

Data Path : Z:\HPCHEM1\BNA F\DATA\BF082416\
 Data File : BF089865.D
 Acq On : 24 Aug 2016 11:03
 Operator : UM/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

umangi
 8/25/2016 6:48:50 PM

Quant Time: Aug 25 13:09:58 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 19 14:02:57 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.67	152	464240	20.00	ng	-0.01
21) Naphthalene-d8	7.95	136	1814249	20.00	ng	-0.02
38) Acenaphthene-d10	9.71	164	862897	20.00	ng	0.00
63) Phenanthrene-d10	11.19	188	1601678	20.00	ng	0.00
75) Chrysene-d12	13.87	240	1301776	20.00	ng	0.00
86) Perylene-d12	15.37	264	1135876	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.23	112	2088575	77.15	ng	-0.02
7) Phenol-d6	6.31	99	2668354	80.30	ng	-0.01
23) Nitrobenzene-d5	7.24	82	2441224	83.59	ng	-0.01
41) 2,4,6-Tribromophenol	10.49	330	626795	76.43	ng	-0.01
44) 2-Fluorobiphenyl	9.04	172	3569575	72.71	ng	-0.01
78) Terphenyl-d14	12.78	244	3376580	58.68	ng	-0.01

Target Compounds						Qvalue
2) 1,4-Dioxane	2.10	88	548631	40.95	ng	93
3) Pyridine	2.74	79	1498825	42.66	ng	91
4) n-Nitrosodimethylamine	2.70	42	539650	41.25	ng	# 87
6) Aniline	6.33	93	1884273	42.27	ng	99
8) 2-Chlorophenol	6.44	128	1242736	38.39	ng	97
9) Benzaldehyde	6.20	77	819542	37.86	ng	91
10) Phenol	6.32	94	1496040	38.41	ng	98
11) bis(2-Chloroethyl)ether	6.41	93	1212324	40.14	ng	99
12) 1,3-Dichlorobenzene	6.60	146	1326189	39.19	ng	99
13) 1,4-Dichlorobenzene	6.68	146	1398091	40.19	ng	99
14) 1,2-Dichlorobenzene	6.84	146	1344929	41.42	ng	99
15) Benzyl Alcohol	6.82	79	954857	43.33	ng	# 87
16) 2,2'-oxybis(1-Chloropropan	6.96	45	1489090	38.52	ng	92
17) 2-Methylphenol	6.94	107	1071517	42.26	ng	96
18) Hexachloroethane	7.19	117	523531	42.19	ng	95
19) n-Nitroso-di-n-propylamine	7.10	70	775475	36.49	ng	91
20) 3+4-Methylphenols	7.10	107	1410771	45.38	ng	# 51
22) Acetophenone	7.08	105	1616859	38.81	ng	# 89
24) Nitrobenzene	7.26	77	1250537	38.17	ng	98
25) Isophorone	7.50	82	2346880	39.85	ng	99
26) 2-Nitrophenol	7.58	139	778182	47.60	ng	# 91
27) 2,4-Dimethylphenol	7.62	122	1221754	41.47	ng	97
28) bis(2-Chloroethoxy)methane	7.72	93	1545350	42.41	ng	# 95
29) 2,4-Dichlorophenol	7.82	162	1059499	42.86	ng	96
30) 1,2,4-Trichlorobenzene	7.90	180	1097954	40.15	ng	100
31) Naphthalene	7.98	128	3218409	37.96	ng	99
32) Benzoic acid	7.75	122	790137	34.41	ng	94
33) 4-Chloroaniline	8.03	127	1664388	45.28	ng	# 94
34) Hexachlorobutadiene	8.10	225	576476	40.26	ng	99
35) Caprolactam	8.41	113	300543	39.92	ng	95
36) 4-Chloro-3-methylphenol	8.51	107	1071952	40.04	ng	89
37) 2-Methylnaphthalene	8.67	142	2438993	44.46	ng	99
39) 1,2,4,5-Tetrachlorobenzene	8.83	216	1109962	39.72	ng	# 97
40) Hexachlorocyclopentadiene	8.83	237	438640	45.21	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	8.95	196	704945	40.10	nq	99
43) 2,4,5-Trichlorophenol	8.99	196	758720	43.40	nq	# 93
45) 1,1'-Biphenyl	9.14	154	2835403	42.46	nq	94
46) 2-Chloronaphthalene	9.15	162	2024345	38.35	nq	99
47) 2-Nitroaniline	9.26	65	682753	45.35	nq	# 76
48) Acenaphthylene	9.56	152	3310224	41.73	nq	99
49) Dimethylphthalate	9.45	163	2610105	42.96	nq	100
50) 2,6-Dinitrotoluene	9.51	165	585872	41.77	nq	# 59
51) Acenaphthene	9.75	154	2232995	42.54	nq	99
52) 3-Nitroaniline	9.67	138	723969	46.59	nq	90
53) 2,4-Dinitrophenol	9.77	184	197085	32.88	nq	# 65
54) Dibenzofuran	9.92	168	2834450	42.99	nq	# 88
55) 4-Nitrophenol	9.83	139	558433	44.02	nq	96
56) 2,4-Dinitrotoluene	9.90	165	651225	35.78	nq	# 69
57) Fluorene	10.25	166	2129198	40.52	nq	98
58) 2,3,4,6-Tetrachlorophenol	10.03	232	538517	42.32	nq	96
59) Diethylphthalate	10.15	149	2498045	40.47	nq	97
60) 4-Chlorophenyl-phenylether	10.25	204	872144	36.39	nq	93
61) 4-Nitroaniline	10.27	138	693936	47.63	nq	95
62) Azobenzene	10.41	77	2219292	38.25	nq	94
64) 4,6-Dinitro-2-methylphenol	10.31	198	364869	39.64	nq	# 72
65) n-Nitrosodiphenylamine	10.38	169	2125385	42.20	nq	100
66) 4-Bromophenyl-phenylether	10.74	248	660168	39.31	nq	# 84
67) Hexachlorobenzene	10.80	284	693751	36.21	nq	# 82
68) Atrazine	10.90	200	565302	36.65	nq	97
69) Pentachlorophenol	10.99	266	429499	39.55	nq	99
70) Phenanthrene	11.21	178	3178607	39.29	nq	97
71) Anthracene	11.26	178	3223889	39.20	nq	99
72) Carbazole	11.42	167	2801880	37.99	nq	98
73) Di-n-butylphthalate	11.77	149	4022131	43.59	nq	# 98
74) Fluoranthene	12.39	202	3164480	41.12	nq	93
76) Benzidine	12.53	184	1777424	42.19	nq	97
77) Pyrene	12.62	202	3189949	30.09	nq	98
79) Butylbenzylphthalate	13.28	149	1886502	39.40	nq	95
80) Benzo(a)anthracene	13.85	228	2961879	39.73	nq	99
81) 3,3'-Dichlorobenzidine	13.83	252	1107396	45.31	nq	96
82) Chrysene	13.90	228	2485689	38.89	nq	99
83) Bis(2-ethylhexyl)phthalate	13.89	149	2355528	35.72	nq	98
84) Di-n-octyl phthalate	14.57	149	4028368	45.97	nq	# 94
85) Indeno(1,2,3-cd)pyrene	16.66	276	2575957	31.59	nq	100
87) Benzo(b)fluoranthene	14.98	252	2528425	40.46	nq	99
88) Benzo(k)fluoranthene	15.02	252	2630673m	46.40	nq	
89) Benzo(a)pyrene	15.33	252	2426620	41.03	nq	# 94
90) Dibenzo(a,h)anthracene	16.69	278	2153395	33.54	nq	97
91) Benzo(g,h,i)perylene	17.05	276	2184893	32.32	nq	100

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

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