

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF020419\
 Data File : BF112378.D
 Acq On : 4 Feb 2019 14:17
 Operator : JU/SJ
 Sample : PB116785BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB116785BS

Manual Integrations
 APPROVED

mohammad
 2/5/2019 8:42:45 AM

Quant Time: Feb 04 16:03:10 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 28 10:38:18 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	161808	20.00	ng	-0.01
21) Naphthalene-d8	8.12	136	630313	20.00	ng	-0.01
39) Acenaphthene-d10	9.87	164	323674	20.00	ng	-0.01
64) Phenanthrene-d10	11.36	188	637471	20.00	ng	-0.01
76) Chrysene-d12	14.00	240	541767	20.00	ng	0.00
87) Perylene-d12	15.47	264	494548	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.48	112	1162062	152.54	ng	0.00
7) Phenol-d6	6.49	99	1465748	138.47	ng	0.00
23) Nitrobenzene-d5	7.40	82	843995	77.15	ng	-0.01
42) 2,4,6-Tribromophenol	10.67	330	498846	132.41	ng	-0.01
45) 2-Fluorobiphenyl	9.19	172	1752230	76.57	ng	-0.01
79) Terphenyl-d14	12.94	244	2053257	71.38	ng	-0.01

Target Compounds				Qvalue		
2) 1,4-Dioxane	2.70	88	147627	48.647	ng	95
3) Pyridine	3.45	79	381425	49.411	ng	97
4) n-Nitrosodimethylamine	3.39	42	176679	54.477	ng	# 90
6) Aniline	6.50	93	363554	28.872	ng	# 35
8) 2-Chlorophenol	6.63	128	414293	40.427	ng	98
9) Benzaldehyde	6.38	77	80024	14.464	ng	98
10) Phenol	6.50	94	474375	40.847	ng	95
11) bis(2-Chloroethyl)ether	6.57	93	361479	39.536	ng	98
12) 1,3-Dichlorobenzene	6.77	146	458299	38.400	ng	100
13) 1,4-Dichlorobenzene	6.85	146	474225	38.609	ng	100
14) 1,2-Dichlorobenzene	7.00	146	441947	38.430	ng	96
15) Benzyl Alcohol	6.98	79	338235	37.905	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.10	45	450112	38.342	ng	# 38
17) 2-Methylphenol	7.10	107	325107	40.019	ng	# 92
18) Hexachloroethane	7.35	117	163489	37.904	ng	94
19) n-Nitroso-di-n-propylamine	7.25	70	272802	37.375	ng	99
20) 3+4-Methylphenols	7.25	107	423029	40.721	ng	# 76
22) Acetophenone	7.24	105	559102	36.428	ng	96
24) Nitrobenzene	7.42	77	415996	38.077	ng	96
25) Isophorone	7.66	82	738256	43.019	ng	98
26) 2-Nitrophenol	7.73	139	227118	38.900	ng	93
27) 2,4-Dimethylphenol	7.77	122	358511	44.568	ng	95
28) bis(2-Chloroethoxy)methane	7.86	93	463123	39.633	ng	99
29) 2,4-Dichlorophenol	7.98	162	360480	40.045	ng	97
30) 1,2,4-Trichlorobenzene	8.06	180	397686	38.433	ng	99
31) Naphthalene	8.14	128	1221150	41.428	ng	99
32) Benzoic acid	7.92	122	188941	41.177	ng	98
33) 4-Chloroaniline	8.19	127	295201	23.000	ng	97
34) Hexachlorobutadiene	8.25	225	220132	36.253	ng	99
35) Caprolactam	8.56	113	107062m	33.714	ng	
36) 4-Chloro-3-methylphenol	8.69	107	368334	38.652	ng	98
37) 2-Methylnaphthalene	8.83	142	861657	42.701	ng	98
38) 1-Methylnaphthalene	8.93	142	806558	41.702	ng	98
40) 1,2,4,5-Tetrachlorobenzene	9.00	216	392768	36.958	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.98	237	242780	62.621	nq	100
43) 2,4,6-Trichlorophenol	9.12	196	242413	37.005	nq	100
44) 2,4,5-Trichlorophenol	9.16	196	257366	37.591	nq	98
46) 1,1'-Biphenyl	9.29	154	1044848	36.925	nq	98
47) 2-Chloronaphthalene	9.32	162	805235	36.696	nq	98
48) 2-Nitroaniline	9.42	65	224303	37.504	nq	89
49) Acenaphthylene	9.73	152	1282743	39.883	nq	99
50) Dimethylphthalate	9.59	163	935933	37.217	nq	99
51) 2,6-Dinitrotoluene	9.66	165	215527	38.267	nq #	85
52) Acenaphthene	9.90	154	744857	38.769	nq	99
53) 3-Nitroaniline	9.83	138	151172	24.775	nq	88
54) 2,4-Dinitrophenol	9.95	184	152002	84.095	nq	95
55) Dibenzofuran	10.08	168	1112443	37.889	nq	97
56) 4-Nitrophenol	10.02	139	242442	75.154	nq	97
57) 2,4-Dinitrotoluene	10.07	165	280904	39.177	nq #	86
58) Fluorene	10.42	166	897608	40.854	nq	98
59) 2,3,4,6-Tetrachlorophenol	10.20	232	207395	40.085	nq	96
60) Diethylphthalate	10.29	149	936096	37.509	nq	100
61) 4-Chlorophenyl-phenylether	10.41	204	429942	35.974	nq	96
62) 4-Nitroaniline	10.45	138	224332	37.483	nq	95
63) Azobenzene	10.57	77	818417	37.288	nq	99
65) 4,6-Dinitro-2-methylphenol	10.49	198	121873	34.099	nq	97
66) n-Nitrosodiphenylamine	10.53	169	794371	36.974	nq	100
67) 4-Bromophenyl-phenylether	10.90	248	270401	36.065	nq	97
68) Hexachlorobenzene	10.97	284	296500	36.860	nq	96
69) Atrazine	11.06	200	261921	37.782	nq	96
70) Pentachlorophenol	11.17	266	218606	69.843	nq	98
71) Phenanthrene	11.39	178	1313334	39.628	nq	98
72) Anthracene	11.44	178	1373667	41.610	nq	99
73) Carbazole	11.59	167	1230385	37.783	nq	99
74) Di-n-butylphthalate	11.91	149	1460953	36.144	nq	99
75) Fluoranthene	12.57	202	1409137	39.643	nq	98
77) Benzdine	12.69	184	533614	35.786	nq	95
78) Pyrene	12.80	202	1437512	38.325	nq	99
80) Butylbenzylphthalate	13.41	149	641876	35.370	nq	96
81) Benzo(a)anthracene	13.99	228	1271231	39.916	nq	100
82) 3,3'-Dichlorobenzidine	13.95	252	307554	25.804	nq	98
83) Chrysene	14.03	228	1195420	39.322	nq	98
84) Bis(2-ethylhexyl)phthalate	13.97	149	888721	34.500	nq	98
85) Di-n-octyl phthalate	14.57	149	1441340	36.367	nq	98
86) Indeno(1,2,3-cd)pyrene	16.95	276	1050349	43.603	nq	97
88) Benzo(b)fluoranthene	15.04	252	1199287	40.549	nq	99
89) Benzo(k)fluoranthene	15.07	252	1117522	39.447	nq	98
90) Benzo(a)pyrene	15.41	252	1102887	41.442	nq	99
91) Dibenzo(a,h)anthracene	16.96	278	876463	40.864	nq	99
92) Benzo(g,h,i)perylene	17.39	276	838236	41.424	nq	98

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

