

Data Path : Z:\HPCHEM1\BNA F\DATA\BF020717\
 Data File : BF092823.D
 Acq On : 7 Feb 2017 12:02
 Operator : SJ/MA
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad
 2/8/2017 10:42:41 AM

Quant Time: Feb 07 15:28:48 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF013017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 30 14:50:05 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.91	152	156688	20.00	ng	-0.06
21) Naphthalene-d8	8.19	136	639787	20.00	ng	-0.07
38) Acenaphthene-d10	9.95	164	374037	20.00	ng	-0.06
63) Phenanthrene-d10	11.44	188	605749	20.00	ng	-0.06
75) Chrysene-d12	14.08	240	520399	20.00	ng	-0.07
86) Perylene-d12	15.53	264	418688	20.00	ng	-0.08

System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	721291	78.98	ng	-0.07
7) Phenol-d6	6.54	99	963296	81.97	ng	-0.06
23) Nitrobenzene-d5	7.47	82	918137	79.92	ng	-0.07
41) 2,4,6-Tribromophenol	10.74	330	198713	73.45	ng	-0.07
44) 2-Fluorobiphenyl	9.26	172	1505231	72.91	ng	-0.07
78) Terphenyl-d14	13.03	244	1665568	75.20	ng	-0.06

Target Compounds						Qvalue
2) 1,4-Dioxane	2.51	88	190463	38.59	ng	86
3) Pyridine	3.25	79	555746	44.29	ng	# 82
4) n-Nitrosodimethylamine	3.22	42	237546	40.32	ng	87
6) Aniline	6.57	93	638399	39.83	ng	# 69
8) 2-Chlorophenol	6.69	128	428947	42.57	ng	98
9) Benzaldehyde	6.45	77	311836	37.04	ng	95
10) Phenol	6.57	94	535386	38.94	ng	85
11) bis(2-Chloroethyl)ether	6.65	93	465479	42.42	ng	98
12) 1,3-Dichlorobenzene	6.84	146	467675m	39.63	ng	
13) 1,4-Dichlorobenzene	6.92	146	482379	40.39	ng	98
14) 1,2-Dichlorobenzene	7.08	146	466533	40.85	ng	95
15) Benzyl Alcohol	7.05	79	345837	39.75	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.18	45	743830	39.01	ng	94
17) 2-Methylphenol	7.17	107	386025	44.66	ng	98
18) Hexachloroethane	7.41	117	162347	39.82	ng	# 85
19) n-Nitroso-di-n-propylamine	7.32	70	285816	38.21	ng	94
20) 3+4-Methylphenols	7.32	107	410814	39.75	ng	# 77
22) Acetophenone	7.32	105	571257	39.11	ng	# 92
24) Nitrobenzene	7.49	77	463626	39.30	ng	97
25) Isophorone	7.73	82	862377	40.28	ng	98
26) 2-Nitrophenol	7.81	139	255462	45.55	ng	# 82
27) 2,4-Dimethylphenol	7.85	122	374421	38.02	ng	98
28) bis(2-Chloroethoxy)methane	7.94	93	529866	37.37	ng	99
29) 2,4-Dichlorophenol	8.05	162	375839	40.92	ng	91
30) 1,2,4-Trichlorobenzene	8.13	180	388626	39.26	ng	95
31) Naphthalene	8.21	128	1209318	37.29	ng	100
32) Benzoic acid	7.98	122	224011	41.32	ng	86
33) 4-Chloroaniline	8.27	127	548162	43.10	ng	# 93
34) Hexachlorobutadiene	8.33	225	198370	36.43	ng	99
35) Caprolactam	8.65	113	130992	45.34	ng	# 63
36) 4-Chloro-3-methylphenol	8.75	107	367056	38.33	ng	99
37) 2-Methylnaphthalene	8.91	142	857827	41.19	ng	96
39) 1,2,4,5-Tetrachlorobenzene	9.07	216	376891	35.90	ng	99
40) Hexachlorocyclopentadiene	9.06	237	156469	33.03	ng	96

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42) 2,4,6-Trichlorophenol	9.18	196	260027	37.69	nq	91
43) 2,4,5-Trichlorophenol	9.23	196	294402	38.95	nq #	87
45) 1,1'-Biphenyl	9.37	154	959981	35.44	nq	96
46) 2-Chloronaphthalene	9.39	162	794056	37.48	nq	95
47) 2-Nitroaniline	9.49	65	266565	37.42	nq	97
48) Acenaphthylene	9.81	152	1401760	38.30	nq	98
49) Dimethylphthalate	9.68	163	985453	39.12	nq	99
50) 2,6-Dinitrotoluene	9.73	165	199260	36.62	nq #	78
51) Acenaphthene	9.98	154	853619	39.01	nq	98
52) 3-Nitroaniline	9.90	138	219030	35.18	nq	97
53) 2,4-Dinitrophenol	10.02	184	108917	41.76	nq	90
54) Dibenzofuran	10.16	168	1107845	37.85	nq	93
55) 4-Nitrophenol	10.08	139	179398	40.00	nq	87
56) 2,4-Dinitrotoluene	10.14	165	309855	42.78	nq #	75
57) Fluorene	10.50	166	915906	40.67	nq	98
58) 2,3,4,6-Tetrachlorophenol	10.27	232	191993	38.07	nq	97
59) Diethylphthalate	10.37	149	955912	40.26	nq	95
60) 4-Chlorophenyl-phenylether	10.49	204	393303	36.36	nq	89
61) 4-Nitroaniline	10.52	138	237796	37.29	nq	92
62) Azobenzene	10.65	77	915071	37.31	nq	96
64) 4,6-Dinitro-2-methylphenol	10.56	198	159463	43.51	nq	84
65) n-Nitrosodiphenylamine	10.61	169	820600	42.41	nq	99
66) 4-Bromophenyl-phenylether	10.98	248	252764	39.94	nq	93
67) Hexachlorobenzene	11.05	284	250617	40.07	nq	99
68) Atrazine	11.14	200	265993	42.83	nq	98
69) Pentachlorophenol	11.24	266	126461	42.85	nq	97
70) Phenanthrene	11.46	178	1236875	35.64	nq	99
71) Anthracene	11.52	178	1436551	41.22	nq	98
72) Carbazole	11.66	167	1246086	37.40	nq	99
73) Di-n-butylphthalate	12.00	149	1532379	42.53	nq #	97
74) Fluoranthene	12.65	202	1327914	37.10	nq	97
76) Benzidine	12.77	184	790231	35.64	nq	96
77) Pyrene	12.88	202	1303151	33.46	nq	99
79) Butylbenzylphthalate	13.49	149	652525	41.26	nq #	82
80) Benzo(a)anthracene	14.07	228	1144365	35.95	nq	99
81) 3,3'-Dichlorobenzidine	14.03	252	471737	41.58	nq #	94
82) Chrysene	14.10	228	958568	32.17	nq	100
83) Bis(2-ethylhexyl)phthalate	14.04	149	834891	38.60	nq	99
84) Di-n-octyl phthalate	14.66	149	1382283	39.25	nq	99
85) Indeno(1,2,3-cd)pyrene	16.97	276	856395	37.10	nq	96
87) Benzo(b)fluoranthene	15.11	252	1013183	36.77	nq	99
88) Benzo(k)fluoranthene	15.14	252	1016624m	42.73	nq	
89) Benzo(a)pyrene	15.47	252	947614	39.75	nq #	97
90) Dibenzo(a,h)anthracene	16.99	278	733134	38.05	nq	97
91) Benzo(g,h,i)perylene	17.41	276	710281	36.49	nq	96

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

