

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF090415\
 Data File : BF081436.D
 Acq On : 4 Sep 2015 23:58
 Operator : UM/IZ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

MMDadoda
 9/8/2015 1:57:18 PM

Quant Time: Sep 07 01:54:45 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 01 17:35:35 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.36	152	71506	20.00	ng	-0.02
21) Naphthalene-d8	7.66	136	292467	20.00	ng	-0.02
38) Acenaphthene-d10	9.39	164	141248	20.00	ng	-0.02
63) Phenanthrene-d10	10.86	188	269901	20.00	ng	-0.02
75) Chrysene-d12	13.46	240	206534	20.00	ng	-0.03
86) Perylene-d12	14.77	264	149024	20.00	ng	-0.03

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	4.90	112	389343	84.48	ng	-0.02
7) Phenol-d6	6.03	99	494851	81.12	ng	-0.02
23) Nitrobenzene-d5	6.95	82	505247	83.35	ng	-0.02
41) 2,4,6-Tribromophenol	10.18	330	95210	75.70	ng	-0.02
44) 2-Fluorobiphenyl	8.73	172	747427	67.81	ng	-0.03
78) Terphenyl-d14	12.44	244	761295	85.81	ng	-0.02

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	1.67	88	87096	42.08	ng	95
3) Pyridine	2.22	79	217710	38.14	ng	# 89
4) n-Nitrosodimethylamine	2.18	42	141731	44.70	ng	# 75
6) Aniline	6.02	93	353204	41.00	ng	# 79
8) 2-Chlorophenol	6.15	128	213831	42.10	ng	88
9) Benzaldehyde	5.90	77	157487	41.20	ng	# 75
10) Phenol	6.04	94	286600	40.51	ng	# 53
11) bis(2-Chloroethyl)ether	6.11	93	204737	40.11	ng	87
12) 1,3-Dichlorobenzene	6.30	146	223632	41.21	ng	# 90
13) 1,4-Dichlorobenzene	6.38	146	222102	40.20	ng	# 88
14) 1,2-Dichlorobenzene	6.54	146	193758	38.95	ng	97
15) Benzyl Alcohol	6.52	79	203376	40.64	ng	# 69
16) 2,2'-oxybis(1-Chloropropan	6.67	45	431210	44.07	ng	84
17) 2-Methylphenol	6.66	107	174432	43.04	ng	# 75
18) Hexachloroethane	6.87	117	85816	39.92	ng	91
19) n-Nitroso-di-n-propylamine	6.81	70	184910	44.54	ng	# 85
20) 3+4-Methylphenols	6.82	107	227657	41.67	ng	90
22) Acetophenone	6.80	105	308225	38.58	ng	# 84
24) Nitrobenzene	6.96	77	226736	36.02	ng	# 61
25) Isophorone	7.21	82	456269	40.47	ng	# 85
26) 2-Nitrophenol	7.28	139	113064	41.21	ng	# 42
27) 2,4-Dimethylphenol	7.35	122	195691	41.29	ng	# 82
28) bis(2-Chloroethoxy)methane	7.44	93	256332	39.36	ng	# 93
29) 2,4-Dichlorophenol	7.53	162	169876	41.34	ng	94
30) 1,2,4-Trichlorobenzene	7.60	180	175510	39.31	ng	97
31) Naphthalene	7.68	128	599949	38.46	ng	99
32) Benzoic acid	7.51	122	113801	34.40	ng	# 54
33) 4-Chloroaniline	7.74	127	228758	35.96	ng	# 78
34) Hexachlorobutadiene	7.80	225	106603	39.24	ng	99
35) Caprolactam	8.14	113	58026	46.04	ng	# 61
36) 4-Chloro-3-methylphenol	8.25	107	196649	42.75	ng	86
37) 2-Methylnaphthalene	8.36	142	408224	40.22	ng	# 89
39) 1,2,4,5-Tetrachlorobenzene	8.54	216	159120	35.62	ng	99
40) Hexachlorocyclopentadiene	8.51	237	71448	32.59	ng	96

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42) 2,4,6-Trichlorophenol	8.65	196	114359	37.10	ng	98
43) 2,4,5-Trichlorophenol	8.70	196	121712	38.70	ng #	87
45) 1,1'-Biphenyl	8.83	154	503410	38.74	ng	95
46) 2-Chloronaphthalene	8.84	162	329900	35.15	ng #	88
47) 2-Nitroaniline	8.96	65	174184	45.54	ng #	65
48) Acenaphthylene	9.26	152	637171	38.27	ng	97
49) Dimethylphthalate	9.15	163	439979	40.09	ng #	97
50) 2,6-Dinitrotoluene	9.21	165	92465	37.12	ng #	94
51) Acenaphthene	9.43	154	366690	38.72	ng	89
52) 3-Nitroaniline	9.37	138	110886	39.11	ng #	47
53) 2,4-Dinitrophenol	9.47	184	40743	38.84	ng #	1
54) Dibenzofuran	9.60	168	534371	38.64	ng #	90
55) 4-Nitrophenol	9.56	139	69877m	39.23	ng	
56) 2,4-Dinitrotoluene	9.60	165	115725	36.49	ng #	7
57) Fluorene	9.93	166	366906	34.96	ng	95
58) 2,3,4,6-Tetrachlorophenol	9.72	232	95943	39.96	ng #	87
59) Diethylphthalate	9.84	149	417092	36.13	ng	94
60) 4-Chlorophenyl-phenylether	9.94	204	196963	40.27	ng	99
61) 4-Nitroaniline	9.98	138	101745	35.52	ng #	25
62) Azobenzene	10.09	77	472657	36.83	ng #	79
64) 4,6-Dinitro-2-methylphenol	10.01	198	67236	44.54	ng	98
65) n-Nitrosodiphenylamine	10.07	169	373470	38.03	ng	89
66) 4-Bromophenyl-phenylether	10.42	248	109891	40.27	ng	99
67) Hexachlorobenzene	10.48	284	112655	39.48	ng #	78
68) Atrazine	10.60	200	120117	42.26	ng	82
69) Pentachlorophenol	10.67	266	57539	45.64	ng	98
70) Phenanthrene	10.88	178	541387	36.08	ng	96
71) Anthracene	10.94	178	608992	39.50	ng	98
72) Carbazole	11.10	167	556788	38.49	ng	95
73) Di-n-butylphthalate	11.45	149	734727	43.01	ng #	98
74) Fluoranthene	12.06	202	658045	40.51	ng	86
76) Benzidine	12.19	184	291650	32.89	ng	98
77) Pyrene	12.27	202	615228	40.21	ng	98
79) Butylbenzylphthalate	12.92	149	312754	47.01	ng #	85
80) Benzo(a)anthracene	13.45	228	507727	38.60	ng	97
81) 3,3'-Dichlorobenzidine	13.44	252	169487	38.20	ng #	92
82) Chrysene	13.50	228	472235	40.05	ng	95
83) Bis(2-ethylhexyl)phthalate	13.50	149	373177	42.50	ng #	94
84) Di-n-octyl phthalate	14.10	149	585541	40.50	ng	97
85) Indeno(1,2,3-cd)pyrene	15.84	276	371788	42.98	ng #	90
87) Benzo(b)fluoranthene	14.45	252	449173m	42.22	ng	
88) Benzo(k)fluoranthene	14.47	252	337851m	38.27	ng	
89) Benzo(a)pyrene	14.72	252	340662	37.79	ng #	90
90) Dibenzo(a,h)anthracene	15.86	278	308503	44.28	ng #	77
91) Benzo(g,h,i)perylene	16.16	276	292183	43.94	ng #	79

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

