

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030220\
 Data File : BF119164.D
 Acq On : 2 Mar 2020 9:41
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad
 3/3/2020 12:25:47 PM

Quant Time: Mar 03 07:28:07 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 26 16:03:05 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.74	152	181587	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	691706	20.00	ng	-0.01
39) Acenaphthene-d10	9.77	164	402623	20.00	ng	-0.01
64) Phenanthrene-d10	11.26	188	758371	20.00	ng	0.00
76) Chrysene-d12	13.90	240	558508	20.00	ng	0.00
87) Perylene-d12	15.32	264	498703	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.36	112	833382	76.70	ng	0.00
7) Phenol-d6	6.39	99	1040673	78.75	ng	-0.01
23) Nitrobenzene-d5	7.31	82	1026022	78.32	ng	-0.01
42) 2,4,6-Tribromophenol	10.57	330	423085	80.33	ng	0.00
45) 2-Fluorobiphenyl	9.10	172	2029243	74.25	ng	-0.01
79) Terphenyl-d14	12.84	244	2584243	79.66	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.42	88	164826	34.070	ng	99
3) Pyridine	3.15	79	473834	35.863	ng	100
4) n-Nitrosodimethylamine	3.10	42	157106	39.253	ng	100
6) Aniline	6.40	93	632935	38.816	ng	93
8) 2-Chlorophenol	6.53	128	463262	39.506	ng	97
9) Benzaldehyde	6.29	77	298371	33.794	ng	99
10) Phenol	6.40	94	548672	38.402	ng	92
11) bis(2-Chloroethyl)ether	6.48	93	440884	40.329	ng	99
12) 1,3-Dichlorobenzene	6.68	146	544717	38.757	ng	98
13) 1,4-Dichlorobenzene	6.76	146	553276	39.339	ng	99
14) 1,2-Dichlorobenzene	6.91	146	507866	39.594	ng	100
15) Benzyl Alcohol	6.89	79	403127	42.209	ng	100
16) 2,2'-oxybis(1-Chloropropan	7.02	45	383664	40.514	ng	73
17) 2-Methylphenol	7.01	107	379245	39.890	ng	97
18) Hexachloroethane	7.25	117	203990	40.541	ng	98
19) n-Nitroso-di-n-propylamine	7.16	70	334143	43.394	ng	99
20) 3+4-Methylphenols	7.16	107	473627	40.663	ng	92
22) Acetophenone	7.16	105	645864	40.356	ng	# 98
24) Nitrobenzene	7.33	77	510256	39.387	ng	97
25) Isophorone	7.57	82	924836	40.649	ng	100
26) 2-Nitrophenol	7.64	139	271158	39.503	ng	97
27) 2,4-Dimethylphenol	7.69	122	360168	38.661	ng	98
28) bis(2-Chloroethoxy)methane	7.77	93	592135	39.985	ng	100
29) 2,4-Dichlorophenol	7.89	162	429904	39.199	ng	98
30) 1,2,4-Trichlorobenzene	7.96	180	494708	38.045	ng	99
31) Naphthalene	8.05	128	1386800	38.659	ng	99
32) Benzoic acid	7.85	122	276156	41.652	ng	99
33) 4-Chloroaniline	8.10	127	604426	39.174	ng	98
34) Hexachlorobutadiene	8.16	225	317539	39.204	ng	99
35) Caprolactam	8.49	113	139183m	41.216	ng	
36) 4-Chloro-3-methylphenol	8.59	107	451639	40.691	ng	98
37) 2-Methylnaphthalene	8.73	142	954246	39.527	ng	99
38) 1-Methylnaphthalene	8.83	142	907862	39.987	ng	97
40) 1,2,4,5-Tetrachlorobenzene	8.90	216	512443	37.750	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.89	237	121051	31.016	ng	100
43) 2,4,6-Trichlorophenol	9.02	196	326699	38.111	ng	99
44) 2,4,5-Trichlorophenol	9.07	196	335177	37.261	ng	99
46) 1,1'-Biphenyl	9.20	154	1240942	38.136	ng	100
47) 2-Chloronaphthalene	9.22	162	997095	38.024	ng	99
48) 2-Nitroaniline	9.33	65	297871	40.727	ng	95
49) Acenaphthylene	9.64	152	1430941	37.783	ng	99
50) Dimethylphthalate	9.50	163	1242070	40.343	ng	99
51) 2,6-Dinitrotoluene	9.57	165	268146	39.202	ng	90
52) Acenaphthene	9.81	154	880300	38.547	ng	100
53) 3-Nitroaniline	9.74	138	292674	38.595	ng	95
54) 2,4-Dinitrophenol	9.86	184	117177	39.429	ng	96
55) Dibenzofuran	9.98	168	1281652	37.600	ng	100
56) 4-Nitrophenol	9.92	139	163502	35.886	ng	98
57) 2,4-Dinitrotoluene	9.97	165	337372	38.052	ng	94
58) Fluorene	10.32	166	1057569	39.928	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.10	232	298096	39.273	ng	98
60) Diethylphthalate	10.20	149	1212335	41.785	ng	100
61) 4-Chlorophenyl-phenylether	10.31	204	564971	40.298	ng	99
62) 4-Nitroaniline	10.36	138	304202	39.209	ng	96
63) Azobenzene	10.47	77	1027460	40.741	ng	98
65) 4,6-Dinitro-2-methylphenol	10.39	198	187089	40.769	ng	99
66) n-Nitrosodiphenylamine	10.43	169	965685	38.889	ng	99
67) 4-Bromophenyl-phenylether	10.80	248	386279	38.967	ng	97
68) Hexachlorobenzene	10.87	284	446328	39.052	ng	96
69) Atrazine	10.96	200	331687	38.230	ng	98
70) Pentachlorophenol	11.07	266	179251	38.935	ng	99
71) Phenanthrene	11.29	178	1565131	37.844	ng	99
72) Anthracene	11.33	178	1598602	38.685	ng	99
73) Carbazole	11.49	167	1422725	38.646	ng	99
74) Di-n-butylphthalate	11.82	149	1878755	42.141	ng	100
75) Fluoranthene	12.47	202	1714289	39.050	ng	99
77) Benzidine	12.59	184	654987	28.836	ng	99
78) Pyrene	12.70	202	1685282	37.578	ng	100
80) Butylbenzylphthalate	13.31	149	819233	43.403	ng	99
81) Benzo(a)anthracene	13.89	228	1491468	39.193	ng	99
82) 3,3'-Dichlorobenzidine	13.84	252	562239	38.049	ng	99
83) Chrysene	13.93	228	1442035	38.030	ng	99
84) Bis(2-ethylhexyl)phthalate	13.87	149	1126701	47.249	ng	99
85) Di-n-octyl phthalate	14.47	149	1750555	46.075	ng	100
86) Indeno(1,2,3-cd)pyrene	16.73	276	1139071	31.024	ng	99
88) Benzo(b)fluoranthene	14.92	252	1357700	41.303	ng	99
89) Benzo(k)fluoranthene	14.94	252	1255319	41.208	ng	99
90) Benzo(a)pyrene	15.26	252	1147275	38.671	ng	99
91) Dibenzo(a,h)anthracene	16.73	278	950265	36.403	ng	99
92) Benzo(g,h,i)perylene	17.14	276	879405	34.096	ng	99

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

