

Data Path : Z:\HPCHEM1\BNA F\DATA\BF030817\
 Data File : BF093430.D
 Acq On : 8 Mar 2017 15:53
 Operator : SJ/MA
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad
 3/10/2017 8:11:33 AM

Quant Time: Mar 08 17:02:39 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF030717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 07 15:55:22 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.77	152	197620	20.00	ng	0.00
21) Naphthalene-d8	8.06	136	793527	20.00	ng	0.00
38) Acenaphthene-d10	9.81	164	381205	20.00	ng	0.00
63) Phenanthrene-d10	11.29	188	700534	20.00	ng	0.01
75) Chrysene-d12	13.93	240	550057	20.00	ng	0.01
86) Perylene-d12	15.35	264	516867	20.00	ng	0.02

System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	981116	82.63	ng	0.00
7) Phenol-d6	6.41	99	1175963	78.54	ng	0.00
23) Nitrobenzene-d5	7.34	82	1209549	84.57	ng	0.00
41) 2,4,6-Tribromophenol	10.60	330	212959	85.75	ng	0.00
44) 2-Fluorobiphenyl	9.13	172	1715078	83.68	ng	0.00
78) Terphenyl-d14	12.87	244	1754545	82.93	ng	0.00

Target Compounds						Qvalue
2) 1,4-Dioxane	2.28	88	271622	41.53	ng	# 85
3) Pyridine	2.97	79	731391	43.86	ng	91
4) n-Nitrosodimethylamine	2.94	42	331668	41.64	ng	88
6) Aniline	6.42	93	786374	38.27	ng	# 66
8) 2-Chlorophenol	6.55	128	511429	38.44	ng	93
9) Benzaldehyde	6.31	77	385837	43.86	ng	88
10) Phenol	6.44	94	696347	37.95	ng	92
11) bis(2-Chloroethyl)ether	6.50	93	561043	37.83	ng	92
12) 1,3-Dichlorobenzene	6.70	146	580175m	38.10	ng	
13) 1,4-Dichlorobenzene	6.78	146	597106	38.30	ng	99
14) 1,2-Dichlorobenzene	6.94	146	560351	40.59	ng	95
15) Benzyl Alcohol	6.93	79	421649	39.50	ng	96
16) 2,2'-oxybis(1-Chloropropan	7.05	45	994406	40.37	ng	83
17) 2-Methylphenol	7.04	107	433781	38.55	ng	# 92
18) Hexachloroethane	7.27	117	208208	39.60	ng	# 88
19) n-Nitroso-di-n-propylamine	7.19	70	382634	37.17	ng	92
20) 3+4-Methylphenols	7.20	107	599164	41.06	ng	# 76
22) Acetophenone	7.19	105	766824	40.16	ng	# 97
24) Nitrobenzene	7.36	77	633842	39.43	ng	100
25) Isophorone	7.60	82	1133656	39.31	ng	94
26) 2-Nitrophenol	7.68	139	306987	41.86	ng	91
27) 2,4-Dimethylphenol	7.73	122	493703	41.38	ng	97
28) bis(2-Chloroethoxy)methane	7.82	93	730762	40.14	ng	99
29) 2,4-Dichlorophenol	7.93	162	437236	38.96	ng	96
30) 1,2,4-Trichlorobenzene	8.00	180	478574	40.10	ng	96
31) Naphthalene	8.08	128	1582339	39.79	ng	99
32) Benzoic acid	7.89	122	209217	33.75	ng	97
33) 4-Chloroaniline	8.14	127	658469	40.80	ng	100
34) Hexachlorobutadiene	8.20	225	243957	39.68	ng	97
35) Caprolactam	8.53	113	150379	39.23	ng	# 29
36) 4-Chloro-3-methylphenol	8.63	107	468525	39.11	ng	96
37) 2-Methylnaphthalene	8.77	142	967952	38.26	ng	98
39) 1,2,4,5-Tetrachlorobenzene	8.94	216	470454	45.20	ng	100
40) Hexachlorocyclopentadiene	8.92	237	88560	40.71	ng	96

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.05	196	287543	44.21	nq	91
43) 2,4,5-Trichlorophenol	9.11	196	310559	44.50	nq #	87
45) 1,1'-Biphenyl	9.24	154	1260810	45.40	nq	99
46) 2-Chloronaphthalene	9.26	162	986646	46.42	nq	99
47) 2-Nitroaniline	9.36	65	335996	44.71	nq	91
48) Acenaphthylene	9.67	152	1464074	41.76	nq	99
49) Dimethylphthalate	9.53	163	1081647	42.80	nq	98
50) 2,6-Dinitrotoluene	9.61	165	261632	47.18	nq #	84
51) Acenaphthene	9.84	154	887936	42.83	nq	97
52) 3-Nitroaniline	9.78	138	309731	46.49	nq #	79
53) 2,4-Dinitrophenol	9.90	184	136733	54.14	nq #	51
54) Dibenzofuran	10.01	168	1169740	45.08	nq	96
55) 4-Nitrophenol	9.97	139	133442	41.24	nq	85
56) 2,4-Dinitrotoluene	10.01	165	282185	41.42	nq #	58
57) Fluorene	10.36	166	996280	45.31	nq	98
58) 2,3,4,6-Tetrachlorophenol	10.14	232	224955	45.68	nq	96
59) Diethylphthalate	10.23	149	978524	40.51	nq	99
60) 4-Chlorophenyl-phenylether	10.34	204	438595	44.50	nq	91
61) 4-Nitroaniline	10.40	138	317456	46.59	nq	82
62) Azobenzene	10.50	77	1221282	44.50	nq	97
64) 4,6-Dinitro-2-methylphenol	10.44	198	176065	38.79	nq	89
65) n-Nitrosodiphenylamine	10.47	169	873627	37.35	nq	99
66) 4-Bromophenyl-phenylether	10.84	248	281584	38.01	nq #	88
67) Hexachlorobenzene	10.90	284	259650	36.71	nq #	88
68) Atrazine	11.00	200	264287	38.60	nq	99
69) Pentachlorophenol	11.11	266	101007	41.36	nq	98
70) Phenanthrene	11.32	178	1411380	35.30	nq	100
71) Anthracene	11.37	178	1587952	39.26	nq	98
72) Carbazole	11.53	167	1559529	41.04	nq	99
73) Di-n-butylphthalate	11.85	149	1694948	38.56	nq #	98
74) Fluoranthene	12.50	202	1424977	36.89	nq	99
76) Benzidine	12.63	184	785023	43.40	nq	99
77) Pyrene	12.73	202	1476172	36.90	nq	99
79) Butylbenzylphthalate	13.35	149	744748	44.22	nq	87
80) Benzo(a)anthracene	13.92	228	1246282	38.37	nq	99
81) 3,3'-Dichlorobenzidine	13.89	252	451684	39.70	nq #	98
82) Chrysene	13.97	228	1253233	41.70	nq	98
83) Bis(2-ethylhexyl)phthalate	13.90	149	957819	40.81	nq #	99
84) Di-n-octyl phthalate	14.52	149	1586085	40.54	nq	98
85) Indeno(1,2,3-cd)pyrene	16.73	276	1096704	43.59	nq #	94
87) Benzo(b)fluoranthene	14.95	252	1230191m	39.08	nq	
88) Benzo(k)fluoranthene	14.97	252	1100699m	36.26	nq	
89) Benzo(a)pyrene	15.29	252	1125416	39.12	nq	99
90) Dibenzo(a,h)anthracene	16.73	278	933476	39.22	nq	98
91) Benzo(g,h,i)perylene	17.13	276	960494	39.32	nq #	91

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

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