

Data Path : Z:\HPCHEM1\BNA F\DATA\BF041918\
 Data File : BF104650.D
 Acq On : 20 Apr 2018 1:09
 Operator : JU/SJ
 Sample : PB108340BS
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB108340BS

Manual Integrations
 APPROVED

Sohil
 4/20/2018 3:05:39 PM

Quant Time: Apr 20 02:28:28 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 19 14:44:03 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	134889	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	540196	20.00	ng	0.00
38) Acenaphthene-d10	9.85	164	231594	20.00	ng	0.00
63) Phenanthrene-d10	11.33	188	371461	20.00	ng	0.00
75) Chrysene-d12	13.98	240	266918	20.00	ng	0.00
86) Perylene-d12	15.44	264	230541	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.43	112	942325	106.82	ng	0.02
7) Phenol-d6	6.45	99	1171974	108.13	ng	0.00
23) Nitrobenzene-d5	7.37	82	745904	90.16	ng	0.00
41) 2,4,6-Tribromophenol	10.64	330	274549	114.28	ng	0.00
44) 2-Fluorobiphenyl	9.17	172	1280952	87.38	ng	0.00
78) Terphenyl-d14	12.92	244	1014132	86.62	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.59	88	131218	31.922	ng	# 89
3) Pyridine	3.32	79	374792	32.430	ng	86
4) n-Nitrosodimethylamine	3.28	42	146235	36.198	ng	# 72
6) Aniline	6.47	93	126207	8.381	ng	# 32
8) 2-Chlorophenol	6.59	128	356922	38.333	ng	98
9) Benzaldehyde	6.36	77	202775	35.690	ng	96
10) Phenol	6.46	94	408780	35.544	ng	97
11) bis(2-Chloroethyl)ether	6.55	93	335129	36.516	ng	95
12) 1,3-Dichlorobenzene	6.75	146	374506	35.504	ng	99
13) 1,4-Dichlorobenzene	6.82	146	381277	36.260	ng	99
14) 1,2-Dichlorobenzene	6.97	146	350090	35.928	ng	100
15) Benzyl Alcohol	6.95	79	289341	35.146	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.08	45	425272	34.505	ng	85
17) 2-Methylphenol	7.07	107	300502	37.128	ng	93
18) Hexachloroethane	7.32	117	114757	30.610	ng	99
19) n-Nitroso-di-n-propylamine	7.22	70	230692	32.570	ng	94
20) 3+4-Methylphenols	7.22	107	364711	36.333	ng	# 74
22) Acetophenone	7.22	105	469679	35.316	ng	99
24) Nitrobenzene	7.39	77	362600	39.699	ng	97
25) Isophorone	7.63	82	661889	37.835	ng	98
26) 2-Nitrophenol	7.70	139	182089	48.970	ng	96
27) 2,4-Dimethylphenol	7.75	122	311911	40.400	ng	99
28) bis(2-Chloroethoxy)methane	7.84	93	430054	40.207	ng	100
29) 2,4-Dichlorophenol	7.95	162	294301	39.951	ng	99
30) 1,2,4-Trichlorobenzene	8.03	180	298593	36.934	ng	99
31) Naphthalene	8.11	128	992110	36.883	ng	100
32) Benzoic acid	7.87	122	211687	34.364	ng	96
33) 4-Chloroaniline	8.16	127	59430	5.157	ng	95
34) Hexachlorobutadiene	8.22	225	162737	36.750	ng	98
35) Caprolactam	8.54	113	98837m	38.460	ng	
36) 4-Chloro-3-methylphenol	8.65	107	307137	38.883	ng	99
37) 2-Methylnaphthalene	8.80	142	649371	37.321	ng	98
39) 1,2,4,5-Tetrachlorobenzene	8.97	216	277786	41.901	ng	99
40) Hexachlorocyclopentadiene	8.95	237	71053	19.144	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.09	196	190696	41.436	nq	97
43) 2,4,5-Trichlorophenol	9.13	196	203262	39.469	nq	97
45) 1,1'-Biphenyl	9.27	154	762080	39.170	nq	99
46) 2-Chloronaphthalene	9.29	162	596417	38.811	nq	99
47) 2-Nitroaniline	9.39	65	181187	47.676	nq	98
48) Acenaphthylene	9.71	152	881159	36.546	nq	99
49) Dimethylphthalate	9.57	163	732065	40.085	nq	100
50) 2,6-Dinitrotoluene	9.63	165	153824	45.519	nq	96
51) Acenaphthene	9.88	154	511398	34.763	nq	99
52) 3-Nitroaniline	9.80	138	64358	16.267	nq	98
53) 2,4-Dinitrophenol	9.91	184	48683	44.944	nq	# 24
54) Dibenzofuran	10.05	168	772892	37.700	nq	98
55) 4-Nitrophenol	9.97	139	234744	75.535	nq	98
56) 2,4-Dinitrotoluene	10.04	165	188335	42.487	nq	# 99
57) Fluorene	10.40	166	583015	39.070	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.17	232	150383	39.141	nq	94
59) Diethylphthalate	10.27	149	693648	38.136	nq	99
60) 4-Chlorophenyl-phenylether	10.39	204	276766	41.646	nq	99
61) 4-Nitroaniline	10.42	138	133664	34.554	nq	90
62) Azobenzene	10.55	77	631912	36.364	nq	96
64) 4,6-Dinitro-2-methylphenol	10.45	198	38348	25.300	nq	89
65) n-Nitrosodiphenylamine	10.51	169	532092	41.313	nq	99
66) 4-Bromophenyl-phenylether	10.87	248	169246	42.835	nq	95
67) Hexachlorobenzene	10.95	284	177055	39.824	nq	# 86
68) Atrazine	11.03	200	172963	44.491	nq	100
69) Pentachlorophenol	11.14	266	163510	59.448	nq	99
70) Phenanthrene	11.36	178	785993	37.996	nq	99
71) Anthracene	11.41	178	800886	38.087	nq	100
72) Carbazole	11.57	167	738121	35.922	nq	100
73) Di-n-butylphthalate	11.89	149	995422	39.657	nq	99
74) Fluoranthene	12.55	202	739799	34.229	nq	97
76) Benzidine	12.67	184	393431	45.017	nq	98
77) Pyrene	12.78	202	741415	37.474	nq	98
79) Butylbenzylphthalate	13.39	149	366891	39.362	nq	99
80) Benzo(a)anthracene	13.97	228	642932	40.319	nq	98
81) 3,3'-Dichlorobenzidine	13.93	252	216589	36.429	nq	97
82) Chrysene	14.01	228	612877	37.419	nq	98
83) Bis(2-ethylhexyl)phthalate	13.95	149	495428	41.323	nq	99
84) Di-n-octyl phthalate	14.57	149	810569	40.462	nq	95
85) Indeno(1,2,3-cd)pyrene	16.90	276	512641	36.844	nq	98
87) Benzo(b)fluoranthene	15.02	252	575168	39.502	nq	98
88) Benzo(k)fluoranthene	15.05	252	559451	40.698	nq	99
89) Benzo(a)pyrene	15.39	252	523229	40.162	nq	99
90) Dibenzo(a,h)anthracene	16.92	278	437126	40.853	nq	98
91) Benzo(g,h,i)perylene	17.34	276	416501	40.178	nq	96

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

