

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040216\
 Data File : BF086024.D
 Acq On : 2 Apr 2016 21:08
 Operator : UM/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 4/4/2016 7:46:06 PM

Quant Time: Apr 04 03:37:08 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040216.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 01 23:17:06 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.77	152	125689	20.00	ng	0.00
21) Naphthalene-d8	8.06	136	522307	20.00	ng	0.00
38) Acenaphthene-d10	9.81	164	242147	20.00	ng	0.00
63) Phenanthrene-d10	11.30	188	450887	20.00	ng	0.00
75) Chrysene-d12	13.93	240	255547	20.00	ng	0.00
86) Perylene-d12	15.33	264	224686	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.37	112	587719	79.93	ng	0.00
7) Phenol-d6	6.42	99	715683	76.63	ng	0.00
23) Nitrobenzene-d5	7.34	82	664270	75.00	ng	0.00
41) 2,4,6-Tribromophenol	10.60	330	166012	73.04	ng	0.00
44) 2-Fluorobiphenyl	9.14	172	1115714	76.61	ng	0.00
78) Terphenyl-d14	12.88	244	973297	89.53	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.32	88	144240	41.38	ng	100
3) Pyridine	3.01	79	370286	47.44	ng	96
4) n-Nitrosodimethylamine	2.97	42	142485	41.52	ng	98
6) Aniline	6.44	93	497693	40.61	ng	# 86
8) 2-Chlorophenol	6.56	128	339952	38.76	ng	90
9) Benzaldehyde	6.33	77	209209	41.63	ng	91
10) Phenol	6.43	94	383746	36.02	ng	# 60
11) bis(2-Chloroethyl)ether	6.51	93	324573	38.47	ng	95
12) 1,3-Dichlorobenzene	6.72	146	374654	40.31	ng	98
13) 1,4-Dichlorobenzene	6.80	146	373073	38.96	ng	92
14) 1,2-Dichlorobenzene	6.94	146	356461	40.45	ng	99
15) Benzyl Alcohol	6.92	79	236871	39.20	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.06	45	396220	38.44	ng	88
17) 2-Methylphenol	7.05	107	284781	41.19	ng	96
18) Hexachloroethane	7.29	117	136609	38.79	ng	93
19) n-Nitroso-di-n-propylamine	7.20	70	207688	35.88	ng	94
20) 3+4-Methylphenols	7.20	107	336120	39.97	ng	# 80
22) Acetophenone	7.20	105	482388	39.29	ng	# 93
24) Nitrobenzene	7.37	77	405746	41.87	ng	93
25) Isophorone	7.61	82	687461	39.32	ng	96
26) 2-Nitrophenol	7.69	139	183954	39.68	ng	# 88
27) 2,4-Dimethylphenol	7.72	122	296204	35.57	ng	98
28) bis(2-Chloroethoxy)methane	7.81	93	373095	36.37	ng	98
29) 2,4-Dichlorophenol	7.93	162	276702	39.31	ng	99
30) 1,2,4-Trichlorobenzene	8.01	180	309764	39.20	ng	95
31) Naphthalene	8.09	128	988050	37.61	ng	99
32) Benzoic acid	7.88	122	247474	40.51	ng	90
33) 4-Chloroaniline	8.14	127	397705	37.48	ng	96
34) Hexachlorobutadiene	8.20	225	167281	39.38	ng	100
35) Caprolactam	8.52	113	91026	40.76	ng	98
36) 4-Chloro-3-methylphenol	8.62	107	297217	37.56	ng	91
37) 2-Methylnaphthalene	8.77	142	613445	35.74	ng	99
39) 1,2,4,5-Tetrachlorobenzene	8.94	216	294377	40.45	ng	99
40) Hexachlorocyclopentadiene	8.92	237	50731	28.40	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.06	196	190098	40.68	nq	99
43) 2,4,5-Trichlorophenol	9.10	196	191127	40.28	nq	# 95
45) 1,1'-Biphenyl	9.24	154	818375	39.20	nq	95
46) 2-Chloronaphthalene	9.26	162	617166	40.78	nq	96
47) 2-Nitroaniline	9.37	65	195408	44.13	nq	95
48) Acenaphthylene	9.68	152	942345	38.87	nq	100
49) Dimethylphthalate	9.54	163	722570	40.26	nq	99
50) 2,6-Dinitrotoluene	9.61	165	144909	38.16	nq	96
51) Acenaphthene	9.85	154	541127	37.53	nq	100
52) 3-Nitroaniline	9.78	138	173912	38.00	nq	100
53) 2,4-Dinitrophenol	9.89	184	48652	29.18	nq	# 82
54) Dibenzofuran	10.02	168	702740	36.90	nq	99
55) 4-Nitrophenol	9.95	139	108971	40.39	nq	83
56) 2,4-Dinitrotoluene	10.02	165	195699	40.79	nq	# 77
57) Fluorene	10.36	166	620851	38.17	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.14	232	144164	40.81	nq	95
59) Diethylphthalate	10.24	149	655248	36.17	nq	97
60) 4-Chlorophenyl-phenylether	10.35	204	281476	37.15	nq	97
61) 4-Nitroaniline	10.40	138	192424	42.69	nq	95
62) Azobenzene	10.51	77	653766	37.68	nq	98
64) 4,6-Dinitro-2-methylphenol	10.43	198	80442	29.44	nq	# 57
65) n-Nitrosodiphenylamine	10.48	169	539929	38.29	nq	99
66) 4-Bromophenyl-phenylether	10.84	248	184660	40.11	nq	99
67) Hexachlorobenzene	10.91	284	192174	40.37	nq	# 94
68) Atrazine	11.00	200	164017	36.00	nq	96
69) Pentachlorophenol	11.10	266	81673	36.61	nq	97
70) Phenanthrene	11.32	178	917017	37.28	nq	100
71) Anthracene	11.37	178	899417	34.91	nq	99
72) Carbazole	11.53	167	739262	33.38	nq	99
73) Di-n-butylphthalate	11.86	149	1041816	37.96	nq	99
74) Fluoranthene	12.51	202	863184	33.74	nq	89
76) Benzidine	12.62	184	294972	32.72	nq	100
77) Pyrene	12.73	202	831865	46.19	nq	99
79) Butylbenzylphthalate	13.36	149	419380	51.72	nq	96
80) Benzo(a)anthracene	13.92	228	587942	39.03	nq	99
81) 3,3'-Dichlorobenzidine	13.88	252	436586	39.68	nq	97
82) Chrysene	13.96	228	598343	41.24	nq	99
83) Bis(2-ethylhexyl)phthalate	13.91	149	529729	49.22	nq	100
84) Di-n-octyl phthalate	14.52	149	855775	49.80	nq	99
85) Indeno(1,2,3-cd)pyrene	16.71	276	544001	46.21	nq	100
87) Benzo(b)fluoranthene	14.95	252	598737m	42.36	nq	
88) Benzo(k)fluoranthene	14.97	252	397388m	31.75	nq	
89) Benzo(a)pyrene	15.28	252	483086	38.58	nq	98
90) Dibenzo(a,h)anthracene	16.71	278	457055	45.36	nq	99
91) Benzo(g,h,i)perylene	17.11	276	444505	45.59	nq	96

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

