

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF110917\
 Data File : BF100370.D
 Acq On : 10 Nov 2017 5:35
 Operator : SJ/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 11/10/2017 1:31:05 PM

Quant Time: Nov 10 06:14:48 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 09 16:04:38 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	158688	20.00	ng	0.00
21) Naphthalene-d8	8.19	136	635085	20.00	ng	0.00
38) Acenaphthene-d10	9.95	164	283085	20.00	ng	0.00
63) Phenanthrene-d10	11.43	188	474428	20.00	ng	0.00
75) Chrysene-d12	14.07	240	298189	20.00	ng	0.00
86) Perylene-d12	15.55	264	319123	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.51	112	792066	80.01	ng	0.00
7) Phenol-d6	6.53	99	976754	80.01	ng	0.00
23) Nitrobenzene-d5	7.47	82	891955	92.85	ng	0.00
41) 2,4,6-Tribromophenol	10.73	330	226311	78.45	ng	0.00
44) 2-Fluorobiphenyl	9.26	172	1528306	82.21	ng	0.00
78) Terphenyl-d14	13.01	244	1132220	87.24	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.69	88	195958	40.619	ng	96
3) Pyridine	3.47	79	526028	39.966	ng	95
4) n-Nitrosodimethylamine	3.42	42	257469	40.290	ng	97
6) Aniline	6.57	93	689208	39.963	ng	100
8) 2-Chlorophenol	6.69	128	432737	41.320	ng	99
9) Benzaldehyde	6.46	77	352134	40.392	ng	96
10) Phenol	6.54	94	519949	40.573	ng	100
11) bis(2-Chloroethyl)ether	6.64	93	439226	40.804	ng	94
12) 1,3-Dichlorobenzene	6.85	146	488852	40.487	ng	97
13) 1,4-Dichlorobenzene	6.92	146	486568	39.815	ng	97
14) 1,2-Dichlorobenzene	7.07	146	452856	40.079	ng	97
15) Benzyl Alcohol	7.04	79	390029	40.938	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.17	45	585175m	33.990	ng	
17) 2-Methylphenol	7.15	107	365326	40.820	ng	99
18) Hexachloroethane	7.42	117	165256	41.013	ng	97
19) n-Nitroso-di-n-propylamine	7.32	70	333194	39.357	ng	93
20) 3+4-Methylphenols	7.30	107	448337	39.588	ng	# 85
22) Acetophenone	7.32	105	601465	38.838	ng	# 98
24) Nitrobenzene	7.49	77	460974	46.624	ng	99
25) Isophorone	7.72	82	809591	40.106	ng	98
26) 2-Nitrophenol	7.80	139	231592	49.740	ng	97
27) 2,4-Dimethylphenol	7.83	122	373855	41.314	ng	99
28) bis(2-Chloroethoxy)methane	7.93	93	538983	40.819	ng	99
29) 2,4-Dichlorophenol	8.04	162	360334	41.712	ng	95
30) 1,2,4-Trichlorobenzene	8.13	180	379983	40.664	ng	96
31) Naphthalene	8.21	128	1211008	39.590	ng	100
32) Benzoic acid	7.94	122	316926m	46.178	ng	
33) 4-Chloroaniline	8.25	127	518081	40.689	ng	98
34) Hexachlorobutadiene	8.32	225	228620	41.149	ng	99
35) Caprolactam	8.63	113	115875	39.086	ng	96
36) 4-Chloro-3-methylphenol	8.73	107	368612	39.730	ng	99
37) 2-Methylnaphthalene	8.90	142	791277	38.964	ng	100
39) 1,2,4,5-Tetrachlorobenzene	9.06	216	409822	43.651	ng	97
40) Hexachlorocyclopentadiene	9.05	237	104722	23.700	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.17	196	257815	43.328	ng	97
43) 2,4,5-Trichlorophenol	9.21	196	270775	43.922	ng	93
45) 1,1'-Biphenyl	9.36	154	1012721	41.363	ng	99
46) 2-Chloronaphthalene	9.39	162	737823	41.539	ng	98
47) 2-Nitroaniline	9.48	65	248111	47.315	ng	92
48) Acenaphthylene	9.80	152	1178268	40.028	ng	99
49) Dimethylphthalate	9.66	163	899952	41.638	ng	100
50) 2,6-Dinitrotoluene	9.72	165	191478	44.755	ng	97
51) Acenaphthene	9.98	154	700395	39.123	ng	99
52) 3-Nitroaniline	9.89	138	221096	43.763	ng	99
53) 2,4-Dinitrophenol	9.99	184	36815	31.123	ng	# 18
54) Dibenzofuran	10.15	168	1000064	39.857	ng	99
55) 4-Nitrophenol	10.04	139	157526	38.400	ng	94
56) 2,4-Dinitrotoluene	10.12	165	246446	45.661	ng	98
57) Fluorene	10.49	166	734150	39.940	ng	99
58) 2,3,4,6-Tetrachlorophenol	10.26	232	200888	39.759	ng	88
59) Diethylphthalate	10.36	149	883077	40.827	ng	95
60) 4-Chlorophenyl-phenylether	10.48	204	371529	41.203	ng	95
61) 4-Nitroaniline	10.50	138	196050	40.925	ng	97
62) Azobenzene	10.64	77	802411	39.389	ng	96
64) 4,6-Dinitro-2-methylphenol	10.53	198	71149	32.969	ng	94
65) n-Nitrosodiphenylamine	10.60	169	740806	44.907	ng	96
66) 4-Bromophenyl-phenylether	10.97	248	232410	45.698	ng	# 83
67) Hexachlorobenzene	11.04	284	233130	43.828	ng	# 85
68) Atrazine	11.12	200	227478	45.555	ng	98
69) Pentachlorophenol	11.23	266	150209	39.382	ng	98
70) Phenanthrene	11.45	178	1005585	40.257	ng	99
71) Anthracene	11.50	178	1018036	40.171	ng	99
72) Carbazole	11.66	167	884360	37.819	ng	98
73) Di-n-butylphthalate	11.99	149	1264200	46.198	ng	99
74) Fluoranthene	12.64	202	887093	33.509	ng	98
76) Benzidine	12.76	184	439232	36.781	ng	99
77) Pyrene	12.87	202	871881	41.018	ng	99
79) Butylbenzylphthalate	13.49	149	406927	46.943	ng	94
80) Benzo(a)anthracene	14.06	228	742474	41.348	ng	99
81) 3,3'-Dichlorobenzidine	14.02	252	281823	43.804	ng	98
82) Chrysene	14.09	228	707336	40.678	ng	100
83) Bis(2-ethylhexyl)phthalate	14.04	149	565567	49.606	ng	99
84) Di-n-octyl phthalate	14.66	149	927124	47.335	ng	100
85) Indeno(1,2,3-cd)pyrene	17.04	276	665562	47.321	ng	96
87) Benzo(b)fluoranthene	15.12	252	741088	39.198	ng	99
88) Benzo(k)fluoranthene	15.14	252	741046	41.483	ng	98
89) Benzo(a)pyrene	15.49	252	695243	41.479	ng	99
90) Dibenzo(a,h)anthracene	17.06	278	573476	41.837	ng	98
91) Benzo(g,h,i)perylene	17.50	276	518708	38.657	ng	95

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

