

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF070220\
 Data File : BF120758.D
 Acq On : 2 Jul 2020 15:07
 Operator : JU/CG
 Sample : SSTDICC005
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC005

Quant Time: Jul 02 19:00:52 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF070220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 02 17:30:37 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	135959	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	528486	20.00	ng	0.00
39) Acenaphthene-d10	9.89	164	270899	20.00	ng	0.00
64) Phenanthrene-d10	11.37	188	499239	20.00	ng	0.00
76) Chrysene-d12	14.00	240	393950	20.00	ng	0.00
86) Perylene-d12	15.47	264	345046	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.45	112	94093	11.58	ng	-0.01
7) Phenol-d6	6.46	99	124624	11.67	ng	-0.02
23) Nitrobenzene-d5	7.40	82	81766	8.88	ng	-0.01
42) 2,4,6-Tribromophenol	10.67	330	23291	6.93	ng	0.00
45) 2-Fluorobiphenyl	9.20	172	206438	10.74	ng	0.00
79) Terphenyl-d14	12.95	244	215273	10.30	ng	-0.01

Target Compounds

					Qvalue
2) 1,4-Dioxane	2.59	88	22229	6.398 ng	96
3) Pyridine	3.35	79	57856	5.871 ng	100
4) n-Nitrosodimethylamine	3.27	42	29668	7.496 ng	# 88
6) Aniline	6.50	93	75356	5.822 ng	99
8) 2-Chlorophenol	6.63	128	49387	5.591 ng	99
9) Benzaldehyde	6.39	77	32499	6.023 ng	98
10) Phenol	6.47	94	61973	6.016 ng	98
11) bis(2-Chloroethyl)ether	6.57	93	50181	5.721 ng	97
12) 1,3-Dichlorobenzene	6.79	146	55330	5.673 ng	97
13) 1,4-Dichlorobenzene	6.86	146	56446	5.767 ng	97
14) 1,2-Dichlorobenzene	7.02	146	54657	5.662 ng	98
15) Benzyl Alcohol	6.98	79	46753	6.958 ng	99
16) 2,2'-oxybis(1-Chloropropan	7.12	45	80166	6.784 ng	100
17) 2-Methylphenol	7.09	107	44836	5.887 ng	97
18) Hexachloroethane	7.36	117	20303	5.757 ng	96
19) n-Nitroso-di-n-propylamine	7.25	70	38354	6.260 ng	99
20) 3+4-Methylphenols	7.24	107	59396	5.767 ng	94
22) Acetophenone	7.25	105	73611	5.489 ng	96
24) Nitrobenzene	7.42	77	47939	4.987 ng	94
25) Isophorone	7.66	82	103194	5.784 ng	97
26) 2-Nitrophenol	7.75	139	12609	2.452 ng	91
27) 2,4-Dimethylphenol	7.77	122	39464	5.119 ng	98
28) bis(2-Chloroethoxy)methane	7.87	93	60527	5.508 ng	97
29) 2,4-Dichlorophenol	7.97	162	39102	5.046 ng	97
30) 1,2,4-Trichlorobenzene	8.07	180	43989	4.847 ng	99
31) Naphthalene	8.15	128	151917	5.761 ng	98
32) Benzoic acid	7.82	122	12817	2.366 ng	94
33) 4-Chloroaniline	8.20	127	61841	5.371 ng	97
34) Hexachlorobutadiene	8.27	225	27382	4.997 ng	98
35) Caprolactam	8.52	113	11247	4.472 ng	100
36) 4-Chloro-3-methylphenol	8.66	107	41254	5.314 ng	98
37) 2-Methylnaphthalene	8.84	142	98518	5.529 ng	99
38) 1-Methylnaphthalene	8.94	142	91971	5.503 ng	96
40) 1,2,4,5-Tetrachlorobenzene	9.00	216	41372	4.784 ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	9.00	237	20416	5.166	ng	99
43) 2,4,6-Trichlorophenol	9.12	196	25592	4.387	ng	99
44) 2,4,5-Trichlorophenol	9.15	196	27325	4.169	ng	# 92
46) 1,1'-Biphenyl	9.30	154	114087	5.197	ng	100
47) 2-Chloronaphthalene	9.33	162	88206	4.993	ng	99
48) 2-Nitroaniline	9.42	65	15452	2.999	ng	# 84
49) Acenaphthylene	9.75	152	142618	5.199	ng	98
50) Dimethylphthalate	9.60	163	102703	4.947	ng	99
51) 2,6-Dinitrotoluene	9.66	165	12442	2.711	ng	# 78
52) Acenaphthene	9.92	154	87714	5.586	ng	99
53) 3-Nitroaniline	9.83	138	16075	2.983	ng	# 90
55) Dibenzofuran	10.09	168	130379	5.447	ng	99
56) 4-Nitrophenol	9.97	139	11214	3.368	ng	87
57) 2,4-Dinitrotoluene	10.06	165	12643	2.253	ng	# 87
58) Fluorene	10.43	166	101548	5.450	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.20	232	20969	4.235	ng	# 89
60) Diethylphthalate	10.30	149	107012	5.244	ng	98
61) 4-Chlorophenyl-phenylether	10.43	204	50697	5.361	ng	94
62) 4-Nitroaniline	10.43	138	15488	2.874	ng	93
63) Azobenzene	10.58	77	115326	6.191	ng	93
65) 4,6-Dinitro-2-methylphenol	10.46	198	2861	1.038	ng	92
66) n-Nitrosodiphenylamine	10.54	169	85073	5.181	ng	99
67) 4-Bromophenyl-phenylether	10.92	248	27516	4.528	ng	93
68) Hexachlorobenzene	10.97	284	29074	4.509	ng	# 91
69) Atrazine	11.06	200	20544	4.648	ng	96
70) Pentachlorophenol	11.17	266	12746	5.905	ng	96
71) Phenanthrene	11.39	178	141756	5.460	ng	98
72) Anthracene	11.44	178	142553	5.474	ng	98
73) Carbazole	11.59	167	137623	5.559	ng	98
74) Di-n-butylphthalate	11.93	149	169975	5.675	ng	98
75) Fluoranthene	12.58	202	148830	5.102	ng	97
78) Pyrene	12.81	202	154059	4.810	ng	99
80) Butylbenzylphthalate	13.43	149	59632	4.415	ng	91
81) Benzo(a)anthracene	13.99	228	131156	5.174	ng	98
82) 3,3'-Dichlorobenzidine	13.96	252	37147	3.926	ng	# 96
83) Chrysene	14.03	228	124712	4.690	ng	98
84) Bis(2-ethylhexyl)phthalate	14.00	149	93939	5.578	ng	100
85) Di-n-octyl phthalate	14.61	149	145018	4.545	ng	96
87) Indeno(1,2,3-cd)pyrene	16.90	276	101975	4.678	ng	98
88) Benzo(b)fluoranthene	15.04	252	108544	4.914	ng	98
89) Benzo(k)fluoranthene	15.07	252	107952	5.578	ng	99
90) Benzo(a)pyrene	15.40	252	97425	5.133	ng	98
91) Dibenzo(a,h)anthracene	16.93	278	82811	4.642	ng	98
92) Benzo(g,h,i)perylene	17.34	276	77253	4.328	ng	98

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

