

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121418\
 Data File : BF111411.D
 Acq On : 14 Dec 2018 11:34
 Operator : JU/SJ
 Sample : SSTDICC050
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC050

Manual Integrations
 APPROVED

Sohil
 12/17/2018 2:36:11 PM

Quant Time: Dec 14 12:11:26 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 14 11:42:38 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	278421	20.00	ng	0.00
21) Naphthalene-d8	8.08	136	1183473	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	559281	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	971447	20.00	ng	0.00
76) Chrysene-d12	13.96	240	610328	20.00	ng	0.00
87) Perylene-d12	15.39	264	479444	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.42	112	1407036	85.46	ng	0.00
7) Phenol-d6	6.45	99	2162080	97.40	ng	0.00
23) Nitrobenzene-d5	7.36	82	2034134	96.88	ng	0.00
42) 2,4,6-Tribromophenol	10.62	330	484395	98.25	ng	0.00
45) 2-Fluorobiphenyl	9.16	172	3271057	93.90	ng	0.00
79) Terphenyl-d14	12.90	244	2507533	90.43	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.50	88	334660	40.692	ng	98
3) Pyridine	3.23	79	882498	43.681	ng	94
4) n-Nitrosodimethylamine	3.19	42	410047	43.507	ng	93
6) Aniline	6.46	93	1357435	50.546	ng	# 76
8) 2-Chlorophenol	6.59	128	996404	49.237	ng	99
9) Benzaldehyde	6.34	77	607132	43.947	ng	98
10) Phenol	6.46	94	1187653	47.229	ng	# 73
11) bis(2-Chloroethyl)ether	6.53	93	989028	49.994	ng	99
12) 1,3-Dichlorobenzene	6.74	146	1068739	48.783	ng	99
13) 1,4-Dichlorobenzene	6.82	146	1083759	48.868	ng	98
14) 1,2-Dichlorobenzene	6.97	146	978558	47.695	ng	99
15) Benzyl Alcohol	6.95	79	826178	48.603	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.07	45	1776459	46.351	ng	94
17) 2-Methylphenol	7.06	107	797269	48.755	ng	95
18) Hexachloroethane	7.31	117	402707	48.911	ng	96
19) n-Nitroso-di-n-propylamine	7.22	70	700579	47.493	ng	98
20) 3+4-Methylphenols	7.22	107	942580	48.335	ng	# 76
22) Acetophenone	7.21	105	1320242	48.026	ng	# 93
24) Nitrobenzene	7.38	77	1036176	47.909	ng	97
25) Isophorone	7.62	82	1734959	48.644	ng	99
26) 2-Nitrophenol	7.69	139	552702	51.990	ng	99
27) 2,4-Dimethylphenol	7.74	122	798234	49.464	ng	99
28) bis(2-Chloroethoxy)methane	7.83	93	1188219	48.537	ng	99
29) 2,4-Dichlorophenol	7.94	162	824602	48.984	ng	98
30) 1,2,4-Trichlorobenzene	8.02	180	835294	48.291	ng	99
31) Naphthalene	8.10	128	2617500	47.471	ng	98
32) Benzoic acid	7.89	122	692025m	51.585	ng	
33) 4-Chloroaniline	8.15	127	1204796	51.163	ng	96
34) Hexachlorobutadiene	8.22	225	449489	46.843	ng	99
35) Caprolactam	8.54	113	311753m	52.851	ng	
36) 4-Chloro-3-methylphenol	8.64	107	883513	49.252	ng	99
37) 2-Methylnaphthalene	8.79	142	1726498	48.438	ng	99
38) 1-Methylnaphthalene	8.89	142	1657438	48.179	ng	98
40) 1,2,4,5-Tetrachlorobenzene	8.96	216	732664	46.605	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.95	237	415252	50.585	nq	100
43) 2,4,6-Trichlorophenol	9.07	196	536305	47.662	nq	99
44) 2,4,5-Trichlorophenol	9.12	196	582785	49.007	nq	98
46) 1,1'-Biphenyl	9.26	154	2236497	46.868	nq	97
47) 2-Chloronaphthalene	9.28	162	1751406	47.925	nq	97
48) 2-Nitroaniline	9.37	65	596822	50.256	nq	96
49) Acenaphthylene	9.70	152	2534751	47.475	nq	97
50) Dimethylphthalate	9.56	163	1955509	48.622	nq	99
51) 2,6-Dinitrotoluene	9.62	165	464381	52.090	nq	98
52) Acenaphthene	9.87	154	1639096	48.672	nq	99
53) 3-Nitroaniline	9.79	138	565233	52.749	nq	92
54) 2,4-Dinitrophenol	9.89	184	191126	54.645	nq	# 83
55) Dibenzofuran	10.04	168	2333932	48.008	nq	98
56) 4-Nitrophenol	9.96	139	410893	53.247	nq	96
57) 2,4-Dinitrotoluene	10.02	165	599357	52.457	nq	98
58) Fluorene	10.38	166	1637952	47.097	nq	99
59) 2,3,4,6-Tetrachlorophenol	10.16	232	441006	48.005	nq	98
60) Diethylphthalate	10.26	149	1915410	47.306	nq	99
61) 4-Chlorophenyl-phenylether	10.37	204	807513	46.837	nq	100
62) 4-Nitroaniline	10.40	138	553234	53.714	nq	98
63) Azobenzene	10.53	77	1940542	47.534	nq	96
65) 4,6-Dinitro-2-methylphenol	10.43	198	285191	51.136	nq	89
66) n-Nitrosodiphenylamine	10.49	169	1627120	47.086	nq	98
67) 4-Bromophenyl-phenylether	10.86	248	515285	47.647	nq	98
68) Hexachlorobenzene	10.93	284	530612	47.648	nq	95
69) Atrazine	11.02	200	467958	46.821	nq	99
70) Pentachlorophenol	11.13	266	345781	46.660	nq	96
71) Phenanthrene	11.34	178	2372867	47.257	nq	95
72) Anthracene	11.39	178	2435259	47.631	nq	97
73) Carbazole	11.55	167	2550493	48.245	nq	96
74) Di-n-butylphthalate	11.88	149	2832582	47.058	nq	98
75) Fluoranthene	12.53	202	2085552	42.546	nq	99
77) Benzidine	12.64	184	981019	85.174	nq	99
78) Pyrene	12.76	202	2166499	45.995	nq	98
80) Butylbenzylphthalate	13.38	149	1045804	46.242	nq	87
81) Benzo(a)anthracene	13.94	228	1589991	46.561	nq	99
82) 3,3'-Dichlorobenzidine	13.91	252	641360	49.660	nq	99
83) Chrysene	13.99	228	1705995	48.966	nq	97
84) Bis(2-ethylhexyl)phthalate	13.94	149	1307990	47.309	nq	99
85) Di-n-octyl phthalate	14.55	149	2218964	48.878	nq	99
86) Indeno(1,2,3-cd)pyrene	16.81	276	1252320	45.958	nq	97
88) Benzo(b)fluoranthene	14.98	252	1399496	51.333	nq	99
89) Benzo(k)fluoranthene	15.01	252	1341022	49.279	nq	99
90) Benzo(a)pyrene	15.33	252	1237970	48.769	nq	97
91) Dibenzo(a,h)anthracene	16.83	278	1048996	49.669	nq	95
92) Benzo(g,h,i)perylene	17.23	276	1118991	52.995	nq	96

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

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