

Data Path : Z:\HPCHEM1\BNA F\DATA\BF031117\
 Data File : BF093583.D
 Acq On : 11 Mar 2017 22:33
 Operator : SJ/MA
 Sample : I2106-01MS
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HY-SB419AMS

Manual Integrations
 APPROVED

mohammad
 3/15/2017 5:33:02 PM

Quant Time: Mar 13 19:21:55 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF030717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 13 15:05:08 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.72	152	163727	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	321234	20.00	ng	0.01
38) Acenaphthene-d10	9.81	164	128764	20.00	ng	0.05
63) Phenanthrene-d10	11.28	188	236695	20.00	ng	0.03
75) Chrysene-d12	13.89	240	321844	20.00	ng	0.00
86) Perylene-d12	15.28	264	266408	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.32	112	1362423	138.49	ng	0.00
7) Phenol-d6	6.39	99	1583094	127.61	ng	0.01
23) Nitrobenzene-d5	7.30	82	822887	142.13	ng	0.00
41) 2,4,6-Tribromophenol	10.61	330	121157	144.42	ng	0.06
44) 2-Fluorobiphenyl	9.11	172	783058	113.12	ng	0.02
78) Terphenyl-d14	12.84	244	766033	61.88	ng	0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.23	88	201140	37.12	ng	# 82
3) Pyridine	2.92	79	472113	34.17	ng	89
4) n-Nitrosodimethylamine	2.87	42	271833	41.20	ng	82
6) Aniline	6.40	93	168243	9.88	ng	# 1
8) 2-Chlorophenol	6.52	128	452350	41.04	ng	91
9) Benzaldehyde	6.28	77	98013	9.59	ng	84
10) Phenol	6.41	94	745627	49.05	ng	# 75
11) bis(2-Chloroethyl)ether	6.47	93	522317	42.51	ng	95
12) 1,3-Dichlorobenzene	6.66	146	478197m	37.90	ng	
13) 1,4-Dichlorobenzene	6.74	146	494788	38.31	ng	99
14) 1,2-Dichlorobenzene	6.89	146	403262m	35.25	ng	
15) Benzyl Alcohol	6.89	79	336655	38.07	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.01	45	721000	35.33	ng	76
17) 2-Methylphenol	7.02	107	337567	36.21	ng	# 88
18) Hexachloroethane	7.22	117	160886	36.93	ng	# 88
19) n-Nitroso-di-n-propylamine	7.16	70	424569	49.78	ng	# 92
20) 3+4-Methylphenols	7.18	107	481055	39.79	ng	# 75
22) Acetophenone	7.14	105	578097	74.78	ng	# 95
24) Nitrobenzene	7.33	77	403071	61.94	ng	# 82
25) Isophorone	7.57	82	1020277	87.39	ng	# 80
26) 2-Nitrophenol	7.65	139	199878	67.33	ng	# 30
27) 2,4-Dimethylphenol	7.69	122	349469	72.36	ng	97
28) bis(2-Chloroethoxy)methane	7.77	93	482968	65.53	ng	# 53
29) 2,4-Dichlorophenol	7.91	162	506673	111.52	ng	# 81
30) 1,2,4-Trichlorobenzene	7.97	180	258314	53.46	ng	# 95
31) Naphthalene	8.05	128	1128289	70.09	ng	# 85
32) Benzoic acid	7.83	122	125997	50.21	ng	# 39
33) 4-Chloroaniline	8.12	127	105890	16.21	ng	# 30
34) Hexachlorobutadiene	8.16	225	150219	60.36	ng	98
35) Caprolactam	8.47	113	356383	229.68	ng	# 1
36) 4-Chloro-3-methylphenol	8.62	107	234003	48.25	ng	94
37) 2-Methylnaphthalene	8.76	142	494026	48.23	ng	# 93
39) 1,2,4,5-Tetrachlorobenzene	8.93	216	170932	48.61	ng	# 87
40) Hexachlorocyclopentadiene	8.90	237	13120	24.64	ng	96

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.04	196	104888	47.75	nq	88
43) 2,4,5-Trichlorophenol	9.10	196	97624	41.42	nq #	1
45) 1,1'-Biphenyl	9.21	154	432254	46.08	nq	93
46) 2-Chloronaphthalene	9.25	162	392906	54.72	nq #	82
47) 2-Nitroaniline	9.35	65	109043	42.96	nq #	70
48) Acenaphthylene	9.67	152	458865	38.75	nq #	68
49) Dimethylphthalate	9.52	163	238030	27.88	nq #	70
50) 2,6-Dinitrotoluene	9.60	165	102830	54.90	nq #	34
51) Acenaphthene	9.84	154	430838	61.53	nq #	76
52) 3-Nitroaniline	9.76	138	111686	49.63	nq #	70
53) 2,4-Dinitrophenol	9.89	184	2465	2.89	nq #	1
54) Dibenzofuran	10.01	168	479691	54.73	nq #	45
55) 4-Nitrophenol	10.01	139	368111	336.78	nq #	39
56) 2,4-Dinitrotoluene	10.04	165	137587	59.79	nq #	81
57) Fluorene	10.36	166	592729	79.81	nq #	86
58) 2,3,4,6-Tetrachlorophenol	10.15	232	48669	29.26	nq #	1
59) Diethylphthalate	10.22	149	256090	31.39	nq #	88
60) 4-Chlorophenyl-phenylether	10.34	204	126525	38.01	nq #	65
61) 4-Nitroaniline	10.39	138	58842m	25.57	nq	
62) Azobenzene	10.50	77	296075	31.94	nq #	34
64) 4,6-Dinitro-2-methylphenol	10.45	198	7475	4.87	nq #	1
65) n-Nitrosodiphenylamine	10.47	169	720861	91.21	nq #	62
66) 4-Bromophenyl-phenylether	10.82	248	100190	40.03	nq #	58
67) Hexachlorobenzene	10.93	284	66788	27.94	nq #	40
68) Atrazine	11.00	200	132171	57.14	nq #	75
69) Pentachlorophenol	11.11	266	63184	76.57	nq	97
70) Phenanthrene	11.31	178	1612313	119.34	nq #	95
71) Anthracene	11.36	178	557751	40.81	nq #	87
72) Carbazole	11.52	167	424842	33.09	nq #	86
73) Di-n-butylphthalate	11.83	149	684183	46.06	nq	99
74) Fluoranthene	12.48	202	832081	63.76	nq	96
76) Benzidine	12.60	184	96199	9.09	nq #	75
77) Pyrene	12.70	202	982846	41.99	nq	98
79) Butylbenzylphthalate	13.31	149	379315	38.50	nq #	85
80) Benzo(a)anthracene	13.88	228	757117	39.83	nq	99
81) 3,3'-Dichlorobenzidine	13.84	252	163705	24.59	nq #	97
82) Chrysene	13.91	228	629967	35.82	nq	99
83) Bis(2-ethylhexyl)phthalate	13.85	149	549621	40.02	nq	99
84) Di-n-octyl phthalate	14.47	149	873293	38.15	nq	99
85) Indeno(1,2,3-cd)pyrene	16.63	276	598928	40.69	nq	95
87) Benzo(b)fluoranthene	14.89	252	653490m	40.28	nq	
88) Benzo(k)fluoranthene	14.92	252	611130m	39.06	nq	
89) Benzo(a)pyrene	15.24	252	617262	41.63	nq	99
90) Dibenzo(a,h)anthracene	16.63	278	508043	41.41	nq #	94
91) Benzo(g,h,i)perylene	17.04	276	500040	39.71	nq #	93

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

