

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF020819\
 Data File : BF112474.D
 Acq On : 8 Feb 2019 21:52
 Operator : JU/SJ
 Sample : K1385-02MS
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 98-SB-01-2-4MS

Quant Time: Feb 09 05:54:05 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 28 10:38:18 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	84932	20.00	ng	-0.01
21) Naphthalene-d8	8.12	136	329499	20.00	ng	-0.01
39) Acenaphthene-d10	9.87	164	163542	20.00	ng	-0.02
64) Phenanthrene-d10	11.36	188	303416	20.00	ng	-0.01
76) Chrysene-d12	14.00	240	205115	20.00	ng	0.00
87) Perylene-d12	15.47	264	200760	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.46	112	465440	116.40	ng	-0.02
7) Phenol-d6	6.48	99	593964	106.90	ng	-0.01
23) Nitrobenzene-d5	7.40	82	368864	64.50	ng	-0.01
42) 2,4,6-Tribromophenol	10.66	330	161844	85.02	ng	-0.02
45) 2-Fluorobiphenyl	9.19	172	784123	67.81	ng	-0.02
79) Terphenyl-d14	12.94	244	765445	70.29	ng	-0.01

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.55	88	54842	34.430	ng	96
3) Pyridine	3.32	79	128835	31.797	ng	96
4) n-Nitrosodimethylamine	3.25	42	67548	39.680	ng	# 87
6) Aniline	6.50	93	68084	10.301	ng	# 1
8) 2-Chlorophenol	6.63	128	186399	34.652	ng	99
9) Benzaldehyde	6.38	77	84451	29.080	ng	99
10) Phenol	6.50	94	220560	36.182	ng	84
11) bis(2-Chloroethyl)ether	6.56	93	159842	33.306	ng	97
12) 1,3-Dichlorobenzene	6.77	146	201511	32.167	ng	98
13) 1,4-Dichlorobenzene	6.85	146	209674	32.522	ng	99
14) 1,2-Dichlorobenzene	7.00	146	197659	32.745	ng	98
15) Benzyl Alcohol	6.98	79	153766	32.830	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.10	45	198260	32.175	ng	# 17
17) 2-Methylphenol	7.10	107	144822	33.963	ng	# 88
18) Hexachloroethane	7.34	117	66815	29.512	ng	99
19) n-Nitroso-di-n-propylamine	7.24	70	123522	32.241	ng	99
20) 3+4-Methylphenols	7.26	107	193221	35.435	ng	# 61
22) Acetophenone	7.24	105	259323	32.321	ng	# 92
24) Nitrobenzene	7.42	77	184123	32.239	ng	97
25) Isophorone	7.65	82	335036	37.346	ng	99
26) 2-Nitrophenol	7.73	139	83656	27.409	ng	95
27) 2,4-Dimethylphenol	7.77	122	160085	38.069	ng	93
28) bis(2-Chloroethoxy)methane	7.86	93	209760	34.339	ng	99
29) 2,4-Dichlorophenol	7.99	162	160835	34.178	ng	96
30) 1,2,4-Trichlorobenzene	8.06	180	178513	33.002	ng	99
31) Naphthalene	8.13	128	562142	36.482	ng	98
33) 4-Chloroaniline	8.19	127	50910	7.588	ng	96
34) Hexachlorobutadiene	8.25	225	98987	31.185	ng	100
35) Caprolactam	8.54	113	46602	28.072	ng	95
36) 4-Chloro-3-methylphenol	8.69	107	161732	32.466	ng	99
37) 2-Methylnaphthalene	8.83	142	396125	37.553	ng	98
38) 1-Methylnaphthalene	8.93	142	370415	36.636	ng	97
40) 1,2,4,5-Tetrachlorobenzene	9.00	216	176291	32.831	ng	99
41) Hexachlorocyclopentadiene	8.98	237	17088	16.903	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) 2,4,6-Trichlorophenol	9.12	196	102666	31.018	nq	98
44) 2,4,5-Trichlorophenol	9.17	196	103365	29.880	nq	98
46) 1,1'-Biphenyl	9.29	154	474589	33.194	nq	98
47) 2-Chloronaphthalene	9.32	162	364456	32.872	nq	96
48) 2-Nitroaniline	9.42	65	102118	33.792	nq	90
49) Acenaphthylene	9.73	152	586440	36.087	nq	98
50) Dimethylphthalate	9.59	163	482820	37.998	nq	99
51) 2,6-Dinitrotoluene	9.66	165	92783	32.604	nq	94
52) Acenaphthene	9.90	154	336740	34.688	nq	98
53) 3-Nitroaniline	9.83	138	58910	19.108	nq	82
54) 2,4-Dinitrophenol	9.99	184	252	0.276	nq #	44
55) Dibenzofuran	10.07	168	510290	34.398	nq	95
56) 4-Nitrophenol	10.02	139	62141	38.124	nq	92
57) 2,4-Dinitrotoluene	10.07	165	122244	33.743	nq #	85
58) Fluorene	10.42	166	414698	37.356	nq	98
59) 2,3,4,6-Tetrachlorophenol	10.20	232	72965	27.911	nq	94
60) Diethylphthalate	10.29	149	442557	35.097	nq	98
61) 4-Chlorophenyl-phenylether	10.40	204	196527	32.544	nq #	89
62) 4-Nitroaniline	10.45	138	94549	31.266	nq	95
63) Azobenzene	10.57	77	369764	33.342	nq	97
65) 4,6-Dinitro-2-methylphenol	10.49	198	5207	3.061	nq	86
66) n-Nitrosodiphenylamine	10.53	169	362737	35.472	nq	99
67) 4-Bromophenyl-phenylether	10.90	248	118688	33.259	nq	95
68) Hexachlorobenzene	10.97	284	126528	33.047	nq	98
69) Atrazine	11.05	200	117050	35.474	nq	97
70) Pentachlorophenol	11.18	266	47672	32.000	nq	99
71) Phenanthrene	11.38	178	629290	39.894	nq	98
72) Anthracene	11.43	178	605364	38.526	nq	97
73) Carbazole	11.59	167	522028	33.680	nq	99
74) Di-n-butylphthalate	11.91	149	698783	36.322	nq	98
75) Fluoranthene	12.57	202	676850	40.006	nq	98
77) Benzidine	12.69	184	86235	15.275	nq	96
78) Pyrene	12.80	202	637742	44.908	nq	99
80) Butylbenzylphthalate	13.40	149	257337	37.454	nq	91
81) Benzo(a)anthracene	13.99	228	491616	40.772	nq	97
82) 3,3'-Dichlorobenzidine	13.95	252	79776	17.679	nq	98
83) Chrysene	14.03	228	456986	39.704	nq	98
84) Bis(2-ethylhexyl)phthalate	13.96	149	372238	38.167	nq	96
85) Di-n-octyl phthalate	14.57	149	565200	37.667	nq	97
86) Indeno(1,2,3-cd)pyrene	16.96	276	404709	44.376	nq	97
88) Benzo(b)fluoranthene	15.04	252	511307	42.586	nq	100
89) Benzo(k)fluoranthene	15.07	252	400001	34.782	nq	98
90) Benzo(a)pyrene	15.42	252	435405	40.303	nq	100
91) Dibenzo(a,h)anthracene	16.96	278	325013	37.329	nq	99
92) Benzo(g,h,i)perylene	17.41	276	318244	38.742	nq	98

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

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