

Data Path : Z:\HPCHEM1\BNA F\DATA\BF030317\
 Data File : BF093386.D
 Acq On : 4 Mar 2017 5:44
 Operator : SJ/MA
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad
 3/6/2017 4:23:41 PM

Quant Time: Mar 05 23:52:20 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF013017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 17 17:31:55 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.78	152	168772	20.00	ng	-0.07
21) Naphthalene-d8	8.08	136	656330	20.00	ng	-0.07
38) Acenaphthene-d10	9.82	164	287060	20.00	ng	-0.07
63) Phenanthrene-d10	11.31	188	449463	20.00	ng	-0.07
75) Chrysene-d12	13.95	240	294577	20.00	ng	-0.07
86) Perylene-d12	15.36	264	282441	20.00	ng	-0.09

System Monitoring Compounds						
5) 2-Fluorophenol	5.37	112	830656	84.44	ng	-0.08
7) Phenol-d6	6.44	99	1022014	80.74	ng	-0.05
23) Nitrobenzene-d5	7.36	82	951400	80.72	ng	-0.07
41) 2,4,6-Tribromophenol	10.61	330	148605	71.58	ng	-0.07
44) 2-Fluorobiphenyl	9.15	172	1415155	89.31	ng	-0.07
78) Terphenyl-d14	12.89	244	1038655	82.85	ng	-0.08

Target Compounds						Qvalue
2) 1,4-Dioxane	2.29	88	218975	41.20	ng	# 85
3) Pyridine	2.99	79	622090	46.02	ng	88
4) n-Nitrosodimethylamine	2.96	42	286246	45.10	ng	86
6) Aniline	6.45	93	671632	38.90	ng	83
8) 2-Chlorophenol	6.57	128	437416	40.30	ng	88
9) Benzaldehyde	6.33	77	324124	35.74	ng	86
10) Phenol	6.45	94	597691	40.36	ng	95
11) bis(2-Chloroethyl)ether	6.52	93	474511	40.14	ng	91
12) 1,3-Dichlorobenzene	6.73	146	503709	39.63	ng	99
13) 1,4-Dichlorobenzene	6.81	146	501660m	38.99	ng	
14) 1,2-Dichlorobenzene	6.96	146	478337	38.88	ng	98
15) Benzyl Alcohol	6.94	79	349505	37.30	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.07	45	798823	38.90	ng	85
17) 2-Methylphenol	7.06	107	362406	38.92	ng	94
18) Hexachloroethane	7.30	117	170702	38.87	ng	# 85
19) n-Nitroso-di-n-propylamine	7.21	70	309697	38.44	ng	93
20) 3+4-Methylphenols	7.22	107	488923	43.92	ng	# 75
22) Acetophenone	7.21	105	619534	41.34	ng	# 97
24) Nitrobenzene	7.38	77	520595	43.02	ng	99
25) Isophorone	7.62	82	907467	41.32	ng	96
26) 2-Nitrophenol	7.70	139	244397	42.48	ng	90
27) 2,4-Dimethylphenol	7.74	122	416683	41.24	ng	94
28) bis(2-Chloroethoxy)methane	7.84	93	574096	39.47	ng	100
29) 2,4-Dichlorophenol	7.95	162	356618	37.85	ng	97
30) 1,2,4-Trichlorobenzene	8.02	180	385678	37.98	ng	97
31) Naphthalene	8.10	128	1293258	38.87	ng	99
32) Benzoic acid	7.88	122	198562	36.67	ng	96
33) 4-Chloroaniline	8.16	127	528381	40.49	ng	98
34) Hexachlorobutadiene	8.21	225	206811	37.02	ng	99
35) Caprolactam	8.53	113	116616	39.35	ng	# 34
36) 4-Chloro-3-methylphenol	8.65	107	388246	39.52	ng	97
37) 2-Methylnaphthalene	8.78	142	769839	36.04	ng	99
39) 1,2,4,5-Tetrachlorobenzene	8.96	216	368866	45.78	ng	99
40) Hexachlorocyclopentadiene	8.93	237	73427	20.20	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.07	196	226290	42.74	nq	93
43) 2,4,5-Trichlorophenol	9.12	196	246310	42.46	nq	96
45) 1,1'-Biphenyl	9.25	154	956129	46.00	nq	99
46) 2-Chloronaphthalene	9.28	162	758068	46.62	nq	98
47) 2-Nitroaniline	9.38	65	261025	47.74	nq	94
48) Acenaphthylene	9.69	152	1179713	42.00	nq	99
49) Dimethylphthalate	9.55	163	817573	42.28	nq	99
50) 2,6-Dinitrotoluene	9.62	165	174668	41.83	nq #	57
51) Acenaphthene	9.86	154	685640	40.83	nq	96
52) 3-Nitroaniline	9.79	138	204706	42.84	nq	96
53) 2,4-Dinitrophenol	9.92	184	33661	19.64	nq #	72
54) Dibenzofuran	10.03	168	891291	39.68	nq	95
55) 4-Nitrophenol	9.97	139	104421	30.34	nq #	85
56) 2,4-Dinitrotoluene	10.03	165	212788	38.28	nq #	86
57) Fluorene	10.37	166	774112	44.79	nq	98
58) 2,3,4,6-Tetrachlorophenol	10.16	232	166246	42.96	nq	95
59) Diethylphthalate	10.25	149	829961	45.55	nq	98
60) 4-Chlorophenyl-phenylether	10.36	204	346310	41.71	nq	89
61) 4-Nitroaniline	10.41	138	223929	45.75	nq	95
62) Azobenzene	10.52	77	907328	48.20	nq	96
64) 4,6-Dinitro-2-methylphenol	10.44	198	80265	30.38	nq #	51
65) n-Nitrosodiphenylamine	10.49	169	673364	46.90	nq	99
66) 4-Bromophenyl-phenylether	10.85	248	201088	42.82	nq #	86
67) Hexachlorobenzene	10.92	284	187505	40.41	nq	96
68) Atrazine	11.01	200	195899	42.52	nq	99
69) Pentachlorophenol	11.13	266	70100	33.83	nq	97
70) Phenanthrene	11.33	178	1034918	40.19	nq	99
71) Anthracene	11.38	178	997892	38.59	nq	99
72) Carbazole	11.54	167	967308	39.13	nq	99
73) Di-n-butylphthalate	11.87	149	1190532	44.53	nq #	97
74) Fluoranthene	12.52	202	897146	33.78	nq	93
76) Benzidine	12.65	184	755090	60.15	nq	98
77) Pyrene	12.75	202	889612	40.35	nq	98
79) Butylbenzylphthalate	13.36	149	380390	42.49	nq	96
80) Benzo(a)anthracene	13.94	228	693105	38.47	nq	99
81) 3,3'-Dichlorobenzidine	13.89	252	253642	39.49	nq	98
82) Chrysene	13.97	228	696973	41.32	nq	99
83) Bis(2-ethylhexyl)phthalate	13.92	149	534505	43.66	nq	99
84) Di-n-octyl phthalate	14.53	149	860031	43.14	nq	98
85) Indeno(1,2,3-cd)pyrene	16.74	276	637526	48.79	nq #	93
87) Benzo(b)fluoranthene	14.96	252	628889m	33.83	nq	
88) Benzo(k)fluoranthene	14.99	252	690756m	43.03	nq	
89) Benzo(a)pyrene	15.31	252	654201	40.68	nq	97
90) Dibenzo(a,h)anthracene	16.75	278	558263	42.95	nq	97
91) Benzo(g,h,i)perylene	17.15	276	545540	41.55	nq #	89

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

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