

Data Path : Z:\HPCHEM1\BNA F\DATA\BF012015\
 Data File : BF076982.D
 Acq On : 21 Jan 2015 3:39
 Operator : IZ / TP
 Sample : G1101-09MS
 Misc : W
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TE-PMW-11MMS

Manual Integrations
 APPROVED

tejaskumar
 1/21/2015 5:30:10 PM

Quant Time: Jan 21 04:59:05 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF011415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 21 04:28:05 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	31126	20.00	ng	0.00
21) Naphthalene-d8	10.82	136	134489	20.00	ng	0.00
38) Acenaphthene-d10	14.95	164	72183	20.00	ng	0.00
63) Phenanthrene-d10	17.74	188	129616	20.00	ng	0.00
75) Chrysene-d12	21.24	240	131379	20.00	ng	0.00
86) Perylene-d12	22.91	264	114077	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	4.96	112	107397	56.73	ng	0.00
7) Phenol-d6	7.32	99	83524	34.97	ng	0.00
23) Nitrobenzene-d5	9.21	82	209108	86.14	ng	0.00
41) 2,4,6-Tribromophenol	16.64	330	78115	133.32	ng	0.00
44) 2-Fluorobiphenyl	13.48	172	442299	93.25	ng	0.00
78) Terphenyl-d14	19.97	244	474728	83.19	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	1.25	88	9859	12.16	ng	96
3) Pyridine	1.76	79	14780	7.05	ng	99
4) n-Nitrosodimethylamine	1.68	42	17698	17.03	ng	99
6) Aniline	7.20	93	17922	5.42	ng	97
8) 2-Chlorophenol	7.44	128	75965	34.81	ng	93
9) Benzaldehyde	6.92	77	28133	18.56	ng	96
10) Phenol	7.35	94	29920	11.80	ng	94
11) bis(2-Chloroethyl)ether	7.44	93	88890	44.80	ng	87
12) 1,3-Dichlorobenzene	7.75	146	103052	41.50	ng	94
13) 1,4-Dichlorobenzene	7.94	146	102684	39.00	ng	99
14) 1,2-Dichlorobenzene	8.26	146	98899	40.76	ng	97
15) Benzyl Alcohol	8.34	79	49667	25.70	ng	99
16) 2,2'-oxybis(1-Chloropropan	8.70	45	116355	43.62	ng	81
17) 2-Methylphenol	8.70	107	44707	25.07	ng	# 86
18) Hexachloroethane	9.01	117	38104	41.61	ng	98
19) n-Nitroso-di-n-propylamine	8.97	70	72238	43.21	ng	94
20) 3+4-Methylphenols	9.09	107	51954	22.00	ng	92
22) Acetophenone	8.89	105	146614	44.51	ng	95
24) Nitrobenzene	9.25	77	109761	44.38	ng	96
25) Isophorone	9.85	82	195316	44.18	ng	97
26) 2-Nitrophenol	9.99	139	50565	40.23	ng	# 82
27) 2,4-Dimethylphenol	10.28	122	55937	29.25	ng	98
28) bis(2-Chloroethoxy)methane	10.47	93	117817	46.89	ng	97
29) 2,4-Dichlorophenol	10.60	162	72967	40.94	ng	99
30) 1,2,4-Trichlorobenzene	10.73	180	83908	42.39	ng	93
31) Naphthalene	10.87	128	289614	41.83	ng	98
32) Benzoic acid	10.57	122	12990	12.26	ng	95
33) 4-Chloroaniline	11.12	127	26734	9.56	ng	93
34) Hexachlorobutadiene	11.24	225	48257	41.08	ng	96
35) Caprolactam	11.94	113	3037	5.79	ng	95
36) 4-Chloro-3-methylphenol	12.40	107	75836	37.16	ng	89
37) 2-Methylnaphthalene	12.53	142	213883	45.23	ng	94
39) 1,2,4,5-Tetrachlorobenzene	12.93	216	82850	46.42	ng	98
40) Hexachlorocyclopentadiene	12.92	237	82250	85.29	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.27	196	53069	40.81	nq	92
43) 2,4,5-Trichlorophenol	13.36	196	58395	44.03	nq	96
45) 1,1'-Biphenyl	13.67	154	246378	46.04	nq	97
46) 2-Chloronaphthalene	13.65	162	199183	45.23	nq	98
47) 2-Nitroaniline	13.99	65	63602	47.61	nq	97
48) Acenaphthylene	14.61	152	322453	43.41	nq	98
49) Dimethylphthalate	14.55	163	253898	49.60	nq	99
50) 2,6-Dinitrotoluene	14.64	165	50910	49.78	nq	92
51) Acenaphthene	15.03	154	184148	43.70	nq	98
52) 3-Nitroaniline	15.00	138	30285	23.63	nq	94
53) 2,4-Dinitrophenol	15.27	184	43502	89.07	nq	92
54) Dibenzofuran	15.45	168	283765	46.22	nq	99
55) 4-Nitrophenol	15.60	139	24322	30.31	nq	93
56) 2,4-Dinitrotoluene	15.58	165	71442	46.66	nq	95
57) Fluorene	16.17	166	237120	45.59	nq	99
58) 2,3,4,6-Tetrachlorophenol	15.80	232	43847	45.41	nq	# 95
59) Diethylphthalate	16.17	149	256291	49.54	nq	99
60) 4-Chlorophenyl-phenylether	16.27	204	101813	47.84	nq	94
61) 4-Nitroaniline	16.32	138	50190	39.20	nq	96
62) Azobenzene	16.55	77	249916	46.36	nq	98
64) 4,6-Dinitro-2-methylphenol	16.39	198	32766	39.93	nq	99
65) n-Nitrosodiphenylamine	16.51	169	134023	29.00	nq	98
66) 4-Bromophenyl-phenylether	17.11	248	56965	43.38	nq	85
67) Hexachlorobenzene	17.12	284	62252	44.99	nq	91
68) Atrazine	17.48	200	69502	49.56	nq	93
69) Pentachlorophenol	17.49	266	48182	76.26	nq	93
70) Phenanthrene	17.77	178	345153	42.70	nq	99
71) Anthracene	17.84	178	337077	42.76	nq	98
72) Carbazole	18.14	167	323929	42.37	nq	99
73) Di-n-butylphthalate	18.72	149	423117	47.81	nq	99
74) Fluoranthene	19.42	202	372586	44.98	nq	98
76) Benzidine	19.66	184	12570	2.99	nq	95
77) Pyrene	19.69	202	377627	41.94	nq	98
79) Butylbenzylphthalate	20.63	149	193925	47.56	nq	99
80) Benzo(a)anthracene	21.23	228	342305	41.53	nq	100
81) 3,3'-Dichlorobenzidine	21.24	252	49541	18.91	nq	94
82) Chrysene	21.27	228	324678	43.19	nq	98
83) Bis(2-ethylhexyl)phthalate	21.37	149	274949	48.73	nq	100
84) Di-n-octyl phthalate	22.14	149	473473	48.35	nq	97
85) Indeno(1,2,3-cd)pyrene	24.04	276	318905	40.03	nq	# 85
87) Benzo(b)fluoranthene	22.48	252	310865m	40.46	nq	
88) Benzo(k)fluoranthene	22.52	252	313616m	46.72	nq	
89) Benzo(a)pyrene	22.84	252	297152	43.85	nq	# 96
90) Dibenzo(a,h)anthracene	24.06	278	255145	41.03	nq	# 95
91) Benzo(g,h,i)perylene	24.32	276	266600	43.13	nq	# 92

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

