

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF022818\
 Data File : BF103310.D
 Acq On : 28 Feb 2018 22:58
 Operator : SJ/JU
 Sample : J1708-01MS
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RO-222-226MS

Manual Integrations
 APPROVED

Sohil
 3/1/2018 11:43:29 AM

Quant Time: Mar 01 01:17:01 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 27 17:05:20 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	146978	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	623910	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	265368	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	370561	20.00	ng	0.00
75) Chrysene-d12	14.06	240	237157	20.00	ng	0.00
86) Perylene-d12	15.55	264	238893	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.50	112	1186013	122.04	ng	0.01
7) Phenol-d6	6.52	99	1392213	117.08	ng	0.00
23) Nitrobenzene-d5	7.45	82	884332	80.95	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	218325	97.20	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	1254833	80.41	ng	0.00
78) Terphenyl-d14	12.99	244	799214	81.01	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.70	88	192750	37.553	ng	91
3) Pyridine	3.49	79	509813	37.205	ng	99
4) n-Nitrosodimethylamine	3.43	42	301172	54.498	ng	# 89
6) Aniline	6.54	93	374953	21.848	ng	# 88
8) 2-Chlorophenol	6.66	128	418669	41.471	ng	92
9) Benzaldehyde	6.43	77	205360	25.620	ng	97
10) Phenol	6.53	94	556041	40.279	ng	88
11) bis(2-Chloroethyl)ether	6.61	93	434129	41.435	ng	90
12) 1,3-Dichlorobenzene	6.82	146	427918	37.092	ng	98
13) 1,4-Dichlorobenzene	6.89	146	432008	37.350	ng	100
14) 1,2-Dichlorobenzene	7.05	146	395546	37.087	ng	100
15) Benzyl Alcohol	7.02	79	353630	39.464	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.15	45	689970	38.349	ng	91
17) 2-Methylphenol	7.13	107	354384	38.628	ng	# 94
18) Hexachloroethane	7.38	117	144160	34.237	ng	94
19) n-Nitroso-di-n-propylamine	7.29	70	284998	38.509	ng	95
20) 3+4-Methylphenols	7.29	107	412404	36.876	ng	# 59
22) Acetophenone	7.29	105	546505	37.527	ng	# 90
24) Nitrobenzene	7.46	77	466852	38.812	ng	99
25) Isophorone	7.70	82	869996	40.433	ng	98
26) 2-Nitrophenol	7.77	139	225541	39.830	ng	94
27) 2,4-Dimethylphenol	7.81	122	362201	41.847	ng	99
28) bis(2-Chloroethoxy)methane	7.90	93	541454	42.800	ng	99
29) 2,4-Dichlorophenol	8.02	162	339570	40.977	ng	98
30) 1,2,4-Trichlorobenzene	8.10	180	327759	37.142	ng	97
31) Naphthalene	8.18	128	1143326	37.430	ng	100
32) Benzoic acid	7.92	122	81163	17.733	ng	98
33) 4-Chloroaniline	8.23	127	159058	12.067	ng	96
34) Hexachlorobutadiene	8.29	225	162126	35.281	ng	99
35) Caprolactam	8.60	113	116860m	40.125	ng	
36) 4-Chloro-3-methylphenol	8.72	107	374117	40.026	ng	100
37) 2-Methylnaphthalene	8.87	142	747606	37.871	ng	100
39) 1,2,4,5-Tetrachlorobenzene	9.04	216	275726	38.007	ng	98
40) Hexachlorocyclopentadiene	9.02	237	20482	16.002	ng	96

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.15	196	200310	41.035	nq	97
43) 2,4,5-Trichlorophenol	9.20	196	209220	37.265	nq	98
45) 1,1'-Biphenyl	9.33	154	829637	37.310	nq	99
46) 2-Chloronaphthalene	9.36	162	660248	37.531	nq	99
47) 2-Nitroaniline	9.46	65	264605	40.606	nq	87
48) Acenaphthylene	9.78	152	977800	36.125	nq	99
49) Dimethylphthalate	9.63	163	905824	42.550	nq	100
50) 2,6-Dinitrotoluene	9.70	165	170745	38.220	nq	96
51) Acenaphthene	9.95	154	561130	35.552	nq	99
52) 3-Nitroaniline	9.87	138	136314	24.050	nq	99
53) 2,4-Dinitrophenol	10.00	184	7208	13.551	nq	# 20
54) Dibenzofuran	10.12	168	820132	37.853	nq	97
55) 4-Nitrophenol	10.05	139	211841	60.305	nq	92
56) 2,4-Dinitrotoluene	10.11	165	206257	36.505	nq	# 90
57) Fluorene	10.47	166	629963	34.132	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.25	232	136936	36.307	nq	97
59) Diethylphthalate	10.33	149	767815	37.711	nq	98
60) 4-Chlorophenyl-phenylether	10.45	204	295292	37.728	nq	98
61) 4-Nitroaniline	10.49	138	168950	29.843	nq	91
62) Azobenzene	10.62	77	757885	36.569	nq	97
64) 4,6-Dinitro-2-methylphenol	10.52	198	12232	9.812	nq	86
65) n-Nitrosodiphenylamine	10.57	169	572681	44.386	nq	100
66) 4-Bromophenyl-phenylether	10.95	248	164930	42.726	nq	93
67) Hexachlorobenzene	11.02	284	160254	38.249	nq	97
68) Atrazine	11.10	200	168275	43.392	nq	97
69) Pentachlorophenol	11.22	266	120949	59.703	nq	98
70) Phenanthrene	11.43	178	750867	36.820	nq	98
71) Anthracene	11.49	178	779679	37.284	nq	99
72) Carbazole	11.64	167	726906	34.906	nq	100
73) Di-n-butylphthalate	11.96	149	1050197	40.302	nq	99
74) Fluoranthene	12.62	202	649997	31.718	nq	98
76) Benzidine	12.75	184	378998	42.747	nq	98
77) Pyrene	12.86	202	667099	36.078	nq	99
79) Butylbenzylphthalate	13.46	149	363148	37.173	nq	93
80) Benzo(a)anthracene	14.05	228	582113	37.830	nq	100
81) 3,3'-Dichlorobenzidine	14.00	252	214759	39.107	nq	99
82) Chrysene	14.08	228	551694	36.925	nq	99
83) Bis(2-ethylhexyl)phthalate	14.01	149	491891	37.312	nq	# 99
84) Di-n-octyl phthalate	14.63	149	903760	42.430	nq	96
85) Indeno(1,2,3-cd)pyrene	17.08	276	460774	38.955	nq	97
87) Benzo(b)fluoranthene	15.12	252	595957	37.942	nq	97
88) Benzo(k)fluoranthene	15.14	252	605260	40.922	nq	98
89) Benzo(a)pyrene	15.49	252	556401	39.668	nq	98
90) Dibenzo(a,h)anthracene	17.09	278	401318	37.028	nq	98
91) Benzo(g,h,i)perylene	17.54	276	364906	34.449	nq	99

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

