

Data Path : Z:\HPCHEM1\BNA F\DATA\BF030218\
 Data File : BF103393.D
 Acq On : 3 Mar 2018 4:38
 Operator : SJ/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 3/5/2018 3:14:22 PM

Quant Time: Mar 05 00:18:17 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 02 13:28:09 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.88	152	146634	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	592502	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	227881	20.00	ng	0.00
63) Phenanthrene-d10	11.42	188	332404	20.00	ng	0.00
75) Chrysene-d12	14.07	240	242106	20.00	ng	0.00
86) Perylene-d12	15.59	264	178893	20.00	ng	0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.50	112	775705	80.01	ng	0.00
7) Phenol-d6	6.52	99	919596	77.51	ng	0.00
23) Nitrobenzene-d5	7.45	82	829720	79.97	ng	0.00
41) 2,4,6-Tribromophenol	10.72	330	115787	60.03	ng	0.00
44) 2-Fluorobiphenyl	9.24	172	1097773	82.36	ng	0.00
78) Terphenyl-d14	13.00	244	737763	73.25	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.61	88	205394	40.111	ng	93
3) Pyridine	3.41	79	554476	40.559	ng	99
4) n-Nitrosodimethylamine	3.36	42	249835	45.314	ng	# 89
6) Aniline	6.55	93	663752	38.766	ng	91
8) 2-Chlorophenol	6.67	128	391509	38.872	ng	93
9) Benzaldehyde	6.43	77	345558	52.836	ng	97
10) Phenol	6.53	94	526450	38.225	ng	83
11) bis(2-Chloroethyl)ether	6.62	93	422495	40.419	ng	91
12) 1,3-Dichlorobenzene	6.82	146	437240	37.989	ng	98
13) 1,4-Dichlorobenzene	6.90	146	445073	38.570	ng	99
14) 1,2-Dichlorobenzene	7.05	146	397591	37.366	ng	98
15) Benzyl Alcohol	7.03	79	322360	36.059	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.15	45	668956	37.268	ng	86
17) 2-Methylphenol	7.14	107	342828	37.456	ng	# 91
18) Hexachloroethane	7.39	117	89531	21.313	ng	93
19) n-Nitroso-di-n-propylamine	7.29	70	280386	37.975	ng	94
20) 3+4-Methylphenols	7.29	107	411989	36.926	ng	# 58
22) Acetophenone	7.29	105	539872	39.036	ng	# 88
24) Nitrobenzene	7.47	77	463476	40.574	ng	95
25) Isophorone	7.70	82	802033	39.250	ng	98
26) 2-Nitrophenol	7.78	139	146872	27.312	ng	92
27) 2,4-Dimethylphenol	7.82	122	318138	38.704	ng	99
28) bis(2-Chloroethoxy)methane	7.91	93	485358	40.400	ng	99
29) 2,4-Dichlorophenol	8.03	162	307191	39.035	ng	99
30) 1,2,4-Trichlorobenzene	8.10	180	316160	37.727	ng	97
31) Naphthalene	8.19	128	1073209	36.997	ng	99
32) Benzoic acid	7.94	122	113066m	22.784	ng	
33) 4-Chloroaniline	8.24	127	477199	38.123	ng	97
34) Hexachlorobutadiene	8.29	225	161614	37.034	ng	99
35) Caprolactam	8.62	113	100698	36.408	ng	# 77
36) 4-Chloro-3-methylphenol	8.72	107	325412	36.661	ng	97
37) 2-Methylnaphthalene	8.88	142	667457	35.603	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.05	216	252227	40.487	ng	98
42) 2,4,6-Trichlorophenol	9.16	196	168629	40.227	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) 2,4,5-Trichlorophenol	9.21	196	176502	36.609	nq	94
45) 1,1'-Biphenyl	9.34	154	766072	40.118	nq	99
46) 2-Chloronaphthalene	9.37	162	587938	38.918	nq	100
47) 2-Nitroaniline	9.47	65	266769	47.672	nq	# 81
48) Acenaphthylene	9.79	152	871969	37.514	nq	100
49) Dimethylphthalate	9.63	163	715774	39.154	nq	99
50) 2,6-Dinitrotoluene	9.71	165	118556	30.904	nq	# 81
51) Acenaphthene	9.96	154	495700	36.573	nq	99
52) 3-Nitroaniline	9.89	138	209549	43.053	nq	99
54) Dibenzofuran	10.13	168	677868	36.434	nq	93
55) 4-Nitrophenol	10.06	139	92286	30.593	nq	88
56) 2,4-Dinitrotoluene	10.12	165	123651	25.485	nq	# 73
57) Fluorene	10.47	166	528245	33.329	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.25	232	107123	33.075	nq	93
59) Diethylphthalate	10.33	149	686891	39.286	nq	100
60) 4-Chlorophenyl-phenylether	10.46	204	248176	36.924	nq	96
61) 4-Nitroaniline	10.50	138	207536	42.690	nq	96
62) Azobenzene	10.62	77	655978	36.859	nq	95
65) n-Nitrosodiphenylamine	10.58	169	473619	40.922	nq	100
66) 4-Bromophenyl-phenylether	10.95	248	135752	39.205	nq	97
67) Hexachlorobenzene	11.03	284	137011	36.455	nq	99
68) Atrazine	11.10	200	111126	31.944	nq	98
69) Pentachlorophenol	11.23	266	63187	34.771	nq	100
70) Phenanthrene	11.44	178	666095	36.412	nq	98
71) Anthracene	11.49	178	696053	37.106	nq	99
72) Carbazole	11.65	167	712157	38.123	nq	99
73) Di-n-butylphthalate	11.96	149	946266	40.482	nq	99
74) Fluoranthene	12.63	202	679503	36.964	nq	98
76) Benzidine	12.76	184	487609	53.872	nq	97
77) Pyrene	12.87	202	724993	38.408	nq	99
79) Butylbenzylphthalate	13.47	149	402268	40.336	nq	99
80) Benzo(a)anthracene	14.06	228	586797	37.355	nq	100
81) 3,3'-Dichlorobenzidine	14.02	252	230164	41.056	nq	98
82) Chrysene	14.10	228	565638	37.084	nq	99
83) Bis(2-ethylhexyl)phthalate	14.02	149	517349	38.441	nq	# 98
84) Di-n-octyl phthalate	14.65	149	918047	42.220	nq	96
85) Indeno(1,2,3-cd)pyrene	17.14	276	387171m	32.063	nq	
87) Benzo(b)fluoranthene	15.14	252	444418	37.784	nq	99
88) Benzo(k)fluoranthene	15.17	252	436395	39.401	nq	99
89) Benzo(a)pyrene	15.53	252	392823	37.399	nq	98
90) Dibenzo(a,h)anthracene	17.14	278	324289	39.956	nq	98
91) Benzo(g,h,i)perylene	17.60	276	322283	40.629	nq	99

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

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