

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF032020\
 Data File : BF119470.D
 Acq On : 20 Mar 2020 14:14
 Operator : CG/JU
 Sample : L1749-01MS
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BU-02-031920MS

Manual Integrations
 APPROVED

mohammad
 3/23/2020 9:40:38 AM

Quant Time: Mar 23 04:23:53 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF031820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 18 15:11:18 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	280370	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	1057972	20.00	ng	0.00
39) Acenaphthene-d10	9.90	164	546375	20.00	ng	0.00
64) Phenanthrene-d10	11.39	188	1037501	20.00	ng	0.00
76) Chrysene-d12	14.03	240	634841	20.00	ng	0.00
87) Perylene-d12	15.50	264	642192	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.47	112	1859480	105.58	ng	0.01
7) Phenol-d6	6.48	99	2392703	106.35	ng	0.00
23) Nitrobenzene-d5	7.42	82	1506509	77.39	ng	0.00
42) 2,4,6-Tribromophenol	10.69	330	659326	116.97	ng	0.00
45) 2-Fluorobiphenyl	9.22	172	2695827	73.10	ng	0.00
79) Terphenyl-d14	12.97	244	2857048	88.17	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.64	88	317798	36.535	ng	99
3) Pyridine	3.39	79	797261	33.269	ng	99
4) n-Nitrosodimethylamine	3.33	42	470737	40.281	ng	95
6) Aniline	6.51	93	974877	31.396	ng	99
8) 2-Chlorophenol	6.63	128	865115	46.768	ng	98
9) Benzaldehyde	6.40	77	554680	38.864	ng	98
10) Phenol	6.49	94	1070157m	44.578	ng	
11) bis(2-Chloroethyl)ether	6.59	93	824331	42.934	ng	97
12) 1,3-Dichlorobenzene	6.79	146	875596	43.735	ng	100
13) 1,4-Dichlorobenzene	6.87	146	885532	44.221	ng	100
14) 1,2-Dichlorobenzene	7.02	146	822587	43.947	ng	99
15) Benzyl Alcohol	6.99	79	754706	44.595	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.13	45	1586146	40.956	ng	99
17) 2-Methylphenol	7.10	107	697784	41.171	ng	99
18) Hexachloroethane	7.37	117	332140	45.259	ng	96
19) n-Nitroso-di-n-propylamine	7.27	70	586763	43.269	ng	97
20) 3+4-Methylphenols	7.26	107	846585m	40.758	ng	
22) Acetophenone	7.26	105	1109373	43.884	ng	97
24) Nitrobenzene	7.43	77	891672	42.861	ng	99
25) Isophorone	7.67	82	1651813	41.830	ng	99
26) 2-Nitrophenol	7.75	139	452976	55.300	ng	98
27) 2,4-Dimethylphenol	7.79	122	797703	47.097	ng	99
28) bis(2-Chloroethoxy)methane	7.89	93	1031112	44.157	ng	99
29) 2,4-Dichlorophenol	7.99	162	694869	44.649	ng	99
30) 1,2,4-Trichlorobenzene	8.08	180	745951	43.341	ng	99
31) Naphthalene	8.16	128	2391994	43.935	ng	100
32) Benzoic acid	7.90	122	502547	46.126	ng	96
33) 4-Chloroaniline	8.20	127	537859	21.560	ng	99
34) Hexachlorobutadiene	8.28	225	419127	43.525	ng	98
35) Caprolactam	8.57	113	221618m	43.511	ng	
36) 4-Chloro-3-methylphenol	8.69	107	754120	44.335	ng	99
37) 2-Methylnaphthalene	8.85	142	1596495	44.681	ng	99
38) 1-Methylnaphthalene	8.95	142	1521802	44.997	ng	100
40) 1,2,4,5-Tetrachlorobenzene	9.02	216	679745	42.445	ng	98

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41) Hexachlorocyclopentadiene	9.01	237	739845	81.261	nq	100
43) 2,4,6-Trichlorophenol	9.13	196	514118	44.860	nq	99
44) 2,4,5-Trichlorophenol	9.17	196	533466	41.552	nq	98
46) 1,1'-Biphenyl	9.32	154	1907312	42.861	nq	98
47) 2-Chloronaphthalene	9.35	162	1486791	42.654	nq	98
48) 2-Nitroaniline	9.43	65	558955	46.458	nq	98
49) Acenaphthylene	9.76	152	2386877	43.605	nq	99
50) Dimethylphthalate	9.62	163	1974700	50.882	nq	99
51) 2,6-Dinitrotoluene	9.68	165	393045	47.249	nq	88
52) Acenaphthene	9.93	154	1572983	46.095	nq	98
53) 3-Nitroaniline	9.85	138	278616	26.530	nq	96
54) 2,4-Dinitrophenol	9.95	184	307799	85.012	nq	94
55) Dibenzofuran	10.10	168	2108477	43.095	nq	98
56) 4-Nitrophenol	10.00	139	731419	88.285	nq	91
57) 2,4-Dinitrotoluene	10.09	165	531558	50.726	nq	96
58) Fluorene	10.45	166	1592097	44.005	nq	98
59) 2,3,4,6-Tetrachlorophenol	10.22	232	451441	47.042	nq	99
60) Diethylphthalate	10.32	149	1772903	45.010	nq	99
61) 4-Chlorophenyl-phenylether	10.44	204	763753	43.168	nq	100
62) 4-Nitroaniline	10.46	138	409612	40.674	nq	98
63) Azobenzene	10.60	77	1702388	42.919	nq	94
65) 4,6-Dinitro-2-methylphenol	10.49	198	222153	51.064	nq	98
66) n-Nitrosodiphenylamine	10.56	169	1474321	43.287	nq	97
67) 4-Bromophenyl-phenylether	10.93	248	499662	42.288	nq	94
68) Hexachlorobenzene	11.00	284	534276	44.180	nq	96
69) Atrazine	11.09	200	437722	46.878	nq	99
70) Pentachlorophenol	11.19	266	604728	85.282	nq	97
71) Phenanthrene	11.41	178	2408422	44.769	nq	99
72) Anthracene	11.46	178	2451202	44.831	nq	100
73) Carbazole	11.62	167	2273058	42.001	nq	99
74) Di-n-butylphthalate	11.95	149	2707648	46.770	nq	99
75) Fluoranthene	12.60	202	2559091	43.955	nq	99
77) Benzidine	12.72	184	1359232	50.592	nq	98
78) Pyrene	12.83	202	2551759	48.977	nq	100
80) Butylbenzylphthalate	13.45	149	1139379	54.070	nq	98
81) Benzo(a)anthracene	14.02	228	1818544	44.051	nq	99
82) 3,3'-Dichlorobenzidine	13.98	252	626657	38.499	nq	98
83) Chrysene	14.06	228	1822663	42.780	nq	99
84) Bis(2-ethylhexyl)phthalate	14.02	149	1396576	55.596	nq	99
85) Di-n-octyl phthalate	14.63	149	2289516	51.824	nq	98
86) Indeno(1,2,3-cd)pyrene	16.98	276	2130525	53.096	nq	98
88) Benzo(b)fluoranthene	15.07	252	1804557	42.066	nq	98
89) Benzo(k)fluoranthene	15.10	252	1663120	40.715	nq	99
90) Benzo(a)pyrene	15.44	252	1617866	41.499	nq	99
91) Dibenzo(a,h)anthracene	17.00	278	1743508	51.130	nq	98
92) Benzo(g,h,i)perylene	17.43	276	1830244	53.427	nq	97

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

