

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051016\
 Data File : BF086876.D
 Acq On : 11 May 2016 1:29
 Operator : UM/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

sohil
 5/11/2016 8:16:01 PM

Quant Time: May 11 05:42:45 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon May 09 16:04:56 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.68	152	200567	20.00	ng	-0.01
21) Naphthalene-d8	7.97	136	746823	20.00	ng	-0.01
38) Acenaphthene-d10	9.72	164	307859	20.00	ng	-0.01
63) Phenanthrene-d10	11.20	188	490432	20.00	ng	-0.01
75) Chrysene-d12	13.83	240	346672	20.00	ng	-0.01
86) Perylene-d12	15.20	264	348961	20.00	ng	-0.01

System Monitoring Compounds						
5) 2-Fluorophenol	5.28	112	881598	74.44	ng	-0.01
7) Phenol-d6	6.35	99	1040337	65.73	ng	-0.01
23) Nitrobenzene-d5	7.26	82	1118841	75.23	ng	0.00
41) 2,4,6-Tribromophenol	10.51	330	208270	63.90	ng	-0.01
44) 2-Fluorobiphenyl	9.06	172	1663423	90.81	ng	0.00
78) Terphenyl-d14	12.78	244	1316502	80.57	ng	-0.01

Target Compounds						Qvalue
2) 1,4-Dioxane	2.17	88	222936	38.34	ng	# 72
3) Pyridine	2.84	79	568155	39.97	ng	# 63
4) n-Nitrosodimethylamine	2.81	42	440153	42.70	ng	# 50
6) Aniline	6.35	93	656046	34.27	ng	# 85
8) 2-Chlorophenol	6.48	128	517373	36.20	ng	# 87
9) Benzaldehyde	6.24	77	345891	37.74	ng	# 96
10) Phenol	6.36	94	542968	28.86	ng	# 28
11) bis(2-Chloroethyl)ether	6.43	93	521452	35.55	ng	# 84
12) 1,3-Dichlorobenzene	6.62	146	585602m	37.87	ng	# 95
13) 1,4-Dichlorobenzene	6.70	146	602659	38.21	ng	# 94
14) 1,2-Dichlorobenzene	6.86	146	475933	34.62	ng	# 93
15) Benzyl Alcohol	6.85	79	406649	37.21	ng	# 85
16) 2,2'-oxybis(1-Chloropropan	6.98	45	1182365	37.07	ng	# 72
17) 2-Methylphenol	6.97	107	380031	33.42	ng	# 97
18) Hexachloroethane	7.20	117	228617	38.53	ng	# 63
19) n-Nitroso-di-n-propylamine	7.12	70	387400	35.02	ng	# 63
20) 3+4-Methylphenols	7.13	107	475977	32.05	ng	# 93
22) Acetophenone	7.10	105	710732	40.29	ng	# 91
24) Nitrobenzene	7.29	77	656495	42.91	ng	# 94
25) Isophorone	7.53	82	1073471	38.90	ng	# 70
26) 2-Nitrophenol	7.60	139	253200	37.65	ng	# 90
27) 2,4-Dimethylphenol	7.65	122	423428	36.96	ng	# 98
28) bis(2-Chloroethoxy)methane	7.73	93	564847	37.95	ng	# 98
29) 2,4-Dichlorophenol	7.85	162	377728	37.46	ng	# 95
30) 1,2,4-Trichlorobenzene	7.92	180	434872	38.57	ng	# 99
31) Naphthalene	8.00	128	1342899	35.90	ng	# 88
32) Benzoic acid	7.80	122	211402	31.42	ng	# 96
33) 4-Chloroaniline	8.06	127	496271	34.14	ng	# 98
34) Hexachlorobutadiene	8.11	225	261243	39.94	ng	# 98
35) Caprolactam	8.45	113	114967	33.51	ng	# 91
36) 4-Chloro-3-methylphenol	8.56	107	376108	30.76	ng	# 96
37) 2-Methylnaphthalene	8.68	142	803001	36.72	ng	# 97
39) 1,2,4,5-Tetrachlorobenzene	8.85	216	384547	42.96	ng	# 93
40) Hexachlorocyclopentadiene	8.84	237	122428	28.82	ng	# 93

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	8.98	196	233361	38.99	nq	92
43) 2,4,5-Trichlorophenol	9.02	196	217361	35.11	nq	95
45) 1,1'-Biphenyl	9.15	154	931850	38.05	nq	96
46) 2-Chloronaphthalene	9.17	162	727050	37.90	nq	93
47) 2-Nitroaniline	9.29	65	256314	36.56	nq	96
48) Acenaphthylene	9.58	152	1062276	37.64	nq	99
49) Dimethylphthalate	9.46	163	839666	35.75	nq	99
50) 2,6-Dinitrotoluene	9.53	165	206897	41.86	nq	87
51) Acenaphthene	9.76	154	659899	37.59	nq	99
52) 3-Nitroaniline	9.70	138	177023	34.02	nq	93
53) 2,4-Dinitrophenol	9.80	184	48477	25.04	nq #	77
54) Dibenzofuran	9.93	168	842425	37.39	nq	94
55) 4-Nitrophenol	9.88	139	84085	27.41	nq #	73
56) 2,4-Dinitrotoluene	9.93	165	226890	37.09	nq #	88
57) Fluorene	10.27	166	747929	38.92	nq	100
58) 2,3,4,6-Tetrachlorophenol	10.05	232	162953	34.97	nq	96
59) Diethylphthalate	10.16	149	842201	36.80	nq	98
60) 4-Chlorophenyl-phenylether	10.27	204	341749	39.68	nq #	78
61) 4-Nitroaniline	10.30	138	156672	31.35	nq #	68
62) Azobenzene	10.42	77	864050	34.99	nq	82
64) 4,6-Dinitro-2-methylphenol	10.34	198	92078	30.22	nq	88
65) n-Nitrosodiphenylamine	10.38	169	576940	38.29	nq	100
66) 4-Bromophenyl-phenylether	10.75	248	201352	40.49	nq	98
67) Hexachlorobenzene	10.82	284	251182	43.22	nq	96
68) Atrazine	10.92	200	206348	41.00	nq	95
69) Pentachlorophenol	11.01	266	104411	38.88	nq	97
70) Phenanthrene	11.23	178	1010786	38.62	nq	98
71) Anthracene	11.28	178	1029588	38.47	nq	100
72) Carbazole	11.44	167	796023	34.59	nq	99
73) Di-n-butylphthalate	11.77	149	1125283	40.05	nq #	98
74) Fluoranthene	12.41	202	1034993	38.45	nq	94
76) Benzidine	12.54	184	215511	24.38	nq	96
77) Pyrene	12.64	202	1036592	39.06	nq	99
79) Butylbenzylphthalate	13.25	149	454269	40.47	nq #	60
80) Benzo(a)anthracene	13.81	228	779792	40.07	nq	100
81) 3,3'-Dichlorobenzidine	13.79	252	473958	38.33	nq #	96
82) Chrysene	13.86	228	866096	43.75	nq	100
83) Bis(2-ethylhexyl)phthalate	13.81	149	598938	44.35	nq #	94
84) Di-n-octyl phthalate	14.43	149	1097187	49.05	nq #	84
85) Indeno(1,2,3-cd)pyrene	16.49	276	785065	47.66	nq	95
87) Benzo(b)fluoranthene	14.82	252	744464m	32.83	nq	
88) Benzo(k)fluoranthene	14.85	252	822817m	44.31	nq	
89) Benzo(a)pyrene	15.15	252	752623	38.47	nq	98
90) Dibenzo(a,h)anthracene	16.50	278	648297	39.32	nq	98
91) Benzo(g,h,i)perylene	16.88	276	642150	38.33	nq #	93

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

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