

Data Path : Z:\HPCHEM1\BNA F\DATA\BF120215\
 Data File : BF083413.D
 Acq On : 2 Dec 2015 12:33
 Operator : UM/IZ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Mmdadoda
 12/3/2015 7:09:22 PM

Quant Time: Dec 03 03:12:48 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF113015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 01 01:17:16 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.54	152	180508m	20.00	ng	0.00
21) Naphthalene-d8	7.84	136	723024	20.00	ng	0.00
38) Acenaphthene-d10	9.58	164	351279	20.00	ng	0.00
63) Phenanthrene-d10	11.12	188	641964	20.00	ng	-0.01
75) Chrysene-d12	14.74	240	542614	20.00	ng	-0.02
86) Perylene-d12	16.80	264	397028	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.13	112	941030	78.23	ng	0.00
7) Phenol-d6	6.22	99	1281944	77.86	ng	0.00
23) Nitrobenzene-d5	7.12	82	1190354	72.37	ng	-0.01
41) 2,4,6-Tribromophenol	10.37	330	267812	75.95	ng	-0.01
44) 2-Fluorobiphenyl	8.91	172	1844696	74.83	ng	-0.01
78) Terphenyl-d14	13.21	244	1738341	76.40	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	1.94	88	210215	32.27	ng	99
3) Pyridine	2.56	79	582451	35.75	ng	96
4) n-Nitrosodimethylamine	2.51	42	322675	40.23	ng	97
6) Aniline	6.21	93	873872	38.79	ng	96
8) 2-Chlorophenol	6.34	128	513623	38.71	ng	92
9) Benzaldehyde	6.08	77	307694	38.52	ng	97
10) Phenol	6.24	94	765106	38.62	ng	99
11) bis(2-Chloroethyl)ether	6.29	93	565967	39.27	ng	89
12) 1,3-Dichlorobenzene	6.48	146	550514m	37.32	ng	
13) 1,4-Dichlorobenzene	6.56	146	559960m	36.73	ng	
14) 1,2-Dichlorobenzene	6.72	146	534034	38.18	ng	96
15) Benzyl Alcohol	6.70	79	485595	38.16	ng	93
16) 2,2'-oxybis(1-Chloropropan	6.84	45	929651	36.76	ng	# 49
17) 2-Methylphenol	6.84	107	443406	40.09	ng	97
18) Hexachloroethane	7.06	117	206445	38.99	ng	# 79
19) n-Nitroso-di-n-propylamine	6.98	70	424191	36.04	ng	92
20) 3+4-Methylphenols	7.00	107	589279	39.34	ng	99
22) Acetophenone	6.97	105	769547	35.57	ng	# 95
24) Nitrobenzene	7.14	77	661055	37.64	ng	90
25) Isophorone	7.38	82	1124152	35.28	ng	96
26) 2-Nitrophenol	7.46	139	292952	40.86	ng	# 78
27) 2,4-Dimethylphenol	7.52	122	437977	36.44	ng	95
28) bis(2-Chloroethoxy)methane	7.61	93	688455	37.94	ng	99
29) 2,4-Dichlorophenol	7.71	162	425678	36.83	ng	95
30) 1,2,4-Trichlorobenzene	7.78	180	457880	36.39	ng	98
31) Naphthalene	7.86	128	1478605	37.69	ng	98
32) Benzoic acid	7.68	122	313088	37.77	ng	94
33) 4-Chloroaniline	7.92	127	567313	33.55	ng	# 90
34) Hexachlorobutadiene	7.98	225	269841	37.77	ng	97
35) Caprolactam	8.30	113	133032	36.42	ng	91
36) 4-Chloro-3-methylphenol	8.42	107	461157	35.57	ng	# 81
37) 2-Methylnaphthalene	8.54	142	941459	35.91	ng	96
39) 1,2,4,5-Tetrachlorobenzene	8.72	216	447764	40.21	ng	98
40) Hexachlorocyclopentadiene	8.70	237	192943	36.48	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	8.84	196	309207	39.59	nq	98
43) 2,4,5-Trichlorophenol	8.89	196	308999	38.18	nq	# 93
45) 1,1'-Biphenyl	9.01	154	1203205	39.06	nq	98
46) 2-Chloronaphthalene	9.04	162	898136	38.26	nq	93
47) 2-Nitroaniline	9.14	65	351219	37.78	nq	# 83
48) Acenaphthylene	9.44	152	1385614	37.01	nq	99
49) Dimethylphthalate	9.33	163	1046944	36.67	nq	98
50) 2,6-Dinitrotoluene	9.38	165	217355	35.65	nq	85
51) Acenaphthene	9.61	154	788514	35.80	nq	98
52) 3-Nitroaniline	9.55	138	244362	34.20	nq	91
53) 2,4-Dinitrophenol	9.66	184	111901	34.62	nq	# 72
54) Dibenzofuran	9.78	168	1167408	36.17	nq	99
55) 4-Nitrophenol	9.74	139	181554m	40.62	nq	
56) 2,4-Dinitrotoluene	9.79	165	287855	37.21	nq	# 58
57) Fluorene	10.12	166	884223	34.69	nq	99
58) 2,3,4,6-Tetrachlorophenol	9.92	232	246162	37.70	nq	96
59) Diethylphthalate	10.02	149	965154	35.47	nq	100
60) 4-Chlorophenyl-phenylether	10.12	204	418143	35.71	nq	96
61) 4-Nitroaniline	10.17	138	254540	35.60	nq	94
62) Azobenzene	10.28	77	1247422	37.95	nq	99
64) 4,6-Dinitro-2-methylphenol	10.19	198	164696	38.21	nq	# 77
65) n-Nitrosodiphenylamine	10.25	169	858400	38.05	nq	100
66) 4-Bromophenyl-phenylether	10.62	248	278897	39.72	nq	98
67) Hexachlorobenzene	10.69	284	304373	39.20	nq	97
68) Atrazine	10.82	200	233181	37.37	nq	93
69) Pentachlorophenol	10.91	266	163870	40.29	nq	97
70) Phenanthrene	11.15	178	1463022	38.51	nq	100
71) Anthracene	11.21	178	1486361	38.86	nq	100
72) Carbazole	11.40	167	1278758	37.21	nq	99
73) Di-n-butylphthalate	11.85	149	1619548	39.77	nq	99
74) Fluoranthene	12.65	202	1446732	36.74	nq	98
76) Benzidine	12.85	184	553253	34.23	nq	98
77) Pyrene	12.96	202	1485833	39.21	nq	98
79) Butylbenzylphthalate	13.95	149	685296	42.63	nq	84
80) Benzo(a)anthracene	14.73	228	1264876	38.05	nq	99
81) 3,3'-Dichlorobenzidine	14.72	252	408756	39.37	nq	# 98
82) Chrysene	14.79	228	1246876	38.82	nq	99
83) Bis(2-ethylhexyl)phthalate	14.85	149	904463	41.94	nq	100
84) Di-n-octyl phthalate	15.83	149	1379972	42.97	nq	100
85) Indeno(1,2,3-cd)pyrene	18.18	276	967795	41.37	nq	100
87) Benzo(b)fluoranthene	16.28	252	956255	37.54	nq	99
88) Benzo(k)fluoranthene	16.33	252	916165	38.09	nq	99
89) Benzo(a)pyrene	16.73	252	839739	38.38	nq	98
90) Dibenzo(a,h)anthracene	18.20	278	813159	42.65	nq	99
91) Benzo(g,h,i)perylene	18.56	276	800414	42.77	nq	97

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

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