38.26	nq	# 88
68) Atrazine	14.65	200	163305	38.62	nq	96
69) Pentachlorophenol	14.75	266	67236	36.71	nq	96
70) Phenanthrene	15.04	178	900113	37.96	nq	98
71) Anthracene	15.12	178	949985	39.22	nq	98
72) Carbazole	15.41	167	881607	40.01	nq	97
73) Di-n-butylphthalate	15.97	149	1107028	40.28	nq	98
74) Fluoranthene	16.78	202	968114	39.80	nq	98
76) Benzidine	17.01	184	432331m	48.19	nq	
77) Pyrene	17.09	202	999368	40.31	nq	98
79) Butylbenzylphthalate	17.98	149	454525	39.41	nq	# 85
80) Benzo(a)anthracene	18.67	228	752543	39.00	nq	99
81) 3,3'-Dichlorobenzidine	18.64	252	262582	39.40	nq	# 96
82) Chrysene	18.71	228	720882	38.64	nq	99
83) Bis(2-ethylhexyl)phthalate	18.72	149	620344	37.75	nq	# 99
84) Di-n-octyl phthalate	19.49	149	1042579	38.70	nq	96
85) Indeno(1,2,3-cd)pyrene	21.68	276	491916	32.47	nq	99
87) Benzo(b)fluoranthene	19.94	252	644621	39.88	nq	96
88) Benzo(k)fluoranthene	19.97	252	624738	40.07	nq	96
89) Benzo(a)pyrene	20.30	252	577359	38.71	nq	99
90) Dibenzo(a,h)anthracene	21.68	278	417605	33.90	nq	97
91) Benzo(g,h,i)perylene	22.07	276	429712	35.55	nq	99

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

