

Data Path : Z:\HPCHEM1\BNA N\DATA\BN041818\
 Data File : BN000731.D
 Acq On : 18 Apr 2018 09:21
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTDCCC040

Quant Time: Apr 18 15:24:14 2018
 Quant Method : Z:\HPCHEM1\BNA N\METHODS\8270-BN041818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 18 14:15:49 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	171526	20.00	ng	0.00
21) Naphthalene-d8	10.40	136	765828	20.00	ng	0.00
38) Acenaphthene-d10	14.26	164	413592	20.00	ng	0.00
63) Phenanthrene-d10	17.01	188	862778	20.00	ng	0.00
75) Chrysene-d12	21.22	240	808145	20.00	ng	0.00
86) Perylene-d12	23.42	264	775847	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.26	112	965314	83.60	ng	0.00
7) Phenol-d6	6.82	99	1356719	86.22	ng	0.00
23) Nitrobenzene-d5	8.78	82	1226833	83.51	ng	0.00
41) 2,4,6-Tribromophenol	15.76	330	368385	88.48	ng	0.00
44) 2-Fluorobiphenyl	12.89	172	2374454	79.25	ng	0.00
78) Terphenyl-d14	19.66	244	2896290	77.27	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.26	88	209620	39.580	ng	95
3) Pyridine	3.63	79	614833	44.071	ng	98
4) n-Nitrosodimethylamine	3.55	42	166329	43.355	ng	# 82
6) Aniline	6.98	93	903243	43.268	ng	92
8) 2-Chlorophenol	7.20	128	502636	42.091	ng	91
9) Benzaldehyde	6.79	77	310521	43.252	ng	93
10) Phenol	6.84	94	713190	43.000	ng	86
11) bis(2-Chloroethyl)ether	7.08	93	570086	42.013	ng	84
12) 1,3-Dichlorobenzene	7.52	146	566826	40.867	ng	99
13) 1,4-Dichlorobenzene	7.67	146	574186	40.775	ng	98
14) 1,2-Dichlorobenzene	7.98	146	551840	41.056	ng	97
15) Benzyl Alcohol	7.87	79	482811	44.612	ng	94
16) 2,2'-oxybis(1-Chloropropan	8.17	45	617829	42.872	ng	79
17) 2-Methylphenol	8.07	107	489405	43.893	ng	96
18) Hexachloroethane	8.70	117	214770	41.389	ng	# 80
19) n-Nitroso-di-n-propylamine	8.44	70	442477	45.204	ng	# 81
20) 3+4-Methylphenols	8.40	107	678988	44.311	ng	95
22) Acetophenone	8.45	105	913288	41.618	ng	# 88
24) Nitrobenzene	8.82	77	657695	41.752	ng	# 96
25) Isophorone	9.34	82	1168865	43.912	ng	98
26) 2-Nitrophenol	9.52	139	260898	40.428	ng	93
27) 2,4-Dimethylphenol	9.59	122	448796	42.600	ng	96
28) bis(2-Chloroethoxy)methane	9.83	93	730079	42.070	ng	97
29) 2,4-Dichlorophenol	10.05	162	444426	42.838	ng	98
30) 1,2,4-Trichlorobenzene	10.26	180	499682	39.960	ng	97
31) Naphthalene	10.45	128	1663482	40.402	ng	98
32) Benzoic acid	9.75	122	413996	42.605	ng	92
33) 4-Chloroaniline	10.56	127	725438	43.150	ng	97
34) Hexachlorobutadiene	10.74	225	268356	39.514	ng	98
35) Caprolactam	11.35	113	167388	46.635	ng	98
36) 4-Chloro-3-methylphenol	11.69	107	548258	44.759	ng	98
37) 2-Methylnaphthalene	12.07	142	1122100	41.603	ng	95
39) 1,2,4,5-Tetrachlorobenzene	12.44	216	503632	39.266	ng	# 97
40) Hexachlorocyclopentadiene	12.42	237	285179	37.914	ng	92

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	12.68	196	327593	42.109	ng	98
43) 2,4,5-Trichlorophenol	12.75	196	379763	42.264	ng	# 92
45) 1,1'-Biphenyl	13.10	154	1472328	39.771	ng	97
46) 2-Chloronaphthalene	13.13	162	1119133	40.319	ng	97
47) 2-Nitroaniline	13.34	65	345714	42.933	ng	85
48) Acenaphthylene	13.99	152	1827097	41.524	ng	98
49) Dimethylphthalate	13.73	163	1397946	42.296	ng	100
50) 2,6-Dinitrotoluene	13.85	165	295149	42.160	ng	93
51) Acenaphthene	14.33	154	1087618	40.551	ng	99
52) 3-Nitroaniline	14.17	138	348140	43.694	ng	# 95
53) 2,4-Dinitrophenol	14.37	184	141259	41.740	ng	97
54) Dibenzofuran	14.67	168	1595513	40.868	ng	97
55) 4-Nitrophenol	14.48	139	266858	45.104	ng	87
56) 2,4-Dinitrotoluene	14.63	165	400477	43.694	ng	95
57) Fluorene	15.32	166	1279450	41.453	ng	97
58) 2,3,4,6-Tetrachlorophenol	14.81	232	306230	44.336	ng	93
59) Diethylphthalate	15.12	149	1433629	43.766	ng	97
60) 4-Chlorophenyl-phenylether	15.32	204	584676	40.524	ng	92
61) 4-Nitroaniline	15.35	138	357370	45.595	ng	94
62) Azobenzene	15.62	77	1533054	43.708	ng	89
64) 4,6-Dinitro-2-methylphenol	15.40	198	221267	41.521	ng	88
65) n-Nitrosodiphenylamine	15.53	169	1160174	40.051	ng	96
66) 4-Bromophenyl-phenylether	16.22	248	343494	39.187	ng	# 89
67) Hexachlorobenzene	16.32	284	394111	38.833	ng	# 92
68) Atrazine	16.50	200	359492	45.810	ng	95
69) Pentachlorophenol	16.66	266	277968	42.387	ng	99
70) Phenanthrene	17.06	178	1988132	39.789	ng	99
71) Anthracene	17.15	178	1995477	40.777	ng	99
72) Carbazole	17.42	167	1954945	42.510	ng	97
73) Di-n-butylphthalate	18.00	149	2408959	45.971	ng	99
74) Fluoranthene	19.08	202	2162312	42.934	ng	93
76) Benzidine	19.27	184	962372	44.828	ng	97
77) Pyrene	19.44	202	2291143	39.357	ng	100
79) Butylbenzylphthalate	20.37	149	1075983	43.874	ng	96
80) Benzo(a)anthracene	21.20	228	2084137	40.921	ng	99
81) 3,3'-Dichlorobenzidine	21.14	252	704041	43.914	ng	# 96
82) Chrysene	21.26	228	1992653	40.414	ng	98
83) Bis(2-ethylhexyl)phthalate	21.17	149	1654861	44.175	ng	# 95
84) Di-n-octyl phthalate	22.04	149	2469964	46.083	ng	# 93
85) Indeno(1,2,3-cd)pyrene	26.27	276	1547958	36.503	ng	# 91
87) Benzo(b)fluoranthene	22.76	252	1961279	41.605	ng	# 95
88) Benzo(k)fluoranthene	22.81	252	1935823	41.232	ng	# 95
89) Benzo(a)pyrene	23.32	252	1802833	41.909	ng	# 94
90) Dibenzo(a,h)anthracene	25.62	278	1623049	38.291	ng	# 91
91) Benzo(g,h,i)perylene	26.27	276	1547958	37.405	ng	# 90

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

