

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030518\
 Data File : BF103433.D
 Acq On : 6 Mar 2018 2:13
 Operator : SJ/JU
 Sample : J1722-01MS
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-24MS

Manual Integrations
 APPROVED

Sohil
 3/6/2018 4:23:57 PM

Quant Time: Mar 06 06:04:10 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 05 13:40:47 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	118621	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	486131	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	183906	20.00	ng	0.00
63) Phenanthrene-d10	11.41	188	283644	20.00	ng	0.00
75) Chrysene-d12	14.07	240	152613	20.00	ng	0.00
86) Perylene-d12	15.59	264	122861	20.00	ng	0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.50	112	1109593	141.47	ng	0.00
7) Phenol-d6	6.53	99	1262052	131.50	ng	0.00
23) Nitrobenzene-d5	7.45	82	793876	93.26	ng	0.00
41) 2,4,6-Tribromophenol	10.72	330	166256	106.80	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	1039853	102.14	ng	0.00
78) Terphenyl-d14	13.00	244	727688	114.62	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.64	88	132015	31.869	ng	91
3) Pyridine	3.48	79	349186	31.575	ng	99
4) n-Nitrosodimethylamine	3.39	42	203587	45.646	ng	# 94
6) Aniline	6.55	93	274238	19.799	ng	# 15
8) 2-Chlorophenol	6.67	128	312550	38.361	ng	93
9) Benzaldehyde	6.43	77	147633	22.278	ng	96
10) Phenol	6.54	94	451081	40.487	ng	81
11) bis(2-Chloroethyl)ether	6.61	93	354214	41.889	ng	95
12) 1,3-Dichlorobenzene	6.82	146	328722	35.305	ng	98
13) 1,4-Dichlorobenzene	6.89	146	337202	36.122	ng	98
14) 1,2-Dichlorobenzene	7.05	146	297297	34.539	ng	98
15) Benzyl Alcohol	7.02	79	246872	34.136	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.15	45	501496	34.537	ng	50
17) 2-Methylphenol	7.14	107	267227	36.091	ng	# 84
18) Hexachloroethane	7.37	117	63884	18.799	ng	# 88
19) n-Nitroso-di-n-propylamine	7.29	70	232422	38.912	ng	95
20) 3+4-Methylphenols	7.30	107	358797	39.752	ng	89
22) Acetophenone	7.29	105	451641	39.802	ng	# 93
24) Nitrobenzene	7.46	77	351618	37.517	ng	97
25) Isophorone	7.69	82	655525	39.100	ng	95
26) 2-Nitrophenol	7.77	139	106544	24.148	ng	91
27) 2,4-Dimethylphenol	7.82	122	288863	42.833	ng	99
28) bis(2-Chloroethoxy)methane	7.90	93	404068	40.993	ng	98
29) 2,4-Dichlorophenol	8.03	162	258417	40.023	ng	100
30) 1,2,4-Trichlorobenzene	8.10	180	237361	34.521	ng	99
31) Naphthalene	8.18	128	925354	38.880	ng	99
32) Benzoic acid	7.96	122	7386m	8.177	ng	
33) 4-Chloroaniline	8.25	127	244523	23.809	ng	96
34) Hexachlorobutadiene	8.29	225	116808	32.623	ng	99
35) Caprolactam	8.60	113	82340	36.285	ng	# 82
36) 4-Chloro-3-methylphenol	8.73	107	274387	37.676	ng	99
37) 2-Methylnaphthalene	8.87	142	569209	37.006	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.04	216	197590	39.301	ng	99
42) 2,4,6-Trichlorophenol	9.16	196	131899	38.989	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) 2,4,5-Trichlorophenol	9.21	196	136156	34.993	nq	93
45) 1,1'-Biphenyl	9.33	154	603615	39.169	nq	99
46) 2-Chloronaphthalene	9.36	162	467514	38.347	nq	99
47) 2-Nitroaniline	9.47	65	208985	46.276	nq	88
48) Acenaphthylene	9.78	152	701526	37.398	nq	100
49) Dimethylphthalate	9.63	163	605166	41.019	nq	100
50) 2,6-Dinitrotoluene	9.70	165	86605	27.973	nq #	82
51) Acenaphthene	9.95	154	394090	36.029	nq	99
52) 3-Nitroaniline	9.88	138	153428	39.060	nq	99
54) Dibenzofuran	10.12	168	576386	38.387	nq	96
55) 4-Nitrophenol	10.06	139	109059	44.798	nq	93
56) 2,4-Dinitrotoluene	10.12	165	89890	22.957	nq #	72
57) Fluorene	10.47	166	455037	35.575	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.25	232	77961	29.827	nq #	95
59) Diethylphthalate	10.33	149	504979	35.788	nq	99
60) 4-Chlorophenyl-phenylether	10.45	204	196961	36.312	nq	99
61) 4-Nitroaniline	10.50	138	156495	39.888	nq	94
62) Azobenzene	10.62	77	476886	33.203	nq	98
64) 4,6-Dinitro-2-methylphenol	10.72	198	375	5.140	nq #	1
65) n-Nitrosodiphenylamine	10.57	169	385174	39.001	nq	100
66) 4-Bromophenyl-phenylether	10.95	248	108589	36.751	nq	94
67) Hexachlorobenzene	11.02	284	110233	34.372	nq	98
68) Atrazine	11.10	200	87217	29.381	nq	96
69) Pentachlorophenol	11.22	266	44932	28.976	nq	97
70) Phenanthrene	11.44	178	847658	54.303	nq	99
71) Anthracene	11.49	178	665709	41.589	nq	100
72) Carbazole	11.65	167	584127	36.645	nq	99
73) Di-n-butylphthalate	11.96	149	736387	36.919	nq	99
74) Fluoranthene	12.64	202	963290	61.410	nq	100
76) Benzidine	12.75	184	111969	19.625	nq	98
77) Pyrene	12.87	202	936800	78.732	nq	99
79) Butylbenzylphthalate	13.46	149	284083	45.189	nq	96
80) Benzo(a)anthracene	14.06	228	555760	56.125	nq	99
81) 3,3'-Dichlorobenzidine	14.02	252	107461	30.409	nq	98
82) Chrysene	14.10	228	497674	51.762	nq	99
83) Bis(2-ethylhexyl)phthalate	14.02	149	363843	42.889	nq	99
84) Di-n-octyl phthalate	14.64	149	532513	38.851	nq	100
85) Indeno(1,2,3-cd)pyrene	17.15	276	347844	45.698	nq	98
87) Benzo(b)fluoranthene	15.14	252	455149	56.344	nq #	94
88) Benzo(k)fluoranthene	15.17	252	351239	46.175	nq	99
89) Benzo(a)pyrene	15.53	252	383672	53.187	nq #	94
90) Dibenzo(a,h)anthracene	17.15	278	243556	43.694	nq	98
91) Benzo(g,h,i)perylene	17.62	276	294938	54.139	nq	95

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

