

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080818\
 Data File : BF108015.D
 Acq On : 8 Aug 2018 13:13
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Aug 08 13:38:53 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF080818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 08 12:24:24 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.02	152	282721	20.00	ng	0.00
21) Naphthalene-d8	8.30	136	1124933	20.00	ng	0.00
38) Acenaphthene-d10	10.07	164	592114	20.00	ng	0.00
63) Phenanthrene-d10	11.56	188	1186411	20.00	ng	0.00
75) Chrysene-d12	14.22	240	923866	20.00	ng	0.00
86) Perylene-d12	15.78	264	764756	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.63	112	1249722	80.03	ng	0.00
7) Phenol-d6	6.63	99	1672292	80.59	ng	0.00
23) Nitrobenzene-d5	7.58	82	1649461	81.82	ng	0.00
41) 2,4,6-Tribromophenol	10.86	330	497148	84.17	ng	0.00
44) 2-Fluorobiphenyl	9.38	172	2845053	77.14	ng	0.00
78) Terphenyl-d14	13.15	244	2840295	78.17	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.92	88	322981	40.252	ng	88
3) Pyridine	3.68	79	835844	40.438	ng	# 83
4) n-Nitrosodimethylamine	3.65	42	468849	40.311	ng	82
6) Aniline	6.68	93	1153036	40.849	ng	94
8) 2-Chlorophenol	6.80	128	711153	40.434	ng	95
9) Benzaldehyde	6.57	77	653994	40.649	ng	97
10) Phenol	6.64	94	851366	39.663	ng	87
11) bis(2-Chloroethyl)ether	6.75	93	700091	40.343	ng	92
12) 1,3-Dichlorobenzene	6.96	146	816130	40.132	ng	99
13) 1,4-Dichlorobenzene	7.04	146	811289	39.706	ng	96
14) 1,2-Dichlorobenzene	7.19	146	758285	39.899	ng	97
15) Benzyl Alcohol	7.15	79	723205	41.398	ng	95
16) 2,2'-oxybis(1-Chloropropan	7.28	45	1438190	40.229	ng	96
17) 2-Methylphenol	7.25	107	616785	40.894	ng	# 94
18) Hexachloroethane	7.53	117	292941	40.271	ng	96
19) n-Nitroso-di-n-propylamine	7.42	70	571179	40.703	ng	# 85
20) 3+4-Methylphenols	7.41	107	791971	40.975	ng	93
22) Acetophenone	7.42	105	1043240	39.661	ng	# 90
24) Nitrobenzene	7.60	77	836261	41.022	ng	97
25) Isophorone	7.84	82	1550803	40.359	ng	98
26) 2-Nitrophenol	7.91	139	327775	42.576	ng	# 88
27) 2,4-Dimethylphenol	7.94	122	689633	40.438	ng	96
28) bis(2-Chloroethoxy)methane	8.04	93	894711	39.886	ng	98
29) 2,4-Dichlorophenol	8.15	162	650140	41.425	ng	98
30) 1,2,4-Trichlorobenzene	8.24	180	690536	39.907	ng	96
31) Naphthalene	8.32	128	2027261	39.626	ng	97
32) Benzoic acid	8.05	122	388513	40.346	ng	91
33) 4-Chloroaniline	8.37	127	899050	40.325	ng	95
34) Hexachlorobutadiene	8.44	225	442080	40.358	ng	99
35) Caprolactam	8.74	113	203280	41.332	ng	# 76
36) 4-Chloro-3-methylphenol	8.84	107	699367	41.307	ng	95
37) 2-Methylnaphthalene	9.01	142	1453506	40.157	ng	96
39) 1,2,4,5-Tetrachlorobenzene	9.18	216	732673	40.294	ng	98
40) Hexachlorocyclopentadiene	9.17	237	486432	41.417	ng	98

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42) 2,4,6-Trichlorophenol	9.29	196	480288	42.565	nq	98
43) 2,4,5-Trichlorophenol	9.32	196	518113	41.712	nq	96
45) 1,1'-Biphenyl	9.48	154	1730177	39.632	nq	97
46) 2-Chloronaphthalene	9.51	162	1389558	40.272	nq	99
47) 2-Nitroaniline	9.60	65	487161	43.635	nq	87
48) Acenaphthylene	9.93	152	2167816	39.741	nq	96
49) Dimethylphthalate	9.78	163	1637731	39.707	nq #	98
50) 2,6-Dinitrotoluene	9.84	165	367313	42.565	nq #	86
51) Acenaphthene	10.10	154	1341411	40.149	nq	98
52) 3-Nitroaniline	10.02	138	399293	42.291	nq	99
53) 2,4-Dinitrophenol	10.11	184	117403	40.184	nq #	37
54) Dibenzofuran	10.27	168	1915778	39.578	nq	91
55) 4-Nitrophenol	10.15	139	302966	42.375	nq #	78
56) 2,4-Dinitrotoluene	10.25	165	485420	43.701	nq #	97
57) Fluorene	10.62	166	1503778	39.244	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.38	232	453980	43.728	nq #	87
59) Diethylphthalate	10.48	149	1601628	40.101	nq	96
60) 4-Chlorophenyl-phenylether	10.61	204	802679	40.201	nq	91
61) 4-Nitroaniline	10.64	138	391613	41.952	nq	88
62) Azobenzene	10.77	77	1652895	40.073	nq	90
64) 4,6-Dinitro-2-methylphenol	10.65	198	177837	40.827	nq	89
65) n-Nitrosodiphenylamine	10.72	169	1407707	40.152	nq	98
66) 4-Bromophenyl-phenylether	11.10	248	523115	40.573	nq #	85
67) Hexachlorobenzene	11.17	284	551402	40.647	nq #	88
68) Atrazine	11.25	200	482455	41.028	nq	97
69) Pentachlorophenol	11.35	266	270487	41.658	nq	95
70) Phenanthrene	11.59	178	2188224	39.104	nq	96
71) Anthracene	11.64	178	2248652	39.207	nq	96
72) Carbazole	11.79	167	2005953	39.365	nq	96
73) Di-n-butylphthalate	12.11	149	2305962	39.952	nq #	97
74) Fluoranthene	12.78	202	2373608	38.866	nq	94
76) Benzidine	12.90	184	1463760	42.406	nq	94
77) Pyrene	13.01	202	2374163	40.591	nq	96
79) Butylbenzylphthalate	13.62	149	1004828	43.264	nq #	91
80) Benzo(a)anthracene	14.21	228	2109962	39.795	nq	96
81) 3,3'-Dichlorobenzidine	14.17	252	789517	41.551	nq #	95
82) Chrysene	14.25	228	1934687	39.801	nq	96
83) Bis(2-ethylhexyl)phthalate	14.18	149	1342561	42.686	nq #	99
84) Di-n-octyl phthalate	14.80	149	2083234	43.431	nq #	93
85) Indeno(1,2,3-cd)pyrene	17.39	276	1667314	39.587	nq #	91
87) Benzo(b)fluoranthene	15.31	252	1903280	41.650	nq	96
88) Benzo(k)fluoranthene	15.34	252	1664308	39.620	nq #	95
89) Benzo(a)pyrene	15.71	252	1656675	40.485	nq #	96
90) Dibenzo(a,h)anthracene	17.42	278	1384446	40.890	nq #	94
91) Benzo(g,h,i)perylene	17.89	276	1360265	40.801	nq #	91

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

