

Data Path : Z:\HPCHEM1\BNA F\DATA\BF012617\
 Data File : BF092532.D
 Acq On : 26 Jan 2017 18:36
 Operator : SJ/MA
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad
 1/27/2017 3:03:01 PM

Quant Time: Jan 27 01:14:45 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 24 13:48:07 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.98	152	184515	20.00	ng	0.01
21) Naphthalene-d8	8.27	136	753680	20.00	ng	0.00
38) Acenaphthene-d10	10.04	164	413682	20.00	ng	0.01
63) Phenanthrene-d10	11.53	188	678857	20.00	ng	0.00
75) Chrysene-d12	14.17	240	501178	20.00	ng	0.00
86) Perylene-d12	15.65	264	430158	20.00	ng	0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.56	112	906585	76.44	ng	0.01
7) Phenol-d6	6.60	99	1146490	75.89	ng	0.01
23) Nitrobenzene-d5	7.55	82	1150374	83.38	ng	0.01
41) 2,4,6-Tribromophenol	10.83	330	239173	84.84	ng	0.01
44) 2-Fluorobiphenyl	9.36	172	1693633	75.71	ng	0.00
78) Terphenyl-d14	13.12	244	1520512	80.79	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.64	88	231850	36.76	ng	# 84
3) Pyridine	3.39	79	655029	36.48	ng	# 85
4) n-Nitrosodimethylamine	3.34	42	309162	35.10	ng	# 83
6) Aniline	6.64	93	878407	38.98	ng	# 45
8) 2-Chlorophenol	6.76	128	513809	41.25	ng	# 93
9) Benzaldehyde	6.52	77	400995	37.41	ng	# 96
10) Phenol	6.61	94	702042	38.50	ng	# 73
11) bis(2-Chloroethyl)ether	6.70	93	481614	35.30	ng	# 88
12) 1,3-Dichlorobenzene	6.92	146	568806	38.61	ng	# 99
13) 1,4-Dichlorobenzene	6.99	146	561424	37.86	ng	# 98
14) 1,2-Dichlorobenzene	7.15	146	557765	39.65	ng	# 97
15) Benzyl Alcohol	7.12	79	436036	38.30	ng	# 97
16) 2,2'-oxybis(1-Chloropropan	7.25	45	1004477	37.22	ng	# 92
17) 2-Methylphenol	7.23	107	421749	39.62	ng	# 95
18) Hexachloroethane	7.49	117	197357	38.64	ng	# 97
19) n-Nitroso-di-n-propylamine	7.39	70	374892	36.44	ng	# 90
20) 3+4-Methylphenols	7.38	107	498961	38.64	ng	# 93
22) Acetophenone	7.39	105	665866	37.09	ng	# 95
24) Nitrobenzene	7.56	77	506257	35.02	ng	# 100
25) Isophorone	7.80	82	1020616	37.54	ng	# 94
26) 2-Nitrophenol	7.88	139	270016	44.58	ng	# 97
27) 2,4-Dimethylphenol	7.92	122	442817	35.15	ng	# 98
28) bis(2-Chloroethoxy)methane	8.02	93	702819	39.38	ng	# 100
29) 2,4-Dichlorophenol	8.12	162	417511	38.23	ng	# 90
30) 1,2,4-Trichlorobenzene	8.21	180	473235	40.09	ng	# 96
31) Naphthalene	8.29	128	1433231	37.15	ng	# 99
32) Benzoic acid	8.03	122	209658	24.78	ng	# 96
33) 4-Chloroaniline	8.34	127	618524	37.90	ng	# 97
34) Hexachlorobutadiene	8.41	225	240631	37.27	ng	# 100
35) Caprolactam	8.73	113	134928	37.00	ng	# 27
36) 4-Chloro-3-methylphenol	8.82	107	455628	38.69	ng	# 94
37) 2-Methylnaphthalene	8.99	142	955404	39.12	ng	# 98
39) 1,2,4,5-Tetrachlorobenzene	9.16	216	434167	41.61	ng	# 99
40) Hexachlorocyclopentadiene	9.14	237	202804	38.81	ng	# 94

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.26	196	314544	39.34	nq	91
43) 2,4,5-Trichlorophenol	9.31	196	318046	43.48	nq #	84
45) 1,1'-Biphenyl	9.46	154	1198173	40.07	nq	99
46) 2-Chloronaphthalene	9.48	162	919671	37.65	nq	97
47) 2-Nitroaniline	9.57	65	300670	36.56	nq	97
48) Acenaphthylene	9.90	152	1437361	40.12	nq	97
49) Dimethylphthalate	9.76	163	966063	36.72	nq	98
50) 2,6-Dinitrotoluene	9.82	165	234601	43.63	nq #	72
51) Acenaphthene	10.08	154	921225	41.38	nq	98
52) 3-Nitroaniline	9.98	138	266434	39.56	nq	98
53) 2,4-Dinitrophenol	10.09	184	115667	45.17	nq #	43
54) Dibenzofuran	10.25	168	1208612	39.99	nq	91
55) 4-Nitrophenol	10.14	139	245002	45.81	nq #	79
56) 2,4-Dinitrotoluene	10.22	165	339066	47.51	nq #	85
57) Fluorene	10.59	166	964229	43.02	nq	98
58) 2,3,4,6-Tetrachlorophenol	10.36	232	236300	43.11	nq	96
59) Diethylphthalate	10.45	149	953066	37.78	nq	99
60) 4-Chlorophenyl-phenylether	10.58	204	446136	41.04	nq #	85
61) 4-Nitroaniline	10.60	138	269652	40.04	nq	94
62) Azobenzene	10.74	77	1101463	38.37	nq	94
64) 4,6-Dinitro-2-methylphenol	10.64	198	185705	50.48	nq	83
65) n-Nitrosodiphenylamine	10.69	169	810191	38.22	nq	100
66) 4-Bromophenyl-phenylether	11.07	248	275307	40.17	nq	92
67) Hexachlorobenzene	11.14	284	281761	40.67	nq	99
68) Atrazine	11.23	200	315363	43.46	nq	95
69) Pentachlorophenol	11.33	266	181599	41.00	nq	98
70) Phenanthrene	11.55	178	1312568	35.11	nq	99
71) Anthracene	11.61	178	1510716	41.35	nq	98
72) Carbazole	11.76	167	1274771	36.54	nq	99
73) Di-n-butylphthalate	12.09	149	1583982	39.71	nq #	98
74) Fluoranthene	12.74	202	1356915	37.11	nq	98
76) Benzidine	12.86	184	837576	45.02	nq	99
77) Pyrene	12.97	202	1456170	40.10	nq	100
79) Butylbenzylphthalate	13.59	149	639368	43.46	nq	97
80) Benzo(a)anthracene	14.16	228	1121646	41.64	nq	99
81) 3,3'-Dichlorobenzidine	14.12	252	437967	42.83	nq #	96
82) Chrysene	14.20	228	1200095	41.72	nq	99
83) Bis(2-ethylhexyl)phthalate	14.15	149	833541	44.99	nq #	98
84) Di-n-octyl phthalate	14.76	149	1494400	44.43	nq	98
85) Indeno(1,2,3-cd)pyrene	17.15	276	898280	42.21	nq	95
87) Benzo(b)fluoranthene	15.22	252	1000756	36.24	nq	99
88) Benzo(k)fluoranthene	15.25	252	1001640m	43.62	nq	
89) Benzo(a)pyrene	15.60	252	940916	40.28	nq	97
90) Dibenzo(a,h)anthracene	17.17	278	774802	41.01	nq	99
91) Benzo(g,h,i)perylene	17.61	276	780678	40.86	nq #	92

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

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