

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF042619\
 Data File : BF114001.D
 Acq On : 26 Apr 2019 11:21
 Operator : JU/SJ
 Sample : PB119139BS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB119139BS

Manual Integrations
 APPROVED

Sohil
 4/27/2019 2:36:37 AM

Quant Time: Apr 27 00:20:00 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF042219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 23 03:32:35 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	217822	20.00	ng	0.00
21) Naphthalene-d8	8.18	136	880277	20.00	ng	0.00
39) Acenaphthene-d10	9.94	164	471534	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	944219	20.00	ng	0.00
76) Chrysene-d12	14.07	240	719081	20.00	ng	0.00
87) Perylene-d12	15.55	264	428531	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.55	112	1545051	121.32	ng	0.03
7) Phenol-d6	6.55	99	1985438	124.26	ng	0.02
23) Nitrobenzene-d5	7.46	82	1143182	80.18	ng	0.00
42) 2,4,6-Tribromophenol	10.73	330	696848	121.31	ng	0.01
45) 2-Fluorobiphenyl	9.26	172	2201843	73.03	ng	0.00
79) Terphenyl-d14	13.01	244	2584566	73.69	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.82	88	237429	39.549	ng	97
3) Pyridine	3.56	79	420882	25.685	ng	98
4) n-Nitrosodimethylamine	3.52	42	205374	36.613	ng	93
6) Aniline	6.57	93	578476	26.365	ng	97
8) 2-Chlorophenol	6.69	128	603833	41.529	ng	99
9) Benzaldehyde	6.45	77	29186	Below Cal		94
10) Phenol	6.56	94	743415	38.909	ng	89
11) bis(2-Chloroethyl)ether	6.64	93	585548	40.726	ng	98
12) 1,3-Dichlorobenzene	6.85	146	639439	38.830	ng	97
13) 1,4-Dichlorobenzene	6.92	146	645708	38.984	ng	98
14) 1,2-Dichlorobenzene	7.07	146	610071	40.078	ng	98
15) Benzyl Alcohol	7.05	79	520551	48.373	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.17	45	667174	39.752	ng	89
17) 2-Methylphenol	7.16	107	554849	43.533	ng	96
18) Hexachloroethane	7.42	117	233944	40.292	ng	97
19) n-Nitroso-di-n-propylamine	7.32	70	435380	42.077	ng	99
20) 3+4-Methylphenols	7.31	107	616450	42.809	ng	# 86
22) Acetophenone	7.31	105	846049	43.190	ng	# 87
24) Nitrobenzene	7.48	77	674424	42.812	ng	99
25) Isophorone	7.72	82	1248251	42.790	ng	99
26) 2-Nitrophenol	7.80	139	350696	48.798	ng	99
27) 2,4-Dimethylphenol	7.83	122	597527	51.031	ng	98
28) bis(2-Chloroethoxy)methane	7.93	93	774341	41.656	ng	99
29) 2,4-Dichlorophenol	8.05	162	580754	43.373	ng	97
30) 1,2,4-Trichlorobenzene	8.12	180	632802	42.385	ng	100
31) Naphthalene	8.20	128	1754789	41.962	ng	100
32) Benzoic acid	7.99	122	421799	53.343	ng	95
33) 4-Chloroaniline	8.25	127	504873	28.533	ng	98
34) Hexachlorobutadiene	8.32	225	376910	40.498	ng	99
35) Caprolactam	8.65	113	192678m	45.236	ng	
36) 4-Chloro-3-methylphenol	8.74	107	578747	43.628	ng	99
37) 2-Methylnaphthalene	8.90	142	1222121	43.338	ng	98
38) 1-Methylnaphthalene	9.00	142	1144226	42.040	ng	98
40) 1,2,4,5-Tetrachlorobenzene	9.06	216	624408	40.766	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	9.05	237	556043	65.100	ng	99
43) 2,4,6-Trichlorophenol	9.17	196	445116	45.765	ng	100
44) 2,4,5-Trichlorophenol	9.23	196	470295	43.109	ng	99
46) 1,1'-Biphenyl	9.36	154	1490712	41.384	ng	98
47) 2-Chloronaphthalene	9.39	162	1220412	41.662	ng	98
48) 2-Nitroaniline	9.48	65	353796	46.072	ng	91
49) Acenaphthylene	9.80	152	1815753	39.594	ng	100
50) Dimethylphthalate	9.66	163	1478604	43.358	ng	100
51) 2,6-Dinitrotoluene	9.72	165	341330	46.654	ng	95
52) Acenaphthene	9.97	154	1123933	41.470	ng	97
53) 3-Nitroaniline	9.89	138	269093	35.002	ng	93
54) 2,4-Dinitrophenol	10.00	184	365369	111.743	ng	# 1
55) Dibenzofuran	10.15	168	1669412	41.121	ng	96
56) 4-Nitrophenol	10.06	139	552462	90.759	ng	93
57) 2,4-Dinitrotoluene	10.13	165	457307	49.521	ng	95
58) Fluorene	10.49	166	1264327	41.366	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.26	232	432148	45.076	ng	100
60) Diethylphthalate	10.36	149	1400554	43.248	ng	100
61) 4-Chlorophenyl-phenylether	10.48	204	656211	40.540	ng	100
62) 4-Nitroaniline	10.52	138	341857	45.566	ng	95
63) Azobenzene	10.64	77	1277783	40.711	ng	96
65) 4,6-Dinitro-2-methylphenol	10.54	198	230371	52.434	ng	96
66) n-Nitrosodiphenylamine	10.60	169	1173941	41.436	ng	99
67) 4-Bromophenyl-phenylether	10.97	248	456039	39.980	ng	97
68) Hexachlorobenzene	11.04	284	500930	39.964	ng	96
69) Atrazine	11.12	200	349046	37.872	ng	98
70) Pentachlorophenol	11.23	266	601982	84.823	ng	96
71) Phenanthrene	11.45	178	1928912	40.649	ng	99
72) Anthracene	11.50	178	1938521	40.454	ng	99
73) Carbazole	11.65	167	1755446	41.062	ng	98
74) Di-n-butylphthalate	11.98	149	2070152	41.165	ng	99
75) Fluoranthene	12.64	202	2017786	38.975	ng	99
77) Benzidine	12.75	184	275441	12.856	ng	97
78) Pyrene	12.87	202	2009730	39.966	ng	99
80) Butylbenzylphthalate	13.47	149	854456	43.187	ng	98
81) Benzo(a)anthracene	14.06	228	1729845	40.830	ng	100
82) 3,3'-Dichlorobenzidine	14.02	252	570770	36.756	ng	# 99
83) Chrysene	14.10	228	1685824	40.857	ng	99
84) Bis(2-ethylhexyl)phthalate	14.04	149	1148616	45.629	ng	99
85) Di-n-octyl phthalate	14.66	149	1755674	44.559	ng	99
86) Indeno(1,2,3-cd)pyrene	17.06	276	1409772	36.070	ng	99
88) Benzo(b)fluoranthene	15.13	252	1551701	57.231	ng	100
89) Benzo(k)fluoranthene	15.16	252	1355101	58.104	ng	99
90) Benzo(a)pyrene	15.50	252	1323907	56.453	ng	99
91) Dibenzo(a,h)anthracene	17.08	278	1172027	53.240	ng	99
92) Benzo(g,h,i)perylene	17.52	276	1211956	54.553	ng	98

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

