

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF041020\
 Data File : BF119893.D
 Acq On : 10 Apr 2020 14:37
 Operator : MA/SJ
 Sample : PB128115BS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB128115BS

Manual Integrations
 APPROVED

mohammad
 4/14/2020 3:39:01 PM

Quant Time: Apr 11 05:32:09 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 10 13:48:19 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.75	152	188687	20.00	ng	0.00
21) Naphthalene-d8	8.04	136	741100	20.00	ng	0.00
39) Acenaphthene-d10	9.80	164	391904	20.00	ng	0.00
64) Phenanthrene-d10	11.28	188	764381	20.00	ng	0.00
76) Chrysene-d12	13.93	240	528996	20.00	ng	0.00
87) Perylene-d12	15.36	264	472046	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.37	112	1324225	121.29	ng	0.02
7) Phenol-d6	6.39	99	1693477	120.37	ng	0.00
23) Nitrobenzene-d5	7.32	82	1061210	74.08	ng	0.00
42) 2,4,6-Tribromophenol	10.59	330	522795	108.96	ng	0.00
45) 2-Fluorobiphenyl	9.12	172	2082714	75.45	ng	0.00
79) Terphenyl-d14	12.87	244	2359421	75.05	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.51	88	160759	33.058	ng	98
3) Pyridine	3.22	79	426920	31.997	ng	98
4) n-Nitrosodimethylamine	3.17	42	255960	37.547	ng	99
6) Aniline	6.42	93	674371	36.521	ng	99
8) 2-Chlorophenol	6.54	128	557923	45.734	ng	98
9) Benzaldehyde	6.30	77	218473	24.146	ng	100
10) Phenol	6.40	94	704610	44.146	ng	97
11) bis(2-Chloroethyl)ether	6.49	93	525484	42.640	ng	98
12) 1,3-Dichlorobenzene	6.69	146	558212	41.796	ng	100
13) 1,4-Dichlorobenzene	6.77	146	559712	41.141	ng	99
14) 1,2-Dichlorobenzene	6.93	146	527714	42.089	ng	97
15) Benzyl Alcohol	6.90	79	414615	37.944	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.03	45	948867	39.568	ng	99
17) 2-Methylphenol	7.02	107	445034	40.878	ng	96
18) Hexachloroethane	7.27	117	212561	41.806	ng	93
19) n-Nitroso-di-n-propylamine	7.17	70	384509	42.502	ng	98
20) 3+4-Methylphenols	7.17	107	553191	41.547	ng	98
22) Acetophenone	7.17	105	722955	41.659	ng	97
24) Nitrobenzene	7.34	77	581948	40.383	ng	99
25) Isophorone	7.58	82	1077383	39.014	ng	100
26) 2-Nitrophenol	7.66	139	302371	41.998	ng	95
27) 2,4-Dimethylphenol	7.70	122	478030	44.024	ng	100
28) bis(2-Chloroethoxy)methane	7.79	93	687409	42.303	ng	99
29) 2,4-Dichlorophenol	7.90	162	470918	42.310	ng	97
30) 1,2,4-Trichlorobenzene	7.98	180	494004	39.172	ng	96
31) Naphthalene	8.06	128	1584065	40.626	ng	100
32) Benzoic acid	7.83	122	387451	40.629	ng	95
33) 4-Chloroaniline	8.11	127	481321	28.356	ng	99
34) Hexachlorobutadiene	8.18	225	286581	37.962	ng	99
35) Caprolactam	8.49	113	149164m	43.812	ng	
36) 4-Chloro-3-methylphenol	8.60	107	518690	43.044	ng	98
37) 2-Methylnaphthalene	8.75	142	1076088	40.707	ng	98
38) 1-Methylnaphthalene	8.85	142	1044074	41.903	ng	99
40) 1,2,4,5-Tetrachlorobenzene	8.92	216	473431	38.248	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.91	237	572145	81.287	nq	99
43) 2,4,6-Trichlorophenol	9.03	196	349619	40.903	nq	98
44) 2,4,5-Trichlorophenol	9.07	196	374751	38.927	nq	98
46) 1,1'-Biphenyl	9.22	154	1317781	39.838	nq	99
47) 2-Chloronaphthalene	9.25	162	1019817	40.021	nq	99
48) 2-Nitroaniline	9.34	65	374069	40.508	nq	96
49) Acenaphthylene	9.66	152	1606997	41.467	nq	99
50) Dimethylphthalate	9.53	163	1208137	39.855	nq	99
51) 2,6-Dinitrotoluene	9.59	165	270451	40.742	nq	89
52) Acenaphthene	9.83	154	1153618m	44.625	nq	
53) 3-Nitroaniline	9.75	138	227400	29.995	nq	97
54) 2,4-Dinitrophenol	9.86	184	338768	85.241	nq	92
55) Dibenzofuran	10.00	168	1458052	41.101	nq	98
56) 4-Nitrophenol	9.92	139	471418	83.572	nq	97
57) 2,4-Dinitrotoluene	9.99	165	359776	41.137	nq	100
58) Fluorene	10.35	166	1139475	40.164	nq	99
59) 2,3,4,6-Tetrachlorophenol	10.12	232	323755	43.971	nq	97
60) Diethylphthalate	10.23	149	1244619	39.615	nq	99
61) 4-Chlorophenyl-phenylether	10.34	204	537603	36.972	nq	95
62) 4-Nitroaniline	10.37	138	299963	41.517	nq	96
63) Azobenzene	10.50	77	1172492	41.643	nq	99
65) 4,6-Dinitro-2-methylphenol	10.40	198	213662	42.429	nq	83
66) n-Nitrosodiphenylamine	10.46	169	1031138	40.562	nq	99
67) 4-Bromophenyl-phenylether	10.83	248	354945	37.340	nq	97
68) Hexachlorobenzene	10.90	284	388523	40.289	nq	93
69) Atrazine	10.99	200	307252	43.440	nq	99
70) Pentachlorophenol	11.09	266	440160	79.205	nq	99
71) Phenanthrene	11.31	178	1630968	40.091	nq	98
72) Anthracene	11.36	178	1708282	41.106	nq	100
73) Carbazole	11.52	167	1576474	40.113	nq	99
74) Di-n-butylphthalate	11.85	149	1932258	37.400	nq	100
75) Fluoranthene	12.50	202	1804613	41.113	nq	99
77) Benzidine	12.62	184	676167	37.814	nq	97
78) Pyrene	12.73	202	1797916	40.316	nq	100
80) Butylbenzylphthalate	13.35	149	791911	36.596	nq	100
81) Benzo(a)anthracene	13.92	228	1383480	39.754	nq	100
82) 3,3'-Dichlorobenzidine	13.88	252	421974	32.497	nq	96
83) Chrysene	13.96	228	1421980	40.214	nq	99
84) Bis(2-ethylhexyl)phthalate	13.92	149	976379	33.319	nq	99
85) Di-n-octyl phthalate	14.52	149	1692735	33.926	nq	98
86) Indeno(1,2,3-cd)pyrene	16.77	276	1194620	38.326	nq	99
88) Benzo(b)fluoranthene	14.95	252	1377238	40.557	nq	100
89) Benzo(k)fluoranthene	14.98	252	1248103	42.598	nq	100
90) Benzo(a)pyrene	15.30	252	1160730	40.654	nq	100
91) Dibenzo(a,h)anthracene	16.79	278	994054	39.305	nq	99
92) Benzo(g,h,i)perylene	17.20	276	974299	39.027	nq	99

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

