

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021116\
 Data File : BF085010.D
 Acq On : 11 Feb 2016 11:09
 Operator : SJ/IZ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 2/12/2016 12:25:57 PM

Quant Time: Feb 11 16:13:02 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 10 13:28:51 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.62	152	199909	20.00	ng	-0.01
21) Naphthalene-d8	7.91	136	827042	20.00	ng	-0.01
38) Acenaphthene-d10	9.66	164	386656	20.00	ng	-0.01
63) Phenanthrene-d10	11.13	188	710219	20.00	ng	-0.01
75) Chrysene-d12	13.76	240	438391	20.00	ng	-0.01
86) Perylene-d12	15.12	264	394496	20.00	ng	-0.01

System Monitoring Compounds						
5) 2-Fluorophenol	5.19	112	940364	73.27	ng	-0.01
7) Phenol-d6	6.27	99	1179288	74.12	ng	-0.01
23) Nitrobenzene-d5	7.19	82	1150791	78.38	ng	-0.01
41) 2,4,6-Tribromophenol	10.44	330	328589	81.47	ng	-0.01
44) 2-Fluorobiphenyl	8.98	172	1712842	70.48	ng	-0.01
78) Terphenyl-d14	12.72	244	1686344	87.17	ng	-0.01

Target Compounds						Qvalue
2) 1,4-Dioxane	2.04	88	206387	36.71	ng	# 74
3) Pyridine	2.68	79	579115	39.54	ng	# 71
4) n-Nitrosodimethylamine	2.65	42	303066	40.00	ng	81
6) Aniline	6.27	93	786451	40.39	ng	98
8) 2-Chlorophenol	6.40	128	543170	37.39	ng	91
9) Benzaldehyde	6.16	77	334431	35.59	ng	90
10) Phenol	6.28	94	635001	35.63	ng	# 62
11) bis(2-Chloroethyl)ether	6.36	93	569218	40.95	ng	92
12) 1,3-Dichlorobenzene	6.55	146	554262m	36.11	ng	
13) 1,4-Dichlorobenzene	6.63	146	561840	36.62	ng	97
14) 1,2-Dichlorobenzene	6.79	146	567497	39.71	ng	96
15) Benzyl Alcohol	6.78	79	453478	41.27	ng	92
16) 2,2'-oxybis(1-Chloropropan	6.90	45	894244	38.25	ng	# 44
17) 2-Methylphenol	6.90	107	451503	39.30	ng	# 86
18) Hexachloroethane	7.12	117	229007	38.76	ng	# 89
19) n-Nitroso-di-n-propylamine	7.05	70	368218	36.92	ng	# 75
20) 3+4-Methylphenols	7.06	107	593354	41.61	ng	88
22) Acetophenone	7.04	105	792719	36.37	ng	# 99
24) Nitrobenzene	7.21	77	614666	39.10	ng	98
25) Isophorone	7.46	82	1090362	38.19	ng	96
26) 2-Nitrophenol	7.53	139	326879	42.16	ng	92
27) 2,4-Dimethylphenol	7.58	122	493995	36.63	ng	96
28) bis(2-Chloroethoxy)methane	7.67	93	592166	34.96	ng	98
29) 2,4-Dichlorophenol	7.77	162	421194	36.50	ng	94
30) 1,2,4-Trichlorobenzene	7.85	180	506838	38.89	ng	96
31) Naphthalene	7.93	128	1582429	38.14	ng	100
32) Benzoic acid	7.75	122	392684	45.52	ng	98
33) 4-Chloroaniline	7.99	127	707656	41.25	ng	95
34) Hexachlorobutadiene	8.04	225	282722	37.62	ng	99
35) Caprolactam	8.39	113	149193	41.95	ng	96
36) 4-Chloro-3-methylphenol	8.48	107	481366	36.56	ng	93
37) 2-Methylnaphthalene	8.62	142	1019826	39.28	ng	98
39) 1,2,4,5-Tetrachlorobenzene	8.79	216	478840	38.99	ng	98
40) Hexachlorocyclopentadiene	8.76	237	180943	41.80	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	8.90	196	311126	39.33	ng	97
43) 2,4,5-Trichlorophenol	8.95	196	323010	40.18	ng #	92
45) 1,1'-Biphenyl	9.08	154	1428509	41.36	ng	98
46) 2-Chloronaphthalene	9.11	162	1005147	39.52	ng #	87
47) 2-Nitroaniline	9.21	65	358483	44.51	ng	88
48) Acenaphthylene	9.52	152	1509640	39.11	ng	99
49) Dimethylphthalate	9.39	163	1096354	37.23	ng #	90
50) 2,6-Dinitrotoluene	9.46	165	253004	39.33	ng #	84
51) Acenaphthene	9.69	154	996624	39.69	ng	97
52) 3-Nitroaniline	9.62	138	253010	35.00	ng	93
53) 2,4-Dinitrophenol	9.74	184	113626	41.29	ng	87
54) Dibenzofuran	9.86	168	1217421	39.67	ng	100
55) 4-Nitrophenol	9.80	139	184224	34.67	ng #	83
56) 2,4-Dinitrotoluene	9.86	165	340978	41.55	ng	96
57) Fluorene	10.20	166	1063882	41.51	ng	99
58) 2,3,4,6-Tetrachlorophenol	9.99	232	257639	42.10	ng #	86
59) Diethylphthalate	10.09	149	1014383	35.27	ng	99
60) 4-Chlorophenyl-phenylether	10.19	204	406869	36.40	ng	89
61) 4-Nitroaniline	10.24	138	262568	37.27	ng	92
62) Azobenzene	10.35	77	1041460	37.66	ng	89
64) 4,6-Dinitro-2-methylphenol	10.27	198	196283	44.78	ng	85
65) n-Nitrosodiphenylamine	10.32	169	816369	36.20	ng	99
66) 4-Bromophenyl-phenylether	10.68	248	315579	40.23	ng #	89
67) Hexachlorobenzene	10.74	284	308662	36.11	ng #	79
68) Atrazine	10.86	200	269374	37.82	ng	98
69) Pentachlorophenol	10.95	266	169083	42.09	ng	98
70) Phenanthrene	11.15	178	1368769	34.75	ng	97
71) Anthracene	11.21	178	1625813	40.96	ng	99
72) Carbazole	11.37	167	1379250	39.85	ng	99
73) Di-n-butylphthalate	11.71	149	1702041	38.63	ng	99
74) Fluoranthene	12.34	202	1582921	39.70	ng	93
76) Benzidine	12.47	184	615114	39.89	ng	99
77) Pyrene	12.56	202	1450137	43.20	ng	99
79) Butylbenzylphthalate	13.20	149	697300	45.80	ng	97
80) Benzo(a)anthracene	13.75	228	1120409	41.18	ng	99
81) 3,3'-Dichlorobenzidine	13.73	252	434028	45.05	ng #	97
82) Chrysene	13.79	228	1109998	42.07	ng	99
83) Bis(2-ethylhexyl)phthalate	13.76	149	736250	38.86	ng	99
84) Di-n-octyl phthalate	14.38	149	1192871	38.16	ng #	90
85) Indeno(1,2,3-cd)pyrene	16.38	276	1015968	48.92	ng	99
87) Benzo(b)fluoranthene	14.75	252	980437m	36.99	ng	
88) Benzo(k)fluoranthene	14.78	252	749012m	34.18	ng	
89) Benzo(a)pyrene	15.06	252	830128	36.76	ng #	95
90) Dibenzo(a,h)anthracene	16.39	278	837004	44.03	ng #	95
91) Benzo(g,h,i)perylene	16.75	276	892134	46.58	ng #	95

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

