

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF073018\
 Data File : BF107793.D
 Acq On : 30 Jul 2018 11:43
 Operator : JU/SJ
 Sample : SSTDICC010
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC010

Manual Integrations
 APPROVED

Sohil
 7/31/2018 2:42:10 PM

Quant Time: Jul 30 17:46:37 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 30 13:05:29 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	161001	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	738142	20.00	ng	0.00
38) Acenaphthene-d10	10.00	164	349273	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	753225	20.00	ng	0.00
75) Chrysene-d12	14.15	240	664483	20.00	ng	0.00
86) Perylene-d12	15.68	264	616774	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.61	112	189659m	18.94	ng	0.01
7) Phenol-d6	6.60	99	264962	22.04	ng	0.00
23) Nitrobenzene-d5	7.53	82	279276	25.53	ng	0.00
41) 2,4,6-Tribromophenol	10.80	330	109498	27.57	ng	0.00
44) 2-Fluorobiphenyl	9.31	172	541384	23.57	ng	0.00
78) Terphenyl-d14	13.08	244	713582	27.49	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.83	88	37166m	7.974	ng	
3) Pyridine	3.67	79	71838m	5.666	ng	
4) n-Nitrosodimethylamine	3.78	42	48176m	9.009	ng	
6) Aniline	6.65	93	118279	7.627	ng	# 71
8) 2-Chlorophenol	6.75	128	129413	12.063	ng	94
9) Benzaldehyde	6.54	77	92162m	11.720	ng	
10) Phenol	6.62	94	170616	12.065	ng	76
11) bis(2-Chloroethyl)ether	6.70	93	157092	13.400	ng	93
12) 1,3-Dichlorobenzene	6.90	146	136190	11.197	ng	99
13) 1,4-Dichlorobenzene	6.98	146	162921	13.068	ng	97
14) 1,2-Dichlorobenzene	7.13	146	152327	13.615	ng	98
15) Benzyl Alcohol	7.12	79	77411m	9.447	ng	
16) 2,2'-oxybis(1-Chloropropan	7.22	45	234510	11.857	ng	75
17) 2-Methylphenol	7.22	107	88575	9.714	ng	# 66
18) Hexachloroethane	7.47	117	59004	13.494	ng	98
19) n-Nitroso-di-n-propylamine	7.36	70	95421	12.284	ng	98
20) 3+4-Methylphenols	7.38	107	104570	9.658	ng	# 65
22) Acetophenone	7.37	105	198612	11.457	ng	94
24) Nitrobenzene	7.55	77	143728	11.473	ng	99
25) Isophorone	7.77	82	248578	10.644	ng	99
26) 2-Nitrophenol	7.87	139	70174	13.265	ng	95
27) 2,4-Dimethylphenol	7.90	122	88713	8.859	ng	98
28) bis(2-Chloroethoxy)methane	7.99	93	145089	9.297	ng	99
29) 2,4-Dichlorophenol	8.12	162	104591	10.219	ng	97
30) 1,2,4-Trichlorobenzene	8.18	180	131336	11.425	ng	98
31) Naphthalene	8.27	128	409459	12.615	ng	98
32) Benzoic acid	8.04	122	37193m	6.063	ng	
33) 4-Chloroaniline	8.34	127	146361	10.316	ng	98
34) Hexachlorobutadiene	8.37	225	85183	12.470	ng	99
35) Caprolactam	8.67	113	28619	8.597	ng	# 70
36) 4-Chloro-3-methylphenol	8.81	107	94528	8.314	ng	97
37) 2-Methylnaphthalene	8.95	142	201311	8.950	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.12	216	107226	9.204	ng	94
40) Hexachlorocyclopentadiene	9.10	237	23966	4.047	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.24	196	57803	7.752	nq	98
43) 2,4,5-Trichlorophenol	9.29	196	76704	9.842	nq	96
45) 1,1'-Biphenyl	9.42	154	320236	10.525	nq	96
46) 2-Chloronaphthalene	9.45	162	236959	10.189	nq	99
47) 2-Nitroaniline	9.57	65	71501	10.546	nq	88
48) Acenaphthylene	9.87	152	448826	12.847	nq	99
49) Dimethylphthalate	9.71	163	268021	10.182	nq	99
50) 2,6-Dinitrotoluene	9.78	165	60752	11.624	nq	# 87
51) Acenaphthene	10.04	154	261975	11.968	nq	98
52) 3-Nitroaniline	9.98	138	78152	12.292	nq	100
53) 2,4-Dinitrophenol	10.12	184	15174	9.005	nq	# 52
54) Dibenzofuran	10.21	168	396966	12.318	nq	99
55) 4-Nitrophenol	10.20	139	10969m	2.304	nq	
56) 2,4-Dinitrotoluene	10.20	165	86877	13.771	nq	# 86
57) Fluorene	10.55	166	309702	13.471	nq	97
58) 2,3,4,6-Tetrachlorophenol	10.33	232	80907	12.474	nq	# 84
59) Diethylphthalate	10.41	149	336200	12.229	nq	98
60) 4-Chlorophenyl-phenylether	10.54	204	169803	13.065	nq	89
61) 4-Nitroaniline	10.60	138	69331	12.979	nq	95
62) Azobenzene	10.70	77	302075	11.718	nq	93
64) 4,6-Dinitro-2-methylphenol	10.61	198	35524	11.267	nq	87
65) n-Nitrosodiphenylamine	10.66	169	294640	11.518	nq	97
66) 4-Bromophenyl-phenylether	11.04	248	104054	11.735	nq	# 85
67) Hexachlorobenzene	11.10	284	113530	11.637	nq	# 84
68) Atrazine	11.18	200	88434	11.322	nq	94
69) Pentachlorophenol	11.31	266	36821	6.042	nq	96
70) Phenanthrene	11.52	178	420993	11.470	nq	98
71) Anthracene	11.58	178	499389	13.513	nq	97
72) Carbazole	11.74	167	484703	12.613	nq	98
73) Di-n-butylphthalate	12.04	149	549184	12.623	nq	98
74) Fluoranthene	12.71	202	506730	12.865	nq	96
76) Benzidine	12.85	184	250461	13.089	nq	97
77) Pyrene	12.95	202	519879	11.438	nq	97
79) Butylbenzylphthalate	13.55	149	224863	11.504	nq	89
80) Benzo(a)anthracene	14.14	228	414403	10.642	nq	99
81) 3,3'-Dichlorobenzidine	14.10	252	176807	14.340	nq	97
82) Chrysene	14.18	228	467133	12.488	nq	99
83) Bis(2-ethylhexyl)phthalate	14.09	149	308314	11.179	nq	99
84) Di-n-octyl phthalate	14.72	149	499792	11.658	nq	98
85) Indeno(1,2,3-cd)pyrene	17.26	276	340707	10.747	nq	# 94
87) Benzo(b)fluoranthene	15.22	252	327739	9.709	nq	98
88) Benzo(k)fluoranthene	15.25	252	435754	13.176	nq	# 95
89) Benzo(a)pyrene	15.61	252	349494	11.319	nq	98
90) Dibenzo(a,h)anthracene	17.27	278	289455	10.845	nq	# 95
91) Benzo(g,h,i)perylene	17.74	276	259344	9.851	nq	# 92

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

