

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF051820\
 Data File : BF120353.D
 Acq On : 18 May 2020 20:28
 Operator : MA/SJ
 Sample : L2506-08MS 2X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TR-05-051520

Manual Integrations
 APPROVED

mohammad
 5/19/2020 3:05:33 PM

Quant Time: May 19 05:34:22 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF051420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon May 18 15:23:07 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.60	152	280270	20.00	ng	0.00
21) Naphthalene-d8	7.89	136	1053236	20.00	ng	0.00
39) Acenaphthene-d10	9.64	164	487444	20.00	ng	0.00
64) Phenanthrene-d10	11.12	188	738213	20.00	ng	0.00
76) Chrysene-d12	13.76	240	513812	20.00	ng	0.00
87) Perylene-d12	15.14	264	591465	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.21	112	1199525	83.27	ng	0.00
7) Phenol-d6	6.26	99	1363869	74.87	ng	0.00
23) Nitrobenzene-d5	7.17	82	777460	50.54	ng	0.00
42) 2,4,6-Tribromophenol	10.43	330	250644	63.07	ng	0.00
45) 2-Fluorobiphenyl	8.97	172	1424735	60.05	ng	0.00
79) Terphenyl-d14	12.72	244	1196543	51.45	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.22	88	134213	23.423	ng	97
3) Pyridine	2.89	79	284135	15.775	ng	97
4) n-Nitrosodimethylamine	2.82	42	162133	22.988	ng	# 91
6) Aniline	6.27	93	295754m	12.695	ng	
8) 2-Chlorophenol	6.39	128	350701	20.881	ng	97
9) Benzaldehyde	6.15	77	224633	20.004	ng	99
10) Phenol	6.27	94	452890	21.298	ng	83
11) bis(2-Chloroethyl)ether	6.34	93	341529	20.107	ng	98
12) 1,3-Dichlorobenzene	6.54	146	356386	19.830	ng	99
13) 1,4-Dichlorobenzene	6.62	146	376630	20.347	ng	99
14) 1,2-Dichlorobenzene	6.77	146	344388	20.661	ng	99
15) Benzyl Alcohol	6.76	79	269257	20.806	ng	97
16) 2,2'-oxybis(1-Chloropropan	6.89	45	554571	20.717	ng	81
17) 2-Methylphenol	6.88	107	263117	18.558	ng	93
18) Hexachloroethane	7.11	117	123172	17.468	ng	99
19) n-Nitroso-di-n-propylamine	7.02	70	232849	19.512	ng	90
20) 3+4-Methylphenols	7.04	107	342384	18.574	ng	96
22) Acetophenone	7.02	105	466243	21.609	ng	# 97
24) Nitrobenzene	7.19	77	344587	20.814	ng	94
25) Isophorone	7.43	82	632105	19.749	ng	97
26) 2-Nitrophenol	7.51	139	176379	20.370	ng	96
27) 2,4-Dimethylphenol	7.56	122	286107	22.715	ng	96
28) bis(2-Chloroethoxy)methane	7.65	93	415619	20.659	ng	98
29) 2,4-Dichlorophenol	7.76	162	264377	20.437	ng	98
30) 1,2,4-Trichlorobenzene	7.83	180	295725	20.629	ng	98
31) Naphthalene	7.91	128	965056	21.710	ng	98
32) Benzoic acid	7.67	122	67421	8.166	ng	95
33) 4-Chloroaniline	7.97	127	202333	9.988	ng	99
34) Hexachlorobutadiene	8.03	225	168148	20.294	ng	96
35) Caprolactam	8.33	113	81052	19.152	ng	# 89
36) 4-Chloro-3-methylphenol	8.47	107	268345	19.176	ng	97
37) 2-Methylnaphthalene	8.60	142	625533	20.564	ng	100
38) 1-Methylnaphthalene	8.70	142	604523	20.490	ng	100
40) 1,2,4,5-Tetrachlorobenzene	8.77	216	283431	23.060	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.75	237	24555	11.069	ng	96
43) 2,4,6-Trichlorophenol	8.89	196	177728	21.542	ng	100
44) 2,4,5-Trichlorophenol	8.94	196	184173	19.438	ng	98
46) 1,1'-Biphenyl	9.07	154	739631	23.158	ng	98
47) 2-Chloronaphthalene	9.09	162	554642	21.650	ng	96
48) 2-Nitroaniline	9.19	65	189728	20.297	ng	94
49) Acenaphthylene	9.50	152	938310	23.977	ng	98
50) Dimethylphthalate	9.37	163	847870	27.516	ng	98
51) 2,6-Dinitrotoluene	9.43	165	132645	19.287	ng	98
52) Acenaphthene	9.67	154	507231	21.624	ng	98
53) 3-Nitroaniline	9.60	138	94736	11.551	ng	92
54) 2,4-Dinitrophenol	9.72	184	34406	10.599	ng	# 88
55) Dibenzofuran	9.85	168	744870	22.051	ng	100
56) 4-Nitrophenol	9.80	139	150157	34.075	ng	94
57) 2,4-Dinitrotoluene	9.84	165	158910	19.698	ng	100
58) Fluorene	10.19	166	588930	22.893	ng	99
59) 2,3,4,6-Tetrachlorophenol	9.97	232	141319	19.819	ng	99
60) Diethylphthalate	10.07	149	643923	21.646	ng	99
61) 4-Chlorophenyl-phenylether	10.19	204	280372	20.962	ng	93
62) 4-Nitroaniline	10.22	138	108261	14.176	ng	97
63) Azobenzene	10.35	77	553528	19.688	ng	98
65) 4,6-Dinitro-2-methylphenol	10.25	198	38893	10.418	ng	91
66) n-Nitrosodiphenylamine	10.30	169	506452	24.798	ng	98
67) 4-Bromophenyl-phenylether	10.67	248	153951	21.491	ng	95
68) Hexachlorobenzene	10.73	284	156325	20.860	ng	98
69) Atrazine	10.83	200	130075	21.285	ng	98
70) Pentachlorophenol	10.94	266	127603	40.034	ng	99
71) Phenanthrene	11.15	178	844138	27.213	ng	99
72) Anthracene	11.20	178	794880	23.936	ng	98
73) Carbazole	11.36	167	728522	23.480	ng	97
74) Di-n-butylphthalate	11.70	149	958248	25.625	ng	97
75) Fluoranthene	12.34	202	926702	27.872	ng	97
77) Benzidine	12.48	184	188886	12.589	ng	97
78) Pyrene	12.57	202	1104252	30.146	ng	98
80) Butylbenzylphthalate	13.19	149	329237	19.088	ng	96
81) Benzo(a)anthracene	13.75	228	777731	28.698	ng	98
82) 3,3'-Dichlorobenzidine	13.72	252	140307	13.739	ng	98
83) Chrysene	13.79	228	755428	25.959	ng	99
84) Bis(2-ethylhexyl)phthalate	13.75	149	441389	20.568	ng	99
85) Di-n-octyl phthalate	14.36	149	792275	20.522	ng	99
86) Indeno(1,2,3-cd)pyrene	16.46	276	662498	25.084	ng	98
88) Benzo(b)fluoranthene	14.76	252	868011	26.872	ng	99
89) Benzo(k)fluoranthene	14.79	252	673284	24.059	ng	100
90) Benzo(a)pyrene	15.09	252	737777	25.468	ng	99
91) Dibenzo(a,h)anthracene	16.47	278	514871	20.063	ng	98
92) Benzo(g,h,i)perylene	16.86	276	566258	21.881	ng	99

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

