

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122718\
 Data File : BF111777.D
 Acq On : 27 Dec 2018 21:55
 Operator : JU/SJ
 Sample : J6188-06MSD
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BU-03-122118MSD

Manual Integrations
 APPROVED

Sohil
 12/28/2018 10:37:31 AM

Quant Time: Dec 28 07:31:00 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 19 17:39:39 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.91	152	132844	20.00	ng	0.01
21) Naphthalene-d8	8.19	136	454841	20.00	ng	0.01
39) Acenaphthene-d10	9.95	164	204301	20.00	ng	0.01
64) Phenanthrene-d10	11.43	188	369721	20.00	ng	0.02
76) Chrysene-d12	14.07	240	289424	20.00	ng	0.02
87) Perylene-d12	15.56	264	226230	20.00	ng	0.04

System Monitoring Compounds

5) 2-Fluorophenol	5.56	112	833945	107.00	ng	0.04
7) Phenol-d6	6.56	99	937576	94.53	ng	0.04
23) Nitrobenzene-d5	7.47	82	748435	95.78	ng	0.01
42) 2,4,6-Tribromophenol	10.74	330	302999	125.52	ng	0.02
45) 2-Fluorobiphenyl	9.26	172	1315834	93.88	ng	0.00
79) Terphenyl-d14	13.02	244	1375784	90.15	ng	0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.86	88	117307	31.992	ng	# 84
3) Pyridine	3.58	79	260777	27.298	ng	98
4) n-Nitrosodimethylamine	3.51	42	196274	41.982	ng	81
6) Aniline	6.57	93	59554	4.710	ng	# 1
8) 2-Chlorophenol	6.70	128	336924	39.026	ng	97
9) Benzaldehyde	6.46	77	200223	29.352	ng	97
10) Phenol	6.57	94	327404	29.256	ng	96
11) bis(2-Chloroethyl)ether	6.64	93	334692	39.910	ng	98
12) 1,3-Dichlorobenzene	6.85	146	394672	39.447	ng	98
13) 1,4-Dichlorobenzene	6.93	146	399193	39.342	ng	98
14) 1,2-Dichlorobenzene	7.08	146	376002	39.718	ng	98
15) Benzyl Alcohol	7.05	79	289098	36.700	ng	90
16) 2,2'-oxybis(1-Chloropropan	7.18	45	552423	35.767	ng	75
17) 2-Methylphenol	7.17	107	253005	36.599	ng	# 83
18) Hexachloroethane	7.42	117	116953	34.204	ng	# 86
19) n-Nitroso-di-n-propylamine	7.32	70	258493	39.243	ng	98
20) 3+4-Methylphenols	7.32	107	325377	37.250	ng	# 72
22) Acetophenone	7.31	105	482332	41.859	ng	# 90
24) Nitrobenzene	7.49	77	373441	45.598	ng	100
25) Isophorone	7.73	82	652066	47.005	ng	98
26) 2-Nitrophenol	7.80	139	157325	47.365	ng	96
27) 2,4-Dimethylphenol	7.85	122	277445	42.373	ng	92
28) bis(2-Chloroethoxy)methane	7.93	93	406834	44.071	ng	99
29) 2,4-Dichlorophenol	8.05	162	280229	39.790	ng	97
30) 1,2,4-Trichlorobenzene	8.13	180	324640	41.366	ng	96
31) Naphthalene	8.22	128	971190	44.439	ng	97
32) Benzoic acid	7.95	122	123864m	28.611	ng	
33) 4-Chloroaniline	8.26	127	41728	4.418	ng	97
34) Hexachlorobutadiene	8.33	225	201791	42.188	ng	98
35) Caprolactam	8.61	113	55332	23.186	ng	93
36) 4-Chloro-3-methylphenol	8.75	107	256461	37.045	ng	98
37) 2-Methylnaphthalene	8.90	142	638943	44.937	ng	97
38) 1-Methylnaphthalene	9.00	142	600062	43.942	ng	98
40) 1,2,4,5-Tetrachlorobenzene	9.07	216	305999	45.581	ng	# 98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	9.06	237	105417	32.359	ng	98
43) 2,4,6-Trichlorophenol	9.19	196	189484	42.275	ng	99
44) 2,4,5-Trichlorophenol	9.23	196	193496	42.081	ng	95
46) 1,1'-Biphenyl	9.36	154	769390	44.612	ng	94
47) 2-Chloronaphthalene	9.39	162	588340	43.058	ng	94
48) 2-Nitroaniline	9.48	65	182991	51.479	ng	97
49) Acenaphthylene	9.81	152	913656	45.797	ng	94
50) Dimethylphthalate	9.66	163	702141	46.998	ng	98
51) 2,6-Dinitrotoluene	9.72	165	138259	46.411	ng	97
52) Acenaphthene	9.98	154	536124	43.571	ng	98
53) 3-Nitroaniline	9.89	138	76439	24.059	ng	98
54) 2,4-Dinitrophenol	9.99	184	21413	27.659	ng	# 73
55) Dibenzofuran	10.15	168	802315	43.160	ng	95
56) 4-Nitrophenol	10.07	139	162505	67.023	ng	93
57) 2,4-Dinitrotoluene	10.12	165	176801	49.381	ng	# 94
58) Fluorene	10.49	166	615127	45.921	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.27	232	162104	41.891	ng	89
60) Diethylphthalate	10.36	149	668439	45.611	ng	99
61) 4-Chlorophenyl-phenylether	10.48	204	316111	42.713	ng	97
62) 4-Nitroaniline	10.50	138	144261	46.718	ng	89
63) Azobenzene	10.65	77	605310	41.142	ng	94
65) 4,6-Dinitro-2-methylphenol	10.53	198	18284	16.990	ng	80
66) n-Nitrosodiphenylamine	10.60	169	534739	43.087	ng	99
67) 4-Bromophenyl-phenylether	10.97	248	193633	42.225	ng	91
68) Hexachlorobenzene	11.05	284	218694	42.772	ng	98
69) Atrazine	11.13	200	172512	40.011	ng	96
70) Pentachlorophenol	11.24	266	163376	51.686	ng	95
71) Phenanthrene	11.46	178	897014	45.748	ng	95
72) Anthracene	11.51	178	922215	46.858	ng	95
73) Carbazole	11.66	167	792264	41.463	ng	95
74) Di-n-butylphthalate	11.99	149	995281	46.457	ng	96
75) Fluoranthene	12.64	202	944045	47.546	ng	95
77) Benzidine	12.76	184	130427	11.976	ng	# 95
78) Pyrene	12.87	202	953969	46.675	ng	96
80) Butylbenzylphthalate	13.49	149	410610	49.286	ng	88
81) Benzo(a)anthracene	14.06	228	785671	47.566	ng	99
82) 3,3'-Dichlorobenzidine	14.02	252	83936	12.862	ng	97
83) Chrysene	14.10	228	731000	45.029	ng	98
84) Bis(2-ethylhexyl)phthalate	14.04	149	572142	54.520	ng	# 98
85) Di-n-octyl phthalate	14.66	149	901483	55.445	ng	96
86) Indeno(1,2,3-cd)pyrene	17.07	276	571677	41.424	ng	# 86
88) Benzo(b)fluoranthene	15.12	252	665390	50.325	ng	# 96
89) Benzo(k)fluoranthene	15.15	252	560536	44.531	ng	# 97
90) Benzo(a)pyrene	15.50	252	568660	47.341	ng	# 93
91) Dibenzo(a,h)anthracene	17.08	278	489018	46.490	ng	# 90
92) Benzo(g,h,i)perylene	17.52	276	468567	45.620	ng	# 88

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

