

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021517\
 Data File : BF092980.D
 Acq On : 16 Feb 2017 6:54
 Operator : SJ/MA
 Sample : I1790-02MSD
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 38-WCS021017-0-15MSD

Manual Integrations
 APPROVED

mohammad
 2/16/2017 11:39:49 AM

Quant Time: Feb 16 07:32:07 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF013017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 13 18:04:38 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.88	152	148825	20.00	ng	-0.01
21) Naphthalene-d8	8.16	136	586829	20.00	ng	-0.01
38) Acenaphthene-d10	9.90	164	307100	20.00	ng	-0.02
63) Phenanthrene-d10	11.39	188	470862	20.00	ng	-0.01
75) Chrysene-d12	14.03	240	342435	20.00	ng	-0.02
86) Perylene-d12	15.46	264	283933	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.49	112	1262790	145.57	ng	0.01
7) Phenol-d6	6.52	99	1421710	127.38	ng	-0.01
23) Nitrobenzene-d5	7.45	82	1086510	103.10	ng	-0.01
41) 2,4,6-Tribromophenol	10.70	330	317782	143.07	ng	-0.01
44) 2-Fluorobiphenyl	9.23	172	1681736	99.21	ng	-0.01
78) Terphenyl-d14	12.98	244	1328218	91.14	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.64	88	187311	39.96	ng	# 71
3) Pyridine	3.34	79	370027m	31.05	ng	
4) n-Nitrosodimethylamine	3.28	42	272940	48.77	ng	89
6) Aniline	6.54	93	219389	14.41	ng	# 60
8) 2-Chlorophenol	6.66	128	456979	47.75	ng	89
9) Benzaldehyde	6.42	77	37044	4.63	ng	96
10) Phenol	6.53	94	590686	45.23	ng	# 76
11) bis(2-Chloroethyl)ether	6.61	93	461438	44.27	ng	92
12) 1,3-Dichlorobenzene	6.82	146	513488	45.81	ng	97
13) 1,4-Dichlorobenzene	6.89	146	512228	45.15	ng	99
14) 1,2-Dichlorobenzene	7.05	146	510254	47.04	ng	95
15) Benzyl Alcohol	7.02	79	412289	49.89	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.15	45	866689	47.86	ng	97
17) 2-Methylphenol	7.14	107	406292	49.49	ng	96
18) Hexachloroethane	7.38	117	169520	43.77	ng	# 85
19) n-Nitroso-di-n-propylamine	7.30	70	363263	51.13	ng	96
20) 3+4-Methylphenols	7.30	107	473443	48.23	ng	# 85
22) Acetophenone	7.29	105	630647	47.07	ng	# 94
24) Nitrobenzene	7.46	77	501022	46.30	ng	98
25) Isophorone	7.70	82	962680	49.02	ng	96
26) 2-Nitrophenol	7.78	139	282115	54.85	ng	89
27) 2,4-Dimethylphenol	7.82	122	455486	50.42	ng	92
28) bis(2-Chloroethoxy)methane	7.92	93	645306	49.62	ng	98
29) 2,4-Dichlorophenol	8.02	162	402860	47.82	ng	90
30) 1,2,4-Trichlorobenzene	8.10	180	430590	47.43	ng	95
31) Naphthalene	8.18	128	1319444	44.35	ng	99
32) Benzoic acid	7.95	122	274152	52.73	ng	96
33) 4-Chloroaniline	8.22	127	95882	8.22	ng	96
34) Hexachlorobutadiene	8.29	225	215297	43.10	ng	99
35) Caprolactam	8.60	113	117261m	44.25	ng	
36) 4-Chloro-3-methylphenol	8.73	107	457586	52.10	ng	90
37) 2-Methylnaphthalene	8.86	142	929443	48.66	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.04	216	413328	47.95	ng	100
40) Hexachlorocyclopentadiene	9.02	237	212281	54.58	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.15	196	281011	49.61	ng	92
43) 2,4,5-Trichlorophenol	9.20	196	287813	46.37	ng	94
45) 1,1'-Biphenyl	9.33	154	1046068	47.04	ng	97
46) 2-Chloronaphthalene	9.36	162	823750	47.35	ng	96
47) 2-Nitroaniline	9.46	65	301311	51.52	ng	95
48) Acenaphthylene	9.78	152	1403196	46.69	ng	99
49) Dimethylphthalate	9.64	163	1108681	53.60	ng	100
50) 2,6-Dinitrotoluene	9.70	165	212701	47.61	ng	# 82
51) Acenaphthene	9.94	154	868212	48.32	ng	98
52) 3-Nitroaniline	9.87	138	68526	13.40	ng	# 77
53) 2,4-Dinitrophenol	9.97	184	158312	70.28	ng	# 56
54) Dibenzofuran	10.12	168	1194046	49.69	ng	# 89
55) 4-Nitrophenol	10.04	139	283689	77.05	ng	93
56) 2,4-Dinitrotoluene	10.11	165	311163	52.32	ng	# 80
57) Fluorene	10.46	166	871901	47.16	ng	98
58) 2,3,4,6-Tetrachlorophenol	10.24	232	223393	53.96	ng	98
59) Diethylphthalate	10.34	149	1037818	53.24	ng	95
60) 4-Chlorophenyl-phenylether	10.45	204	423778	47.71	ng	# 79
61) 4-Nitroaniline	10.49	138	232650	44.43	ng	85
62) Azobenzene	10.61	77	1013301m	50.31	ng	
64) 4,6-Dinitro-2-methylphenol	10.51	198	108975	38.58	ng	# 49
65) n-Nitrosodiphenylamine	10.57	169	754180	50.14	ng	100
66) 4-Bromophenyl-phenylether	10.94	248	251607	51.15	ng	# 76
67) Hexachlorobenzene	11.00	284	226837	46.66	ng	# 78
68) Atrazine	11.09	200	218310	45.23	ng	99
69) Pentachlorophenol	11.21	266	261087	101.87	ng	99
70) Phenanthrene	11.41	178	1271215	47.12	ng	98
71) Anthracene	11.47	178	1277157	47.14	ng	98
72) Carbazole	11.63	167	1218634	47.06	ng	98
73) Di-n-butylphthalate	11.95	149	1438978	51.38	ng	98
74) Fluoranthene	12.60	202	1072787	38.55	ng	99
76) Benzidine	12.73	184	36403	2.49	ng	98
77) Pyrene	12.83	202	1059941	41.36	ng	99
79) Butylbenzylphthalate	13.45	149	536975	51.60	ng	96
80) Benzo(a)anthracene	14.02	228	923267	44.08	ng	99
81) 3,3'-Dichlorobenzidine	13.99	252	209284	28.03	ng	# 95
82) Chrysene	14.05	228	747228	38.11	ng	99
83) Bis(2-ethylhexyl)phthalate	14.01	149	760541	53.44	ng	# 97
84) Di-n-octyl phthalate	14.61	149	1216033	52.47	ng	100
85) Indeno(1,2,3-cd)pyrene	16.88	276	796308	52.43	ng	# 94
87) Benzo(b)fluoranthene	15.05	252	814421m	43.58	ng	
88) Benzo(k)fluoranthene	15.08	252	854213m	52.94	ng	
89) Benzo(a)pyrene	15.40	252	786052	48.63	ng	# 96
90) Dibenzo(a,h)anthracene	16.90	278	673447	51.55	ng	99
91) Benzo(g,h,i)perylene	17.31	276	663110	50.24	ng	# 92

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

