

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF021420\
 Data File : BF118932.D
 Acq On : 14 Feb 2020 14:16
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Feb 15 03:56:48 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF020320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Feb 15 03:51:59 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.78	152	195923	20.00	ng	-0.04
21) Naphthalene-d8	8.06	136	716783	20.00	ng	-0.04
39) Acenaphthene-d10	9.82	164	392038	20.00	ng	-0.04
64) Phenanthrene-d10	11.30	188	729966	20.00	ng	-0.04
76) Chrysene-d12	13.93	240	826645	20.00	ng	-0.05
87) Perylene-d12	15.36	264	731877	20.00	ng	-0.06

System Monitoring Compounds						
5) 2-Fluorophenol	5.40	112	1092910	105.59	ng	-0.04
7) Phenol-d6	6.42	99	1299558	101.15	ng	-0.03
23) Nitrobenzene-d5	7.35	82	1001899	83.31	ng	-0.04
42) 2,4,6-Tribromophenol	10.60	330	383979	81.86	ng	-0.04
45) 2-Fluorobiphenyl	9.14	172	1921206	87.61	ng	-0.04
79) Terphenyl-d14	12.88	244	3002190	74.60	ng	-0.05

Target Compounds				Qvalue		
2) 1,4-Dioxane	2.49	88	198961	42.18	ng	99
3) Pyridine	3.23	79	737705	56.32	ng	100
4) n-Nitrosodimethylamine	3.17	42	265452	55.66	ng	99
6) Aniline	6.45	93	860638	52.01	ng	96
8) 2-Chlorophenol	6.57	128	623291	54.07	ng	99
9) Benzaldehyde	6.33	77	371269	60.41	ng	99
10) Phenol	6.43	94	735843	51.51	ng	84
11) bis(2-Chloroethyl)ether	6.52	93	600450	54.94	ng	98
12) 1,3-Dichlorobenzene	6.72	146	581010	42.34	ng	98
13) 1,4-Dichlorobenzene	6.80	146	583672	42.64	ng	99
14) 1,2-Dichlorobenzene	6.95	146	534395	40.87	ng	98
15) Benzyl Alcohol	6.92	79	411208	41.41	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.06	45	528897	39.51	ng	96
17) 2-Methylphenol	7.04	107	398846	41.61	ng	99
18) Hexachloroethane	7.29	117	205364	41.80	ng	99
19) n-Nitroso-di-n-propylamine	7.20	70	310708	39.52	ng	99
20) 3+4-Methylphenols	7.19	107	462818	38.52	ng	# 87
22) Acetophenone	7.19	105	649496	41.71	ng	98
24) Nitrobenzene	7.36	77	502201	41.80	ng	98
25) Isophorone	7.60	82	866561	40.10	ng	100
26) 2-Nitrophenol	7.68	139	266668	41.66	ng	96
27) 2,4-Dimethylphenol	7.72	122	355538	40.89	ng	99
28) bis(2-Chloroethoxy)methane	7.81	93	569446	40.63	ng	98
29) 2,4-Dichlorophenol	7.92	162	428717	41.17	ng	99
30) 1,2,4-Trichlorobenzene	8.00	180	500279	41.10	ng	98
31) Naphthalene	8.09	128	1398117	42.65	ng	98
32) Benzoic acid	7.86	122	265293	36.85	ng	98
33) 4-Chloroaniline	8.13	127	600876	40.97	ng	98
34) Hexachlorobutadiene	8.20	225	312357	41.88	ng	99
35) Caprolactam	8.51	113	133180	39.45	ng	97
36) 4-Chloro-3-methylphenol	8.62	107	432309	41.25	ng	98
37) 2-Methylnaphthalene	8.77	142	937751	41.24	ng	98
38) 1-Methylnaphthalene	8.87	142	885417	41.40	ng	99
40) 1,2,4,5-Tetrachlorobenzene	8.94	216	504730	42.66	ng	100

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41) Hexachlorocyclopentadiene	8.93	237	163241	35.18	ng	99
43) 2,4,6-Trichlorophenol	9.05	196	318918	40.31	ng	99
44) 2,4,5-Trichlorophenol	9.10	196	332784	41.16	ng	99
46) 1,1'-Biphenyl	9.24	154	1171349	43.19	ng	98
47) 2-Chloronaphthalene	9.26	162	966507	42.87	ng	99
48) 2-Nitroaniline	9.36	65	271025	41.68	ng	94
49) Acenaphthylene	9.67	152	1402035	43.07	ng	100
50) Dimethylphthalate	9.54	163	1125856	42.18	ng	99
51) 2,6-Dinitrotoluene	9.60	165	256918	42.06	ng	99
52) Acenaphthene	9.85	154	846035	39.44	ng	99
53) 3-Nitroaniline	9.77	138	282988	41.78	ng	96
54) 2,4-Dinitrophenol	9.88	184	103324	35.19	ng	91
55) Dibenzofuran	10.02	168	1266830	40.87	ng	98
56) 4-Nitrophenol	9.94	139	181364	36.99	ng	93
57) 2,4-Dinitrotoluene	10.00	165	332753	41.17	ng	# 87
58) Fluorene	10.36	166	977862	41.93	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.14	232	286239	40.54	ng	99
60) Diethylphthalate	10.23	149	1065209	41.04	ng	98
61) 4-Chlorophenyl-phenylether	10.35	204	510565	40.37	ng	97
62) 4-Nitroaniline	10.38	138	291608	42.99	ng	96
63) Azobenzene	10.52	77	927620	41.86	ng	97
65) 4,6-Dinitro-2-methylphenol	10.42	198	164396	40.08	ng	98
66) n-Nitrosodiphenylamine	10.47	169	893060	42.24	ng	100
67) 4-Bromophenyl-phenylether	10.84	248	354369	41.04	ng	96
68) Hexachlorobenzene	10.92	284	412036	41.60	ng	95
69) Atrazine	11.00	200	274739	39.00	ng	99
70) Pentachlorophenol	11.10	266	172924	33.07	ng	98
71) Phenanthrene	11.32	178	1486261	41.38	ng	99
72) Anthracene	11.37	178	1481969	41.48	ng	98
73) Carbazole	11.53	167	1322501	42.18	ng	99
74) Di-n-butylphthalate	11.86	149	1594897	39.93	ng	99
75) Fluoranthene	12.51	202	1595137	40.68	ng	98
77) Benzidine	12.63	184	891371	35.65	ng	98
78) Pyrene	12.73	202	2129925	37.56	ng	99
80) Butylbenzylphthalate	13.35	149	847197	32.73	ng	98
81) Benzo(a)anthracene	13.92	228	1930504	39.25	ng	100
82) 3,3'-Dichlorobenzidine	13.88	252	748719	37.21	ng	# 98
83) Chrysene	13.96	228	1929895	39.09	ng	100
84) Bis(2-ethylhexyl)phthalate	13.92	149	1164155	35.45	ng	99
85) Di-n-octyl phthalate	14.52	149	1916254	34.06	ng	99
86) Indeno(1,2,3-cd)pyrene	16.79	276	1606523	32.39	ng	99
88) Benzo(b)fluoranthene	14.95	252	1891171	41.64	ng	100
89) Benzo(k)fluoranthene	14.99	252	1703005	43.06	ng	99
90) Benzo(a)pyrene	15.30	252	1618356	41.64	ng	100
91) Dibenzo(a,h)anthracene	16.79	278	1324115	38.38	ng	99
92) Benzo(g,h,i)perylene	17.20	276	1291719	38.93	ng	97

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

