

Data Path : Z:\HPCHEM1\BNA F\DATA\BF101116\
 Data File : BF090515.D
 Acq On : 11 Oct 2016 21:19
 Operator : UM/SJ
 Sample : H5146-03MSD
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S8-0617MSD

Manual Integrations
 APPROVED

sohil
 10/12/2016 4:09:17 PM

Quant Time: Oct 12 01:27:13 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF101116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 11 16:14:43 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	233537	20.00	ng	0.00
21) Naphthalene-d8	8.25	136	962803	20.00	ng	0.00
38) Acenaphthene-d10	10.01	164	500240	20.00	ng	0.00
63) Phenanthrene-d10	11.49	188	888180	20.00	ng	0.01
75) Chrysene-d12	14.13	240	794885	20.00	ng	0.00
86) Perylene-d12	15.61	264	519006	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.56	112	770928	51.36	ng	0.00
7) Phenol-d6	6.58	99	607620	32.08	ng	-0.01
23) Nitrobenzene-d5	7.51	82	1477035	83.21	ng	0.00
41) 2,4,6-Tribromophenol	10.79	330	695757	140.31	ng	0.00
44) 2-Fluorobiphenyl	9.32	172	2457849	82.57	ng	0.00
78) Terphenyl-d14	13.08	244	3054590	93.91	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.66	88	93590	12.71	ng	# 85
3) Pyridine	3.41	79	240042m	11.80	ng	
4) n-Nitrosodimethylamine	3.34	42	103217	15.48	ng	# 48
6) Aniline	6.61	93	525999	21.89	ng	96
8) 2-Chlorophenol	6.74	128	524398	32.84	ng	94
9) Benzaldehyde	6.50	77	237706	19.92	ng	96
10) Phenol	6.59	94	246274	11.53	ng	94
11) bis(2-Chloroethyl)ether	6.68	93	584549	36.55	ng	94
12) 1,3-Dichlorobenzene	6.90	146	589714	34.23	ng	98
13) 1,4-Dichlorobenzene	6.97	146	559725	31.84	ng	98
14) 1,2-Dichlorobenzene	7.13	146	610149	38.20	ng	97
15) Benzyl Alcohol	7.09	79	374596	25.75	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.23	45	847641	36.98	ng	88
17) 2-Methylphenol	7.21	107	339416	26.13	ng	98
18) Hexachloroethane	7.47	117	219124	33.78	ng	100
19) n-Nitroso-di-n-propylamine	7.37	70	483668	40.03	ng	# 80
20) 3+4-Methylphenols	7.35	107	371142	22.57	ng	# 45
22) Acetophenone	7.37	105	910269	42.94	ng	# 84
24) Nitrobenzene	7.54	77	777550	40.84	ng	91
25) Isophorone	7.78	82	1353082	41.88	ng	99
26) 2-Nitrophenol	7.86	139	356239	41.01	ng	98
27) 2,4-Dimethylphenol	7.89	122	508296	34.42	ng	98
28) bis(2-Chloroethoxy)methane	7.98	93	822661	41.23	ng	# 96
29) 2,4-Dichlorophenol	8.10	162	489552	39.64	ng	95
30) 1,2,4-Trichlorobenzene	8.18	180	505567	35.64	ng	100
31) Naphthalene	8.27	128	1752339	36.53	ng	98
32) Benzoic acid	7.96	122	79038	6.78	ng	97
33) 4-Chloroaniline	8.30	127	502424	24.74	ng	98
34) Hexachlorobutadiene	8.38	225	272539	34.36	ng	99
35) Caprolactam	8.67	113	22513	5.62	ng	# 84
36) 4-Chloro-3-methylphenol	8.78	107	473441	32.83	ng	97
37) 2-Methylnaphthalene	8.95	142	1164307	36.16	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.13	216	596565	42.60	ng	98
40) Hexachlorocyclopentadiene	9.11	237	617330	88.81	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.23	196	365769	38.74	nq	96
43) 2,4,5-Trichlorophenol	9.27	196	375919	40.23	nq #	91
45) 1,1'-Biphenyl	9.42	154	1608307	42.10	nq	96
46) 2-Chloronaphthalene	9.45	162	1170122	41.16	nq	98
47) 2-Nitroaniline	9.54	65	407207	40.16	nq	97
48) Acenaphthylene	9.87	152	1832855	39.96	nq	97
49) Dimethylphthalate	9.72	163	1350842	40.15	nq	99
50) 2,6-Dinitrotoluene	9.78	165	304485	40.99	nq	96
51) Acenaphthene	10.04	154	1323622	43.06	nq	100
52) 3-Nitroaniline	9.95	138	246023	26.69	nq	97
53) 2,4-Dinitrophenol	10.05	184	346432	84.04	nq #	82
54) Dibenzofuran	10.21	168	1788591	43.73	nq	98
55) 4-Nitrophenol	10.11	139	218193	30.46	nq #	71
56) 2,4-Dinitrotoluene	10.19	165	502818	51.18	nq #	89
57) Fluorene	10.55	166	1429246	42.80	nq	100
58) 2,3,4,6-Tetrachlorophenol	10.33	232	367365	47.13	nq	98
59) Diethylphthalate	10.42	149	1490487	43.56	nq #	92
60) 4-Chlorophenyl-phenylether	10.54	204	720400	46.76	nq	91
61) 4-Nitroaniline	10.57	138	340046	36.16	nq	94
62) Azobenzene	10.70	77	1660468	44.44	nq	88
64) 4,6-Dinitro-2-methylphenol	10.60	198	236377	39.85	nq #	33
65) n-Nitrosodiphenylamine	10.66	169	1132430	39.21	nq	99
66) 4-Bromophenyl-phenylether	11.03	248	462789	47.02	nq #	89
67) Hexachlorobenzene	11.10	284	492569	45.31	nq #	90
68) Atrazine	11.18	200	390184	43.65	nq	98
69) Pentachlorophenol	11.30	266	591084	91.37	nq	98
70) Phenanthrene	11.51	178	2038586	41.65	nq	99
71) Anthracene	11.56	178	1996683	40.26	nq	99
72) Carbazole	11.72	167	2043343	43.63	nq	98
73) Di-n-butylphthalate	12.05	149	2530326	44.86	nq #	98
74) Fluoranthene	12.70	202	2152461	43.08	nq	96
76) Benzidine	12.82	184	678840	27.17	nq	99
77) Pyrene	12.93	202	2073145	41.67	nq	100
79) Butylbenzylphthalate	13.55	149	1043556	44.80	nq	91
80) Benzo(a)anthracene	14.12	228	1900655	40.12	nq	99
81) 3,3'-Dichlorobenzidine	14.09	252	481810	28.26	nq #	98
82) Chrysene	14.16	228	1725884	39.14	nq	99
83) Bis(2-ethylhexyl)phthalate	14.11	149	1404573	41.32	nq	99
84) Di-n-octyl phthalate	14.73	149	2283323	43.37	nq #	87
85) Indeno(1,2,3-cd)pyrene	17.08	276	1163470	31.69	nq	97
87) Benzo(b)fluoranthene	15.17	252	1750851	51.13	nq #	96
88) Benzo(k)fluoranthene	15.21	252	1189136m	38.98	nq	
89) Benzo(a)pyrene	15.55	252	1354934	45.06	nq	97
90) Dibenzo(a,h)anthracene	17.10	278	1000103	39.15	nq	98
91) Benzo(g,h,i)perylene	17.53	276	933423	39.51	nq	96

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

