

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101319\
 Data File : BP000673.D
 Acq On : 12 Oct 2019 17:33
 Operator : JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDCCC040

Quant Time: Oct 12 22:20:36 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP100119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 01 18:03:42 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.75	152	227745	20.00	ng	-0.02
21) Naphthalene-d8	8.03	136	866576	20.00	ng	-0.02
39) Acenaphthene-d10	9.80	164	483755	20.00	ng	-0.02
64) Phenanthrene-d10	11.30	188	887003	20.00	ng	-0.01
76) Chrysene-d12	13.97	240	724748	20.00	ng	0.00
87) Perylene-d12	15.45	264	860682	20.00	ng	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.37	112	1138984	80.39	ng	-0.02
7) Phenol-d6	6.39	99	1390070	81.42	ng	-0.02
23) Nitrobenzene-d5	7.32	82	1113131	78.45	ng	-0.02
42) 2,4,6-Tribromophenol	10.59	330	406941	78.94	ng	-0.02
45) 2-Fluorobiphenyl	9.12	172	2324146	77.73	ng	-0.02
79) Terphenyl-d14	12.90	244	2788359	77.90	ng	-0.01

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.42	88	229209	40.512	ng	99
3) Pyridine	3.15	79	598899	38.743	ng	99
4) n-Nitrosodimethylamine	3.09	42	173117	39.798	ng	98
6) Aniline	6.41	93	858084	41.370	ng	# 90
8) 2-Chlorophenol	6.54	128	565235	40.423	ng	96
9) Benzaldehyde	6.30	77	392267	44.109	ng	97
10) Phenol	6.40	94	711500	40.702	ng	92
11) bis(2-Chloroethyl)ether	6.49	93	540203	40.167	ng	99
12) 1,3-Dichlorobenzene	6.69	146	670907	39.581	ng	98
13) 1,4-Dichlorobenzene	6.77	146	684368	40.125	ng	98
14) 1,2-Dichlorobenzene	6.92	146	630117	40.014	ng	99
15) Benzyl Alcohol	6.89	79	437219	39.653	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.03	45	505064	40.713	ng	100
17) 2-Methylphenol	7.01	107	456019	39.462	ng	98
18) Hexachloroethane	7.26	117	236224	38.915	ng	100
19) n-Nitroso-di-n-propylamine	7.16	70	306519	38.949	ng	99
20) 3+4-Methylphenols	7.16	107	552495	40.859	ng	93
22) Acetophenone	7.16	105	862463	43.072	ng	98
24) Nitrobenzene	7.33	77	537888	39.307	ng	98
25) Isophorone	7.57	82	985480	39.134	ng	96
26) 2-Nitrophenol	7.65	139	288842	40.137	ng	95
27) 2,4-Dimethylphenol	7.69	122	438064	39.711	ng	98
28) bis(2-Chloroethoxy)methane	7.78	93	679263	39.692	ng	99
29) 2,4-Dichlorophenol	7.89	162	493477	39.564	ng	96
30) 1,2,4-Trichlorobenzene	7.98	180	578777	39.511	ng	99
31) Naphthalene	8.06	128	1608940	39.754	ng	99
32) Benzoic acid	7.82	122	306083	43.943	ng	99
33) 4-Chloroaniline	8.10	127	713988	40.179	ng	100
34) Hexachlorobutadiene	8.18	225	340945	38.513	ng	98
35) Caprolactam	8.48	113	167307	40.048	ng	97
36) 4-Chloro-3-methylphenol	8.59	107	482408	39.489	ng	98
37) 2-Methylnaphthalene	8.75	142	1100894	40.520	ng	100
38) 1-Methylnaphthalene	8.85	142	1030980	40.160	ng	98
40) 1,2,4,5-Tetrachlorobenzene	8.92	216	686301	44.822	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.91	237	186064	34.887	ng	100
43) 2,4,6-Trichlorophenol	9.03	196	356929	37.869	ng	99
44) 2,4,5-Trichlorophenol	9.08	196	369985	38.019	ng	98
46) 1,1'-Biphenyl	9.22	154	1593024	43.627	ng	99
47) 2-Chloronaphthalene	9.25	162	1098942	38.920	ng	99
48) 2-Nitroaniline	9.34	65	272614	39.880	ng	88
49) Acenaphthylene	9.66	152	1658388	38.930	ng	100
50) Dimethylphthalate	9.52	163	1283353	38.180	ng	100
51) 2,6-Dinitrotoluene	9.58	165	284500	38.812	ng	96
52) Acenaphthene	9.83	154	972930	37.651	ng	99
53) 3-Nitroaniline	9.75	138	329096	39.182	ng	95
54) 2,4-Dinitrophenol	9.86	184	84136	34.398	ng	94
55) Dibenzofuran	10.00	168	1520654	39.367	ng	100
56) 4-Nitrophenol	9.92	139	197541	36.839	ng	94
57) 2,4-Dinitrotoluene	9.99	165	383519	39.460	ng	95
58) Fluorene	10.35	166	1141312	41.456	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.13	232	320425	38.950	ng	98
60) Diethylphthalate	10.23	149	1234245	38.681	ng	99
61) 4-Chlorophenyl-phenylether	10.35	204	603811	40.786	ng	99
62) 4-Nitroaniline	10.37	138	341277	42.248	ng	96
63) Azobenzene	10.50	77	1030396	42.060	ng	99
65) 4,6-Dinitro-2-methylphenol	10.40	198	148715	37.821	ng	99
66) n-Nitrosodiphenylamine	10.46	169	1039079	40.221	ng	98
67) 4-Bromophenyl-phenylether	10.84	248	384217	38.480	ng	96
68) Hexachlorobenzene	10.91	284	425182	37.473	ng	95
69) Atrazine	11.00	200	325620	38.408	ng	99
70) Pentachlorophenol	11.10	266	162582	35.676	ng	96
71) Phenanthrene	11.32	178	1779864	39.954	ng	99
72) Anthracene	11.37	178	1782139	40.347	ng	100
73) Carbazole	11.53	167	1656890	40.466	ng	98
74) Di-n-butylphthalate	11.87	149	1969609	39.488	ng	99
75) Fluoranthene	12.52	202	1979931	39.573	ng	98
77) Benzidine	12.65	184	851768	40.586	ng	99
78) Pyrene	12.76	202	2006582	40.143	ng	99
80) Butylbenzylphthalate	13.39	149	872264	40.046	ng	100
81) Benzo(a)anthracene	13.96	228	1778979	40.230	ng	97
82) 3,3'-Dichlorobenzidine	13.92	252	717188	40.832	ng	98
83) Chrysene	14.00	228	1752896	40.388	ng	100
84) Bis(2-ethylhexyl)phthalate	13.96	149	1097966	40.080	ng	98
85) Di-n-octyl phthalate	14.58	149	2045517	41.196	ng	100
86) Indeno(1,2,3-cd)pyrene	16.92	276	2142191	39.513	ng	99
88) Benzo(b)fluoranthene	15.02	252	1852894	37.947	ng	98
89) Benzo(k)fluoranthene	15.05	252	1847828	39.749	ng	98
90) Benzo(a)pyrene	15.39	252	1773129	38.955	ng	98
91) Dibenzo(a,h)anthracene	16.94	278	1720840	38.980	ng	98
92) Benzo(g,h,i)perylene	17.36	276	1739711	38.782	ng	100

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

