

Data Path : Z:\HPCHEM1\BNA F\DATA\BF031115\
 Data File : BF077697.D
 Acq On : 11 Mar 2015 20:00
 Operator : TP/IZ
 Sample : G1475-01MS
 Misc : S
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-GF-01-012-AMS

Manual Integrations
 APPROVED

mohammad
 3/13/2015 9:40:27 AM

Quant Time: Mar 12 05:46:16 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 12 02:24:12 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.17	152	57956	20.00	ng	0.00
21) Naphthalene-d8	8.75	136	218473	20.00	ng	0.00
38) Acenaphthene-d10	10.91	164	114329	20.00	ng	0.00
63) Phenanthrene-d10	12.75	188	225195	20.00	ng	0.00
75) Chrysene-d12	16.03	240	234768	20.00	ng	0.00
86) Perylene-d12	17.79	264	226528	20.00	ng	0.07

System Monitoring Compounds

5) 2-Fluorophenol	5.46	112	315951	95.16	ng	0.00
7) Phenol-d6	6.74	99	423481	103.23	ng	0.00
23) Nitrobenzene-d5	7.86	82	205476	61.65	ng	0.00
41) 2,4,6-Tribromophenol	11.89	330	104058	95.13	ng	0.00
44) 2-Fluorobiphenyl	10.09	172	455900	58.60	ng	0.00
78) Terphenyl-d14	14.75	244	564929	58.78	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.06	88	37722	25.02	ng	# 100
3) Pyridine	2.76	79	107859	26.63	ng	96
4) n-Nitrosodimethylamine	2.68	42	51138	29.00	ng	84
6) Aniline	6.75	93	132226	22.53	ng	# 1
8) 2-Chlorophenol	6.90	128	121787	30.82	ng	97
9) Benzaldehyde	6.61	77	12931	7.69	ng	99
10) Phenol	6.75	94	157546	32.49	ng	91
11) bis(2-Chloroethyl)ether	6.86	93	109754	30.22	ng	99
12) 1,3-Dichlorobenzene	7.09	146	127320	28.96	ng	99
13) 1,4-Dichlorobenzene	7.20	146	123393	27.02	ng	98
14) 1,2-Dichlorobenzene	7.38	146	110575	26.41	ng	98
15) Benzyl Alcohol	7.36	79	88692	27.70	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.54	45	128766	26.31	ng	99
17) 2-Methylphenol	7.50	107	88902	27.63	ng	99
18) Hexachloroethane	7.79	117	39246	26.20	ng	97
19) n-Nitroso-di-n-propylamine	7.69	70	75852	26.73	ng	97
20) 3+4-Methylphenols	7.70	107	119450	28.29	ng	98
22) Acetophenone	7.68	105	149542	30.92	ng	# 98
24) Nitrobenzene	7.88	77	104023	28.97	ng	98
25) Isophorone	8.18	82	195636	29.07	ng	# 90
26) 2-Nitrophenol	8.28	139	54484	30.99	ng	97
27) 2,4-Dimethylphenol	8.35	122	93497	29.20	ng	99
28) bis(2-Chloroethoxy)methane	8.46	93	124836	30.56	ng	99
29) 2,4-Dichlorophenol	8.58	162	98230	32.18	ng	97
30) 1,2,4-Trichlorobenzene	8.68	180	100390	29.39	ng	98
31) Naphthalene	8.77	128	322872	28.77	ng	100
32) Benzoic acid	8.44	122	34659	21.24	ng	89
33) 4-Chloroaniline	8.85	127	101684	22.33	ng	98
34) Hexachlorobutadiene	8.93	225	52859	27.31	ng	98
35) Caprolactam	9.25	113	28124	31.52	ng	98
36) 4-Chloro-3-methylphenol	9.46	107	96636	31.46	ng	97
37) 2-Methylnaphthalene	9.63	142	226069	29.99	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.84	216	97918	28.71	ng	# 100
40) Hexachlorocyclopentadiene	9.82	237	91667	54.74	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.98	196	67666	28.65	nq	98
43) 2,4,5-Trichlorophenol	10.03	196	74079	29.03	nq	98
45) 1,1'-Biphenyl	10.21	154	262616	28.73	nq	99
46) 2-Chloronaphthalene	10.22	162	203959	27.46	nq	# 92
47) 2-Nitroaniline	10.36	65	59446	32.23	nq	96
48) Acenaphthylene	10.74	152	334412	27.08	nq	99
49) Dimethylphthalate	10.60	163	264573	31.20	nq	100
50) 2,6-Dinitrotoluene	10.67	165	54873	31.22	nq	92
51) Acenaphthene	10.96	154	193282	27.29	nq	98
52) 3-Nitroaniline	10.88	138	52014	25.48	nq	91
53) 2,4-Dinitrophenol	11.00	184	33432	54.62	nq	97
54) Dibenzofuran	11.17	168	302936	28.84	nq	99
55) 4-Nitrophenol	11.09	139	98852	65.37	nq	99
56) 2,4-Dinitrotoluene	11.16	165	69525	30.33	nq	# 91
57) Fluorene	11.60	166	246016	28.21	nq	99
58) 2,3,4,6-Tetrachlorophenol	11.32	232	61199	31.15	nq	# 100
59) Diethylphthalate	11.48	149	241993	29.29	nq	98
60) 4-Chlorophenyl-phenylether	11.61	204	112879	28.36	nq	96
61) 4-Nitroaniline	11.62	138	56570	27.81	nq	79
62) Azobenzene	11.80	77	219959	27.86	nq	98
64) 4,6-Dinitro-2-methylphenol	11.67	198	29473	29.53	nq	92
65) n-Nitrosodiphenylamine	11.76	169	212565	28.35	nq	99
66) 4-Bromophenyl-phenylether	12.21	248	67917	28.26	nq	95
67) Hexachlorobenzene	12.27	284	74282	28.13	nq	93
68) Atrazine	12.43	200	69241	32.95	nq	98
69) Pentachlorophenol	12.52	266	80024	58.71	nq	98
70) Phenanthrene	12.79	178	367465	28.42	nq	100
71) Anthracene	12.84	178	358924	28.10	nq	99
72) Carbazole	13.05	167	337950	27.82	nq	99
73) Di-n-butylphthalate	13.51	149	392584	28.82	nq	99
74) Fluoranthene	14.26	202	386963	27.14	nq	99
76) Benzidine	14.43	184	174105	29.86	nq	99
77) Pyrene	14.53	202	414513	28.08	nq	100
79) Butylbenzylphthalate	15.37	149	170578	29.31	nq	98
80) Benzo(a)anthracene	16.02	228	377195	27.10	nq	99
81) 3,3'-Dichlorobenzidine	16.01	252	110229	24.01	nq	# 98
82) Chrysene	16.07	228	357833	28.60	nq	99
83) Bis(2-ethylhexyl)phthalate	16.09	149	257078	29.16	nq	# 95
84) Di-n-octyl phthalate	16.89	149	428032	28.39	nq	99
85) Indeno(1,2,3-cd)pyrene	19.17	276	412205m	26.39	nq	
87) Benzo(b)fluoranthene	17.33	252	451866	30.56	nq	97
88) Benzo(k)fluoranthene	17.37	252	273566m	23.68	nq	
89) Benzo(a)pyrene	17.72	252	358967	29.09	nq	99
90) Dibenzo(a,h)anthracene	19.21	278	340116	27.79	nq	99
91) Benzo(g,h,i)perylene	19.59	276	340030	27.29	nq	99

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

