

Data Path : Z:\HPCHEM1\BNA F\DATA\BF080416\
 Data File : BF089605.D
 Acq On : 5 Aug 2016 00:45
 Operator : UM/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040EC

Manual Integrations
 APPROVED

sohil
 8/5/2016 6:58:55 PM

Quant Time: Aug 05 02:06:52 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF080416.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 05 01:39:23 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.45	152	47981	20.00	ng	0.00
21) Naphthalene-d8	9.47	136	202919	20.00	ng	0.00
38) Acenaphthene-d10	12.30	164	94942	20.00	ng	0.00
63) Phenanthrene-d10	14.69	188	179743	20.00	ng	-0.01
75) Chrysene-d12	18.39	240	103550	20.00	ng	0.00
86) Perylene-d12	20.05	264	80095	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.31	112	224432	78.40	ng	0.00
7) Phenol-d6	6.98	99	301318	77.59	ng	0.00
23) Nitrobenzene-d5	8.36	82	292547	79.74	ng	0.00
41) 2,4,6-Tribromophenol	13.59	330	62843	80.21	ng	0.00
44) 2-Fluorobiphenyl	11.26	172	556337	76.16	ng	0.00
78) Terphenyl-d14	17.05	244	417721	79.64	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	1.74	88	58287	35.52	ng	96
3) Pyridine	2.31	79	116158	38.52	ng	98
4) n-Nitrosodimethylamine	2.28	42	46106	36.62	ng	97
6) Aniline	6.94	93	202997	40.34	ng	99
8) 2-Chlorophenol	7.12	128	137775	39.20	ng	99
9) Benzaldehyde	6.75	77	59624	42.32	ng	97
10) Phenol	7.00	94	137939	34.60	ng	100
11) bis(2-Chloroethyl)ether	7.10	93	130434	38.08	ng	98
12) 1,3-Dichlorobenzene	7.35	146	145874	38.22	ng	98
13) 1,4-Dichlorobenzene	7.47	146	145918	38.33	ng	100
14) 1,2-Dichlorobenzene	7.70	146	147118	36.66	ng	99
15) Benzyl Alcohol	7.74	79	96502	37.90	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.94	45	162353	37.22	ng	95
17) 2-Methylphenol	7.94	107	98075	38.64	ng	95
18) Hexachloroethane	8.23	117	59386	37.01	ng	97
19) n-Nitroso-di-n-propylamine	8.17	70	108212	38.57	ng	99
20) 3+4-Methylphenols	8.20	107	126242	39.40	ng	99
22) Acetophenone	8.14	105	207596	42.91	ng	99
24) Nitrobenzene	8.39	77	143751	41.25	ng	99
25) Isophorone	8.79	82	268723	38.25	ng	99
26) 2-Nitrophenol	8.89	139	63686	39.83	ng	99
27) 2,4-Dimethylphenol	9.03	122	118651	41.27	ng	98
28) bis(2-Chloroethoxy)methane	9.18	93	170639	40.15	ng	100
29) 2,4-Dichlorophenol	9.30	162	106900	39.30	ng	98
30) 1,2,4-Trichlorobenzene	9.40	180	121106	38.98	ng	98
31) Naphthalene	9.51	128	419848	38.93	ng	99
32) Benzoic acid	9.36	122	66967	37.07	ng	99
33) 4-Chloroaniline	9.64	127	161195	39.80	ng	100
34) Hexachlorobutadiene	9.72	225	69813	39.02	ng	99
35) Caprolactam	10.28	113	31895	40.23	ng	93
36) 4-Chloro-3-methylphenol	10.50	107	127705	40.39	ng	98
37) 2-Methylnaphthalene	10.63	142	253882	38.28	ng	99
39) 1,2,4,5-Tetrachlorobenzene	10.91	216	139695	39.15	ng	99
40) Hexachlorocyclopentadiene	10.89	237	45953	42.05	ng	100

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42) 2,4,6-Trichlorophenol	11.13	196	70317	39.95	nq	98
43) 2,4,5-Trichlorophenol	11.20	196	74161	39.42	nq	96
45) 1,1'-Biphenyl	11.40	154	333766	38.98	nq	99
46) 2-Chloronaphthalene	11.42	162	251707	39.21	nq	98
47) 2-Nitroaniline	11.62	65	80700	39.00	nq	99
48) Acenaphthylene	12.07	152	405785	38.37	nq	100
49) Dimethylphthalate	11.96	163	285976	38.43	nq	100
50) 2,6-Dinitrotoluene	12.04	165	60366	40.80	nq	99
51) Acenaphthene	12.35	154	236040	38.65	nq	99
52) 3-Nitroaniline	12.31	138	67141	37.94	nq	97
53) 2,4-Dinitrophenol	12.51	184	17523	36.08	nq	97
54) Dibenzofuran	12.64	168	327675	39.02	nq	99
55) 4-Nitrophenol	12.68	139	32391m	31.88	nq	
56) 2,4-Dinitrotoluene	12.70	165	79375	38.14	nq	99
57) Fluorene	13.19	166	273193	38.06	nq	99
58) 2,3,4,6-Tetrachlorophenol	12.87	232	52649	36.87	nq	96
59) Diethylphthalate	13.11	149	291548	37.88	nq	98
60) 4-Chlorophenyl-phenylether	13.22	204	128024	38.90	nq	98
61) 4-Nitroaniline	13.31	138	62701	39.41	nq	97
62) Azobenzene	13.47	77	298132	40.05	nq	92
64) 4,6-Dinitro-2-methylphenol	13.36	198	40469	38.05	nq	95
65) n-Nitrosodiphenylamine	13.44	169	236843	39.01	nq	100
66) 4-Bromophenyl-phenylether	14.01	248	70839	40.75	nq	97
67) Hexachlorobenzene	14.07	284	73065	38.20	nq	96
68) Atrazine	14.35	200	73116	43.13	nq	98
69) Pentachlorophenol	14.42	266	33046	42.40	nq	99
70) Phenanthrene	14.73	178	372298	38.09	nq	100
71) Anthracene	14.81	178	391399	37.35	nq	100
72) Carbazole	15.10	167	334710	38.33	nq	99
73) Di-n-butylphthalate	15.69	149	451151	38.20	nq	99
74) Fluoranthene	16.49	202	371871	37.01	nq	99
76) Benzidine	16.73	184	138701	44.19	nq	99
77) Pyrene	16.79	202	362206	39.56	nq	99
79) Butylbenzylphthalate	17.71	149	154577	39.68	nq	# 80
80) Benzo(a)anthracene	18.38	228	231446	38.90	nq	98
81) 3,3'-Dichlorobenzidine	18.37	252	77346	41.26	nq	98
82) Chrysene	18.42	228	236483	37.89	nq	99
83) Bis(2-ethylhexyl)phthalate	18.47	149	209990	38.93	nq	99
84) Di-n-octyl phthalate	19.23	149	300583	38.72	nq	100
85) Indeno(1,2,3-cd)pyrene	21.22	276	207589	43.80	nq	99
87) Benzo(b)fluoranthene	19.63	252	186294	37.82	nq	# 95
88) Benzo(k)fluoranthene	19.67	252	184306m	37.09	nq	
89) Benzo(a)pyrene	19.99	252	178051	38.60	nq	99
90) Dibenzo(a,h)anthracene	21.23	278	181357	42.09	nq	100
91) Benzo(g,h,i)perylene	21.55	276	189465	42.95	nq	98

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

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