

Data Path : Z:\HPCHEM1\BNA F\DATA\BF122116\
 Data File : BF091835.D
 Acq On : 21 Dec 2016 15:00
 Operator : UM/SJ
 Sample : SSTDICC080
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC080

Manual Integrations
 APPROVED

sohil
 12/22/2016 7:08:38 PM

Quant Time: Dec 21 16:04:21 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF122116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 21 14:14:16 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.75	152	277045	20.00	ng	0.00
21) Naphthalene-d8	8.04	136	1136252	20.00	ng	0.00
38) Acenaphthene-d10	9.79	164	574846	20.00	ng	0.00
63) Phenanthrene-d10	11.28	188	920300	20.00	ng	0.01
75) Chrysene-d12	13.91	240	527377	20.00	ng	0.00
86) Perylene-d12	15.30	264	531996	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.36	112	2259588	136.65	ng	0.01
7) Phenol-d6	6.42	99	2714748	134.61	ng	0.01
23) Nitrobenzene-d5	7.33	82	2648095	134.40	ng	0.01
41) 2,4,6-Tribromophenol	10.59	330	669628	137.37	ng	0.01
44) 2-Fluorobiphenyl	9.13	172	4125551	144.18	ng	0.01
78) Terphenyl-d14	12.86	244	3354390	160.60	ng	0.01

Target Compounds						Qvalue
2) 1,4-Dioxane	2.29	88	617873	75.87	ng	99
3) Pyridine	2.99	79	2228626	84.95	ng	96
4) n-Nitrosodimethylamine	2.97	42	863676	85.06	ng	98
6) Aniline	6.43	93	1753218	67.56	ng	94
8) 2-Chlorophenol	6.54	128	1212047	66.59	ng	97
10) Phenol	6.43	94	1589438	69.92	ng	98
11) bis(2-Chloroethyl)ether	6.50	93	1244388	68.54	ng	99
12) 1,3-Dichlorobenzene	6.69	146	1323125m	66.19	ng	
13) 1,4-Dichlorobenzene	6.77	146	1400475	69.24	ng	98
15) Benzyl Alcohol	6.91	79	1030726	66.01	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.05	45	1940582	69.95	ng	94
17) 2-Methylphenol	7.04	107	1070035	70.35	ng	# 87
18) Hexachloroethane	7.26	117	534038	71.07	ng	91
19) n-Nitroso-di-n-propylamine	7.20	70	881200	65.69	ng	98
20) 3+4-Methylphenols	7.20	107	1289090	75.98	ng	# 71
22) Acetophenone	7.18	105	1943030	70.82	ng	98
24) Nitrobenzene	7.36	77	1627249	76.91	ng	# 85
25) Isophorone	7.60	82	2745700	73.04	ng	98
26) 2-Nitrophenol	7.66	139	811167	79.47	ng	93
27) 2,4-Dimethylphenol	7.71	122	1157617	69.52	ng	99
28) bis(2-Chloroethoxy)methane	7.80	93	1480087	67.29	ng	99
29) 2,4-Dichlorophenol	7.92	162	1069117	69.62	ng	99
30) 1,2,4-Trichlorobenzene	7.98	180	1073188	64.48	ng	# 96
31) Naphthalene	8.06	128	3263942	60.13	ng	98
32) Benzoic acid	7.89	122	1006183	84.79	ng	99
33) 4-Chloroaniline	8.12	127	1649605	72.14	ng	99
34) Hexachlorobutadiene	8.18	225	595619	63.65	ng	99
35) Caprolactam	8.53	113	374715m	75.83	ng	
36) 4-Chloro-3-methylphenol	8.62	107	1250384	73.81	ng	88
39) 1,2,4,5-Tetrachlorobenzene	8.92	216	896685	62.58	ng	# 97
42) 2,4,6-Trichlorophenol	9.04	196	705733	68.93	ng	97
43) 2,4,5-Trichlorophenol	9.08	196	700789	69.68	ng	97
45) 1,1'-Biphenyl	9.22	154	2536507	61.56	ng	97
46) 2-Chloronaphthalene	9.24	162	1981634	64.36	ng	94

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
47) 2-Nitroaniline	9.34	65	715745	68.08	nq	99
48) Acenaphthylene	9.66	152	3347337	70.16	nq	100
49) Dimethylphthalate	9.54	163	2588048	68.92	nq	98
50) 2,6-Dinitrotoluene	9.60	165	580892	70.56	nq	# 81
51) Acenaphthene	9.84	154	2272970	69.42	nq	99
52) 3-Nitroaniline	9.77	138	746472	72.83	nq	95
53) 2,4-Dinitrophenol	9.87	184	387327	87.09	nq	# 78
55) 4-Nitrophenol	9.94	139	473272	66.49	nq	98
56) 2,4-Dinitrotoluene	10.00	165	600947	58.43	nq	# 77
58) 2,3,4,6-Tetrachlorophenol	10.12	232	513943	61.81	nq	96
59) Diethylphthalate	10.22	149	2291689	62.09	nq	99
60) 4-Chlorophenyl-phenylether	10.34	204	975444	70.81	nq	91
61) 4-Nitroaniline	10.38	138	677690	69.78	nq	99
62) Azobenzene	10.50	77	2826128	71.96	nq	94
64) 4,6-Dinitro-2-methylphenol	10.41	198	412188	67.91	nq	# 51
65) n-Nitrosodiphenylamine	10.46	169	2014855	70.51	nq	100
66) 4-Bromophenyl-phenylether	10.82	248	652162	67.58	nq	92
67) Hexachlorobenzene	10.89	284	642013	63.49	nq	# 84
69) Pentachlorophenol	11.08	266	366504	67.16	nq	98
73) Di-n-butylphthalate	11.85	149	3665713	67.13	nq	98
74) Fluoranthene	12.49	202	3343940	69.62	nq	95
76) Benzidine	12.60	184	924978	60.93	nq	# 94
77) Pyrene	12.72	202	3397106	89.20	nq	99
79) Butylbenzylphthalate	13.33	149	1371410	76.94	nq	99
80) Benzo(a)anthracene	13.89	228	2261440	74.97	nq	99
81) 3,3'-Dichlorobenzidine	13.87	252	873811	72.39	nq	# 95
82) Chrysene	13.94	228	2290263	73.25	nq	99
83) Bis(2-ethylhexyl)phthalate	13.89	149	1595018	70.54	nq	99
84) Di-n-octyl phthalate	14.51	149	3138560	75.33	nq	99
85) Indeno(1,2,3-cd)pyrene	16.65	276	2143169	69.71	nq	100
87) Benzo(b)fluoranthene	14.91	252	2177179m	64.54	nq	
88) Benzo(k)fluoranthene	14.95	252	2043359m	86.33	nq	
89) Benzo(a)pyrene	15.25	252	2074358	71.65	nq	98
90) Dibenzo(a,h)anthracene	16.67	278	1735201	68.75	nq	98
91) Benzo(a,h,i)perylene	17.06	276	1891488	73.74	nq	97

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

