

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062316\
 Data File : BF088318.D
 Acq On : 24 Jun 2016 6:37
 Operator : UM/SJ
 Sample : H3683-05MS
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-03-COMPMS

Manual Integrations
 APPROVED

sohil
 6/24/2016 1:32:47 PM

Quant Time: Jun 24 12:58:37 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 24 01:47:17 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.59	152	78958m	20.00	ng	0.00
21) Naphthalene-d8	7.88	136	318909	20.00	ng	0.00
38) Acenaphthene-d10	9.63	164	142740	20.00	ng	0.00
63) Phenanthrene-d10	11.11	188	247035	20.00	ng	0.00
75) Chrysene-d12	13.74	240	145705	20.00	ng	0.00
86) Perylene-d12	15.09	264	158856	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.19	112	435056	94.20	ng	0.00
7) Phenol-d6	6.27	99	625460	111.74	ng	-0.01
23) Nitrobenzene-d5	7.18	82	390267	71.46	ng	0.00
41) 2,4,6-Tribromophenol	10.42	330	165762	109.98	ng	-0.01
44) 2-Fluorobiphenyl	8.96	172	725688	79.51	ng	-0.01
78) Terphenyl-d14	12.69	244	573364	89.55	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.03	88	32764	14.60	ng	# 28
3) Pyridine	2.71	79	96883	18.28	ng	93
4) n-Nitrosodimethylamine	2.64	42	44140	19.67	ng	80
6) Aniline	6.26	93	196244m	24.63	ng	
8) 2-Chlorophenol	6.39	128	186082	34.04	ng	89
10) Phenol	6.28	94	244700	42.19	ng	92
11) bis(2-Chloroethyl)ether	6.34	93	177195	36.10	ng	85
12) 1,3-Dichlorobenzene	6.54	146	148130m	25.25	ng	
13) 1,4-Dichlorobenzene	6.62	146	159504m	27.00	ng	
14) 1,2-Dichlorobenzene	6.76	146	157701	29.35	ng	96
15) Benzyl Alcohol	6.76	79	151896	34.69	ng	95
16) 2,2'-oxybis(1-Chloropropan	6.89	45	177177	26.72	ng	# 1
17) 2-Methylphenol	6.89	107	156931m	35.40	ng	
18) Hexachloroethane	7.10	117	54098	25.80	ng	# 85
19) n-Nitroso-di-n-propylamine	7.03	70	135347m	37.80	ng	
20) 3+4-Methylphenols	7.05	107	221151	42.75	ng	# 78
22) Acetophenone	7.02	105	274212	38.48	ng	# 96
24) Nitrobenzene	7.19	77	184717	34.92	ng	89
25) Isophorone	7.43	82	379054	37.47	ng	98
26) 2-Nitrophenol	7.51	139	115015	40.59	ng	# 87
27) 2,4-Dimethylphenol	7.56	122	157133	30.12	ng	95
28) bis(2-Chloroethoxy)methane	7.64	93	232666	37.08	ng	# 96
29) 2,4-Dichlorophenol	7.76	162	189751	43.56	ng	99
30) 1,2,4-Trichlorobenzene	7.83	180	160647	33.87	ng	99
31) Naphthalene	7.91	128	571114	36.19	ng	98
32) Benzoic acid	7.74	122	99481	34.63	ng	92
33) 4-Chloroaniline	7.98	127	205740	29.19	ng	# 92
34) Hexachlorobutadiene	8.02	225	83063	33.20	ng	97
35) Caprolactam	8.35	113	65892m	45.66	ng	
36) 4-Chloro-3-methylphenol	8.48	107	203989	43.78	ng	94
37) 2-Methylnaphthalene	8.59	142	397065m	38.17	ng	
39) 1,2,4,5-Tetrachlorobenzene	8.76	216	172486	47.22	ng	# 1
40) Hexachlorocyclopentadiene	8.74	237	23693	10.29	ng	99
42) 2,4,6-Trichlorophenol	8.89	196	118397	41.48	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) 2,4,5-Trichlorophenol	8.95	196	119261	40.16	ng	99
45) 1,1'-Biphenyl	9.06	154	519662	49.06	ng	97
46) 2-Chloronaphthalene	9.08	162	403086	43.64	ng	97
47) 2-Nitroaniline	9.20	65	122838	45.08	ng	# 69
48) Acenaphthylene	9.50	152	629104	42.82	ng	99
49) Dimethylphthalate	9.37	163	648741	62.74	ng	99
50) 2,6-Dinitrotoluene	9.44	165	115783	48.90	ng	# 83
51) Acenaphthene	9.67	154	377804	41.22	ng	99
52) 3-Nitroaniline	9.61	138	98807	36.16	ng	89
53) 2,4-Dinitrophenol	9.72	184	53107	46.44	ng	# 89
54) Dibenzofuran	9.84	168	544398	43.13	ng	# 90
55) 4-Nitrophenol	9.83	139	156480m	73.78	ng	
56) 2,4-Dinitrotoluene	9.84	165	134332	44.14	ng	# 25
57) Fluorene	10.18	166	433403	51.51	ng	98
58) 2,3,4,6-Tetrachlorophenol	9.96	232	101555	43.51	ng	# 55
59) Diethylphthalate	10.07	149	492057	47.01	ng	99
60) 4-Chlorophenyl-phenylether	10.17	204	180593	43.02	ng	95
61) 4-Nitroaniline	10.23	138	105721	40.79	ng	99
62) Azobenzene	10.33	77	457266	44.18	ng	96
64) 4,6-Dinitro-2-methylphenol	10.25	198	51439	34.01	ng	95
65) n-Nitrosodiphenylamine	10.30	169	381506	46.54	ng	100
66) 4-Bromophenyl-phenylether	10.66	248	123832	46.72	ng	93
67) Hexachlorobenzene	10.72	284	125198	45.47	ng	97
68) Atrazine	10.83	200	125377	50.51	ng	93
69) Pentachlorophenol	10.92	266	89748	61.83	ng	97
70) Phenanthrene	11.13	178	597199	46.07	ng	99
71) Anthracene	11.19	178	641524	49.06	ng	99
72) Carbazole	11.35	167	566112	46.26	ng	100
73) Di-n-butylphthalate	11.68	149	752545	48.58	ng	99
74) Fluoranthene	12.31	202	552007	40.35	ng	98
76) Benzidine	12.45	184	238025	59.00	ng	98
77) Pyrene	12.54	202	539290	49.88	ng	99
79) Butylbenzylphthalate	13.17	149	275466	57.63	ng	96
80) Benzo(a)anthracene	13.73	228	395626	47.65	ng	99
81) 3,3'-Dichlorobenzidine	13.69	252	122446	54.84	ng	# 97
82) Chrysene	13.76	228	429061	54.93	ng	98
83) Bis(2-ethylhexyl)phthalate	13.73	149	315866	54.28	ng	# 96
84) Di-n-octyl phthalate	14.33	149	557653	58.98	ng	# 100
85) Indeno(1,2,3-cd)pyrene	16.34	276	406796	56.05	ng	# 100
87) Benzo(b)fluoranthene	14.72	252	584304m	52.77	ng	
88) Benzo(k)fluoranthene	14.75	252	262785m	31.46	ng	
89) Benzo(a)pyrene	15.04	252	420310	46.92	ng	# 95
90) Dibenzo(a,h)anthracene	16.34	278	347306	41.74	ng	99
91) Benzo(g,h,i)perylene	16.71	276	342371	40.00	ng	97

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

