

Data Path : Z:\HPCHEM1\BNA F\DATA\BF122316\
 Data File : BF091936.D
 Acq On : 23 Dec 2016 20:55
 Operator : UM/SJ
 Sample : H6068-05MS
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 H6058-05MS

Quant Time: Dec 23 23:09:20 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 16:17:51 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.74	152	230116	20.00	ng	-0.01
21) Naphthalene-d8	8.03	136	851811	20.00	ng	-0.01
38) Acenaphthene-d10	9.79	164	407851	20.00	ng	0.00
63) Phenanthrene-d10	11.26	188	583227	20.00	ng	0.00
75) Chrysene-d12	13.89	240	455224	20.00	ng	-0.01
86) Perylene-d12	15.29	264	457032	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.39	112	1724219	125.11	ng	0.05
7) Phenol-d6	6.44	99	2170338	128.99	ng	0.03
23) Nitrobenzene-d5	7.32	82	1433692	93.59	ng	0.00
41) 2,4,6-Tribromophenol	10.58	330	369002	109.33	ng	0.00
44) 2-Fluorobiphenyl	9.12	172	1936640	99.49	ng	0.00
78) Terphenyl-d14	12.84	244	1282271	68.31	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.33	88	235709	34.74	ng	99
3) Pyridine	3.01	79	533210	22.52	ng	97
4) n-Nitrosodimethylamine	2.98	42	373214	41.78	ng	96
6) Aniline	6.41	93	327446	14.98	ng	# 45
8) 2-Chlorophenol	6.54	128	619177	39.50	ng	97
9) Benzaldehyde	6.28	77	96867	7.67	ng	94
10) Phenol	6.45	94	1013149	52.84	ng	# 72
11) bis(2-Chloroethyl)ether	6.49	93	676505	44.34	ng	99
12) 1,3-Dichlorobenzene	6.68	146	669691	40.02	ng	# 92
13) 1,4-Dichlorobenzene	6.76	146	686372	40.29	ng	97
14) 1,2-Dichlorobenzene	6.91	146	563235	38.11	ng	95
15) Benzyl Alcohol	6.91	79	583622	44.64	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.04	45	927900	40.84	ng	67
17) 2-Methylphenol	7.05	107	545704	43.28	ng	98
18) Hexachloroethane	7.25	117	254696	40.33	ng	93
19) n-Nitroso-di-n-propylamine	7.17	70	516861	46.17	ng	97
20) 3+4-Methylphenols	7.21	107	747223	49.69	ng	# 74
22) Acetophenone	7.16	105	955246	45.47	ng	94
24) Nitrobenzene	7.33	77	673393	41.27	ng	100
25) Isophorone	7.57	82	1270038	44.94	ng	# 93
26) 2-Nitrophenol	7.65	139	362486	45.05	ng	95
27) 2,4-Dimethylphenol	7.71	122	535259	40.11	ng	97
28) bis(2-Chloroethoxy)methane	7.79	93	811716	47.36	ng	96
29) 2,4-Dichlorophenol	7.92	162	529517	46.86	ng	94
30) 1,2,4-Trichlorobenzene	7.97	180	511138	41.28	ng	# 94
31) Naphthalene	8.05	128	1630077	41.81	ng	100
32) Benzoic acid	7.85	122	174613	17.64	ng	95
33) 4-Chloroaniline	8.11	127	199579	11.35	ng	96
34) Hexachlorobutadiene	8.17	225	292987	44.82	ng	98
35) Caprolactam	8.46	113	276093m	74.35	ng	
36) 4-Chloro-3-methylphenol	8.62	107	516801	39.74	ng	96
37) 2-Methylnaphthalene	8.75	142	1123598	49.22	ng	99
39) 1,2,4,5-Tetrachlorobenzene	8.91	216	426897	41.43	ng	98
40) Hexachlorocyclopentadiene	8.90	237	185738	42.44	ng	96

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.04	196	298338	41.40	nq	96
43) 2,4,5-Trichlorophenol	9.09	196	306961	41.39	nq	94
45) 1,1'-Biphenyl	9.21	154	1184200	42.41	nq	96
46) 2-Chloronaphthalene	9.24	162	920431	41.62	nq	99
47) 2-Nitroaniline	9.34	65	352859	44.81	nq	# 76
48) Acenaphthylene	9.65	152	1423341	41.85	nq	99
49) Dimethylphthalate	9.52	163	1179224	45.21	nq	98
50) 2,6-Dinitrotoluene	9.58	165	262885	46.04	nq	# 87
51) Acenaphthene	9.82	154	932401	40.44	nq	100
52) 3-Nitroaniline	9.76	138	151100	21.38	nq	# 70
53) 2,4-Dinitrophenol	9.86	184	81432	25.63	nq	# 62
54) Dibenzofuran	10.00	168	1254716	40.29	nq	99
55) 4-Nitrophenol	9.96	139	382628	79.76	nq	# 74
56) 2,4-Dinitrotoluene	9.98	165	322674	45.03	nq	# 96
57) Fluorene	10.34	166	998781	44.09	nq	98
58) 2,3,4,6-Tetrachlorophenol	10.12	232	229287	39.09	nq	# 87
59) Diethylphthalate	10.21	149	1000617	39.50	nq	100
60) 4-Chlorophenyl-phenylether	10.33	204	406805	41.01	nq	99
61) 4-Nitroaniline	10.36	138	195998	26.77	nq	84
62) Azobenzene	10.49	77	1160311	41.60	nq	97
64) 4,6-Dinitro-2-methylphenol	10.40	198	78245	22.19	nq	# 61
65) n-Nitrosodiphenylamine	10.45	169	847781	48.99	nq	96
66) 4-Bromophenyl-phenylether	10.82	248	280497	46.23	nq	90
67) Hexachlorobenzene	10.89	284	270830	41.76	nq	96
68) Atrazine	10.98	200	261420	46.83	nq	98
69) Pentachlorophenol	11.08	266	249459	78.40	nq	99
70) Phenanthrene	11.29	178	1299023	41.08	nq	98
71) Anthracene	11.34	178	1337577	49.32	nq	99
72) Carbazole	11.50	167	1144432	43.15	nq	99
73) Di-n-butylphthalate	11.84	149	1576585	53.38	nq	98
74) Fluoranthene	12.48	202	1308286	47.35	nq	94
76) Benzidine	12.60	184	391271	31.30	nq	96
77) Pyrene	12.71	202	1262772	37.81	nq	98
79) Butylbenzylphthalate	13.32	149	620106	40.82	nq	91
80) Benzo(a)anthracene	13.88	228	989623	38.51	nq	100
81) 3,3'-Dichlorobenzidine	13.85	252	271007	27.81	nq	97
82) Chrysene	13.92	228	863385	33.50	nq	98
83) Bis(2-ethylhexyl)phthalate	13.88	149	753082	42.10	nq	# 98
84) Di-n-octyl phthalate	14.49	149	1345853	40.61	nq	100
85) Indeno(1,2,3-cd)pyrene	16.63	276	962719	43.11	nq	99
87) Benzo(b)fluoranthene	14.90	252	1134112m	39.86	nq	
88) Benzo(k)fluoranthene	14.92	252	868816m	35.81	nq	
89) Benzo(a)pyrene	15.23	252	979427	38.78	nq	99
90) Dibenzo(a,h)anthracene	16.64	278	798450	38.46	nq	97
91) Benzo(g,h,i)perylene	17.03	276	778965	36.09	nq	98

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

