

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP020320\
 Data File : BP001715.D
 Acq On : 03 Feb 2020 11:31
 Operator : CG/JU
 Sample : SSTDICC020
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDICC020

Quant Time: Feb 03 13:55:41 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP020320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 03 13:47:55 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	277130	20.00	ng	0.00
21) Naphthalene-d8	10.50	136	1113766	20.00	ng	0.00
39) Acenaphthene-d10	14.35	164	656006	20.00	ng	0.00
64) Phenanthrene-d10	17.10	188	1340392	20.00	ng	0.00
76) Chrysene-d12	21.19	240	1289533	20.00	ng	0.00
87) Perylene-d12	23.43	264	1435801	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.33	112	643698	45.50	ng	0.00
7) Phenol-d6	6.90	99	852497	37.70	ng	0.00
23) Nitrobenzene-d5	8.86	82	700589	27.49	ng	0.00
42) 2,4,6-Tribromophenol	15.84	330	251630	25.20	ng	0.00
45) 2-Fluorobiphenyl	12.97	172	1926204	43.17	ng	0.00
79) Terphenyl-d14	19.67	244	2798116	39.69	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.30	88	136503	30.234	ng	99
3) Pyridine	3.68	79	382054	23.418	ng	97
4) n-Nitrosodimethylamine	3.59	42	145474	12.581	ng	94
6) Aniline	7.06	93	529245	18.720	ng	100
8) 2-Chlorophenol	7.29	128	357570	21.649	ng	100
9) Benzaldehyde	6.87	77	249872	18.502	ng	98
10) Phenol	6.92	94	436019	18.642	ng	96
11) bis(2-Chloroethyl)ether	7.16	93	352075	19.193	ng	99
12) 1,3-Dichlorobenzene	7.62	146	443895	21.381	ng	99
13) 1,4-Dichlorobenzene	7.76	146	441525	20.567	ng	99
14) 1,2-Dichlorobenzene	8.08	146	428327	20.646	ng	100
15) Benzyl Alcohol	7.96	79	279825	13.142	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.26	45	451460	13.248	ng	100
17) 2-Methylphenol	8.16	107	290828	17.909	ng	97
18) Hexachloroethane	8.80	117	149440	17.831	ng	99
19) n-Nitroso-di-n-propylamine	8.53	70	261235	15.155	ng	97
20) 3+4-Methylphenols	8.49	107	376594	16.579	ng	98
22) Acetophenone	8.53	105	566555	18.146	ng	97
24) Nitrobenzene	8.90	77	365936	14.192	ng	98
25) Isophorone	9.43	82	680721	14.893	ng	99
26) 2-Nitrophenol	9.62	139	91183	9.357	ng	100
27) 2,4-Dimethylphenol	9.69	122	265580	18.243	ng	100
28) bis(2-Chloroethoxy)methane	9.92	93	469237	17.525	ng	99
29) 2,4-Dichlorophenol	10.15	162	302058	15.158	ng	99
30) 1,2,4-Trichlorobenzene	10.36	180	391510	16.275	ng	96
31) Naphthalene	10.55	128	1171184	19.925	ng	99
32) Benzoic acid	9.78	122	66180	5.351	ng	99
33) 4-Chloroaniline	10.65	127	490131	18.110	ng	99
34) Hexachlorobutadiene	10.85	225	229548	13.480	ng	100
35) Caprolactam	11.39	113	85202	11.774	ng	99
36) 4-Chloro-3-methylphenol	11.78	107	316359	13.228	ng	100
37) 2-Methylnaphthalene	12.16	142	807964	17.247	ng	100
38) 1-Methylnaphthalene	12.38	142	772510	17.136	ng	95
40) 1,2,4,5-Tetrachlorobenzene	12.54	216	421402	18.752	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.53	237	158897	13.940	nq	97
43) 2,4,6-Trichlorophenol	12.77	196	218136	15.051	nq	95
44) 2,4,5-Trichlorophenol	12.84	196	242011	15.714	nq	98
46) 1,1'-Biphenyl	13.19	154	1059941	22.548	nq	99
47) 2-Chloronaphthalene	13.22	162	830723	21.263	nq	99
48) 2-Nitroaniline	13.42	65	151075	9.842	nq	99
49) Acenaphthylene	14.07	152	1234046	20.481	nq	99
50) Dimethylphthalate	13.82	163	984122	18.214	nq	99
51) 2,6-Dinitrotoluene	13.92	165	169276	14.774	nq	90
52) Acenaphthene	14.42	154	784323	21.018	nq	95
53) 3-Nitroaniline	14.24	138	202490	16.738	nq	# 95
54) 2,4-Dinitrophenol	14.45	184	34005	5.278	nq	88
55) Dibenzofuran	14.75	168	1191949	18.888	nq	99
56) 4-Nitrophenol	14.55	139	133515	13.821	nq	99
57) 2,4-Dinitrotoluene	14.70	165	207117	12.354	nq	97
58) Fluorene	15.40	166	981727	19.226	nq	98
59) 2,3,4,6-Tetrachlorophenol	14.97	232	198218	12.464	nq	97
60) Diethylphthalate	15.19	149	978568	17.609	nq	98
61) 4-Chlorophenyl-phenylether	15.40	204	500862	17.138	nq	99
62) 4-Nitroaniline	15.40	138	223257	16.187	nq	97
63) Azobenzene	15.70	77	930144	16.751	nq	98
65) 4,6-Dinitro-2-methylphenol	15.47	198	50730	7.000	nq	99
66) n-Nitrosodiphenylamine	15.62	169	823431	23.189	nq	98
67) 4-Bromophenyl-phenylether	16.30	248	294144	19.452	nq	96
68) Hexachlorobenzene	16.41	284	346477	19.721	nq	96
69) Atrazine	16.57	200	256107	19.638	nq	99
70) Pentachlorophenol	16.75	266	130426	13.443	nq	99
71) Phenanthrene	17.14	178	1511945	20.742	nq	100
72) Anthracene	17.23	178	1466843	20.670	nq	98
73) Carbazole	17.50	167	1384681	20.784	nq	99
74) Di-n-butylphthalate	18.09	149	1536179	19.489	nq	99
75) Fluoranthene	19.12	202	1695027	17.672	nq	99
77) Benzidine	19.29	184	556144	18.888	nq	97
78) Pyrene	19.46	202	1792541	18.923	nq	98
80) Butylbenzylphthalate	20.36	149	571124	15.863	nq	98
81) Benzo(a)anthracene	21.17	228	1777808	19.796	nq	96
82) 3,3'-Dichlorobenzidine	21.11	252	537888	15.972	nq	99
83) Chrysene	21.22	228	1705978	19.492	nq	100
84) Bis(2-ethylhexyl)phthalate	21.14	149	1005573	20.374	nq	96
85) Di-n-octyl phthalate	22.02	149	1484883	18.869	nq	99
86) Indeno(1,2,3-cd)pyrene	25.70	276	1959791	19.129	nq	99
88) Benzo(b)fluoranthene	22.76	252	1732968	19.154	nq	99
89) Benzo(k)fluoranthene	22.80	252	1719443	19.491	nq	100
90) Benzo(a)pyrene	23.33	252	1559683	18.156	nq	97
91) Dibenzo(a,h)anthracene	25.72	278	1634281	18.438	nq	99
92) Benzo(g,h,i)perylene	26.40	276	1558237	17.685	nq	100

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

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