

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062518\
 Data File : BF106791.D
 Acq On : 26 Jun 2018 00:06
 Operator : JU/SJ
 Sample : J3626-02MSD
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 A-1MSD

Quant Time: Jun 26 05:00:27 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 25 18:50:17 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.96	152	55347	20.00	ng	0.00
21) Naphthalene-d8	8.25	136	197152	20.00	ng	0.00
38) Acenaphthene-d10	10.01	164	88559	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	157020	20.00	ng	0.00
75) Chrysene-d12	14.15	240	134494	20.00	ng	-0.01
86) Perylene-d12	15.69	264	100595	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.61	112	277370	84.86	ng	0.00
7) Phenol-d6	6.61	99	395593	96.92	ng	0.00
23) Nitrobenzene-d5	7.53	82	251338	70.85	ng	0.00
41) 2,4,6-Tribromophenol	10.80	330	74696	72.77	ng	0.00
44) 2-Fluorobiphenyl	9.32	172	442194	84.41	ng	0.00
78) Terphenyl-d14	13.08	244	472127	85.03	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.86	88	41681	24.804	ng	98
3) Pyridine	3.64	79	127962	28.078	ng	95
4) n-Nitrosodimethylamine	3.60	42	51429	26.570	ng	84
6) Aniline	6.63	93	98673	17.821	ng	# 76
8) 2-Chlorophenol	6.76	128	109392	30.514	ng	94
9) Benzaldehyde	6.52	77	55431	17.098	ng	95
10) Phenol	6.62	94	147182	31.596	ng	82
11) bis(2-Chloroethyl)ether	6.70	93	116476	30.288	ng	99
12) 1,3-Dichlorobenzene	6.91	146	137468	32.739	ng	96
13) 1,4-Dichlorobenzene	6.98	146	137787	32.458	ng	98
14) 1,2-Dichlorobenzene	7.14	146	126886	32.615	ng	95
15) Benzyl Alcohol	7.11	79	99936	30.491	ng	96
16) 2,2'-oxybis(1-Chloropropan	7.23	45	160563	31.822	ng	69
17) 2-Methylphenol	7.22	107	98131	31.986	ng	# 90
18) Hexachloroethane	7.47	117	35350	23.680	ng	# 85
19) n-Nitroso-di-n-propylamine	7.37	70	77214	28.698	ng	94
20) 3+4-Methylphenols	7.37	107	123227	34.968	ng	# 61
22) Acetophenone	7.37	105	155413	35.235	ng	# 96
24) Nitrobenzene	7.55	77	120859	33.369	ng	99
25) Isophorone	7.78	82	217367	34.771	ng	99
26) 2-Nitrophenol	7.86	139	42384	24.789	ng	97
27) 2,4-Dimethylphenol	7.90	122	102921	38.944	ng	98
28) bis(2-Chloroethoxy)methane	7.98	93	141908	33.809	ng	98
29) 2,4-Dichlorophenol	8.11	162	92715	33.510	ng	96
30) 1,2,4-Trichlorobenzene	8.18	180	113904	37.371	ng	98
31) Naphthalene	8.27	128	386006	41.878	ng	99
33) 4-Chloroaniline	8.32	127	78475	20.290	ng	98
34) Hexachlorobutadiene	8.37	225	74593	37.559	ng	98
35) Caprolactam	8.67	113	28683	31.803	ng	89
36) 4-Chloro-3-methylphenol	8.81	107	97137	33.355	ng	98
37) 2-Methylnaphthalene	8.96	142	281633	46.097	ng	97
39) 1,2,4,5-Tetrachlorobenzene	9.12	216	114891	38.523	ng	# 95
40) Hexachlorocyclopentadiene	9.10	237	3220	2.098	ng	95
42) 2,4,6-Trichlorophenol	9.24	196	20945	10.565	ng	93

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43) 2,4,5-Trichlorophenol	9.29	196	50856	24.584	nq	89
45) 1,1'-Biphenyl	9.42	154	288165	40.830	nq	97
46) 2-Chloronaphthalene	9.45	162	188678	33.963	nq	94
47) 2-Nitroaniline	9.55	65	61433	32.885	nq	94
48) Acenaphthylene	9.87	152	296519	33.807	nq	99
49) Dimethylphthalate	9.72	163	277318	41.752	nq	99
50) 2,6-Dinitrotoluene	9.79	165	45091	32.060	nq #	83
51) Acenaphthene	10.04	154	292453	52.950	nq	98
52) 3-Nitroaniline	9.97	138	52380	32.364	nq	92
54) Dibenzofuran	10.21	168	409621	54.162	nq	97
55) 4-Nitrophenol	10.14	139	5483	4.853	nq	93
56) 2,4-Dinitrotoluene	10.20	165	52613	28.659	nq #	84
57) Fluorene	10.56	166	320547	53.412	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.33	232	7707	4.687	nq #	63
59) Diethylphthalate	10.41	149	204615	31.918	nq #	93
60) 4-Chlorophenyl-phenylether	10.54	204	105285	33.529	nq #	84
61) 4-Nitroaniline	10.58	138	54917	34.190	nq	94
62) Azobenzene	10.70	77	177822	28.168	nq	90
65) n-Nitrosodiphenylamine	10.66	169	177389	33.587	nq	98
66) 4-Bromophenyl-phenylether	11.04	248	69038	36.841	nq #	76
67) Hexachlorobenzene	11.11	284	69115	35.239	nq #	80
68) Atrazine	11.19	200	59812	35.624	nq	93
70) Phenanthrene	11.54	178	970347	128.126	nq	98
71) Anthracene	11.58	178	407067	52.659	nq	100
72) Carbazole	11.73	167	294862	40.742	nq	98
73) Di-n-butylphthalate	12.04	149	290484	34.069	nq	99
74) Fluoranthene	12.72	202	866360	103.788	nq	96
76) Benzidine	12.84	184	157892	41.221	nq	97
77) Pvrene	12.95	202	720212	81.206	nq	98
79) Butylbenzylphthalate	13.55	149	120586	33.069	nq #	93
80) Benzo(a)anthracene	14.14	228	399313	48.714	nq	100
81) 3,3'-Dichlorobenzidine	14.10	252	83904	29.124	nq #	95
82) Chrysene	14.18	228	364036	48.273	nq	99
83) Bis(2-ethylhexyl)phthalate	14.10	149	165157	35.427	nq	98
84) Di-n-octyl phthalate	14.73	149	284367	37.421	nq #	94
85) Indeno(1,2,3-cd)pyrene	17.27	276	214878	33.576	nq #	89
87) Benzo(b)fluoranthene	15.24	252	284858	45.585	nq #	95
88) Benzo(k)fluoranthene	15.27	252	220745	37.095	nq #	95
89) Benzo(a)pyrene	15.63	252	221792	39.926	nq #	96
90) Dibenzo(a,h)anthracene	17.28	278	169989	36.107	nq #	93
91) Benzo(g,h,i)perylene	17.75	276	176127	38.275	nq #	90

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

