

Data Path : Z:\HPCHEM1\BNA F\DATA\BF013117\
 Data File : BF092679.D
 Acq On : 31 Jan 2017 14:11
 Operator : SJ/MA
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad
 2/1/2017 2:27:25 PM

Quant Time: Feb 01 10:44:54 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	6.94	152	160423	20.00	ng	-0.02
21) Naphthalene-d8	8.24	136	622079	20.00	ng	-0.02
38) Acenaphthene-d10	10.01	164	365082	20.00	ng	0.00
63) Phenanthrene-d10	11.49	188	652556	20.00	ng	0.00
75) Chrysene-d12	14.15	240	461517	20.00	ng	0.00
86) Perylene-d12	15.61	264	380264	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.56	112	767439	82.07	ng	-0.01
7) Phenol-d6	6.59	99	948838	78.86	ng	-0.01
23) Nitrobenzene-d5	7.52	82	881405	78.90	ng	-0.02
41) 2,4,6-Tribromophenol	10.80	330	206523	78.21	ng	-0.01
44) 2-Fluorobiphenyl	9.32	172	1576584	78.24	ng	-0.01
78) Terphenyl-d14	13.08	244	1641844	83.59	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.60	88	203607	40.30	ng	# 85
3) Pyridine	3.36	79	617152m	48.04	ng	
4) n-Nitrosodimethylamine	3.31	42	289164	47.93	ng	81
6) Aniline	6.61	93	689926	42.04	ng	# 50
8) 2-Chlorophenol	6.74	128	438309	42.49	ng	93
9) Benzaldehyde	6.50	77	355298	41.22	ng	97
10) Phenol	6.60	94	501499	35.63	ng	97
11) bis(2-Chloroethyl)ether	6.68	93	427842	38.08	ng	92
12) 1,3-Dichlorobenzene	6.89	146	510574m	42.26	ng	
13) 1,4-Dichlorobenzene	6.97	146	537490	43.95	ng	95
14) 1,2-Dichlorobenzene	7.12	146	452443	38.69	ng	97
15) Benzyl Alcohol	7.09	79	357907	40.18	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.23	45	838937	42.98	ng	94
17) 2-Methylphenol	7.22	107	383246	43.31	ng	95
18) Hexachloroethane	7.46	117	168688	40.41	ng	88
19) n-Nitroso-di-n-propylamine	7.37	70	312531	40.81	ng	95
20) 3+4-Methylphenols	7.37	107	421535	39.84	ng	94
22) Acetophenone	7.37	105	595168	41.90	ng	# 91
24) Nitrobenzene	7.54	77	537813	46.89	ng	93
25) Isophorone	7.78	82	905186	43.48	ng	99
26) 2-Nitrophenol	7.86	139	240208	44.05	ng	# 76
27) 2,4-Dimethylphenol	7.89	122	380243	39.71	ng	98
28) bis(2-Chloroethoxy)methane	7.98	93	525437	38.11	ng	98
29) 2,4-Dichlorophenol	8.10	162	366079	40.99	ng	90
30) 1,2,4-Trichlorobenzene	8.18	180	392802	40.81	ng	95
31) Naphthalene	8.26	128	1220589	38.71	ng	99
32) Benzoic acid	8.02	122	225707	42.56	ng	99
33) 4-Chloroaniline	8.32	127	550655	44.52	ng	95
34) Hexachlorobutadiene	8.37	225	206781	39.05	ng	99
35) Caprolactam	8.69	113	129960	46.26	ng	# 51
36) 4-Chloro-3-methylphenol	8.81	107	392918	42.20	ng	# 83
37) 2-Methylnaphthalene	8.96	142	883707	43.65	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.12	216	388344	37.89	ng	99
40) Hexachlorocyclopentadiene	9.10	237	153737	33.25	ng	95

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42) 2,4,6-Trichlorophenol	9.24	196	268066	39.81	nq	94
43) 2,4,5-Trichlorophenol	9.28	196	292648	39.66	nq	96
45) 1,1'-Biphenyl	9.42	154	1095414	41.44	nq	100
46) 2-Chloronaphthalene	9.45	162	826644	39.97	nq	97
47) 2-Nitroaniline	9.55	65	281723	40.52	nq	# 70
48) Acenaphthylene	9.87	152	1444986	40.45	nq	98
49) Dimethylphthalate	9.72	163	902785	36.71	nq	98
50) 2,6-Dinitrotoluene	9.79	165	234100	44.08	nq	96
51) Acenaphthene	10.04	154	862456	40.38	nq	97
52) 3-Nitroaniline	9.96	138	261268	42.99	nq	92
53) 2,4-Dinitrophenol	10.06	184	102062	40.28	nq	# 88
54) Dibenzofuran	10.21	168	1110440	38.87	nq	90
55) 4-Nitrophenol	10.13	139	196238	44.83	nq	# 73
56) 2,4-Dinitrotoluene	10.19	165	253454	35.85	nq	98
57) Fluorene	10.56	166	921341	41.92	nq	98
58) 2,3,4,6-Tetrachlorophenol	10.33	232	195821	39.79	nq	97
59) Diethylphthalate	10.42	149	910663	39.30	nq	100
60) 4-Chlorophenyl-phenylether	10.54	204	416321	39.43	nq	# 86
61) 4-Nitroaniline	10.58	138	245169	39.39	nq	90
62) Azobenzene	10.70	77	1098999	45.90	nq	96
64) 4,6-Dinitro-2-methylphenol	10.60	198	135536	34.89	nq	# 61
65) n-Nitrosodiphenylamine	10.67	169	851709	40.86	nq	99
66) 4-Bromophenyl-phenylether	11.04	248	269630	39.55	nq	94
67) Hexachlorobenzene	11.10	284	273206	40.55	nq	97
68) Atrazine	11.20	200	288961	43.19	nq	96
69) Pentachlorophenol	11.30	266	120852	38.82	nq	99
70) Phenanthrene	11.52	178	1311106	35.07	nq	99
71) Anthracene	11.57	178	1537794	40.96	nq	99
72) Carbazole	11.73	167	1476119	41.13	nq	97
73) Di-n-butylphthalate	12.05	149	1589293	40.95	nq	# 97
74) Fluoranthene	12.70	202	1351270	35.04	nq	98
76) Benzidine	12.83	184	786806	40.01	nq	99
77) Pyrene	12.94	202	1499153	43.41	nq	98
79) Butylbenzylphthalate	13.55	149	619505	44.17	nq	100
80) Benzo(a)anthracene	14.12	228	1196539	42.39	nq	100
81) 3,3'-Dichlorobenzidine	14.09	252	413606	41.11	nq	98
82) Chrysene	14.17	228	1195469	45.23	nq	99
83) Bis(2-ethylhexyl)phthalate	14.11	149	872233	45.47	nq	# 97
84) Di-n-octyl phthalate	14.73	149	1395570	44.68	nq	100
85) Indeno(1,2,3-cd)pyrene	17.09	276	803575	39.25	nq	95
87) Benzo(b)fluoranthene	15.19	252	1033130	41.28	nq	97
88) Benzo(k)fluoranthene	15.21	252	886759m	41.03	nq	
89) Benzo(a)pyrene	15.55	252	865831	39.99	nq	99
90) Dibenzo(a,h)anthracene	17.12	278	691560	39.52	nq	98
91) Benzo(g,h,i)perylene	17.55	276	687272	38.88	nq	# 94

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

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