

Data Path : S:\HPCHEM1\BNA_F\DATA\BF031016\
 Data File : BF085329.D
 Acq On : 10 Mar 2016 13:08
 Operator : UM/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 3/11/2016 4:58:12 PM

Quant Time: Mar 10 16:28:07 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF030316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 10 16:13:07 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.94	152	185960	20.00	ng	0.00
21) Naphthalene-d8	8.23	136	749656	20.00	ng	0.00
38) Acenaphthene-d10	9.98	164	335511	20.00	ng	0.00
63) Phenanthrene-d10	11.46	188	623050	20.00	ng	0.00
75) Chrysene-d12	14.11	240	375413	20.00	ng	0.00
86) Perylene-d12	15.57	264	311283	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.53	112	875207	78.95	ng	0.00
7) Phenol-d6	6.56	99	1159921	79.86	ng	0.00
23) Nitrobenzene-d5	7.51	82	1025811	73.22	ng	0.00
41) 2,4,6-Tribromophenol	10.78	330	257938	89.85	ng	0.00
44) 2-Fluorobiphenyl	9.30	172	1633192	74.03	ng	0.00
78) Terphenyl-d14	13.05	244	1297883	80.26	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.60	88	219465	38.71	ng	97
3) Pyridine	3.36	79	582804	41.07	ng	97
4) n-Nitrosodimethylamine	3.32	42	236755	38.99	ng	95
6) Aniline	6.61	93	788729	38.74	ng	94
8) 2-Chlorophenol	6.72	128	494854	37.08	ng	98
9) Benzaldehyde	6.48	77	263907	38.17	ng	96
10) Phenol	6.58	94	701596	45.10	ng	92
11) bis(2-Chloroethyl)ether	6.68	93	488806	37.83	ng	98
12) 1,3-Dichlorobenzene	6.88	146	563000m	39.29	ng	
13) 1,4-Dichlorobenzene	6.96	146	582200	40.54	ng	94
14) 1,2-Dichlorobenzene	7.11	146	495089	38.00	ng	97
15) Benzyl Alcohol	7.08	79	406585	38.13	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.21	45	679271	36.47	ng	99
17) 2-Methylphenol	7.19	107	407580	37.43	ng	98
18) Hexachloroethane	7.45	117	201789	38.09	ng	91
19) n-Nitroso-di-n-propylamine	7.36	70	363175	38.36	ng	98
20) 3+4-Methylphenols	7.35	107	515149	39.93	ng	# 79
22) Acetophenone	7.35	105	632284	34.55	ng	# 93
24) Nitrobenzene	7.52	77	506012	35.56	ng	98
25) Isophorone	7.76	82	1017148	39.18	ng	99
26) 2-Nitrophenol	7.84	139	294486	42.52	ng	# 79
27) 2,4-Dimethylphenol	7.88	122	472606	37.34	ng	94
28) bis(2-Chloroethoxy)methane	7.98	93	601113	41.58	ng	97
29) 2,4-Dichlorophenol	8.08	162	408407	40.01	ng	92
30) 1,2,4-Trichlorobenzene	8.17	180	446414	40.45	ng	95
31) Naphthalene	8.25	128	1476563	40.06	ng	99
32) Benzoic acid	7.99	122	224093	26.66	ng	98
33) 4-Chloroaniline	8.30	127	654361	43.89	ng	# 92
34) Hexachlorobutadiene	8.37	225	239581	41.72	ng	98
35) Caprolactam	8.68	113	137240	41.17	ng	91
36) 4-Chloro-3-methylphenol	8.78	107	479799	41.43	ng	87
37) 2-Methylnaphthalene	8.94	142	885573	37.44	ng	97
39) 1,2,4,5-Tetrachlorobenzene	9.11	216	425092	44.30	ng	97
40) Hexachlorocyclopentadiene	9.10	237	218412	42.20	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.21	196	269804	37.77	ng	99
43) 2,4,5-Trichlorophenol	9.26	196	285145	42.10	ng #	87
45) 1,1'-Biphenyl	9.41	154	1222940	42.66	ng	95
46) 2-Chloronaphthalene	9.43	162	857437	39.24	ng	95
47) 2-Nitroaniline	9.52	65	259201	38.28	ng	89
48) Acenaphthylene	9.84	152	1350802	39.72	ng	99
49) Dimethylphthalate	9.70	163	937191	36.39	ng	99
50) 2,6-Dinitrotoluene	9.76	165	208834	37.91	ng #	82
51) Acenaphthene	10.01	154	833597	40.37	ng	99
52) 3-Nitroaniline	9.93	138	239504	36.34	ng	84
53) 2,4-Dinitrophenol	10.04	184	101051	40.55	ng #	50
54) Dibenzofuran	10.18	168	1010962	37.37	ng	99
55) 4-Nitrophenol	10.08	139	204308	39.44	ng #	68
56) 2,4-Dinitrotoluene	10.17	165	305692	43.36	ng #	67
57) Fluorene	10.53	166	827559	38.95	ng	99
58) 2,3,4,6-Tetrachlorophenol	10.30	232	190750	40.09	ng	94
59) Diethylphthalate	10.40	149	944798	37.81	ng	96
60) 4-Chlorophenyl-phenylether	10.53	204	443950	42.94	ng #	83
61) 4-Nitroaniline	10.55	138	263626	42.76	ng	98
62) Azobenzene	10.69	77	981117	39.85	ng	85
64) 4,6-Dinitro-2-methylphenol	10.57	198	129076	34.60	ng #	56
65) n-Nitrosodiphenylamine	10.64	169	755642	38.99	ng	99
66) 4-Bromophenyl-phenylether	11.02	248	261984	43.30	ng #	81
67) Hexachlorobenzene	11.08	284	255875	39.65	ng #	74
68) Atrazine	11.17	200	236494	43.88	ng	98
69) Pentachlorophenol	11.27	266	153439	42.83	ng	99
70) Phenanthrene	11.50	178	1356312	41.54	ng	98
71) Anthracene	11.54	178	1300023	37.50	ng	99
72) Carbazole	11.69	167	1103498	36.31	ng	99
73) Di-n-butylphthalate	12.02	149	1249765	32.53	ng	99
74) Fluoranthene	12.68	202	1148141	35.04	ng	97
76) Benzidine	12.80	184	509465	45.60	ng	99
77) Pyrene	12.92	202	1203767	42.60	ng	97
79) Butylbenzylphthalate	13.52	149	521900	40.71	ng	98
80) Benzo(a)anthracene	14.09	228	904345	40.24	ng	98
81) 3,3'-Dichlorobenzidine	14.06	252	314650	44.38	ng	99
82) Chrysene	14.14	228	919834	44.31	ng	100
83) Bis(2-ethylhexyl)phthalate	14.08	149	660876	39.17	ng	100
84) Di-n-octyl phthalate	14.70	149	1032052	40.17	ng	96
85) Indeno(1,2,3-cd)pyrene	17.03	276	784911	48.44	ng	96
87) Benzo(b)fluoranthene	15.15	252	768174	37.02	ng #	96
88) Benzo(k)fluoranthene	15.18	252	718414m	42.93	ng	
89) Benzo(a)pyrene	15.51	252	691419	40.21	ng	99
90) Dibenzo(a,h)anthracene	17.05	278	658729	44.88	ng #	95
91) Benzo(g,h,i)perylene	17.48	276	664596	45.03	ng	97

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

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