

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011720\
 Data File : BF118555.D
 Acq On : 17 Jan 2020 12:02
 Operator : JU
 Sample : SSTDICC050
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC050

Manual Integrations
 APPROVED

mohammad
 1/21/2020 7:58:20 AM

Quant Time: Jan 17 13:33:27 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF011720.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 17 12:09:51 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	432815	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	1569865	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	889736	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	1557145	20.00	ng	0.00
76) Chrysene-d12	14.05	240	1026640	20.00	ng	0.00
87) Perylene-d12	15.52	264	872902	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.51	112	2427917	96.03	ng	0.00
7) Phenol-d6	6.52	99	3165347	104.06	ng	0.00
23) Nitrobenzene-d5	7.46	82	2890924	112.77	ng	0.00
42) 2,4,6-Tribromophenol	10.72	330	1073486	99.47	ng	0.00
45) 2-Fluorobiphenyl	9.26	172	5033538	91.26	ng	0.00
79) Terphenyl-d14	13.00	244	5360239	86.53	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.68	88	578756	56.273	ng	99
3) Pyridine	3.45	79	1586863	58.464	ng	99
4) n-Nitrosodimethylamine	3.39	42	739832	78.549	ng	99
6) Aniline	6.56	93	2195599	57.207	ng	98
8) 2-Chlorophenol	6.68	128	1406074	49.613	ng	97
9) Benzaldehyde	6.44	77	749372	52.897	ng	97
10) Phenol	6.53	94	1815169	53.070	ng	97
11) bis(2-Chloroethyl)ether	6.63	93	1403608	54.583	ng	98
12) 1,3-Dichlorobenzene	6.83	146	1652110	50.296	ng	100
13) 1,4-Dichlorobenzene	6.91	146	1653820	50.247	ng	99
14) 1,2-Dichlorobenzene	7.07	146	1559268	51.135	ng	97
15) Benzyl Alcohol	7.03	79	1330364	61.362	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.17	45	2066836	77.621	ng	100
17) 2-Methylphenol	7.14	107	1184784	53.576	ng	99
18) Hexachloroethane	7.41	117	615018	52.659	ng	98
19) n-Nitroso-di-n-propylamine	7.32	70	1159897	65.915	ng	98
20) 3+4-Methylphenols	7.30	107	1421767	49.384	ng	90
22) Acetophenone	7.30	105	1992365	54.706	ng	96
24) Nitrobenzene	7.47	77	1512442	59.355	ng	98
25) Isophorone	7.72	82	2865563	62.031	ng	99
26) 2-Nitrophenol	7.79	139	628725	42.052	ng	98
27) 2,4-Dimethylphenol	7.82	122	1164149	58.331	ng	99
28) bis(2-Chloroethoxy)methane	7.92	93	1867788	58.652	ng	100
29) 2,4-Dichlorophenol	8.03	162	1293726	55.725	ng	99
30) 1,2,4-Trichlorobenzene	8.12	180	1493814	56.530	ng	100
31) Naphthalene	8.20	128	3852791	48.921	ng	98
32) Benzoic acid	7.96	122	791771	54.942	ng	99
33) 4-Chloroaniline	8.25	127	1769674	52.809	ng	98
34) Hexachlorobutadiene	8.32	225	928375	59.800	ng	98
35) Caprolactam	8.63	113	390016m	54.512	ng	
36) 4-Chloro-3-methylphenol	8.72	107	1258896	52.539	ng	97
37) 2-Methylnaphthalene	8.89	142	2730237	51.859	ng	99
38) 1-Methylnaphthalene	8.99	142	2605556	52.194	ng	99
40) 1,2,4,5-Tetrachlorobenzene	9.06	216	1510066	55.347	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	9.05	237	819089	198.644	nq	98
43) 2,4,6-Trichlorophenol	9.16	196	1009662	57.147	nq	100
44) 2,4,5-Trichlorophenol	9.20	196	992992	56.279	nq	97
46) 1,1'-Biphenyl	9.36	154	3264910	46.311	nq	97
47) 2-Chloronaphthalene	9.38	162	2794947	49.378	nq	96
48) 2-Nitroaniline	9.47	65	804721	52.936	nq	99
49) Acenaphthylene	9.80	152	3887290	47.131	nq	99
50) Dimethylphthalate	9.66	163	3147720	47.932	nq	99
51) 2,6-Dinitrotoluene	9.71	165	683573	47.341	nq	91
52) Acenaphthene	9.97	154	2593995	51.118	nq	98
53) 3-Nitroaniline	9.88	138	743021	44.488	nq	98
54) 2,4-Dinitrophenol	9.98	184	224674	41.661	nq	# 5
55) Dibenzofuran	10.14	168	3614170	49.084	nq	95
56) 4-Nitrophenol	10.03	139	539295	70.447	nq	97
57) 2,4-Dinitrotoluene	10.12	165	821623	44.516	nq	93
58) Fluorene	10.48	166	2779518	46.470	nq	99
59) 2,3,4,6-Tetrachlorophenol	10.25	232	873375	60.530	nq	99
60) Diethylphthalate	10.36	149	3082461	47.107	nq	98
61) 4-Chlorophenyl-phenylether	10.47	204	1537796	51.103	nq	97
62) 4-Nitroaniline	10.50	138	729037	44.204	nq	99
63) Azobenzene	10.63	77	2979748	54.978	nq	99
65) 4,6-Dinitro-2-methylphenol	10.52	198	323903	39.421	nq	81
66) n-Nitrosodiphenylamine	10.59	169	2607987	51.012	nq	99
67) 4-Bromophenyl-phenylether	10.96	248	1106868	58.967	nq	97
68) Hexachlorobenzene	11.03	284	1221563	54.886	nq	96
69) Atrazine	11.12	200	670473	54.687	nq	98
70) Pentachlorophenol	11.22	266	641645	104.463	nq	99
71) Phenanthrene	11.44	178	3968143	47.339	nq	99
72) Anthracene	11.49	178	4032900	48.272	nq	99
73) Carbazole	11.64	167	3591870	48.124	nq	99
74) Di-n-butylphthalate	11.98	149	4244945	43.973	nq	100
75) Fluoranthene	12.63	202	4087828	48.818	nq	97
77) Benzidine	12.74	184	1609765	70.250	nq	98
78) Pyrene	12.86	202	4110796	47.387	nq	98
80) Butylbenzylphthalate	13.47	149	1807546	46.548	nq	98
81) Benzo(a)anthracene	14.04	228	3391132	47.653	nq	98
82) 3,3'-Dichlorobenzidine	14.00	252	1212165	50.168	nq	98
83) Chrysene	14.08	228	3241119	47.519	nq	99
84) Bis(2-ethylhexyl)phthalate	14.03	149	2202811	42.459	nq	100
85) Di-n-octyl phthalate	14.64	149	3358177	43.261	nq	100
86) Indeno(1,2,3-cd)pyrene	17.01	276	3056311	47.685	nq	100
88) Benzo(b)fluoranthene	15.09	252	3085484	55.154	nq	100
89) Benzo(k)fluoranthene	15.13	252	2722965	53.734	nq	99
90) Benzo(a)pyrene	15.46	252	2625946	52.675	nq	99
91) Dibenzo(a,h)anthracene	17.03	278	2513086	55.518	nq	99
92) Benzo(g,h,i)perylene	17.46	276	2412422	54.782	nq	99

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

