

Data Path : U:\HPCHEM1\BNA F\DATA\BF101915\
 Data File : BF082368.D
 Acq On : 19 Oct 2015 21:16
 Operator : UM/IZ
 Sample : G3993-02MS
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 STOCKPILEMS

Manual Integrations
 APPROVED

MMdadoda
 10/20/2015 6:25:48 PM

Quant Time: Oct 20 03:03:10 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF101015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 20 01:57:16 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.81	152	88380m	20.00	ng	0.01
21) Naphthalene-d8	8.09	136	352507	20.00	ng	0.00
38) Acenaphthene-d10	9.85	164	178433	20.00	ng	0.00
63) Phenanthrene-d10	11.34	188	349374	20.00	ng	-0.01
75) Chrysene-d12	14.02	240	315669	20.00	ng	-0.07
86) Perylene-d12	15.53	264	260614	20.00	ng	-0.14

System Monitoring Compounds

5) 2-Fluorophenol	5.43	112	544547	124.35	ng	0.03
7) Phenol-d6	6.47	99	626958	110.02	ng	0.02
23) Nitrobenzene-d5	7.38	82	507134	96.61	ng	0.00
41) 2,4,6-Tribromophenol	10.65	330	183230	109.50	ng	0.00
44) 2-Fluorobiphenyl	9.18	172	1016928	94.51	ng	0.00
78) Terphenyl-d14	12.94	244	1006336	84.48	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.50	88	80268	36.48	ng	# 85
3) Pyridine	3.54	79	132679m	25.93	ng	
4) n-Nitrosodimethylamine	3.13	42	95203	43.87	ng	# 78
6) Aniline	6.55	93	259326	35.30	ng	# 55
8) 2-Chlorophenol	6.61	128	209277	39.15	ng	95
9) Benzaldehyde	6.35	77	49323	16.95	ng	91
10) Phenol	6.48	94	237845	36.20	ng	92
11) bis(2-Chloroethyl)ether	6.55	93	259326	46.67	ng	94
12) 1,3-Dichlorobenzene	6.74	146	260221m	41.80	ng	
13) 1,4-Dichlorobenzene	6.82	146	270106m	41.54	ng	
14) 1,2-Dichlorobenzene	6.98	146	240364	38.93	ng	# 90
15) Benzyl Alcohol	6.98	79	358538	91.14	ng	96
16) 2,2'-oxybis(1-Chloropropan	7.10	45	362856	46.90	ng	65
17) 2-Methylphenol	7.09	107	169725m	40.15	ng	
18) Hexachloroethane	7.31	117	93020	41.80	ng	94
19) n-Nitroso-di-n-propylamine	7.23	70	175140m	43.76	ng	
20) 3+4-Methylphenols	7.25	107	298106	59.18	ng	# 62
22) Acetophenone	7.23	105	339970	48.75	ng	# 88
24) Nitrobenzene	7.41	77	244061	42.95	ng	# 81
25) Isophorone	7.65	82	465077	43.90	ng	99
26) 2-Nitrophenol	7.71	139	118738	41.85	ng	# 89
27) 2,4-Dimethylphenol	7.76	122	205819	45.03	ng	97
28) bis(2-Chloroethoxy)methane	7.85	93	293853	47.67	ng	99
29) 2,4-Dichlorophenol	7.97	162	202440	44.30	ng	99
30) 1,2,4-Trichlorobenzene	8.03	180	228377	43.36	ng	98
31) Naphthalene	8.11	128	674015	44.11	ng	100
32) Benzoic acid	7.94	122	276362	90.09	ng	98
33) 4-Chloroaniline	8.19	127	20876	3.29	ng	# 91
34) Hexachlorobutadiene	8.23	225	138364	42.16	ng	97
35) Caprolactam	8.61	113	54205	40.88	ng	# 74
36) 4-Chloro-3-methylphenol	8.67	107	197778	44.26	ng	# 83
37) 2-Methylnaphthalene	8.81	142	505481	47.72	ng	97
39) 1,2,4,5-Tetrachlorobenzene	8.98	216	225349	42.46	ng	98
40) Hexachlorocyclopentadiene	8.96	237	164271	63.12	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.10	196	149543	42.42	nq	99
43) 2,4,5-Trichlorophenol	9.14	196	153365	40.16	nq	# 91
45) 1,1'-Biphenyl	9.28	154	655794	48.20	nq	96
46) 2-Chloronaphthalene	9.30	162	499170	46.60	nq	96
47) 2-Nitroaniline	9.41	65	137571	46.78	nq	92
48) Acenaphthylene	9.71	152	715498	42.88	nq	100
49) Dimethylphthalate	9.59	163	695167	55.33	nq	99
50) 2,6-Dinitrotoluene	9.65	165	121896	45.98	nq	# 70
51) Acenaphthene	9.89	154	433059	43.23	nq	99
52) 3-Nitroaniline	9.82	138	34352	11.47	nq	# 80
53) 2,4-Dinitrophenol	9.93	184	128581	86.25	nq	98
54) Dibenzofuran	10.06	168	632282	45.98	nq	96
55) 4-Nitrophenol	10.00	139	132786	68.21	nq	93
56) 2,4-Dinitrotoluene	10.06	165	179808	54.26	nq	94
57) Fluorene	10.40	166	511917	45.32	nq	98
58) 2,3,4,6-Tetrachlorophenol	10.18	232	128771	41.27	nq	97
59) Diethylphthalate	10.29	149	617874	52.76	nq	98
60) 4-Chlorophenyl-phenylether	10.40	204	272025	52.57	nq	# 84
61) 4-Nitroaniline	10.43	138	113953	39.23	nq	91
62) Azobenzene	10.56	77	571740	49.88	nq	89
64) 4,6-Dinitro-2-methylphenol	10.47	198	83683	42.68	nq	88
65) n-Nitrosodiphenylamine	10.53	169	487398	46.59	nq	99
66) 4-Bromophenyl-phenylether	10.89	248	159854	45.72	nq	# 85
67) Hexachlorobenzene	10.95	284	164242	43.43	nq	# 73
68) Atrazine	11.05	200	127855	38.58	nq	99
69) Pentachlorophenol	11.15	266	167211	85.93	nq	100
70) Phenanthrene	11.37	178	857113	50.05	nq	99
71) Anthracene	11.42	178	803299	45.52	nq	99
72) Carbazole	11.58	167	704920	43.79	nq	100
73) Di-n-butylphthalate	11.91	149	943068	47.49	nq	100
74) Fluoranthene	12.56	202	935632	50.87	nq	99
77) Pyrene	12.79	202	929474	49.62	nq	99
79) Butylbenzylphthalate	13.42	149	397029	46.44	nq	95
80) Benzo(a)anthracene	14.01	228	726024	44.20	nq	100
81) 3,3'-Dichlorobenzidine	13.98	252	92854	14.89	nq	99
82) Chrysene	14.06	228	759493	43.89	nq	99
83) Bis(2-ethylhexyl)phthalate	14.00	149	577957	47.38	nq	# 98
84) Di-n-octyl phthalate	14.68	149	971772	46.39	nq	100
85) Indeno(1,2,3-cd)pyrene	16.97	276	617409	44.89	nq	98
87) Benzo(b)fluoranthene	15.12	252	630773m	39.78	nq	
88) Benzo(k)fluoranthene	15.16	252	686892m	51.54	nq	
89) Benzo(a)pyrene	15.49	252	620329	45.52	nq	# 95
90) Dibenzo(a,h)anthracene	16.98	278	517598	46.35	nq	98
91) Benzo(g,h,i)perylene	17.40	276	508863	46.91	nq	99

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

