

Data Path : Z:\HPCHEM1\BNA F\DATA\BF031218\
 Data File : BF103587.D
 Acq On : 12 Mar 2018 17:41
 Operator : SJ/JU
 Sample : J1867-13MS
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 J1867-09MSMS

Manual Integrations
 APPROVED

Sohil
 3/13/2018 4:15:42 PM

Quant Time: Mar 13 09:09:59 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 09 11:47:13 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	117561	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	501352	20.00	ng	-0.01
38) Acenaphthene-d10	9.90	164	228681	20.00	ng	-0.01
63) Phenanthrene-d10	11.40	188	402592	20.00	ng	0.00
75) Chrysene-d12	14.06	240	296848	20.00	ng	0.00
86) Perylene-d12	15.56	264	285512	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.49	112	586971	75.51	ng	0.00
7) Phenol-d6	6.51	99	493776	51.91	ng	0.00
23) Nitrobenzene-d5	7.43	82	896791	102.15	ng	0.00
41) 2,4,6-Tribromophenol	10.70	330	254627	131.55	ng	0.00
44) 2-Fluorobiphenyl	9.22	172	1278914	100.59	ng	0.00
78) Terphenyl-d14	12.99	244	1211890	98.14	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.63	88	72726	17.71	ng	94
3) Pyridine	3.52	79	160014m	14.60	ng	
4) n-Nitrosodimethylamine	3.46	42	98700m	22.33	ng	
6) Aniline	6.53	93	241009	17.56	ng	# 64
8) 2-Chlorophenol	6.65	128	341770	42.33	ng	92
9) Benzaldehyde	6.42	77	218780	37.03	ng	98
10) Phenol	6.53	94	186387	16.88	ng	# 44
11) bis(2-Chloroethyl)ether	6.59	93	375666	44.83	ng	93
12) 1,3-Dichlorobenzene	6.80	146	376368	40.79	ng	97
13) 1,4-Dichlorobenzene	6.87	146	402510	43.51	ng	99
14) 1,2-Dichlorobenzene	7.03	146	346472	40.61	ng	100
15) Benzyl Alcohol	7.02	79	219918	30.68	ng	93
16) 2,2'-oxybis(1-Chloropropan	7.13	45	667916	46.41	ng	64
17) 2-Methylphenol	7.13	107	245356	33.44	ng	# 72
18) Hexachloroethane	7.36	117	139744	41.49	ng	# 88
19) n-Nitroso-di-n-propylamine	7.27	70	312798	52.84	ng	92
20) 3+4-Methylphenols	7.29	107	306156	34.23	ng	# 67
22) Acetophenone	7.27	105	569758	48.69	ng	# 95
24) Nitrobenzene	7.45	77	458281	47.41	ng	97
25) Isophorone	7.68	82	872509	50.46	ng	97
26) 2-Nitrophenol	7.76	139	225952	49.66	ng	91
27) 2,4-Dimethylphenol	7.80	122	342508	49.25	ng	96
28) bis(2-Chloroethoxy)methane	7.89	93	527447	51.89	ng	99
29) 2,4-Dichlorophenol	8.02	162	324420	48.72	ng	99
30) 1,2,4-Trichlorobenzene	8.08	180	303470	42.80	ng	97
31) Naphthalene	8.16	128	1144280	46.62	ng	99
32) Benzoic acid	7.93	122	52741	15.66	ng	94
33) 4-Chloroaniline	8.23	127	171979	16.24	ng	97
34) Hexachlorobutadiene	8.27	225	142399	38.56	ng	98
35) Caprolactam	8.60	113	58013	24.79	ng	# 78
36) 4-Chloro-3-methylphenol	8.72	107	352668	46.96	ng	98
37) 2-Methylnaphthalene	8.86	142	746808	47.08	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.02	216	284106	45.44	ng	99
40) Hexachlorocyclopentadiene	9.00	237	90566	45.42	ng	95

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.15	196	192098	45.67	nq	98
43) 2,4,5-Trichlorophenol	9.20	196	197970	40.92	nq	98
45) 1,1'-Biphenyl	9.32	154	844865	44.09	nq	99
46) 2-Chloronaphthalene	9.35	162	683871	45.11	nq	98
47) 2-Nitroaniline	9.46	65	288130	51.31	nq	89
48) Acenaphthylene	9.76	152	998198	42.79	nq	98
49) Dimethylphthalate	9.62	163	833197	45.42	nq	100
50) 2,6-Dinitrotoluene	9.70	165	179469	46.62	nq	# 74
51) Acenaphthene	9.93	154	614986	45.22	nq	100
52) 3-Nitroaniline	9.88	138	112085	22.95	nq	97
53) 2,4-Dinitrophenol	10.00	184	115193	82.50	nq	# 17
54) Dibenzofuran	10.11	168	907157	48.59	nq	95
55) 4-Nitrophenol	10.07	139	47481m	15.68	nq	
56) 2,4-Dinitrotoluene	10.12	165	233979	48.06	nq	# 81
57) Fluorene	10.46	166	754459	47.43	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.24	232	149041	45.86	nq	97
59) Diethylphthalate	10.32	149	822780	46.89	nq	99
60) 4-Chlorophenyl-phenylether	10.44	204	315703	46.81	nq	98
61) 4-Nitroaniline	10.50	138	212296	43.52	nq	97
62) Azobenzene	10.60	77	816460	45.72	nq	97
64) 4,6-Dinitro-2-methylphenol	10.53	198	93376	39.14	nq	95
65) n-Nitrosodiphenylamine	10.56	169	624918	44.58	nq	99
66) 4-Bromophenyl-phenylether	10.93	248	188941	45.05	nq	91
67) Hexachlorobenzene	11.01	284	203206	44.64	nq	# 90
68) Atrazine	11.09	200	196649	46.67	nq	97
69) Pentachlorophenol	11.21	266	135722	61.67	nq	97
70) Phenanthrene	11.43	178	1013937	45.76	nq	99
71) Anthracene	11.48	178	1068894	47.05	nq	100
72) Carbazole	11.64	167	1068123	47.21	nq	98
73) Di-n-butylphthalate	11.94	149	1289291	45.54	nq	100
74) Fluoranthene	12.62	202	1064600	47.82	nq	99
76) Benzidine	12.74	184	255027	22.98	nq	99
77) Pyrene	12.85	202	1078048	46.58	nq	99
79) Butylbenzylphthalate	13.45	149	585940	47.92	nq	100
80) Benzo(a)anthracene	14.05	228	927456	48.15	nq	99
81) 3,3'-Dichlorobenzidine	14.00	252	178572	25.98	nq	96
82) Chrysene	14.09	228	915831	48.97	nq	99
83) Bis(2-ethylhexyl)phthalate	14.00	149	733634	44.46	nq	# 98
84) Di-n-octyl phthalate	14.62	149	1328104	49.81	nq	100
85) Indeno(1,2,3-cd)pyrene	17.12	276	783262	52.90	nq	97
87) Benzo(b)fluoranthene	15.12	252	908217	48.38	nq	# 96
88) Benzo(k)fluoranthene	15.15	252	770106	43.57	nq	98
89) Benzo(a)pyrene	15.50	252	795034	47.43	nq	# 96
90) Dibenzo(a,h)anthracene	17.12	278	652225	50.35	nq	98
91) Benzo(g,h,i)perylene	17.59	276	685855	54.18	nq	98

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

