

Data Path : Z:\HPCHEM1\BNA F\DATA\BF110716\
 Data File : BF091050.D
 Acq On : 7 Nov 2016 12:43
 Operator : UM/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

sohil
 11/8/2016 4:30:54 PM

Quant Time: Nov 07 17:42:21 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 04 11:03:12 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.67	152	174379	20.00	ng	-0.03
21) Naphthalene-d8	7.96	136	747844	20.00	ng	-0.03
38) Acenaphthene-d10	9.71	164	347461	20.00	ng	-0.03
63) Phenanthrene-d10	11.20	188	612090	20.00	ng	-0.03
75) Chrysene-d12	13.82	240	509098	20.00	ng	-0.03
86) Perylene-d12	15.21	264	391363	20.00	ng	-0.03

System Monitoring Compounds						
5) 2-Fluorophenol	5.25	112	888744	84.17	ng	-0.03
7) Phenol-d6	6.33	99	1092009	76.20	ng	-0.02
23) Nitrobenzene-d5	7.24	82	1096167	79.96	ng	-0.03
41) 2,4,6-Tribromophenol	10.51	330	313085	99.67	ng	-0.02
44) 2-Fluorobiphenyl	9.05	172	1849083	91.62	ng	-0.02
78) Terphenyl-d14	12.78	244	1873527	91.83	ng	-0.02

Target Compounds						Qvalue
2) 1,4-Dioxane	2.14	88	229043	37.92	ng	97
3) Pyridine	2.80	79	662617	46.90	ng	98
4) n-Nitrosodimethylamine	2.76	42	223421	39.98	ng	91
6) Aniline	6.34	93	675061	39.93	ng	# 69
8) 2-Chlorophenol	6.45	128	499673	39.67	ng	98
9) Benzaldehyde	6.21	77	301547	37.91	ng	99
10) Phenol	6.34	94	641531	40.45	ng	# 80
11) bis(2-Chloroethyl)ether	6.42	93	552287	41.32	ng	90
12) 1,3-Dichlorobenzene	6.61	146	534793	41.98	ng	100
13) 1,4-Dichlorobenzene	6.69	146	531662	38.90	ng	97
14) 1,2-Dichlorobenzene	6.84	146	522389	41.65	ng	98
15) Benzyl Alcohol	6.83	79	418746	41.42	ng	91
16) 2,2'-oxybis(1-Chloropropan	6.96	45	673223	35.18	ng	73
17) 2-Methylphenol	6.94	107	380065	38.12	ng	99
18) Hexachloroethane	7.18	117	211585	40.00	ng	98
19) n-Nitroso-di-n-propylamine	7.10	70	406811	40.96	ng	# 89
20) 3+4-Methylphenols	7.10	107	508340	42.46	ng	95
22) Acetophenone	7.09	105	710448	41.26	ng	# 96
24) Nitrobenzene	7.26	77	638020	42.14	ng	97
25) Isophorone	7.50	82	1006871	38.12	ng	98
26) 2-Nitrophenol	7.58	139	282755	44.13	ng	# 71
27) 2,4-Dimethylphenol	7.63	122	432432	40.81	ng	98
28) bis(2-Chloroethoxy)methane	7.72	93	595822	41.29	ng	100
29) 2,4-Dichlorophenol	7.82	162	372127	40.24	ng	95
30) 1,2,4-Trichlorobenzene	7.90	180	434280	41.76	ng	99
31) Naphthalene	7.98	128	1493673	41.76	ng	100
32) Benzoic acid	7.78	122	312134	38.84	ng	100
33) 4-Chloroaniline	8.04	127	620836	41.74	ng	98
34) Hexachlorobutadiene	8.10	225	217084	38.68	ng	100
35) Caprolactam	8.43	113	131565	45.28	ng	# 66
36) 4-Chloro-3-methylphenol	8.53	107	427662	38.20	ng	89
37) 2-Methylnaphthalene	8.67	142	875025	38.06	ng	99
39) 1,2,4,5-Tetrachlorobenzene	8.84	216	407547	46.00	ng	99
40) Hexachlorocyclopentadiene	8.83	237	126641	40.56	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	8.96	196	248673	43.48	nq	99
43) 2,4,5-Trichlorophenol	9.00	196	270171	45.00	nq #	90
45) 1,1'-Biphenyl	9.14	154	1052542	41.49	nq	99
46) 2-Chloronaphthalene	9.16	162	812382	42.17	nq	99
47) 2-Nitroaniline	9.26	65	272429	38.11	nq	97
48) Acenaphthylene	9.57	152	1167610	38.90	nq	100
49) Dimethylphthalate	9.46	163	939342	41.23	nq	97
50) 2,6-Dinitrotoluene	9.52	165	234895	48.11	nq #	58
51) Acenaphthene	9.74	154	715411	37.96	nq	99
52) 3-Nitroaniline	9.69	138	246503	42.57	nq #	71
53) 2,4-Dinitrophenol	9.79	184	110196	44.45	nq #	90
54) Dibenzofuran	9.92	168	1037389	40.89	nq	98
55) 4-Nitrophenol	9.86	139	141007	39.68	nq #	77
56) 2,4-Dinitrotoluene	9.92	165	252938	40.71	nq #	86
57) Fluorene	10.26	166	803309	38.74	nq	100
58) 2,3,4,6-Tetrachlorophenol	10.04	232	206867	43.98	nq	93
59) Diethylphthalate	10.14	149	885172	37.89	nq	98
60) 4-Chlorophenyl-phenylether	10.26	204	431848	45.76	nq	95
61) 4-Nitroaniline	10.30	138	249273	42.56	nq	84
62) Azobenzene	10.42	77	954994	34.73	nq	98
64) 4,6-Dinitro-2-methylphenol	10.33	198	157185	41.94	nq	73
65) n-Nitrosodiphenylamine	10.38	169	790942	40.65	nq	99
66) 4-Bromophenyl-phenylether	10.75	248	253773	39.86	nq #	63
67) Hexachlorobenzene	10.81	284	288896	41.14	nq	99
68) Atrazine	10.91	200	192191	34.24	nq	97
69) Pentachlorophenol	11.01	266	134646	43.29	nq	98
70) Phenanthrene	11.22	178	1268467	38.98	nq	100
71) Anthracene	11.28	178	1405404	40.63	nq	99
72) Carbazole	11.44	167	1249023	40.07	nq #	96
73) Di-n-butylphthalate	11.77	149	1444391	36.84	nq #	98
74) Fluoranthene	12.41	202	1405412	41.65	nq	88
76) Benzidine	12.53	184	554421	48.75	nq	99
77) Pyrene	12.62	202	1361812	40.85	nq	99
79) Butylbenzylphthalate	13.25	149	653728	40.99	nq	89
80) Benzo(a)anthracene	13.81	228	1242948	43.00	nq	99
81) 3,3'-Dichlorobenzidine	13.79	252	485750	47.83	nq #	99
82) Chrysene	13.86	228	1355296	47.55	nq	99
83) Bis(2-ethylhexyl)phthalate	13.82	149	852033	39.48	nq #	96
84) Di-n-octyl phthalate	14.43	149	1350860	38.07	nq	97
85) Indeno(1,2,3-cd)pyrene	16.51	276	922023	38.99	nq	99
87) Benzo(b)fluoranthene	14.83	252	1258977m	47.43	nq	
88) Benzo(k)fluoranthene	14.85	252	898296m	38.58	nq	
89) Benzo(a)pyrene	15.15	252	949827	41.10	nq #	97
90) Dibenzo(a,h)anthracene	16.52	278	805111	40.30	nq #	96
91) Benzo(g,h,i)perylene	16.89	276	729968	40.41	nq	100

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

