

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061118\
 Data File : BF106286.D
 Acq On : 11 Jun 2018 11:21
 Operator : JU/SJ
 Sample : SSTDICC2.5
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC2.5

Manual Integrations
 APPROVED

Sohil
 6/12/2018 12:57:54 PM

Quant Time: Jun 11 14:50:34 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 11 13:10:46 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	187406	20.00	ng	0.00
21) Naphthalene-d8	8.25	136	756823	20.00	ng	0.00
38) Acenaphthene-d10	10.01	164	382975	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	776591	20.00	ng	0.00
75) Chrysene-d12	14.20	240	715439	20.00	ng	0.04
86) Perylene-d12	15.79	264	627735	20.00	ng	0.09

System Monitoring Compounds

5) 2-Fluorophenol	5.57	112	61062	4.90	ng	0.00
7) Phenol-d6	6.57	99	76414	5.16	ng	-0.02
23) Nitrobenzene-d5	7.52	82	60989	4.69	ng	-0.01
41) 2,4,6-Tribromophenol	10.79	330	13805	2.79	ng	-0.01
44) 2-Fluorobiphenyl	0.00	172	0d	0.00	ng	
78) Terphenyl-d14	13.08	244	183078	5.50	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.83	88	15146	2.683	ng	95
3) Pyridine	3.62	79	40911m	2.793	ng	
4) n-Nitrosodimethylamine	3.55	42	17394	2.426	ng	# 91
6) Aniline	6.63	93	52727	2.746	ng	98
8) 2-Chlorophenol	6.75	128	31583	2.386	ng	94
9) Benzaldehyde	6.52	77	32263	3.205	ng	91
10) Phenol	6.58	94	41861	2.478	ng	96
11) bis(2-Chloroethyl)ether	6.69	93	35557	2.864	ng	92
12) 1,3-Dichlorobenzene	6.91	146	38585	2.648	ng	95
13) 1,4-Dichlorobenzene	6.99	146	39400	2.648	ng	100
14) 1,2-Dichlorobenzene	7.14	146	38794	2.830	ng	96
15) Benzyl Alcohol	7.10	79	30689	2.672	ng	95
16) 2,2'-oxybis(1-Chloropropan	7.23	45	49572	2.273	ng	99
17) 2-Methylphenol	7.20	107	25814	2.311	ng	# 94
18) Hexachloroethane	7.48	117	13658	2.541	ng	90
19) n-Nitroso-di-n-propylamine	7.36	70	30170	3.224	ng	# 96
20) 3+4-Methylphenols	7.35	107	36314	2.730	ng	93
22) Acetophenone	7.37	105	56760	3.264	ng	# 97
24) Nitrobenzene	7.55	77	35000	2.556	ng	96
25) Isophorone	7.78	82	67484	2.742	ng	98
27) 2,4-Dimethylphenol	7.89	122	25869	2.280	ng	93
28) bis(2-Chloroethoxy)methane	7.99	93	44991	2.980	ng	95
29) 2,4-Dichlorophenol	8.09	162	24517	2.306	ng	91
30) 1,2,4-Trichlorobenzene	8.19	180	31147	2.769	ng	89
31) Naphthalene	8.27	128	105495	2.863	ng	98
33) 4-Chloroaniline	8.32	127	41017	2.681	ng	97
34) Hexachlorobutadiene	8.39	225	20199	3.334	ng	94
36) 4-Chloro-3-methylphenol	8.78	107	28995	2.533	ng	96
37) 2-Methylnaphthalene	8.96	142	68712	2.772	ng	96
39) 1,2,4,5-Tetrachlorobenzene	9.13	216	34474	2.984	ng	97
40) Hexachlorocyclopentadiene	9.12	237	13491	1.725	ng	92
42) 2,4,6-Trichlorophenol	9.23	196	17441	2.220	ng	99
43) 2,4,5-Trichlorophenol	9.27	196	18814	2.216	ng	# 91
45) 1,1'-Biphenyl	9.43	154	90528	2.656	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
46) 2-Chloronaphthalene	9.45	162	69594	2.726	nq	95
47) 2-Nitroaniline	9.54	65	11806	1.376	nq #	75
48) Acenaphthylene	9.87	152	105457	2.695	nq	100
49) Dimethylphthalate	9.72	163	78113	2.657	nq	97
50) 2,6-Dinitrotoluene	9.78	165	10487	1.620	nq #	71
51) Acenaphthene	10.04	154	69050	2.666	nq	98
52) 3-Nitroaniline	9.95	138	12723	1.706	nq #	87
54) Dibenzofuran	10.22	168	95400	2.742	nq	100
55) 4-Nitrophenol	10.09	139	8941	1.466	nq #	77
56) 2,4-Dinitrotoluene	10.18	165	10962	1.302	nq #	80
57) Fluorene	10.56	166	78661	2.987	nq	95
58) 2,3,4,6-Tetrachlorophenol	10.33	232	14659	2.090	nq #	94
59) Diethylphthalate	10.42	149	77898	2.656	nq	98
60) 4-Chlorophenyl-phenylether	10.55	204	39210	3.215	nq	89
61) 4-Nitroaniline	10.56	138	12757	1.711	nq	91
62) Azobenzene	10.70	77	85899	2.992	nq	96
65) n-Nitrosodiphenylamine	10.66	169	74300	2.853	nq	93
66) 4-Bromophenyl-phenylether	11.04	248	22220	2.572	nq	91
67) Hexachlorobenzene	11.10	284	23313	2.361	nq #	89
68) Atrazine	11.18	200	18681	2.257	nq	97
69) Pentachlorophenol	11.29	266	9379	1.574	nq	95
70) Phenanthrene	11.52	178	118799	2.714	nq	98
71) Anthracene	11.57	178	114423	2.526	nq	99
72) Carbazole	11.73	167	103974	2.622	nq	97
73) Di-n-butylphthalate	12.05	149	103880	2.122	nq	98
74) Fluoranthene	12.72	202	116406	2.676	nq	95
76) Benzidine	12.83	184	59788	2.078	nq	99
77) Pvrene	12.94	202	124586	2.257	nq	99
79) Butylbenzylphthalate	13.57	149	23705	0.975	nq	98
80) Benzo(a)anthracene	14.19	228	117681	2.817	nq	97
81) 3,3'-Dichlorobenzidine	14.14	252	31415	1.954	nq #	98
82) Chrysene	14.23	228	112062	2.723	nq	97
83) Bis(2-ethylhexyl)phthalate	14.17	149	46601	1.582	nq #	95
84) Di-n-octyl phthalate	14.85	149	55847	1.155	nq #	90
85) Indeno(1,2,3-cd)pvrene	17.34	276	70657	1.885	nq	95
87) Benzo(b)fluoranthene	15.32	252	88481	2.365	nq	99
88) Benzo(k)fluoranthene	15.35	252	93445	2.583	nq	97
89) Benzo(a)pyrene	15.72	252	78210	2.245	nq	97
90) Dibenzo(a,h)anthracene	17.37	278	59294	1.815	nq	97
91) Benzo(g,h,i)perylene	17.82	276	61113	1.946	nq	95

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

