

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF010820\
 Data File : BF118502.D
 Acq On : 8 Jan 2020 16:27
 Operator : JU
 Sample : SSTDICV040
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ICVBF010820

Quant Time: Jan 08 17:09:08 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF010820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 08 16:06:11 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	79029	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	319652	20.00	ng	0.00
39) Acenaphthene-d10	9.88	164	174210	20.00	ng	0.00
64) Phenanthrene-d10	11.37	188	323202	20.00	ng	0.00
76) Chrysene-d12	14.03	240	260260	20.00	ng	0.00
87) Perylene-d12	15.52	264	203705	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.45	112	369331	77.15	ng	0.00
7) Phenol-d6	6.47	99	440522	76.96	ng	0.00
23) Nitrobenzene-d5	7.40	82	458497	75.35	ng	0.00
42) 2,4,6-Tribromophenol	10.68	330	169522	82.47	ng	0.00
45) 2-Fluorobiphenyl	9.20	172	922643	75.96	ng	0.00
79) Terphenyl-d14	12.97	244	1256767	76.52	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.53	88	67748	42.447	ng	98
3) Pyridine	3.29	79	182104	42.386	ng	97
4) n-Nitrosodimethylamine	3.25	42	86898	40.363	ng	95
6) Aniline	6.50	93	278222	39.095	ng	98
8) 2-Chlorophenol	6.62	128	208698	39.871	ng	100
9) Benzaldehyde	6.39	77	112557	35.323	ng	98
10) Phenol	6.49	94	251582	39.667	ng	99
11) bis(2-Chloroethyl)ether	6.57	93	181349	40.106	ng	100
12) 1,3-Dichlorobenzene	6.77	146	244839	40.013	ng	99
13) 1,4-Dichlorobenzene	6.85	146	253036	40.528	ng	98
14) 1,2-Dichlorobenzene	7.00	146	239808	39.747	ng	100
15) Benzyl Alcohol	6.98	79	183223	39.463	ng	100
16) 2,2'-oxybis(1-Chloropropan	7.11	45	202154	40.302	ng	98
17) 2-Methylphenol	7.09	107	168292	39.122	ng	98
18) Hexachloroethane	7.35	117	92814	39.234	ng	# 89
19) n-Nitroso-di-n-propylamine	7.25	70	146680	39.788	ng	98
20) 3+4-Methylphenols	7.25	107	225081	38.573	ng	98
22) Acetophenone	7.25	105	319975	39.491	ng	98
24) Nitrobenzene	7.42	77	226040	37.358	ng	98
25) Isophorone	7.66	82	383483	39.339	ng	100
26) 2-Nitrophenol	7.74	139	118693	39.731	ng	97
27) 2,4-Dimethylphenol	7.77	122	156522	38.960	ng	97
28) bis(2-Chloroethoxy)methane	7.87	93	245092	38.242	ng	99
29) 2,4-Dichlorophenol	7.99	162	180240	38.424	ng	98
30) 1,2,4-Trichlorobenzene	8.06	180	209570	38.466	ng	99
31) Naphthalene	8.15	128	645672	38.616	ng	99
32) Benzoic acid	7.92	122	84457	32.886	ng	96
33) 4-Chloroaniline	8.20	127	270209	39.254	ng	99
34) Hexachlorobutadiene	8.26	225	128069	38.295	ng	99
35) Caprolactam	8.57	113	54634	38.665	ng	90
36) 4-Chloro-3-methylphenol	8.69	107	188579	37.395	ng	98
37) 2-Methylnaphthalene	8.84	142	431670	37.617	ng	98
38) 1-Methylnaphthalene	8.93	142	406601	37.738	ng	98
40) 1,2,4,5-Tetrachlorobenzene	9.00	216	200955	37.543	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.99	237	32615	32.244	ng	98
43) 2,4,6-Trichlorophenol	9.12	196	123586	37.364	ng	96
44) 2,4,5-Trichlorophenol	9.17	196	126402	36.929	ng	99
46) 1,1'-Biphenyl	9.30	154	543506	38.183	ng	98
47) 2-Chloronaphthalene	9.33	162	422021	37.459	ng	99
48) 2-Nitroaniline	9.43	65	124532	38.252	ng	96
49) Acenaphthylene	9.75	152	648225	38.653	ng	99
50) Dimethylphthalate	9.60	163	503513	37.658	ng	99
51) 2,6-Dinitrotoluene	9.67	165	108517	37.687	ng	89
52) Acenaphthene	9.92	154	392069	38.285	ng	99
53) 3-Nitroaniline	9.85	138	128014	37.954	ng	94
54) 2,4-Dinitrophenol	9.96	184	40883	35.438	ng	99
55) Dibenzofuran	10.09	168	591257	38.439	ng	99
56) 4-Nitrophenol	10.02	139	61108	37.314	ng	96
57) 2,4-Dinitrotoluene	10.09	165	147566	38.275	ng	99
58) Fluorene	10.43	166	475907	38.493	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.22	232	109159	37.712	ng	97
60) Diethylphthalate	10.30	149	509700	37.447	ng	99
61) 4-Chlorophenyl-phenylether	10.42	204	236528	38.289	ng	92
62) 4-Nitroaniline	10.47	138	135286	40.964	ng	96
63) Azobenzene	10.59	77	455506	39.381	ng	97
65) 4,6-Dinitro-2-methylphenol	10.50	198	68568	40.132	ng	87
66) n-Nitrosodiphenylamine	10.55	169	419312	39.839	ng	99
67) 4-Bromophenyl-phenylether	10.92	248	148917	39.337	ng	97
68) Hexachlorobenzene	10.99	284	174837	38.675	ng	# 92
69) Atrazine	11.07	200	112498	35.971	ng	99
70) Pentachlorophenol	11.19	266	49066	35.531	ng	99
71) Phenanthrene	11.40	178	699134	38.615	ng	99
72) Anthracene	11.46	178	711151	39.482	ng	99
73) Carbazole	11.61	167	618222	39.622	ng	99
74) Di-n-butylphthalate	11.93	149	820808	39.432	ng	99
75) Fluoranthene	12.60	202	786638	42.747	ng	98
77) Benzidine	12.72	184	252387	40.869	ng	97
78) Pyrene	12.83	202	800876	37.401	ng	99
80) Butylbenzylphthalate	13.44	149	338311	34.499	ng	98
81) Benzo(a)anthracene	14.02	228	709718	38.959	ng	99
82) 3,3'-Dichlorobenzidine	13.98	252	220189	37.170	ng	99
83) Chrysene	14.06	228	681291	38.486	ng	100
84) Bis(2-ethylhexyl)phthalate	14.00	149	507150	38.089	ng	99
85) Di-n-octyl phthalate	14.61	149	809299	42.375	ng	100
86) Indeno(1,2,3-cd)pyrene	17.03	276	458645	32.182	ng	100
88) Benzo(b)fluoranthene	15.08	252	556656	37.797	ng	99
89) Benzo(k)fluoranthene	15.11	252	508973	36.206	ng	98
90) Benzo(a)pyrene	15.46	252	446262	36.976	ng	97
91) Dibenzo(a,h)anthracene	17.03	278	380314	33.606	ng	99
92) Benzo(g,h,i)perylene	17.47	276	412493	38.468	ng	99

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

