

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052920\
 Data File : BF120432.D
 Acq On : 29 May 2020 12:06
 Operator : MA/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad
 6/1/2020 12:20:53 PM

Quant Time: May 29 12:55:41 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 29 12:39:43 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.84	152	298567	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	1143206	20.00	ng	0.00
39) Acenaphthene-d10	9.87	164	589623	20.00	ng	0.00
64) Phenanthrene-d10	11.36	188	1072857	20.00	ng	0.00
76) Chrysene-d12	14.00	240	742812	20.00	ng	0.00
86) Perylene-d12	15.45	264	711224	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.45	112	1235047	78.83	ng	0.00
7) Phenol-d6	6.47	99	1559695	80.77	ng	0.00
23) Nitrobenzene-d5	7.40	82	1374250	78.63	ng	0.00
42) 2,4,6-Tribromophenol	10.66	330	455509	83.38	ng	0.00
45) 2-Fluorobiphenyl	9.20	172	2609801	75.10	ng	0.00
79) Terphenyl-d14	12.94	244	2678945	80.25	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.61	88	303441	39.100	ng	99
3) Pyridine	3.36	79	841857	39.168	ng	96
4) n-Nitrosodimethylamine	3.31	42	358498	39.414	ng	100
6) Aniline	6.50	93	1095252	40.678	ng	99
8) 2-Chlorophenol	6.63	128	718858	40.693	ng	99
9) Benzaldehyde	6.39	77	451169	38.335	ng	99
10) Phenol	6.48	94	869024m	41.153	ng	
11) bis(2-Chloroethyl)ether	6.57	93	725148	41.094	ng	97
12) 1,3-Dichlorobenzene	6.78	146	805412	41.500	ng	98
13) 1,4-Dichlorobenzene	6.86	146	800186	41.325	ng	99
14) 1,2-Dichlorobenzene	7.01	146	742308	41.505	ng	98
15) Benzyl Alcohol	6.98	79	616770	42.371	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.12	45	1147261	41.240	ng	100
17) 2-Methylphenol	7.09	107	642306	40.661	ng	99
18) Hexachloroethane	7.35	117	298431	42.105	ng	97
19) n-Nitroso-di-n-propylamine	7.26	70	498244	40.590	ng	100
20) 3+4-Methylphenols	7.25	107	724262	36.293	ng	98
22) Acetophenone	7.25	105	923485	39.775	ng	99
24) Nitrobenzene	7.42	77	737294	39.163	ng	98
25) Isophorone	7.66	82	1412709	40.219	ng	99
26) 2-Nitrophenol	7.74	139	347201	40.904	ng	98
27) 2,4-Dimethylphenol	7.77	122	594601	40.338	ng	99
28) bis(2-Chloroethoxy)methane	7.87	93	873254	41.282	ng	99
29) 2,4-Dichlorophenol	7.97	162	617830	41.496	ng	99
30) 1,2,4-Trichlorobenzene	8.06	180	696847	41.706	ng	98
31) Naphthalene	8.15	128	2034836	40.529	ng	99
32) Benzoic acid	7.89	122	433694	39.648	ng	95
33) 4-Chloroaniline	8.19	127	904447	40.646	ng	98
34) Hexachlorobutadiene	8.26	225	422394	42.649	ng	99
35) Caprolactam	8.57	113	187737	38.935	ng	98
36) 4-Chloro-3-methylphenol	8.67	107	619111	40.986	ng	95
37) 2-Methylnaphthalene	8.83	142	1381566	41.544	ng	100
38) 1-Methylnaphthalene	8.93	142	1298081	41.353	ng	99
40) 1,2,4,5-Tetrachlorobenzene	9.00	216	637495	42.161	ng	100

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41) Hexachlorocyclopentadiene	8.99	237	404147	47.891	nq	98
43) 2,4,6-Trichlorophenol	9.11	196	470596	43.033	nq	99
44) 2,4,5-Trichlorophenol	9.15	196	510075	41.148	nq	97
46) 1,1'-Biphenyl	9.30	154	1614972	41.695	nq	99
47) 2-Chloronaphthalene	9.32	162	1321352	41.428	nq	98
48) 2-Nitroaniline	9.42	65	377789	38.889	nq	98
49) Acenaphthylene	9.74	152	2008946	40.760	nq	100
50) Dimethylphthalate	9.60	163	1483891	40.256	nq	99
51) 2,6-Dinitrotoluene	9.66	165	318849	38.923	nq #	85
52) Acenaphthene	9.91	154	1252954	40.732	nq	100
53) 3-Nitroaniline	9.83	138	377553	39.248	nq	99
54) 2,4-Dinitrophenol	9.93	184	127395	42.553	nq	94
55) Dibenzofuran	10.09	168	1815906	40.623	nq	97
56) 4-Nitrophenol	9.98	139	287752	39.072	nq	99
57) 2,4-Dinitrotoluene	10.06	165	401881	38.332	nq	92
58) Fluorene	10.43	166	1282242	38.224	nq	99
59) 2,3,4,6-Tetrachlorophenol	10.20	232	401843	41.601	nq	99
60) Diethylphthalate	10.30	149	1502525	40.336	nq	100
61) 4-Chlorophenyl-phenylether	10.42	204	678565	37.310	nq	99
62) 4-Nitroaniline	10.45	138	356125	39.214	nq	95
63) Azobenzene	10.58	77	1395678	39.834	nq	99
65) 4,6-Dinitro-2-methylphenol	10.47	198	184575	47.167	nq	92
66) n-Nitrosodiphenylamine	10.53	169	1251546	41.995	nq	98
67) 4-Bromophenyl-phenylether	10.91	248	463296	42.781	nq	98
68) Hexachlorobenzene	10.97	284	492295	42.686	nq #	90
69) Atrazine	11.06	200	355601	42.334	nq	98
70) Pentachlorophenol	11.16	266	304917	43.242	nq	98
71) Phenanthrene	11.39	178	1919650	40.608	nq	100
72) Anthracene	11.44	178	1958568	41.293	nq	99
73) Carbazole	11.59	167	1863905	40.629	nq	99
74) Di-n-butylphthalate	11.92	149	2211851	41.446	nq	98
75) Fluoranthene	12.57	202	2071398	40.722	nq	97
77) Benzidine	12.69	184	880259	45.102	nq	98
78) Pyrene	12.80	202	2112964	40.903	nq	99
80) Butylbenzylphthalate	13.42	149	935582	43.382	nq	99
81) Benzo(a)anthracene	13.99	228	1619027	40.668	nq	99
82) 3,3'-Dichlorobenzidine	13.95	252	649744	44.955	nq	99
83) Chrysene	14.03	228	1706842	40.896	nq	100
84) Bis(2-ethylhexyl)phthalate	13.98	149	1150680	45.453	nq	99
85) Di-n-octyl phthalate	14.59	149	2081125	47.509	nq	100
87) Indeno(1,2,3-cd)pyrene	16.89	276	1544759	39.711	nq	99
88) Benzo(b)fluoranthene	15.03	252	1690227	42.787	nq	99
89) Benzo(k)fluoranthene	15.06	252	1395032	37.828	nq	99
90) Benzo(a)pyrene	15.39	252	1417047	41.010	nq	99
91) Dibenzo(a,h)anthracene	16.91	278	1259642	39.718	nq	99
92) Benzo(g,h,i)perylene	17.33	276	1231458	39.369	nq	98

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

