

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072518\
 Data File : BF107653.D
 Acq On : 25 Jul 2018 13:32
 Operator : JU/SJ
 Sample : J4101-02MSD
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-1MSD

Manual Integrations
 APPROVED

Sohil
 7/26/2018 2:02:44 PM

Quant Time: Jul 25 14:24:08 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF072018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 20 16:29:27 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	237065	20.00	ng	0.00
21) Naphthalene-d8	8.25	136	939893	20.00	ng	-0.01
38) Acenaphthene-d10	10.01	164	421226	20.00	ng	0.00
63) Phenanthrene-d10	11.51	188	845975	20.00	ng	0.00
75) Chrysene-d12	14.16	240	595429	20.00	ng	0.00
86) Perylene-d12	15.69	264	520532	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.62	112	1591467	107.94	ng	0.02
7) Phenol-d6	6.61	99	1774999	100.26	ng	0.00
23) Nitrobenzene-d5	7.54	82	1385724	99.50	ng	-0.01
41) 2,4,6-Tribromophenol	10.81	330	697232	145.56	ng	0.00
44) 2-Fluorobiphenyl	9.32	172	2565197	114.52	ng	-0.01
78) Terphenyl-d14	13.09	244	2501067	107.54	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.08	88	188084	27.406	ng	# 69
3) Pyridine	3.77	79	374028	20.033	ng	# 84
4) n-Nitrosodimethylamine	3.72	42	252476	32.064	ng	94
6) Aniline	6.65	93	255933	11.208	ng	# 72
8) 2-Chlorophenol	6.77	128	659223	41.731	ng	99
9) Benzaldehyde	6.53	77	148115	10.677	ng	96
10) Phenol	6.62	94	643092	30.886	ng	91
11) bis(2-Chloroethyl)ether	6.71	93	638088	36.964	ng	92
12) 1,3-Dichlorobenzene	6.91	146	730521	40.788	ng	97
13) 1,4-Dichlorobenzene	6.98	146	767810	41.827	ng	98
14) 1,2-Dichlorobenzene	7.14	146	667925	40.545	ng	98
15) Benzyl Alcohol	7.11	79	528509	43.801	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.24	45	1086173	37.297	ng	77
17) 2-Methylphenol	7.22	107	555561	41.379	ng	# 89
18) Hexachloroethane	7.48	117	271255	42.130	ng	94
19) n-Nitroso-di-n-propylamine	7.38	70	465141	40.669	ng	98
20) 3+4-Methylphenols	7.38	107	689232	43.231	ng	# 57
22) Acetophenone	7.38	105	981282	44.455	ng	# 89
24) Nitrobenzene	7.55	77	702343	44.030	ng	99
25) Isophorone	7.79	82	1312099	44.125	ng	98
26) 2-Nitrophenol	7.87	139	393753	58.454	ng	96
27) 2,4-Dimethylphenol	7.90	122	627447	49.209	ng	99
28) bis(2-Chloroethoxy)methane	8.00	93	849568	42.751	ng	100
29) 2,4-Dichlorophenol	8.11	162	604529	46.386	ng	97
30) 1,2,4-Trichlorobenzene	8.19	180	636082	43.455	ng	99
31) Naphthalene	8.27	128	1986180	48.055	ng	98
32) Benzoic acid	8.04	122	345408	40.316	ng	97
33) 4-Chloroaniline	8.34	127	138775	7.682	ng	98
34) Hexachlorobutadiene	8.38	225	378282	43.491	ng	98
35) Caprolactam	8.71	113	163038m	38.461	ng	
36) 4-Chloro-3-methylphenol	8.81	107	656607	45.355	ng	96
37) 2-Methylnaphthalene	8.97	142	1400026	48.882	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.13	216	661559	47.087	ng	99
40) Hexachlorocyclopentadiene	9.11	237	732800	102.604	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.25	196	419258	46.621	nq	99
43) 2,4,5-Trichlorophenol	9.30	196	460450	48.990	nq	95
45) 1,1'-Biphenyl	9.43	154	1695444	46.205	nq	97
46) 2-Chloronaphthalene	9.46	162	1283862	45.775	nq	99
47) 2-Nitroaniline	9.56	65	438440	53.620	nq	92
48) Acenaphthylene	9.88	152	2028480	48.145	nq	97
49) Dimethylphthalate	9.73	163	1466582	46.197	nq	99
50) 2,6-Dinitrotoluene	9.80	165	330678	52.464	nq	93
51) Acenaphthene	10.05	154	1143558	43.319	nq	99
52) 3-Nitroaniline	9.98	138	141682	18.477	nq	95
53) 2,4-Dinitrophenol	10.08	184	383328	109.214	nq #	23
54) Dibenzofuran	10.22	168	1749594	45.018	nq	100
55) 4-Nitrophenol	10.15	139	429708	74.850	nq	96
56) 2,4-Dinitrotoluene	10.21	165	406359	53.410	nq #	96
57) Fluorene	10.57	166	1392716	50.231	nq	95
58) 2,3,4,6-Tetrachlorophenol	10.34	232	396569	50.697	nq #	87
59) Diethylphthalate	10.43	149	1480285	44.647	nq	96
60) 4-Chlorophenyl-phenylether	10.55	204	690259	44.037	nq	96
61) 4-Nitroaniline	10.60	138	320539	49.755	nq	97
62) Azobenzene	10.71	77	1361636	43.797	nq	92
64) 4,6-Dinitro-2-methylphenol	10.62	198	231090	55.032	nq	86
65) n-Nitrosodiphenylamine	10.68	169	1229653	42.798	nq	99
66) 4-Bromophenyl-phenylether	11.04	248	437879	43.970	nq #	90
67) Hexachlorobenzene	11.11	284	476952	43.528	nq #	84
68) Atrazine	11.20	200	378996	43.204	nq	96
69) Pentachlorophenol	11.31	266	506157	73.954	nq	98
70) Phenanthrene	11.53	178	1878040	45.556	nq	98
71) Anthracene	11.58	178	2012783	48.493	nq	98
72) Carbazole	11.74	167	1894565	43.894	nq	99
73) Di-n-butylphthalate	12.05	149	2133214	49.471	nq	98
74) Fluoranthene	12.72	202	2067936	46.747	nq	96
76) Benzidine	12.85	184	157923	9.210	nq	97
77) Pyrene	12.95	202	2053708	50.425	nq	98
79) Butylbenzylphthalate	13.56	149	924291	52.771	nq #	90
80) Benzo(a)anthracene	14.15	228	1657665	47.507	nq	98
81) 3,3'-Dichlorobenzidine	14.11	252	251742	20.707	nq #	95
82) Chrysene	14.18	228	1610441	48.046	nq	98
83) Bis(2-ethylhexyl)phthalate	14.11	149	1162218	47.027	nq	99
84) Di-n-octyl phthalate	14.73	149	1793973	46.700	nq	98
85) Indeno(1,2,3-cd)pyrene	17.27	276	1454198	51.190	nq	94
87) Benzo(b)fluoranthene	15.23	252	1390360	48.805	nq	98
88) Benzo(k)fluoranthene	15.27	252	1297832	46.500	nq	97
89) Benzo(a)pyrene	15.62	252	1285040	49.313	nq	98
90) Dibenzo(a,h)anthracene	17.28	278	1196124	53.099	nq	96
91) Benzo(g,h,i)perylene	17.75	276	1240508	55.830	nq #	91

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

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