

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP090819\
 Data File : BP000104.D
 Acq On : 08 Sep 2019 21:25
 Operator : HP/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDCCC040

Quant Time: Sep 09 01:24:37 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP083019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 30 17:47:17 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	332391	20.00	ng	0.00
21) Naphthalene-d8	8.18	136	1278408	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	688667	20.00	ng	-0.01
64) Phenanthrene-d10	11.43	188	1253161	20.00	ng	-0.01
76) Chrysene-d12	14.10	240	1070361	20.00	ng	-0.01
87) Perylene-d12	15.62	264	1175879	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.50	112	1632816	78.33	ng	0.00
7) Phenol-d6	6.51	99	2030525	77.16	ng	0.00
23) Nitrobenzene-d5	7.45	82	1628855	78.18	ng	0.00
42) 2,4,6-Tribromophenol	10.73	330	480035	82.70	ng	0.00
45) 2-Fluorobiphenyl	9.25	172	3425753	83.07	ng	0.00
79) Terphenyl-d14	13.03	244	3957864	80.60	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.66	88	375492	39.163	ng	96
3) Pyridine	3.44	79	1010889	39.133	ng	97
4) n-Nitrosodimethylamine	3.38	42	275985	33.087	ng	87
6) Aniline	6.55	93	1414940	37.154	ng	98
8) 2-Chlorophenol	6.68	128	855764	39.705	ng	98
9) Benzaldehyde	6.44	77	529048	32.603	ng	97
10) Phenol	6.52	94	1101679	37.682	ng	97
11) bis(2-Chloroethyl)ether	6.62	93	878693	38.046	ng	98
12) 1,3-Dichlorobenzene	6.83	146	1032540	41.071	ng	98
13) 1,4-Dichlorobenzene	6.91	146	1033850	41.298	ng	99
14) 1,2-Dichlorobenzene	7.06	146	963020	41.868	ng	98
15) Benzyl Alcohol	7.02	79	699798	36.272	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.17	45	875684	35.258	ng	95
17) 2-Methylphenol	7.13	107	716894	38.728	ng	97
18) Hexachloroethane	7.41	117	368661	39.705	ng	98
19) n-Nitroso-di-n-propylamine	7.30	70	537832	35.913	ng	97
20) 3+4-Methylphenols	7.29	107	932261	40.381	ng	# 83
22) Acetophenone	7.29	105	1229418	40.150	ng	98
24) Nitrobenzene	7.47	77	866450	38.388	ng	90
25) Isophorone	7.70	82	1614449	36.674	ng	99
26) 2-Nitrophenol	7.79	139	343504	39.080	ng	97
27) 2,4-Dimethylphenol	7.82	122	693491	39.098	ng	99
28) bis(2-Chloroethoxy)methane	7.92	93	1113715	38.906	ng	99
29) 2,4-Dichlorophenol	8.02	162	725404	39.773	ng	99
30) 1,2,4-Trichlorobenzene	8.12	180	859296	41.328	ng	99
31) Naphthalene	8.20	128	2478568	42.257	ng	99
32) Benzoic acid	7.90	122	371146	34.062	ng	97
33) 4-Chloroaniline	8.24	127	1112519	40.216	ng	# 93
34) Hexachlorobutadiene	8.32	225	485793	41.465	ng	98
35) Caprolactam	8.59	113	237955	36.687	ng	85
36) 4-Chloro-3-methylphenol	8.71	107	752095	37.794	ng	98
37) 2-Methylnaphthalene	8.89	142	1734908	42.325	ng	98
38) 1-Methylnaphthalene	8.99	142	1633173	42.286	ng	95
40) 1,2,4,5-Tetrachlorobenzene	9.06	216	824794	42.313	ng	98

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41) Hexachlorocyclopentadiene	9.05	237	402125	37.467	nq	100
43) 2,4,6-Trichlorophenol	9.16	196	522467	38.867	nq	98
44) 2,4,5-Trichlorophenol	9.20	196	549341	39.058	nq	98
46) 1,1'-Biphenyl	9.36	154	2093122	43.053	nq	97
47) 2-Chloronaphthalene	9.38	162	1667600	41.321	nq	96
48) 2-Nitroaniline	9.46	65	405139	37.166	nq	90
49) Acenaphthylene	9.80	152	2561876	42.781	nq	98
50) Dimethylphthalate	9.65	163	1927924	41.082	nq	99
51) 2,6-Dinitrotoluene	9.71	165	410290	41.190	nq	93
52) Acenaphthene	9.97	154	1571311	41.817	nq	98
53) 3-Nitroaniline	9.87	138	483698	40.648	nq	95
54) 2,4-Dinitrophenol	9.97	184	122210	37.838	nq	# 1
55) Dibenzofuran	10.14	168	2333616	43.735	nq	99
56) 4-Nitrophenol	10.02	139	339008	39.857	nq	92
57) 2,4-Dinitrotoluene	10.11	165	533113	41.922	nq	90
58) Fluorene	10.49	166	1767073	44.004	nq	99
59) 2,3,4,6-Tetrachlorophenol	10.26	232	442084	41.514	nq	96
60) Diethylphthalate	10.36	149	1913386	40.733	nq	99
61) 4-Chlorophenyl-phenylether	10.48	204	912144	43.701	nq	97
62) 4-Nitroaniline	10.49	138	478914	42.193	nq	95
63) Azobenzene	10.64	77	1750597	38.917	nq	95
65) 4,6-Dinitro-2-methylphenol	10.52	198	170711	37.426	nq	88
66) n-Nitrosodiphenylamine	10.59	169	1550237	41.638	nq	97
67) 4-Bromophenyl-phenylether	10.97	248	547514	42.991	nq	94
68) Hexachlorobenzene	11.05	284	613011	44.176	nq	94
69) Atrazine	11.12	200	379653	37.826	nq	96
70) Pentachlorophenol	11.23	266	299859	39.880	nq	97
71) Phenanthrene	11.46	178	2720529	43.556	nq	98
72) Anthracene	11.51	178	2662180	43.147	nq	99
73) Carbazole	11.66	167	2484367	42.658	nq	99
74) Di-n-butylphthalate	12.00	149	2957390	41.878	nq	99
75) Fluoranthene	12.66	202	2963348	43.659	nq	97
77) Benzidine	12.77	184	1078738	29.210	nq	96
78) Pyrene	12.89	202	3103793	39.417	nq	98
80) Butylbenzylphthalate	13.52	149	1239175	35.320	nq	# 87
81) Benzo(a)anthracene	14.09	228	2780804	40.093	nq	99
82) 3,3'-Dichlorobenzidine	14.05	252	955877	37.358	nq	# 98
83) Chrysene	14.13	228	2702850	41.199	nq	99
84) Bis(2-ethylhexyl)phthalate	14.09	149	1756153	38.463	nq	98
85) Di-n-octyl phthalate	14.71	149	2920331	36.297	nq	97
86) Indeno(1,2,3-cd)pyrene	17.17	276	3075915	37.417	nq	97
88) Benzo(b)fluoranthene	15.17	252	2890527	40.244	nq	97
89) Benzo(k)fluoranthene	15.20	252	2666752	40.573	nq	99
90) Benzo(a)pyrene	15.56	252	2576233	39.734	nq	97
91) Dibenzo(a,h)anthracene	17.19	278	2550530	39.336	nq	99
92) Benzo(g,h,i)perylene	17.63	276	2466330	38.431	nq	95

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

