

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080218\
 Data File : BF107923.D
 Acq On : 2 Aug 2018 21:33
 Operator : JU/SJ
 Sample : J4285-05MSD
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 KG-PIT-2MSD

Quant Time: Aug 03 00:57:20 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 30 17:48:53 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.96	152	133135	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	513931	20.00	ng	-0.01
38) Acenaphthene-d10	10.00	164	188846	20.00	ng	-0.01
63) Phenanthrene-d10	11.49	188	334296	20.00	ng	-0.01
75) Chrysene-d12	14.14	240	333614	20.00	ng	-0.01
86) Perylene-d12	15.68	264	287092	20.00	ng	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.60	112	724699	92.51	ng	0.00
7) Phenol-d6	6.61	99	926646	89.75	ng	0.00
23) Nitrobenzene-d5	7.52	82	593514	66.53	ng	-0.01
41) 2,4,6-Tribromophenol	10.80	330	215951	84.16	ng	0.00
44) 2-Fluorobiphenyl	9.31	172	1094891	80.79	ng	-0.01
78) Terphenyl-d14	13.07	244	1005086	65.33	ng	-0.01

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.85	88	90048	27.337	ng	# 81
3) Pyridine	3.66	79	203934	25.542	ng	87
4) n-Nitrosodimethylamine	3.64	42	33529	8.412	ng	# 96
6) Aniline	6.64	93	264679	24.648	ng	# 54
8) 2-Chlorophenol	6.75	128	312941	31.803	ng	92
9) Benzaldehyde	6.52	77	70778	11.363	ng	89
10) Phenol	6.62	94	410177	33.316	ng	88
11) bis(2-Chloroethyl)ether	6.69	93	339922	31.049	ng	97
12) 1,3-Dichlorobenzene	6.90	146	316445	30.662	ng	97
13) 1,4-Dichlorobenzene	6.97	146	369698	32.180	ng	98
14) 1,2-Dichlorobenzene	7.12	146	322393	30.231	ng	98
15) Benzyl Alcohol	7.11	79	188027	29.301	ng	93
16) 2,2'-oxybis(1-Chloropropan	7.22	45	498690	31.683	ng	65
17) 2-Methylphenol	7.22	107	232251	31.973	ng	# 72
18) Hexachloroethane	7.45	117	118780	28.764	ng	94
19) n-Nitroso-di-n-propylamine	7.36	70	209733	32.024	ng	100
20) 3+4-Methylphenols	7.38	107	271815	30.479	ng	# 65
22) Acetophenone	7.37	105	446078	36.077	ng	# 95
24) Nitrobenzene	7.54	77	276641	29.574	ng	97
25) Isophorone	7.77	82	571977	39.060	ng	99
26) 2-Nitrophenol	7.86	139	105836	22.545	ng	98
27) 2,4-Dimethylphenol	7.90	122	252692	40.204	ng	98
28) bis(2-Chloroethoxy)methane	7.98	93	356632	36.956	ng	99
29) 2,4-Dichlorophenol	8.11	162	238300	33.519	ng	98
30) 1,2,4-Trichlorobenzene	8.18	180	271229	33.442	ng	98
31) Naphthalene	8.26	128	887925	37.168	ng	100
32) Benzoic acid	7.90	122	252692	59.047	ng	# 14
33) 4-Chloroaniline	8.32	127	220280	21.587	ng	97
34) Hexachlorobutadiene	8.36	225	163820	32.686	ng	98
35) Caprolactam	8.67	113	61350	28.665	ng	# 74
36) 4-Chloro-3-methylphenol	8.81	107	229934	29.780	ng	97
37) 2-Methylnaphthalene	8.95	142	534028	35.587	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.11	216	258273	40.113	ng	# 96
40) Hexachlorocyclopentadiene	9.09	237	38839	20.129	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.24	196	131731	37.008	nq	99
43) 2,4,5-Trichlorophenol	9.29	196	151387	35.328	nq	95
45) 1,1'-Biphenyl	9.41	154	671821	39.091	nq	97
46) 2-Chloronaphthalene	9.44	162	482802	37.253	nq	97
47) 2-Nitroaniline	9.55	65	143281	33.989	nq	88
48) Acenaphthylene	9.86	152	752665	37.461	nq	99
49) Dimethylphthalate	9.71	163	683224	48.174	nq	100
50) 2,6-Dinitrotoluene	9.78	165	108055	35.059	nq	89
51) Acenaphthene	10.02	154	433090	37.107	nq	100
52) 3-Nitroaniline	9.98	138	100445	25.863	nq	95
53) 2,4-Dinitrophenol	10.15	184	6260	11.887	nq #	23
54) Dibenzofuran	10.20	168	607629	33.683	nq	97
55) 4-Nitrophenol	10.20	139	362690	184.636	nq #	19
56) 2,4-Dinitrotoluene	10.20	165	124153	30.648	nq #	82
57) Fluorene	10.55	166	478648	34.141	nq	95
58) 2,3,4,6-Tetrachlorophenol	10.32	232	126065	32.029	nq #	77
59) Diethylphthalate	10.40	149	540437	35.442	nq	100
60) 4-Chlorophenyl-phenylether	10.53	204	239941	31.417	nq	93
61) 4-Nitroaniline	10.62	138	90038	26.123	nq	89
62) Azobenzene	10.70	77	452154	31.838	nq	94
64) 4,6-Dinitro-2-methylphenol	10.61	198	11221	6.170	nq #	44
65) n-Nitrosodiphenylamine	10.65	169	411052	37.343	nq	99
66) 4-Bromophenyl-phenylether	11.02	248	140319	36.013	nq	89
67) Hexachlorobenzene	11.09	284	143667	33.251	nq #	90
68) Atrazine	11.18	200	116091	35.634	nq	95
69) Pentachlorophenol	11.30	266	119669	56.262	nq	99
70) Phenanthrene	11.51	178	576689	36.331	nq	100
71) Anthracene	11.57	178	713866	38.990	nq	99
72) Carbazole	11.74	167	644463	35.619	nq	97
73) Di-n-butylphthalate	12.03	149	738099	37.393	nq	99
74) Fluoranthene	12.71	202	745276	39.965	nq	97
76) Benzidine	12.85	184	519724	48.182	nq	98
77) Pyrene	12.94	202	751557	31.285	nq	97
79) Butylbenzylphthalate	13.54	149	335341	31.877	nq	95
80) Benzo(a)anthracene	14.13	228	678791	35.919	nq	99
81) 3,3'-Dichlorobenzidine	14.10	252	257942	35.211	nq #	97
82) Chrysene	14.17	228	764055	37.484	nq	98
83) Bis(2-ethylhexyl)phthalate	14.09	149	463918	35.079	nq	98
84) Di-n-octyl phthalate	14.71	149	860526	40.538	nq	97
85) Indeno(1,2,3-cd)pyrene	17.25	276	498837	32.174	nq #	92
87) Benzo(b)fluoranthene	15.21	252	530933	37.497	nq	98
88) Benzo(k)fluoranthene	15.25	252	681447	39.199	nq	98
89) Benzo(a)pyrene	15.61	252	542637	37.123	nq	99
90) Dibenzo(a,h)anthracene	17.26	278	427556	33.532	nq	96
91) Benzo(g,h,i)perylene	17.73	276	400381	32.453	nq #	91

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

