

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062519\
 Data File : BF115178.D
 Acq On : 25 Jun 2019 9:44
 Operator : HP/JU
 Sample : PB120877BS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB120877BS

Manual Integrations
 APPROVED

mohammad
 6/26/2019 2:03:04 PM

Quant Time: Jun 26 06:33:22 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF062119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 24 04:21:29 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.61	152	270863	20.00	ng	0.00
21) Naphthalene-d8	7.90	136	1072494	20.00	ng	0.00
39) Acenaphthene-d10	9.64	164	534430	20.00	ng	0.00
64) Phenanthrene-d10	11.11	188	1060499	20.00	ng	0.00
76) Chrysene-d12	13.74	240	716126	20.00	ng	-0.01
87) Perylene-d12	15.09	264	785761	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.22	112	1979472	112.78	ng	0.02
7) Phenol-d6	6.26	99	2409416	113.10	ng	0.00
23) Nitrobenzene-d5	7.18	82	1500611	86.07	ng	0.00
42) 2,4,6-Tribromophenol	10.43	330	699134	124.05	ng	0.00
45) 2-Fluorobiphenyl	8.98	172	2775780	88.22	ng	0.00
79) Terphenyl-d14	12.70	244	2880767	85.53	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.28	88	305993	36.939	ng	99
3) Pyridine	2.94	79	834801	35.755	ng	99
4) n-Nitrosodimethylamine	2.88	42	344335	37.620	ng	96
6) Aniline	6.27	93	721940	24.656	ng	# 1
8) 2-Chlorophenol	6.40	128	828485	41.977	ng	99
9) Benzaldehyde	6.15	77	118557	8.530	ng	98
10) Phenol	6.27	94	948861	38.023	ng	91
11) bis(2-Chloroethyl)ether	6.36	93	747733	40.648	ng	98
12) 1,3-Dichlorobenzene	6.55	146	874717	39.934	ng	99
13) 1,4-Dichlorobenzene	6.63	146	881616	40.138	ng	99
14) 1,2-Dichlorobenzene	6.78	146	820206	40.323	ng	98
15) Benzyl Alcohol	6.76	79	578292	37.278	ng	98
16) 2,2'-oxybis(1-Chloropropan	6.90	45	1067598	40.014	ng	95
17) 2-Methylphenol	6.88	107	737561	45.274	ng	98
18) Hexachloroethane	7.12	117	316648	39.987	ng	96
19) n-Nitroso-di-n-propylamine	7.04	70	531975	39.003	ng	99
20) 3+4-Methylphenols	7.03	107	819155	43.546	ng	96
22) Acetophenone	7.02	105	1087916	44.309	ng	# 97
24) Nitrobenzene	7.20	77	841000	43.785	ng	99
25) Isophorone	7.44	82	1551850	43.450	ng	99
26) 2-Nitrophenol	7.51	139	474748	50.050	ng	100
27) 2,4-Dimethylphenol	7.56	122	782685	50.397	ng	98
28) bis(2-Chloroethoxy)methane	7.65	93	962958	42.658	ng	99
29) 2,4-Dichlorophenol	7.75	162	733143	45.744	ng	97
30) 1,2,4-Trichlorobenzene	7.84	180	776031	43.835	ng	100
31) Naphthalene	7.92	128	2310931	43.896	ng	100
32) Benzoic acid	7.70	122	627187	56.576	ng	99
33) 4-Chloroaniline	7.97	127	557305	23.163	ng	98
34) Hexachlorobutadiene	8.04	225	419581	43.249	ng	98
35) Caprolactam	8.34	113	243538m	45.196	ng	
36) 4-Chloro-3-methylphenol	8.46	107	746693	46.004	ng	97
37) 2-Methylnaphthalene	8.61	142	1608834	45.323	ng	99
38) 1-Methylnaphthalene	8.71	142	1515973	44.717	ng	99
40) 1,2,4,5-Tetrachlorobenzene	8.78	216	682626	45.201	ng	99

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41) Hexachlorocyclopentadiene	8.77	237	714065	79.739	nq	99
43) 2,4,6-Trichlorophenol	8.88	196	530604	45.509	nq	100
44) 2,4,5-Trichlorophenol	8.93	196	580386	48.338	nq	99
46) 1,1'-Biphenyl	9.07	154	1869302	43.714	nq	99
47) 2-Chloronaphthalene	9.09	162	1496647	43.148	nq	97
48) 2-Nitroaniline	9.18	65	455355	45.029	nq	98
49) Acenaphthylene	9.51	152	2307644	42.700	nq	98
50) Dimethylphthalate	9.38	163	1734105	44.103	nq	100
51) 2,6-Dinitrotoluene	9.43	165	419090	46.168	nq	99
52) Acenaphthene	9.68	154	1420681	42.011	nq	99
53) 3-Nitroaniline	9.60	138	318964	28.794	nq	97
54) 2,4-Dinitrophenol	9.70	184	469859	98.700	nq	93
55) Dibenzofuran	9.85	168	2039923	42.028	nq	98
56) 4-Nitrophenol	9.77	139	778672	87.679	nq	97
57) 2,4-Dinitrotoluene	9.83	165	558887	47.682	nq	92
58) Fluorene	10.19	166	1402240	43.116	nq	98
59) 2,3,4,6-Tetrachlorophenol	9.97	232	493102	48.977	nq	98
60) Diethylphthalate	10.08	149	1701757	43.056	nq	100
61) 4-Chlorophenyl-phenylether	10.18	204	680728	44.207	nq	97
62) 4-Nitroaniline	10.21	138	472721	42.538	nq	97
63) Azobenzene	10.34	77	1557638	42.193	nq	97
65) 4,6-Dinitro-2-methylphenol	10.24	198	275354	48.715	nq	84
66) n-Nitrosodiphenylamine	10.31	169	1427117	43.194	nq	99
67) 4-Bromophenyl-phenylether	10.67	248	490619	42.670	nq	96
68) Hexachlorobenzene	10.74	284	536971	43.025	nq	99
69) Atrazine	10.84	200	458395	43.130	nq	100
70) Pentachlorophenol	10.92	266	679538	85.467	nq	98
71) Phenanthrene	11.14	178	2277861	39.893	nq	99
72) Anthracene	11.20	178	2361336	40.969	nq	99
73) Carbazole	11.34	167	2176101	42.281	nq	98
74) Di-n-butylphthalate	11.70	149	2464624	41.440	nq	99
75) Fluoranthene	12.32	202	2380191	41.780	nq	99
77) Benzidine	12.44	184	836217	26.297	nq	100
78) Pyrene	12.54	202	2435213	44.249	nq	98
80) Butylbenzylphthalate	13.18	149	1136955	46.813	nq	97
81) Benzo(a)anthracene	13.72	228	2254401	43.873	nq	99
82) 3,3'-Dichlorobenzidine	13.69	252	730904	39.390	nq	99
83) Chrysene	13.77	228	2103701	44.694	nq	98
84) Bis(2-ethylhexyl)phthalate	13.75	149	1458043	45.816	nq	# 98
85) Di-n-octyl phthalate	14.35	149	2520347	47.136	nq	100
86) Indeno(1,2,3-cd)pyrene	16.38	276	2617693	46.999	nq	99
88) Benzo(b)fluoranthene	14.73	252	2573519	52.490	nq	99
89) Benzo(k)fluoranthene	14.75	252	1991525	45.294	nq	100
90) Benzo(a)pyrene	15.05	252	2275550	51.494	nq	99
91) Dibenzo(a,h)anthracene	16.40	278	2132177	51.683	nq	98
92) Benzo(g,h,i)perylene	16.76	276	2209156	52.908	nq	99

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

