

Data Path : Z:\HPCHEM1\BNA F\DATA\BF031317\
 Data File : BF093650.D
 Acq On : 14 Mar 2017 11:15
 Operator : SJ/MA
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad
 3/15/2017 5:38:49 PM

Quant Time: Mar 15 04:41:43 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF030717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 13 15:05:08 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.71	152	189128	20.00	ng	-0.01
21) Naphthalene-d8	8.00	136	795989	20.00	ng	-0.01
38) Acenaphthene-d10	9.76	164	356046	20.00	ng	0.00
63) Phenanthrene-d10	11.25	188	630894	20.00	ng	0.00
75) Chrysene-d12	13.90	240	485880	20.00	ng	0.01
86) Perylene-d12	15.31	264	344065	20.00	ng	0.02

System Monitoring Compounds						
5) 2-Fluorophenol	5.29	112	995823	87.63	ng	-0.02
7) Phenol-d6	6.37	99	1160659	81.00	ng	-0.01
23) Nitrobenzene-d5	7.29	82	1190758	83.00	ng	-0.01
41) 2,4,6-Tribromophenol	10.56	330	195418	84.24	ng	0.01
44) 2-Fluorobiphenyl	9.09	172	1658429	86.64	ng	0.00
78) Terphenyl-d14	12.84	244	1500348	80.28	ng	0.01

Target Compounds						Qvalue
2) 1,4-Dioxane	2.18	88	249653	39.89	ng	# 81
3) Pyridine	2.86	79	568421	35.62	ng	89
4) n-Nitrosodimethylamine	2.84	42	343950	45.12	ng	85
6) Aniline	6.38	93	689412	35.05	ng	# 83
8) 2-Chlorophenol	6.50	128	513447	40.33	ng	96
9) Benzaldehyde	6.26	77	286265	29.42	ng	90
10) Phenol	6.38	94	751977	42.82	ng	# 76
11) bis(2-Chloroethyl)ether	6.45	93	679867	47.90	ng	91
12) 1,3-Dichlorobenzene	6.64	146	596569	40.94	ng	# 91
13) 1,4-Dichlorobenzene	6.72	146	615834	41.27	ng	99
14) 1,2-Dichlorobenzene	6.88	146	539889	40.86	ng	97
15) Benzyl Alcohol	6.88	79	341477	33.43	ng	96
16) 2,2'-oxybis(1-Chloropropan	7.00	45	981185	41.62	ng	# 40
17) 2-Methylphenol	7.00	107	455877	42.33	ng	# 89
18) Hexachloroethane	7.21	117	211744	42.08	ng	95
19) n-Nitroso-di-n-propylamine	7.14	70	422656	42.90	ng	97
20) 3+4-Methylphenols	7.16	107	639106	45.76	ng	# 74
22) Acetophenone	7.13	105	777641	40.60	ng	# 98
24) Nitrobenzene	7.30	77	649537	40.28	ng	94
25) Isophorone	7.54	82	1163207	40.21	ng	# 92
26) 2-Nitrophenol	7.62	139	286634	38.96	ng	# 81
27) 2,4-Dimethylphenol	7.68	122	443979	37.10	ng	97
28) bis(2-Chloroethoxy)methane	7.76	93	731129	40.03	ng	99
29) 2,4-Dichlorophenol	7.89	162	433421	38.50	ng	93
30) 1,2,4-Trichlorobenzene	7.94	180	439484	36.71	ng	95
31) Naphthalene	8.02	128	1564085	39.21	ng	99
32) Benzoic acid	7.85	122	242749	39.04	ng	97
33) 4-Chloroaniline	8.10	127	600079m	37.07	ng	
34) Hexachlorobutadiene	8.14	225	239616	38.85	ng	98
35) Caprolactam	8.48	113	147024	38.24	ng	# 36
36) 4-Chloro-3-methylphenol	8.60	107	503087	41.87	ng	97
37) 2-Methylnaphthalene	8.72	142	978654	38.56	ng	98
39) 1,2,4,5-Tetrachlorobenzene	8.89	216	425942	43.81	ng	99
40) Hexachlorocyclopentadiene	8.87	237	72545	37.19	ng	98

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42) 2,4,6-Trichlorophenol	9.02	196	259069	42.65	nq	95
43) 2,4,5-Trichlorophenol	9.08	196	264235	40.54	nq #	90
45) 1,1'-Biphenyl	9.18	154	1131858	43.64	nq	95
46) 2-Chloronaphthalene	9.21	162	926510	46.67	nq	97
47) 2-Nitroaniline	9.33	65	359321	51.20	nq	98
48) Acenaphthylene	9.62	152	1457079	44.50	nq	99
49) Dimethylphthalate	9.50	163	1034852	43.84	nq	100
50) 2,6-Dinitrotoluene	9.57	165	225955	43.63	nq #	63
51) Acenaphthene	9.80	154	861514	44.49	nq	96
52) 3-Nitroaniline	9.75	138	265184	42.62	nq	88
53) 2,4-Dinitrophenol	9.88	184	107100	45.41	nq #	63
54) Dibenzofuran	9.97	168	1123565	46.36	nq	99
55) 4-Nitrophenol	9.96	139	72705m	24.06	nq	
56) 2,4-Dinitrotoluene	9.98	165	315377	49.56	nq #	82
57) Fluorene	10.31	166	889173	43.30	nq	97
58) 2,3,4,6-Tetrachlorophenol	10.10	232	198520	43.16	nq	96
59) Diethylphthalate	10.20	149	1046531	46.39	nq	96
60) 4-Chlorophenyl-phenylether	10.30	204	403368	43.82	nq	97
61) 4-Nitroaniline	10.37	138	234547	36.86	nq	95
62) Azobenzene	10.46	77	1083039m	42.25	nq	
64) 4,6-Dinitro-2-methylphenol	10.40	198	128668	31.48	nq	81
65) n-Nitrosodiphenylamine	10.42	169	782220	37.13	nq	99
66) 4-Bromophenyl-phenylether	10.79	248	242277	36.32	nq	98
67) Hexachlorobenzene	10.86	284	234321	36.78	nq #	73
68) Atrazine	10.96	200	218860	35.50	nq	99
69) Pentachlorophenol	11.08	266	90912	41.33	nq	98
70) Phenanthrene	11.27	178	1322395	36.72	nq	99
71) Anthracene	11.33	178	1395253	38.30	nq	99
72) Carbazole	11.50	167	1413783	41.31	nq	97
73) Di-n-butylphthalate	11.82	149	1610012	40.67	nq #	96
74) Fluoranthene	12.47	202	1388489	39.92	nq	93
76) Benzidine	12.62	184	224835	14.07	nq	98
77) Pyrene	12.70	202	1434662	40.60	nq	99
79) Butylbenzylphthalate	13.30	149	640087	43.03	nq	90
80) Benzo(a)anthracene	13.89	228	1091536	38.04	nq	99
81) 3,3'-Dichlorobenzidine	13.85	252	334115	33.25	nq #	97
82) Chrysene	13.92	228	981751	36.98	nq	100
83) Bis(2-ethylhexyl)phthalate	13.87	149	867343	41.83	nq	99
84) Di-n-octyl phthalate	14.48	149	1430821	41.40	nq	97
85) Indeno(1,2,3-cd)pyrene	16.65	276	760454	34.22	nq #	93
87) Benzo(b)fluoranthene	14.91	252	805566m	38.44	nq	
88) Benzo(k)fluoranthene	14.93	252	879525m	43.52	nq	
89) Benzo(a)pyrene	15.25	252	789451	41.22	nq	98
90) Dibenzo(a,h)anthracene	16.65	278	655707	41.39	nq	97
91) Benzo(g,h,i)perylene	17.05	276	681716	41.92	nq #	92

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

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