

Data Path : Z:\HPCHEM1\BNA F\DATA\BF120815\
 Data File : BF083611.D
 Acq On : 8 Dec 2015 23:20
 Operator : UM/IZ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Mmdadoda
 12/9/2015 6:13:27 PM

Quant Time: Dec 17 10:26:30 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 17 09:56:43 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	5.70	152	167627	20.00	ng	0.00
21) Naphthalene-d8	7.60	136	659267	20.00	ng	-0.01
38) Acenaphthene-d10	10.38	164	321160	20.00	ng	0.00
63) Phenanthrene-d10	12.79	188	597489	20.00	ng	0.00
75) Chrysene-d12	16.99	240	496683	20.00	ng	0.00
86) Perylene-d12	18.64	264	384639	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	3.97	112	890816	80.60	ng	0.00
7) Phenol-d6	5.26	99	1171307	79.32	ng	-0.01
23) Nitrobenzene-d5	6.53	82	1115542	78.96	ng	0.00
41) 2,4,6-Tribromophenol	11.68	330	236261	82.57	ng	-0.01
44) 2-Fluorobiphenyl	9.34	172	1776588	77.86	ng	0.00
78) Terphenyl-d14	15.41	244	1652388	78.28	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	1.81	88	217178	39.77	ng	95
3) Pyridine	2.26	79	567223	42.76	ng	100
4) n-Nitrosodimethylamine	2.22	42	293414	42.01	ng	95
6) Aniline	5.25	93	773617	42.51	ng	89
8) 2-Chlorophenol	5.42	128	491586	40.67	ng	92
9) Benzaldehyde	5.10	77	316559	39.31	ng	97
10) Phenol	5.29	94	736755	41.19	ng	79
11) bis(2-Chloroethyl)ether	5.36	93	543486	41.30	ng	97
12) 1,3-Dichlorobenzene	5.61	146	535520m	39.75	ng	
13) 1,4-Dichlorobenzene	5.72	146	558947	40.44	ng	98
14) 1,2-Dichlorobenzene	5.94	146	519155	40.48	ng	98
15) Benzyl Alcohol	5.95	79	441258	42.00	ng	98
16) 2,2'-oxybis(1-Chloropropan	6.12	45	977515	40.17	ng	# 61
17) 2-Methylphenol	6.14	107	411025	41.80	ng	93
18) Hexachloroethane	6.42	117	195732	40.11	ng	99
19) n-Nitroso-di-n-propylamine	6.34	70	413045	40.07	ng	95
20) 3+4-Methylphenols	6.38	107	520299	40.18	ng	96
22) Acetophenone	6.32	105	747906	40.44	ng	# 96
24) Nitrobenzene	6.57	77	601577	38.52	ng	93
25) Isophorone	6.93	82	1090296	39.77	ng	98
26) 2-Nitrophenol	7.05	139	254622	41.06	ng	90
27) 2,4-Dimethylphenol	7.16	122	439253	40.50	ng	95
28) bis(2-Chloroethoxy)methane	7.29	93	617713	40.57	ng	99
29) 2,4-Dichlorophenol	7.45	162	401513	40.70	ng	99
30) 1,2,4-Trichlorobenzene	7.53	180	434598	39.88	ng	97
31) Naphthalene	7.63	128	1383302	39.62	ng	99
32) Benzoic acid	7.49	122	243276	35.81	ng	94
33) 4-Chloroaniline	7.77	127	522215	41.01	ng	96
34) Hexachlorobutadiene	7.84	225	255073	40.07	ng	97
35) Caprolactam	8.38	113	123426	41.69	ng	93
36) 4-Chloro-3-methylphenol	8.61	107	447020	40.34	ng	99
37) 2-Methylnaphthalene	8.74	142	867982	39.57	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.00	216	415984	39.45	ng	# 100
40) Hexachlorocyclopentadiene	8.98	237	65223	43.38	ng	96

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.23	196	251418	40.07	nq	99
43) 2,4,5-Trichlorophenol	9.31	196	259898	41.84	nq #	91
45) 1,1'-Biphenyl	9.49	154	1105400	39.60	nq	99
46) 2-Chloronaphthalene	9.50	162	842691	39.98	nq	94
47) 2-Nitroaniline	9.72	65	316478	41.64	nq	98
48) Acenaphthylene	10.16	152	1378213	39.66	nq	100
49) Dimethylphthalate	10.04	163	1014423	39.68	nq	99
50) 2,6-Dinitrotoluene	10.13	165	219673	41.62	nq	92
51) Acenaphthene	10.44	154	797634	40.09	nq	98
52) 3-Nitroaniline	10.41	138	234190	41.75	nq	83
53) 2,4-Dinitrophenol	10.62	184	97601	40.35	nq	96
54) Dibenzofuran	10.73	168	1093479	39.63	nq	98
55) 4-Nitrophenol	10.82	139	104724	42.40	nq	83
56) 2,4-Dinitrotoluene	10.80	165	297625	42.52	nq #	77
57) Fluorene	11.28	166	951736	39.53	nq	98
58) 2,3,4,6-Tetrachlorophenol	10.97	232	206373	41.37	nq #	100
59) Diethylphthalate	11.18	149	986680	39.70	nq	99
60) 4-Chlorophenyl-phenylether	11.31	204	465372	40.60	nq	91
61) 4-Nitroaniline	11.40	138	218657	42.90	nq	84
62) Azobenzene	11.56	77	1111430	40.63	nq	97
64) 4,6-Dinitro-2-methylphenol	11.46	198	166124	41.10	nq	89
65) n-Nitrosodiphenylamine	11.52	169	784135	39.77	nq	98
66) 4-Bromophenyl-phenylether	12.09	248	261552	39.25	nq	95
67) Hexachlorobenzene	12.17	284	284825	40.29	nq	94
68) Atrazine	12.43	200	239159	40.76	nq	98
69) Pentachlorophenol	12.53	266	79864	45.34	nq	97
70) Phenanthrene	12.82	178	1372813	39.99	nq	99
71) Anthracene	12.91	178	1423395	40.03	nq	100
72) Carbazole	13.21	167	1240155	40.67	nq	99
73) Di-n-butylphthalate	13.84	149	1585638	41.34	nq	99
74) Fluoranthene	14.75	202	1427598	40.65	nq	99
76) Benzidine	15.03	184	687576	44.51	nq	99
77) Pyrene	15.11	202	1438592	40.23	nq	98
79) Butylbenzylphthalate	16.24	149	651215	41.03	nq	89
80) Benzo(a)anthracene	16.98	228	1160946	39.93	nq	98
81) 3,3'-Dichlorobenzidine	16.97	252	402835	41.02	nq	98
82) Chrysene	17.03	228	1165305	39.99	nq	99
83) Bis(2-ethylhexyl)phthalate	17.09	149	908677	41.25	nq	99
84) Di-n-octyl phthalate	17.86	149	1411619	41.94	nq #	100
85) Indeno(1,2,3-cd)pyrene	19.86	276	763619	38.82	nq #	100
87) Benzo(b)fluoranthene	18.25	252	855755m	39.07	nq	
88) Benzo(k)fluoranthene	18.28	252	998722m	41.38	nq	
89) Benzo(a)pyrene	18.58	252	840175	39.31	nq #	96
90) Dibenzo(a,h)anthracene	19.86	278	639135	39.11	nq	97
91) Benzo(g,h,i)perylene	20.23	276	599427	37.62	nq	96

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

