

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090518\
 Data File : BF108780.D
 Acq On : 5 Sep 2018 17:58
 Operator : JU/SJ
 Sample : SSTDICV040
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ICVBF090518

Quant Time: Sep 05 18:20:19 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF090518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 05 18:03:31 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.73	152	279791	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	1139352	20.00	ng	0.00
38) Acenaphthene-d10	9.77	164	552245	20.00	ng	0.00
63) Phenanthrene-d10	11.24	188	1051269	20.00	ng	0.00
75) Chrysene-d12	13.87	240	673123	20.00	ng	0.00
86) Perylene-d12	15.27	264	684830	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.34	112	1321867	78.27	ng	0.00
7) Phenol-d6	6.37	99	1702384	78.34	ng	0.00
23) Nitrobenzene-d5	7.30	82	1544260	84.82	ng	0.00
41) 2,4,6-Tribromophenol	10.55	330	439823	83.43	ng	0.00
44) 2-Fluorobiphenyl	9.10	172	2614621	80.77	ng	0.00
78) Terphenyl-d14	12.82	244	2215960	76.77	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.42	88	326241	40.195	ng	93
3) Pyridine	3.12	79	910970	40.983	ng	90
4) n-Nitrosodimethylamine	3.06	42	345607	40.481	ng	94
6) Aniline	6.40	93	1152875	39.446	ng	98
8) 2-Chlorophenol	6.52	128	747243	39.973	ng	97
9) Benzaldehyde	6.28	77	600509	38.954	ng	98
10) Phenol	6.38	94	986786	39.648	ng	93
11) bis(2-Chloroethyl)ether	6.47	93	765499	38.394	ng	98
12) 1,3-Dichlorobenzene	6.68	146	843339	39.533	ng	98
13) 1,4-Dichlorobenzene	6.75	146	844345	39.579	ng	97
14) 1,2-Dichlorobenzene	6.91	146	773741	39.349	ng	99
15) Benzyl Alcohol	6.88	79	668729	40.100	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.02	45	1155054	38.756	ng	96
17) 2-Methylphenol	6.99	107	623633	39.475	ng	98
18) Hexachloroethane	7.25	117	309702	40.262	ng	96
19) n-Nitroso-di-n-propylamine	7.15	70	582879	38.589	ng	98
20) 3+4-Methylphenols	7.14	107	713955	38.579	ng	# 65
22) Acetophenone	7.15	105	979836	37.893	ng	95
24) Nitrobenzene	7.31	77	818269	41.889	ng	97
25) Isophorone	7.56	82	1383945	38.109	ng	99
26) 2-Nitrophenol	7.63	139	350726	45.572	ng	98
27) 2,4-Dimethylphenol	7.67	122	617767	39.529	ng	98
28) bis(2-Chloroethoxy)methane	7.77	93	932427	38.403	ng	99
29) 2,4-Dichlorophenol	7.87	162	634474	39.372	ng	98
30) 1,2,4-Trichlorobenzene	7.96	180	676993	38.765	ng	98
31) Naphthalene	8.04	128	1986657	38.757	ng	99
32) Benzoic acid	7.80	122	454941	45.017	ng	98
33) 4-Chloroaniline	8.08	127	905648	38.469	ng	98
34) Hexachlorobutadiene	8.17	225	405628	38.879	ng	99
35) Caprolactam	8.46	113	217979	39.918	ng	# 83
36) 4-Chloro-3-methylphenol	8.57	107	666968	38.852	ng	98
37) 2-Methylnaphthalene	8.73	142	1299032	38.360	ng	98
39) 1,2,4,5-Tetrachlorobenzene	8.90	216	646059	39.136	ng	99
40) Hexachlorocyclopentadiene	8.89	237	340570	41.008	ng	99

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42) 2,4,6-Trichlorophenol	9.00	196	460921	40.992	nq	100
43) 2,4,5-Trichlorophenol	9.04	196	476502	40.883	nq	97
45) 1,1'-Biphenyl	9.20	154	1646974	38.738	nq	98
46) 2-Chloronaphthalene	9.21	162	1348278	39.212	nq	99
47) 2-Nitroaniline	9.31	65	421782	44.902	nq	98
48) Acenaphthylene	9.63	152	2006107	39.142	nq	98
49) Dimethylphthalate	9.50	163	1509909	39.091	nq	98
50) 2,6-Dinitrotoluene	9.55	165	340875	46.314	nq	92
51) Acenaphthene	9.80	154	1229951	38.780	nq	99
52) 3-Nitroaniline	9.71	138	401622	45.841	nq	92
53) 2,4-Dinitrophenol	9.81	184	103709	43.199	nq	# 28
54) Dibenzofuran	9.97	168	1776563	38.427	nq	96
55) 4-Nitrophenol	9.87	139	303710	46.370	nq	94
56) 2,4-Dinitrotoluene	9.95	165	433145	46.018	nq	# 95
57) Fluorene	10.31	166	1192608	40.198	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.08	232	396249	42.157	nq	88
59) Diethylphthalate	10.20	149	1549559	38.874	nq	96
60) 4-Chlorophenyl-phenylether	10.31	204	646390	40.795	nq	94
61) 4-Nitroaniline	10.32	138	362836	45.715	nq	98
62) Azobenzene	10.47	77	1586877	38.568	nq	97
64) 4,6-Dinitro-2-methylphenol	10.35	198	173477	42.826	nq	92
65) n-Nitrosodiphenylamine	10.42	169	1338104	38.283	nq	96
66) 4-Bromophenyl-phenylether	10.80	248	455777	38.386	nq	# 88
67) Hexachlorobenzene	10.86	284	488574	38.546	nq	93
68) Atrazine	10.95	200	434841	38.862	nq	98
69) Pentachlorophenol	11.05	266	338598	43.694	nq	99
70) Phenanthrene	11.27	178	1923482	38.314	nq	98
71) Anthracene	11.32	178	1929187	40.136	nq	96
72) Carbazole	11.47	167	1939414	38.313	nq	98
73) Di-n-butylphthalate	11.81	149	2257395	40.158	nq	# 98
74) Fluoranthene	12.45	202	2007180	40.381	nq	97
76) Benzidine	12.57	184	1221158	38.277	nq	96
77) Pyrene	12.68	202	2052863	39.634	nq	98
79) Butylbenzylphthalate	13.30	149	967300	41.470	nq	96
80) Benzo(a)anthracene	13.85	228	1705855	39.538	nq	98
81) 3,3'-Dichlorobenzidine	13.82	252	657552	39.118	nq	# 96
82) Chrysene	13.89	228	1629864	39.022	nq	98
83) Bis(2-ethylhexyl)phthalate	13.87	149	1186036	40.462	nq	99
84) Di-n-octyl phthalate	14.47	149	1930964	39.513	nq	100
85) Indeno(1,2,3-cd)pyrene	16.62	276	1428625	38.649	nq	96
87) Benzo(b)fluoranthene	14.87	252	1441568	38.712	nq	98
88) Benzo(k)fluoranthene	14.90	252	1358981	39.629	nq	97
89) Benzo(a)pyrene	15.21	252	1305348	38.849	nq	97
90) Dibenzo(a,h)anthracene	16.64	278	1188006	40.126	nq	97
91) Benzo(g,h,i)perylene	17.02	276	1197898	40.409	nq	95

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

