

Data Path : Z:\HPCHEM1\BNA F\DATA\BF072517\
 Data File : BF097019.D
 Acq On : 25 Jul 2017 10:42
 Operator : SJ/JU
 Sample : SSTDICC025
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC025

Quant Time: Jul 25 13:20:15 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 25 13:17:51 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.50	152	138998	20.00	ng	0.00
21) Naphthalene-d8	7.79	136	593761	20.00	ng	0.00
38) Acenaphthene-d10	9.53	164	253968	20.00	ng	0.00
63) Phenanthrene-d10	11.00	188	410032	20.00	ng	0.00
75) Chrysene-d12	13.63	240	304003	20.00	ng	0.00
86) Perylene-d12	14.97	264	267340	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.10	112	482546	52.20	ng	0.00
7) Phenol-d6	6.16	99	543238	48.97	ng	0.00
23) Nitrobenzene-d5	7.07	82	524530	54.72	ng	0.00
41) 2,4,6-Tribromophenol	10.32	330	105467	48.65	ng	0.00
44) 2-Fluorobiphenyl	8.86	172	814626	50.68	ng	0.00
78) Terphenyl-d14	12.59	244	697421	39.79	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.05	88	96996	20.391	ng	# 74
3) Pyridine	2.66	79	280709	28.080	ng	# 64
4) n-Nitrosodimethylamine	2.61	42	155363	27.791	ng	# 83
6) Aniline	6.16	93	340782	24.939	ng	# 76
8) 2-Chlorophenol	6.29	128	267845	26.857	ng	# 92
9) Benzaldehyde	6.04	77	173757	27.099	ng	# 94
10) Phenol	6.17	94	327148	26.087	ng	# 96
11) bis(2-Chloroethyl)ether	6.25	93	269423	28.569	ng	# 94
12) 1,3-Dichlorobenzene	6.44	146	279760	26.390	ng	# 99
13) 1,4-Dichlorobenzene	6.52	146	278131	25.907	ng	# 96
14) 1,2-Dichlorobenzene	6.67	146	255716	26.188	ng	# 98
15) Benzyl Alcohol	6.66	79	181693	25.016	ng	# 98
16) 2,2'-oxybis(1-Chloropropan	6.79	45	528907	27.176	ng	# 97
17) 2-Methylphenol	6.78	107	198923	25.650	ng	# 93
18) Hexachloroethane	7.01	117	102385	26.895	ng	# 81
19) n-Nitroso-di-n-propylamine	6.93	70	194272	29.068	ng	# 83
20) 3+4-Methylphenols	6.94	107	279933	29.746	ng	# 63
22) Acetophenone	6.92	105	374650	28.231	ng	# 98
24) Nitrobenzene	7.09	77	273819	27.619	ng	# 95
25) Isophorone	7.33	82	510583	28.632	ng	# 97
26) 2-Nitrophenol	7.40	139	144917	26.288	ng	# 86
27) 2,4-Dimethylphenol	7.46	122	234458	25.853	ng	# 96
28) bis(2-Chloroethoxy)methane	7.55	93	323166	26.819	ng	# 100
29) 2,4-Dichlorophenol	7.66	162	208716	25.425	ng	# 98
30) 1,2,4-Trichlorobenzene	7.73	180	196693	25.200	ng	# 98
31) Naphthalene	7.80	128	716201	25.462	ng	# 100
32) Benzoic acid	7.61	122	180060	31.064	ng	# 95
33) 4-Chloroaniline	7.86	127	286775	25.733	ng	# 98
34) Hexachlorobutadiene	7.93	225	100768	24.184	ng	# 99
35) Caprolactam	8.23	113	62741	23.853	ng	# 59
36) 4-Chloro-3-methylphenol	8.36	107	214159	24.262	ng	# 90
37) 2-Methylnaphthalene	8.50	142	453510	25.290	ng	# 98
39) 1,2,4,5-Tetrachlorobenzene	8.66	216	176651	25.731	ng	# 99
40) Hexachlorocyclopentadiene	8.65	237	43721	19.162	ng	# 99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	8.79	196	125654	24.813	nq	99
43) 2,4,5-Trichlorophenol	8.83	196	127246	24.914	nq	96
45) 1,1'-Biphenyl	8.96	154	559438	25.684	nq	98
46) 2-Chloronaphthalene	8.98	162	399351	24.899	nq	# 93
47) 2-Nitroaniline	9.08	65	140508	25.486	nq	98
48) Acenaphthylene	9.39	152	650518	25.455	nq	99
49) Dimethylphthalate	9.27	163	475367	24.858	nq	100
50) 2,6-Dinitrotoluene	9.33	165	103172	24.909	nq	# 80
51) Acenaphthene	9.56	154	367733	24.359	nq	100
52) 3-Nitroaniline	9.49	138	126146	25.708	nq	94
53) 2,4-Dinitrophenol	9.61	184	43808	25.142	nq	# 76
54) Dibenzofuran	9.73	168	500813	23.673	nq	98
55) 4-Nitrophenol	9.68	139	80789	22.536	nq	# 73
56) 2,4-Dinitrotoluene	9.73	165	122547	23.024	nq	# 64
57) Fluorene	10.07	166	401762	25.700	nq	98
58) 2,3,4,6-Tetrachlorophenol	9.86	232	89502	25.268	nq	# 99
59) Diethylphthalate	9.96	149	453965	24.385	nq	99
60) 4-Chlorophenyl-phenylether	10.07	204	167869	25.540	nq	98
61) 4-Nitroaniline	10.10	138	129005	27.997	nq	92
62) Azobenzene	10.23	77	423857	25.809	nq	94
64) 4,6-Dinitro-2-methylphenol	10.14	198	70406	27.348	nq	91
65) n-Nitrosodiphenylamine	10.19	169	366362	24.996	nq	98
66) 4-Bromophenyl-phenylether	10.56	248	108055	25.812	nq	97
67) Hexachlorobenzene	10.62	284	118046	25.318	nq	# 86
68) Atrazine	10.72	200	107582	25.763	nq	92
69) Pentachlorophenol	10.82	266	47195	20.953	nq	99
70) Phenanthrene	11.03	178	589420	26.438	nq	99
71) Anthracene	11.07	178	599892	26.612	nq	99
72) Carbazole	11.24	167	567975	26.019	nq	98
73) Di-n-butylphthalate	11.58	149	730171	26.858	nq	98
74) Fluoranthene	12.20	202	570728	26.744	nq	96
76) Benzidine	12.33	184	285329	28.650	nq	# 94
77) Pyrene	12.43	202	585335	21.216	nq	97
79) Butylbenzylphthalate	13.06	149	305255	23.286	nq	# 79
80) Benzo(a)anthracene	13.62	228	474471	25.049	nq	98
81) 3,3'-Dichlorobenzidine	13.59	252	182099	26.734	nq	# 96
82) Chrysene	13.65	228	449770	24.120	nq	99
83) Bis(2-ethylhexyl)phthalate	13.63	149	393081	24.623	nq	99
84) Di-n-octyl phthalate	14.24	149	726816	27.543	nq	# 92
85) Indeno(1,2,3-cd)pyrene	16.20	276	370574	21.896	nq	97
87) Benzo(b)fluoranthene	14.62	252	449407	27.873	nq	99
88) Benzo(k)fluoranthene	14.64	252	396547	25.972	nq	97
89) Benzo(a)pyrene	14.92	252	385169	26.046	nq	99
90) Dibenzo(a,h)anthracene	16.21	278	295786	21.737	nq	97
91) Benzo(g,h,i)perylene	16.56	276	306591	21.040	nq	99

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

