

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120216\
 Data File : BF091416.D
 Acq On : 3 Dec 2016 3:43
 Operator : UM/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040EC

Manual Integrations
 APPROVED

sohil

12/5/2016 11:21:14 AM

Quant Time: Dec 03 04:14:29 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 29 13:33:46 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	159750	20.00	ng	-0.02
21) Naphthalene-d8	8.15	136	644942	20.00	ng	-0.02
38) Acenaphthene-d10	9.90	164	321186	20.00	ng	-0.02
63) Phenanthrene-d10	11.38	188	507064	20.00	ng	-0.02
75) Chrysene-d12	14.01	240	363064	20.00	ng	-0.03
86) Perylene-d12	15.45	264	309047	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.45	112	743634	76.04	ng	-0.02
7) Phenol-d6	6.49	99	948264	78.77	ng	0.00
23) Nitrobenzene-d5	7.43	82	900141	78.22	ng	-0.02
41) 2,4,6-Tribromophenol	10.69	330	200355	65.89	ng	-0.02
44) 2-Fluorobiphenyl	9.22	172	1456768	82.12	ng	-0.02
78) Terphenyl-d14	12.97	244	1336409	80.00	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.42	88	157246	35.55	ng	# 85
3) Pyridine	3.16	79	489981	39.14	ng	88
4) n-Nitrosodimethylamine	3.10	42	271868	41.39	ng	98
6) Aniline	6.53	93	666034	41.65	ng	# 68
8) 2-Chlorophenol	6.64	128	410530	38.29	ng	# 76
9) Benzaldehyde	6.40	77	307942	39.78	ng	92
10) Phenol	6.50	94	488684	34.50	ng	98
11) bis(2-Chloroethyl)ether	6.60	93	381343	37.68	ng	94
12) 1,3-Dichlorobenzene	6.80	146	466816	39.14	ng	96
13) 1,4-Dichlorobenzene	6.88	146	510363	43.02	ng	100
14) 1,2-Dichlorobenzene	7.03	146	414320	37.49	ng	97
15) Benzyl Alcohol	7.01	79	425915	44.21	ng	95
16) 2,2'-oxybis(1-Chloropropan	7.14	45	862802	39.65	ng	72
17) 2-Methylphenol	7.12	107	358725	42.33	ng	# 76
18) Hexachloroethane	7.37	117	136357	32.37	ng	# 78
19) n-Nitroso-di-n-propylamine	7.28	70	297344	39.21	ng	# 96
20) 3+4-Methylphenols	7.27	107	393431	38.03	ng	# 65
22) Acetophenone	7.28	105	613218	40.12	ng	98
24) Nitrobenzene	7.45	77	470996	39.97	ng	# 72
25) Isophorone	7.69	82	803454	39.41	ng	95
26) 2-Nitrophenol	7.76	139	158984	29.61	ng	95
27) 2,4-Dimethylphenol	7.81	122	404752	40.29	ng	94
28) bis(2-Chloroethoxy)methane	7.90	93	457905	37.31	ng	99
29) 2,4-Dichlorophenol	8.00	162	317761	35.96	ng	96
30) 1,2,4-Trichlorobenzene	8.09	180	408693	41.25	ng	97
31) Naphthalene	8.17	128	1269462	41.42	ng	100
32) Benzoic acid	7.91	122	179399	24.23	ng	94
33) 4-Chloroaniline	8.22	127	531205	40.54	ng	97
34) Hexachlorobutadiene	8.30	225	230331	37.81	ng	99
35) Caprolactam	8.60	113	108122	42.57	ng	# 86
36) 4-Chloro-3-methylphenol	8.70	107	362461	37.99	ng	97
37) 2-Methylnaphthalene	8.86	142	732492	35.17	ng	93
39) 1,2,4,5-Tetrachlorobenzene	9.03	216	407095	44.96	ng	# 97
40) Hexachlorocyclopentadiene	9.02	237	33954	8.09	ng	98

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42) 2,4,6-Trichlorophenol	9.14	196	230079	39.10	ng	94
43) 2,4,5-Trichlorophenol	9.18	196	236165	40.05	ng	95
45) 1,1'-Biphenyl	9.33	154	997993	43.21	ng	97
46) 2-Chloronaphthalene	9.35	162	707426	41.13	ng	99
47) 2-Nitroaniline	9.44	65	268915	43.53	ng	# 77
48) Acenaphthylene	9.76	152	1026603	37.92	ng	98
49) Dimethylphthalate	9.62	163	802681	41.27	ng	97
50) 2,6-Dinitrotoluene	9.68	165	167459	38.52	ng	89
51) Acenaphthene	9.93	154	655312	35.88	ng	97
52) 3-Nitroaniline	9.85	138	200646	39.88	ng	90
54) Dibenzofuran	10.10	168	911394	37.27	ng	98
55) 4-Nitrophenol	10.00	139	115383	31.75	ng	88
56) 2,4-Dinitrotoluene	10.08	165	198166	35.14	ng	# 55
57) Fluorene	10.45	166	700848	37.33	ng	97
58) 2,3,4,6-Tetrachlorophenol	10.23	232	170052	33.77	ng	96
59) Diethylphthalate	10.32	149	834293	43.46	ng	99
60) 4-Chlorophenyl-phenylether	10.45	204	372473	39.56	ng	98
61) 4-Nitroaniline	10.46	138	196299	41.55	ng	85
62) Azobenzene	10.61	77	815966	41.65	ng	96
64) 4,6-Dinitro-2-methylphenol	10.49	198	7959	2.85	ng	# 55
65) n-Nitrosodiphenylamine	10.56	169	675073	43.91	ng	99
66) 4-Bromophenyl-phenylether	10.93	248	226361	41.53	ng	95
67) Hexachlorobenzene	11.00	284	225101	38.22	ng	# 83
68) Atrazine	11.09	200	178498	35.85	ng	96
69) Pentachlorophenol	11.19	266	117279	33.83	ng	93
70) Phenanthrene	11.41	178	967980	36.74	ng	100
71) Anthracene	11.46	178	1054088	40.31	ng	97
72) Carbazole	11.61	167	917901	38.68	ng	98
73) Di-n-butylphthalate	11.96	149	1176333	43.74	ng	# 98
74) Fluoranthene	12.60	202	1049237	42.04	ng	98
76) Benzidine	12.71	184	377083	35.38	ng	98
77) Pyrene	12.82	202	1063595	38.60	ng	99
79) Butylbenzylphthalate	13.44	149	414924	40.22	ng	89
80) Benzo(a)anthracene	14.00	228	852385	40.15	ng	99
81) 3,3'-Dichlorobenzidine	13.97	252	333165	44.54	ng	# 98
82) Chrysene	14.05	228	912909	43.45	ng	99
83) Bis(2-ethylhexyl)phthalate	14.01	149	575199	43.98	ng	# 91
84) Di-n-octyl phthalate	14.62	149	1017127	51.40	ng	99
85) Indeno(1,2,3-cd)pyrene	16.85	276	623592	32.41	ng	95
87) Benzo(b)fluoranthene	15.04	252	839358m	43.70	ng	
88) Benzo(k)fluoranthene	15.07	252	665255m	41.18	ng	
89) Benzo(a)pyrene	15.39	252	697954	40.46	ng	98
90) Dibenzo(a,h)anthracene	16.87	278	549283	34.43	ng	# 94
91) Benzo(g,h,i)perylene	17.27	276	508442	30.91	ng	99

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

