

Data Path : Z:\HPCHEM1\BNA F\DATA\BF052616\
 Data File : BF087430.D
 Acq On : 27 May 2016 4:21
 Operator : UM/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

sohil
 5/27/2016 7:38:11 PM

Quant Time: May 27 07:44:42 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF052316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 25 16:26:52 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.51	152	152602	20.00	ng	-0.01
21) Naphthalene-d8	9.54	136	631392	20.00	ng	-0.01
38) Acenaphthene-d10	12.38	164	289379	20.00	ng	-0.01
63) Phenanthrene-d10	14.78	188	549198	20.00	ng	0.00
75) Chrysene-d12	18.47	240	387526	20.00	ng	-0.01
86) Perylene-d12	20.14	264	314621	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.40	112	747269	86.05	ng	-0.01
7) Phenol-d6	7.07	99	998982	85.13	ng	0.00
23) Nitrobenzene-d5	8.43	82	1059549	84.57	ng	-0.01
41) 2,4,6-Tribromophenol	13.67	330	215542	77.51	ng	-0.01
44) 2-Fluorobiphenyl	11.32	172	1538869	74.97	ng	-0.01
78) Terphenyl-d14	17.12	244	1450127	79.56	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	1.82	88	186249	40.64	ng	# 73
3) Pyridine	2.40	79	400645	39.99	ng	# 68
4) n-Nitrosodimethylamine	2.38	42	344590	44.64	ng	# 45
6) Aniline	7.02	93	639049	42.10	ng	93
8) 2-Chlorophenol	7.20	128	444795	41.07	ng	96
9) Benzaldehyde	6.82	77	247051	38.13	ng	94
10) Phenol	7.08	94	543558	42.42	ng	78
11) bis(2-Chloroethyl)ether	7.15	93	417009	40.21	ng	85
12) 1,3-Dichlorobenzene	7.40	146	473298m	40.07	ng	
13) 1,4-Dichlorobenzene	7.54	146	488911	40.31	ng	96
14) 1,2-Dichlorobenzene	7.77	146	456035	40.65	ng	98
15) Benzyl Alcohol	7.82	79	367764	42.99	ng	90
16) 2,2'-oxybis(1-Chloropropan	8.00	45	893644	38.29	ng	86
17) 2-Methylphenol	8.02	107	362640	41.78	ng	# 83
18) Hexachloroethane	8.28	117	184684	40.82	ng	96
19) n-Nitroso-di-n-propylamine	8.24	70	343653	39.27	ng	# 70
20) 3+4-Methylphenols	8.28	107	447987	42.70	ng	# 59
22) Acetophenone	8.20	105	632515	41.93	ng	# 97
24) Nitrobenzene	8.47	77	557896	42.19	ng	96
25) Isophorone	8.86	82	892265	39.71	ng	99
26) 2-Nitrophenol	8.97	139	232336	42.98	ng	# 80
27) 2,4-Dimethylphenol	9.11	122	375535	40.02	ng	# 84
28) bis(2-Chloroethoxy)methane	9.23	93	486221	38.42	ng	98
29) 2,4-Dichlorophenol	9.38	162	326523	38.61	ng	97
30) 1,2,4-Trichlorobenzene	9.47	180	386833	39.96	ng	97
31) Naphthalene	9.58	128	1240953	39.22	ng	99
32) Benzoic acid	9.46	122	166940	36.18	ng	# 81
33) 4-Chloroaniline	9.72	127	480025	41.19	ng	# 94
34) Hexachlorobutadiene	9.79	225	228677	38.86	ng	98
35) Caprolactam	10.38	113	109332m	44.30	ng	
36) 4-Chloro-3-methylphenol	10.59	107	417675	43.68	ng	92
37) 2-Methylnaphthalene	10.71	142	788442	39.89	ng	98
39) 1,2,4,5-Tetrachlorobenzene	10.98	216	360793	38.95	ng	99
40) Hexachlorocyclopentadiene	10.95	237	97894	32.10	ng	95

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	11.21	196	210574	39.38	nq	93
43) 2,4,5-Trichlorophenol	11.29	196	212645	39.05	nq	# 90
45) 1,1'-Biphenyl	11.47	154	958698	39.35	nq	97
46) 2-Chloronaphthalene	11.48	162	730428	39.19	nq	# 91
47) 2-Nitroaniline	11.71	65	300788	44.79	nq	85
48) Acenaphthylene	12.15	152	1188347	40.27	nq	99
49) Dimethylphthalate	12.03	163	913540	39.41	nq	99
50) 2,6-Dinitrotoluene	12.12	165	188753	40.41	nq	# 80
51) Acenaphthene	12.42	154	722448	39.83	nq	98
52) 3-Nitroaniline	12.39	138	201857	42.11	nq	# 77
53) 2,4-Dinitrophenol	12.60	184	65701	31.78	nq	# 81
54) Dibenzofuran	12.71	168	958478	39.03	nq	92
55) 4-Nitrophenol	12.80	139	75428	32.98	nq	# 49
56) 2,4-Dinitrotoluene	12.79	165	269087	43.10	nq	93
57) Fluorene	13.27	166	835986	39.26	nq	99
58) 2,3,4,6-Tetrachlorophenol	12.95	232	152687	36.45	nq	95
59) Diethylphthalate	13.18	149	871565	38.67	nq	99
60) 4-Chlorophenyl-phenylether	13.29	204	390199	37.57	nq	99
61) 4-Nitroaniline	13.41	138	187256	41.85	nq	# 71
62) Azobenzene	13.55	77	969191	41.45	nq	86
64) 4,6-Dinitro-2-methylphenol	13.45	198	133795	38.31	nq	# 69
65) n-Nitrosodiphenylamine	13.51	169	672843	37.88	nq	100
66) 4-Bromophenyl-phenylether	14.08	248	229121	38.13	nq	93
67) Hexachlorobenzene	14.15	284	240931	38.34	nq	# 73
68) Atrazine	14.42	200	175983	31.08	nq	93
69) Pentachlorophenol	14.51	266	58112	33.64	nq	98
70) Phenanthrene	14.81	178	1166548	38.35	nq	100
71) Anthracene	14.89	178	1214794	39.53	nq	100
72) Carbazole	15.19	167	1060005	40.44	nq	99
73) Di-n-butylphthalate	15.76	149	1225537	36.16	nq	# 98
74) Fluoranthene	16.57	202	1200277	39.35	nq	96
76) Benzidine	16.80	184	423972	42.52	nq	97
77) Pyrene	16.88	202	1204215	43.17	nq	99
79) Butylbenzylphthalate	17.78	149	474430	38.35	nq	# 71
80) Benzo(a)anthracene	18.46	228	858826	38.96	nq	99
81) 3,3'-Dichlorobenzidine	18.45	252	490052	40.24	nq	# 98
82) Chrysene	18.50	228	844393	38.60	nq	99
83) Bis(2-ethylhexyl)phthalate	18.53	149	604978	33.52	nq	# 95
84) Di-n-octyl phthalate	19.30	149	883903	32.11	nq	# 85
85) Indeno(1,2,3-cd)pyrene	21.36	276	802151	45.74	nq	94
87) Benzo(b)fluoranthene	19.73	252	767020	38.27	nq	98
88) Benzo(k)fluoranthene	19.76	252	652734m	37.84	nq	
89) Benzo(a)pyrene	20.08	252	674795	38.93	nq	99
90) Dibenzo(a,h)anthracene	21.37	278	664226	44.93	nq	# 92
91) Benzo(g,h,i)perylene	21.71	276	676421	46.37	nq	# 91

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

