

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011916\
 Data File : BF084551.D
 Acq On : 19 Jan 2016 14:47
 Operator : SJ/UM
 Sample : PB87864BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB87864BS

Manual Integrations
 APPROVED

mohammad
 1/20/2016 2:22:26 PM

Quant Time: Jan 20 02:58:12 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF123115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 14 14:15:52 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	209276	20.00	ng	-0.03
21) Naphthalene-d8	8.08	136	851954	20.00	ng	-0.05
38) Acenaphthene-d10	9.84	164	424941	20.00	ng	-0.03
63) Phenanthrene-d10	11.32	188	798959	20.00	ng	-0.03
75) Chrysene-d12	13.95	240	576280	20.00	ng	-0.05
86) Perylene-d12	15.36	264	468440	20.00	ng	-0.06

System Monitoring Compounds

5) 2-Fluorophenol	5.40	112	1332811	103.01	ng	-0.03
7) Phenol-d6	6.44	99	1838595	113.15	ng	-0.03
23) Nitrobenzene-d5	7.37	82	1279060	83.56	ng	-0.03
41) 2,4,6-Tribromophenol	10.62	330	475842	110.92	ng	-0.05
44) 2-Fluorobiphenyl	9.16	172	2007111	82.94	ng	-0.03
78) Terphenyl-d14	12.91	244	2173296	84.18	ng	-0.03

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.44	88	214142	34.94	ng	# 71
3) Pyridine	3.14	79	492808	28.84	ng	# 71
4) n-Nitrosodimethylamine	3.09	42	284485	32.43	ng	# 75
6) Aniline	6.45	93	449672	18.92	ng	# 1
8) 2-Chlorophenol	6.58	128	600449	40.90	ng	89
9) Benzaldehyde	6.34	77	60052	5.68	ng	88
10) Phenol	6.45	94	713690	38.63	ng	91
11) bis(2-Chloroethyl)ether	6.53	93	575328	40.20	ng	# 83
12) 1,3-Dichlorobenzene	6.73	146	611790m	40.40	ng	
13) 1,4-Dichlorobenzene	6.81	146	607681	40.02	ng	97
14) 1,2-Dichlorobenzene	6.97	146	611096	42.02	ng	98
15) Benzyl Alcohol	6.94	79	510147	37.32	ng	92
16) 2,2'-oxybis(1-Chloropropan	7.08	45	974395	37.39	ng	88
17) 2-Methylphenol	7.06	107	467391	37.99	ng	94
18) Hexachloroethane	7.31	117	250316	41.22	ng	# 87
19) n-Nitroso-di-n-propylamine	7.22	70	438116	39.83	ng	# 80
20) 3+4-Methylphenols	7.22	107	615513	42.56	ng	# 85
22) Acetophenone	7.21	105	833560	40.69	ng	# 97
24) Nitrobenzene	7.38	77	599388	38.61	ng	96
25) Isophorone	7.62	82	1144499	39.09	ng	99
26) 2-Nitrophenol	7.70	139	330259	46.74	ng	# 87
27) 2,4-Dimethylphenol	7.74	122	584842	40.89	ng	94
28) bis(2-Chloroethoxy)methane	7.84	93	726183	40.61	ng	100
29) 2,4-Dichlorophenol	7.95	162	505578	42.44	ng	92
30) 1,2,4-Trichlorobenzene	8.02	180	518991	42.20	ng	96
31) Naphthalene	8.10	128	1639072	42.40	ng	99
32) Benzoic acid	7.88	122	370806	41.30	ng	98
33) 4-Chloroaniline	8.16	127	224006	12.64	ng	95
34) Hexachlorobutadiene	8.22	225	313894	44.40	ng	98
35) Caprolactam	8.52	113	174690m	43.49	ng	
36) 4-Chloro-3-methylphenol	8.65	107	556511	41.70	ng	95
37) 2-Methylnaphthalene	8.80	142	1190574	42.91	ng	98
39) 1,2,4,5-Tetrachlorobenzene	8.97	216	470193	38.49	ng	99
40) Hexachlorocyclopentadiene	8.96	237	549889	74.56	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.08	196	352153	38.53	nq	94
43) 2,4,5-Trichlorophenol	9.13	196	374644	42.76	nq	97
45) 1,1'-Biphenyl	9.26	154	1446700	42.84	nq	97
46) 2-Chloronaphthalene	9.29	162	1086336	40.95	nq	# 89
47) 2-Nitroaniline	9.38	65	322233	36.80	nq	97
48) Acenaphthylene	9.70	152	1609461	41.07	nq	97
49) Dimethylphthalate	9.57	163	1298363	42.74	nq	# 91
50) 2,6-Dinitrotoluene	9.63	165	308074	46.24	nq	# 89
51) Acenaphthene	9.87	154	1031733m	39.24	nq	
52) 3-Nitroaniline	9.79	138	139098	16.85	nq	# 95
53) 2,4-Dinitrophenol	9.90	184	336697	116.91	nq	# 89
54) Dibenzofuran	10.04	168	1354622	39.42	nq	92
55) 4-Nitrophenol	9.97	139	451352	75.31	nq	# 83
56) 2,4-Dinitrotoluene	10.03	165	359793	42.67	nq	# 91
57) Fluorene	10.38	166	1108981	41.96	nq	100
58) 2,3,4,6-Tetrachlorophenol	10.17	232	307531	41.11	nq	88
59) Diethylphthalate	10.27	149	1304221	43.54	nq	97
60) 4-Chlorophenyl-phenylether	10.37	204	503515	45.39	nq	# 88
61) 4-Nitroaniline	10.41	138	271774	36.93	nq	92
62) Azobenzene	10.53	77	1299219	39.55	nq	86
64) 4,6-Dinitro-2-methylphenol	10.44	198	234908	48.60	nq	79
65) n-Nitrosodiphenylamine	10.50	169	976124	38.21	nq	99
66) 4-Bromophenyl-phenylether	10.86	248	330037	38.38	nq	# 78
67) Hexachlorobenzene	10.93	284	392718	40.57	nq	95
68) Atrazine	11.04	200	333435	39.89	nq	90
69) Pentachlorophenol	11.13	266	368892	73.50	nq	98
70) Phenanthrene	11.34	178	1757545	42.05	nq	96
71) Anthracene	11.39	178	1687986	41.20	nq	99
72) Carbazole	11.55	167	1441058	35.98	nq	98
73) Di-n-butylphthalate	11.89	149	1956151	46.29	nq	99
74) Fluoranthene	12.53	202	1832180	41.97	nq	93
76) Benzidine	12.65	184	446941	21.63	nq	99
77) Pyrene	12.76	202	1857431	41.07	nq	99
79) Butylbenzylphthalate	13.38	149	773700	39.54	nq	88
80) Benzo(a)anthracene	13.94	228	1420099	41.97	nq	98
81) 3,3'-Dichlorobenzidine	13.91	252	345558	29.26	nq	# 94
82) Chrysene	13.99	228	1473942	45.33	nq	99
83) Bis(2-ethylhexyl)phthalate	13.94	149	982870	39.76	nq	# 100
84) Di-n-octyl phthalate	14.56	149	1708359	40.93	nq	# 90
85) Indeno(1,2,3-cd)pyrene	16.72	276	1047485	40.64	nq	100
87) Benzo(b)fluoranthene	14.97	252	1432894m	46.76	nq	
88) Benzo(k)fluoranthene	14.99	252	976645m	38.97	nq	
89) Benzo(a)pyrene	15.30	252	1093676	42.08	nq	# 94
90) Dibenzo(a,h)anthracene	16.74	278	864477	38.65	nq	# 95
91) Benzo(g,h,i)perylene	17.13	276	891127	38.48	nq	99

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

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