

Data Path : Z:\HPCHEM1\BNA F\DATA\BF122015\
 Data File : BF083949.D
 Acq On : 19 Dec 2015 12:42
 Operator : UM/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

apatel
 12/22/2015 11:59:23 AM

Quant Time: Dec 21 02:59:21 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 18 14:23:22 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	69037	20.00	ng	-0.02
21) Naphthalene-d8	8.13	136	265871	20.00	ng	-0.02
38) Acenaphthene-d10	9.89	164	131772	20.00	ng	-0.02
63) Phenanthrene-d10	11.37	188	236482	20.00	ng	-0.02
75) Chrysene-d12	14.01	240	193529	20.00	ng	-0.02
86) Perylene-d12	15.44	264	148137	20.00	ng	-0.03

System Monitoring Compounds

5) 2-Fluorophenol	5.46	112	306437	69.46	ng	-0.02
7) Phenol-d6	6.50	99	375180	69.48	ng	-0.01
23) Nitrobenzene-d5	7.42	82	383408	82.30	ng	-0.01
41) 2,4,6-Tribromophenol	10.68	330	116746	80.42	ng	-0.02
44) 2-Fluorobiphenyl	9.22	172	727209	76.24	ng	-0.02
78) Terphenyl-d14	12.96	244	754962	91.70	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.42	88	64856	33.89	ng	98
3) Pyridine	3.15	79	172516	36.46	ng	99
4) n-Nitrosodimethylamine	3.10	42	69534	38.52	ng	96
6) Aniline	6.51	93	243514	35.02	ng	# 63
8) 2-Chlorophenol	6.64	128	185043	35.24	ng	95
9) Benzaldehyde	6.40	77	115744	35.58	ng	96
10) Phenol	6.51	94	195624	32.47	ng	# 65
11) bis(2-Chloroethyl)ether	6.59	93	163920	35.81	ng	97
12) 1,3-Dichlorobenzene	6.80	146	215480	38.16	ng	99
13) 1,4-Dichlorobenzene	6.86	146	200858	34.84	ng	96
14) 1,2-Dichlorobenzene	7.02	146	200901	43.09	ng	99
15) Benzyl Alcohol	7.00	79	159083	38.93	ng	94
16) 2,2'-oxybis(1-Chloropropan	7.13	45	172112	36.47	ng	57
17) 2-Methylphenol	7.12	107	139389	33.85	ng	96
18) Hexachloroethane	7.37	117	76778	36.30	ng	93
19) n-Nitroso-di-n-propylamine	7.28	70	121222	38.51	ng	91
20) 3+4-Methylphenols	7.28	107	186876	38.61	ng	# 54
22) Acetophenone	7.26	105	243525	38.05	ng	95
24) Nitrobenzene	7.44	77	172446	35.79	ng	97
25) Isophorone	7.68	82	334817	38.20	ng	99
26) 2-Nitrophenol	7.76	139	106168	42.57	ng	94
27) 2,4-Dimethylphenol	7.80	122	179058	41.68	ng	95
28) bis(2-Chloroethoxy)methane	7.89	93	208971	40.31	ng	98
29) 2,4-Dichlorophenol	8.01	162	150171	38.53	ng	93
30) 1,2,4-Trichlorobenzene	8.08	180	155302	37.10	ng	99
31) Naphthalene	8.16	128	501854	37.08	ng	100
32) Benzoic acid	7.93	122	88524	53.52	ng	93
33) 4-Chloroaniline	8.21	127	231879	38.42	ng	# 90
34) Hexachlorobutadiene	8.28	225	99468	38.05	ng	98
35) Caprolactam	8.59	113	49784	38.66	ng	94
36) 4-Chloro-3-methylphenol	8.70	107	180206	40.80	ng	92
37) 2-Methylnaphthalene	8.85	142	369452	39.98	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.01	216	148208	35.67	ng	97
40) Hexachlorocyclopentadiene	9.00	237	31157	34.65	ng	96

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.13	196	100349	36.58	nq	98
43) 2,4,5-Trichlorophenol	9.17	196	104575	38.52	nq #	84
45) 1,1'-Biphenyl	9.31	154	410702	36.83	nq	99
46) 2-Chloronaphthalene	9.33	162	333780	38.29	nq	99
47) 2-Nitroaniline	9.44	65	104093	39.94	nq #	75
48) Acenaphthylene	9.76	152	583770	38.00	nq	100
49) Dimethylphthalate	9.62	163	416547	37.67	nq #	91
50) 2,6-Dinitrotoluene	9.68	165	84414	35.00	nq #	66
51) Acenaphthene	9.93	154	362745	38.40	nq	100
52) 3-Nitroaniline	9.85	138	99159	39.00	nq	86
53) 2,4-Dinitrophenol	9.95	184	28993	39.47	nq #	65
54) Dibenzofuran	10.10	168	460407	37.53	nq	99
55) 4-Nitrophenol	10.02	139	63213	39.83	nq	91
56) 2,4-Dinitrotoluene	10.09	165	117904	38.13	nq #	78
57) Fluorene	10.44	166	388797	41.43	nq	100
58) 2,3,4,6-Tetrachlorophenol	10.21	232	73213	36.45	nq	97
59) Diethylphthalate	10.32	149	439024	38.41	nq	99
60) 4-Chlorophenyl-phenylether	10.43	204	178707	38.01	nq	96
61) 4-Nitroaniline	10.46	138	100783	38.31	nq	86
62) Azobenzene	10.59	77	377190	39.37	nq	97
64) 4,6-Dinitro-2-methylphenol	10.50	198	57172	40.05	nq	73
65) n-Nitrosodiphenylamine	10.54	169	305219	35.47	nq	100
66) 4-Bromophenyl-phenylether	10.92	248	110161	38.10	nq	96
67) Hexachlorobenzene	10.99	284	117280	38.53	nq #	90
68) Atrazine	11.08	200	111199	40.79	nq	94
69) Pentachlorophenol	11.18	266	51967	42.57	nq	97
70) Phenanthrene	11.40	178	531431	37.19	nq	99
71) Anthracene	11.45	178	500049	36.21	nq	100
72) Carbazole	11.61	167	513448	37.63	nq	98
73) Di-n-butylphthalate	11.94	149	674449	42.06	nq	100
74) Fluoranthene	12.58	202	535517	37.56	nq	99
76) Benzidine	12.71	184	216763	27.97	nq	99
77) Pyrene	12.81	202	523999	42.76	nq	100
79) Butylbenzylphthalate	13.43	149	261129	40.53	nq	97
80) Benzo(a)anthracene	14.00	228	446977	37.57	nq	100
81) 3,3'-Dichlorobenzidine	13.96	252	170444	38.33	nq	97
82) Chrysene	14.03	228	425758	37.04	nq	100
83) Bis(2-ethylhexyl)phthalate	13.99	149	363336	42.24	nq	100
84) Di-n-octyl phthalate	14.60	149	616392	45.79	nq	100
85) Indeno(1,2,3-cd)pyrene	16.87	276	359805	41.86	nq	99
87) Benzo(b)fluoranthene	15.03	252	453615m	42.00	nq	
88) Benzo(k)fluoranthene	15.06	252	295780m	33.20	nq	
89) Benzo(a)pyrene	15.38	252	333820	37.43	nq #	96
90) Dibenzo(a,h)anthracene	16.87	278	297807	42.92	nq	97
91) Benzo(g,h,i)perylene	17.28	276	300145	44.14	nq	98

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

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