

Data Path : Z:\HPCHEM1\BNA F\DATA\BF010318\
 Data File : BF101873.D
 Acq On : 3 Jan 2018 21:12
 Operator : SJ/JU
 Sample : J1004-02MSD
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-08-COMPMSD

Manual Integrations
 APPROVED

Sohil
 1/4/2018 1:56:41 PM

Quant Time: Jan 04 10:23:03 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 29 13:21:45 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.77	152	98457	20.00	ng	-0.01
21) Naphthalene-d8	8.05	136	336873	20.00	ng	-0.02
38) Acenaphthene-d10	9.80	164	126139	20.00	ng	-0.02
63) Phenanthrene-d10	11.30	188	249006	20.00	ng	-0.02
75) Chrysene-d12	13.95	240	240669	20.00	ng	-0.02
86) Perylene-d12	15.42	264	190101	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.40	112	577976	87.44	ng	0.00
7) Phenol-d6	6.42	99	611091	74.29	ng	-0.01
23) Nitrobenzene-d5	7.35	82	344534	56.93	ng	-0.02
41) 2,4,6-Tribromophenol	10.60	330	102464	84.30	ng	-0.02
44) 2-Fluorobiphenyl	9.12	172	530059	65.19	ng	-0.02
78) Terphenyl-d14	12.87	244	574681	57.57	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.50	88	91436	26.922	ng	94
3) Pyridine	3.30	79	167070	17.705	ng	97
4) n-Nitrosodimethylamine	3.23	42	84791	19.516	ng	95
6) Aniline	6.45	93	174888	15.299	ng	# 65
8) 2-Chlorophenol	6.56	128	182114	26.192	ng	97
9) Benzaldehyde	6.35	77	51147	8.419	ng	98
10) Phenol	6.44	94	234817	25.045	ng	# 66
11) bis(2-Chloroethyl)ether	6.51	93	191015	25.973	ng	99
12) 1,3-Dichlorobenzene	6.71	146	210893	26.875	ng	98
13) 1,4-Dichlorobenzene	6.79	146	215612	26.906	ng	100
14) 1,2-Dichlorobenzene	6.94	146	194877	27.017	ng	97
15) Benzyl Alcohol	6.93	79	141125	24.925	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.04	45	312937	27.803	ng	71
17) 2-Methylphenol	7.05	107	145954	22.820	ng	# 85
18) Hexachloroethane	7.27	117	53286	19.126	ng	97
19) n-Nitroso-di-n-propylamine	7.18	70	131507	24.343	ng	95
20) 3+4-Methylphenols	7.21	107	191247	28.218	ng	97
22) Acetophenone	7.19	105	261190	33.980	ng	# 98
24) Nitrobenzene	7.36	77	181903	27.159	ng	99
25) Isophorone	7.60	82	325620	26.822	ng	99
26) 2-Nitrophenol	7.69	139	64970	22.579	ng	95
27) 2,4-Dimethylphenol	7.72	122	147767	30.228	ng	97
28) bis(2-Chloroethoxy)methane	7.80	93	215155	27.900	ng	# 97
29) 2,4-Dichlorophenol	7.95	162	128474	26.602	ng	98
30) 1,2,4-Trichlorobenzene	8.00	180	139696	27.410	ng	97
31) Naphthalene	8.07	128	456778	26.934	ng	100
32) Benzoic acid	7.86	122	47077	19.883	ng	98
33) 4-Chloroaniline	8.15	127	124450	16.883	ng	98
34) Hexachlorobutadiene	8.17	225	81639	27.696	ng	99
35) Caprolactam	8.50	113	36946	22.031	ng	94
36) 4-Chloro-3-methylphenol	8.63	107	127153	23.954	ng	98
37) 2-Methylnaphthalene	8.76	142	283524	25.660	ng	99
39) 1,2,4,5-Tetrachlorobenzene	8.93	216	130085	30.545	ng	97
42) 2,4,6-Trichlorophenol	9.06	196	77394	30.186	ng	97

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43) 2,4,5-Trichlorophenol	9.13	196	73553	26.201	nq	96
45) 1,1'-Biphenyl	9.23	154	330938	29.583	nq	99
46) 2-Chloronaphthalene	9.26	162	242519	28.333	nq	97
47) 2-Nitroaniline	9.37	65	88656	29.983	nq	96
48) Acenaphthylene	9.67	152	361045	26.705	nq	99
49) Dimethylphthalate	9.52	163	320930	31.631	nq	100
50) 2,6-Dinitrotoluene	9.60	165	47840	23.199	nq	91
51) Acenaphthene	9.84	154	211571	26.047	nq	98
52) 3-Nitroaniline	9.79	138	75366	29.314	nq	95
54) Dibenzofuran	10.02	168	319181	30.921	nq	97
55) 4-Nitrophenol	10.00	139	31591m	28.180	nq	
56) 2,4-Dinitrotoluene	10.03	165	58632	25.236	nq	# 97
57) Fluorene	10.36	166	251289	28.777	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.15	232	62375m	30.926	nq	
59) Diethylphthalate	10.22	149	265173	26.392	nq	97
60) 4-Chlorophenyl-phenylether	10.34	204	120139	28.707	nq	96
61) 4-Nitroaniline	10.40	138	76018	29.416	nq	95
62) Azobenzene	10.50	77	247658	23.794	nq	98
65) n-Nitrosodiphenylamine	10.46	169	226571	24.642	nq	95
66) 4-Bromophenyl-phenylether	10.83	248	71383	25.513	nq	# 89
67) Hexachlorobenzene	10.92	284	76668	25.891	nq	# 84
68) Atrazine	10.99	200	64316	23.460	nq	97
69) Pentachlorophenol	11.12	266	47964	44.795	nq	98
70) Phenanthrene	11.32	178	383394	27.687	nq	98
71) Anthracene	11.38	178	394146	27.865	nq	99
72) Carbazole	11.54	167	378600	28.588	nq	98
73) Di-n-butylphthalate	11.84	149	430246	26.767	nq	99
74) Fluoranthene	12.52	202	459534	31.616	nq	96
76) Benzidine	12.64	184	354155	34.842	nq	97
77) Pyrene	12.75	202	477278	26.424	nq	98
79) Butylbenzylphthalate	13.34	149	209556	25.641	nq	96
80) Benzo(a)anthracene	13.94	228	412453	25.954	nq	99
81) 3,3'-Dichlorobenzidine	13.90	252	150741	29.091	nq	95
82) Chrysene	13.98	228	394382	26.284	nq	98
83) Bis(2-ethylhexyl)phthalate	13.89	149	274466	26.514	nq	99
84) Di-n-octyl phthalate	14.50	149	502004	27.192	nq	98
85) Indeno(1,2,3-cd)pyrene	16.90	276	271030	20.284	nq	98
87) Benzo(b)fluoranthene	14.99	252	317289	27.383	nq	98
88) Benzo(k)fluoranthene	15.02	252	318980	28.314	nq	97
89) Benzo(a)pyrene	15.36	252	278867	26.107	nq	99
90) Dibenzo(a,h)anthracene	16.89	278	228653	24.932	nq	99
91) Benzo(g,h,i)perylene	17.35	276	221065	24.343	nq	94

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

