

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030916\
 Data File : BF085327.D
 Acq On : 10 Mar 2016 4:43
 Operator : UM/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040EC

Manual Integrations
 APPROVED

UMANGI
 3/10/2016 3:04:59 PM

Quant Time: Mar 10 11:56:31 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF030316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 04 00:54:58 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	195192	20.00	ng	-0.03
21) Naphthalene-d8	8.23	136	790098	20.00	ng	-0.03
38) Acenaphthene-d10	9.99	164	357224	20.00	ng	-0.03
63) Phenanthrene-d10	11.48	188	637384	20.00	ng	-0.03
75) Chrysene-d12	14.12	240	332970	20.00	ng	-0.03
86) Perylene-d12	15.58	264	307570	20.00	ng	-0.03

System Monitoring Compounds

5) 2-Fluorophenol	5.54	112	868991	74.68	ng	-0.03
7) Phenol-d6	6.57	99	1164744	76.40	ng	-0.03
23) Nitrobenzene-d5	7.51	82	1132717	76.72	ng	-0.03
41) 2,4,6-Tribromophenol	10.78	330	248632	81.34	ng	-0.03
44) 2-Fluorobiphenyl	9.32	172	1896879	80.76	ng	-0.03
78) Terphenyl-d14	13.07	244	1442712	100.58	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.62	88	231396	38.88	ng	96
3) Pyridine	3.37	79	662679	44.49	ng	99
4) n-Nitrosodimethylamine	3.33	42	242902	38.11	ng	96
6) Aniline	6.61	93	815612	38.17	ng	99
8) 2-Chlorophenol	6.73	128	562672	40.16	ng	96
9) Benzaldehyde	6.49	77	291136	40.81	ng	98
10) Phenol	6.58	94	611727m	37.46	ng	
11) bis(2-Chloroethyl)ether	6.69	93	546319	40.28	ng	95
12) 1,3-Dichlorobenzene	6.89	146	592505	39.39	ng	99
13) 1,4-Dichlorobenzene	6.96	146	557215	36.97	ng	98
14) 1,2-Dichlorobenzene	7.12	146	562494	41.13	ng	99
15) Benzyl Alcohol	7.09	79	415323	37.11	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.22	45	670747	34.31	ng	98
17) 2-Methylphenol	7.20	107	455671	39.86	ng	98
18) Hexachloroethane	7.46	117	215197	38.70	ng	95
19) n-Nitroso-di-n-propylamine	7.36	70	351449	35.37	ng	# 90
20) 3+4-Methylphenols	7.35	107	522925	38.62	ng	95
22) Acetophenone	7.36	105	714753	37.06	ng	99
24) Nitrobenzene	7.53	77	566234	37.75	ng	99
25) Isophorone	7.77	82	1041016	38.05	ng	98
26) 2-Nitrophenol	7.85	139	297406	40.75	ng	100
27) 2,4-Dimethylphenol	7.89	122	548358	41.11	ng	97
28) bis(2-Chloroethoxy)methane	7.98	93	563379	36.97	ng	100
29) 2,4-Dichlorophenol	8.09	162	446629	41.52	ng	97
30) 1,2,4-Trichlorobenzene	8.17	180	455059	39.12	ng	99
31) Naphthalene	8.25	128	1451059	37.35	ng	99
32) Benzoic acid	8.00	122	336617	37.99	ng	97
33) 4-Chloroaniline	8.30	127	610206	38.83	ng	98
34) Hexachlorobutadiene	8.38	225	252431	41.71	ng	99
35) Caprolactam	8.69	113	140984m	40.13	ng	
36) 4-Chloro-3-methylphenol	8.78	107	459567	37.66	ng	97
37) 2-Methylnaphthalene	8.95	142	1007403	40.41	ng	100
39) 1,2,4,5-Tetrachlorobenzene	9.11	216	410715	40.20	ng	98
40) Hexachlorocyclopentadiene	9.10	237	133951	24.31	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.22	196	326295	42.91	ng	97
43) 2,4,5-Trichlorophenol	9.26	196	301051	41.75	ng	93
45) 1,1'-Biphenyl	9.42	154	1206414	39.53	ng	92
46) 2-Chloronaphthalene	9.44	162	956040	41.10	ng	99
47) 2-Nitroaniline	9.53	65	315689	43.79	ng	# 81
48) Acenaphthylene	9.85	152	1400333	38.68	ng	99
49) Dimethylphthalate	9.72	163	1069343	39.00	ng	100
50) 2,6-Dinitrotoluene	9.77	165	216864	36.97	ng	89
51) Acenaphthene	10.02	154	801447	36.46	ng	99
52) 3-Nitroaniline	9.94	138	280380	39.95	ng	88
53) 2,4-Dinitrophenol	10.05	184	48383	22.65	ng	# 40
54) Dibenzofuran	10.20	168	1090430	37.86	ng	100
55) 4-Nitrophenol	10.09	139	186184	33.76	ng	# 69
56) 2,4-Dinitrotoluene	10.17	165	306330	40.81	ng	# 87
57) Fluorene	10.54	166	858583	37.95	ng	99
58) 2,3,4,6-Tetrachlorophenol	10.31	232	198857	39.26	ng	97
59) Diethylphthalate	10.41	149	1060968	39.88	ng	98
60) 4-Chlorophenyl-phenylether	10.53	204	430138	39.08	ng	94
61) 4-Nitroaniline	10.56	138	302543	46.09	ng	91
62) Azobenzene	10.69	77	993786	37.91	ng	97
64) 4,6-Dinitro-2-methylphenol	10.58	198	90871	23.81	ng	76
65) n-Nitrosodiphenylamine	10.65	169	869137	43.84	ng	100
66) 4-Bromophenyl-phenylether	11.02	248	256587	41.46	ng	# 89
67) Hexachlorobenzene	11.09	284	262471	39.75	ng	# 83
68) Atrazine	11.18	200	275723	50.01	ng	96
69) Pentachlorophenol	11.28	266	154007	42.02	ng	98
70) Phenanthrene	11.50	178	1261200	37.76	ng	99
71) Anthracene	11.56	178	1382099	38.98	ng	100
72) Carbazole	11.70	167	1065682	34.27	ng	98
73) Di-n-butylphthalate	12.04	149	1482817	37.73	ng	99
74) Fluoranthene	12.69	202	1113965	33.23	ng	97
76) Benzidine	12.81	184	486939	49.14	ng	99
77) Pyrene	12.92	202	1076618	42.96	ng	99
79) Butylbenzylphthalate	13.53	149	480965	42.30	ng	87
80) Benzo(a)anthracene	14.11	228	735075	36.88	ng	99
81) 3,3'-Dichlorobenzidine	14.07	252	282457	44.92	ng	# 96
82) Chrysene	14.14	228	614038	33.35	ng	99
83) Bis(2-ethylhexyl)phthalate	14.09	149	605797	40.48	ng	# 98
84) Di-n-octyl phthalate	14.70	149	889476	39.03	ng	95
85) Indeno(1,2,3-cd)pyrene	17.05	276	807021	56.16	ng	97
87) Benzo(b)fluoranthene	15.16	252	759534	37.05	ng	# 94
88) Benzo(k)fluoranthene	15.18	252	583174m	35.27	ng	
89) Benzo(a)pyrene	15.52	252	684669	40.30	ng	99
90) Dibenzo(a,h)anthracene	17.07	278	677688	46.73	ng	98
91) Benzo(g,h,i)perylene	17.49	276	647545	44.41	ng	99

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

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