

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122419\
 Data File : BF118305.D
 Acq On : 24 Dec 2019 13:20
 Operator : JU
 Sample : SSTDICC080
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC080

Manual Integrations
 APPROVED

mohammad
 12/26/2019 4:03:22 PM

Quant Time: Dec 24 15:59:48 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF122419.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 24 14:40:48 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	147724	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	557072	20.00	ng	0.00
39) Acenaphthene-d10	9.90	164	295593	20.00	ng	0.00
64) Phenanthrene-d10	11.40	188	542206	20.00	ng	0.00
76) Chrysene-d12	14.05	240	337303	20.00	ng	0.00
87) Perylene-d12	15.53	264	318014	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	1298085	153.91	ng	0.00
7) Phenol-d6	6.50	99	1610800	157.90	ng	0.01
23) Nitrobenzene-d5	7.43	82	1527088	154.22	ng	0.01
42) 2,4,6-Tribromophenol	10.70	330	566768	157.83	ng	0.01
45) 2-Fluorobiphenyl	9.23	172	2864996	147.48	ng	0.00
79) Terphenyl-d14	12.99	244	3146374	160.41	ng	0.00

Target Compounds					Qvalue
2) 1,4-Dioxane	2.58	88	264653	74.417 ng	99
3) Pyridine	3.35	79	727783	79.319 ng	97
4) n-Nitrosodimethylamine	3.32	42	297864	75.463 ng	98
6) Aniline	6.52	93	1031024	79.009 ng	96
8) 2-Chlorophenol	6.65	128	735499	77.336 ng	98
10) Phenol	6.51	94	911714	78.365 ng	82
11) bis(2-Chloroethyl)ether	6.60	93	649711	77.215 ng	98
12) 1,3-Dichlorobenzene	6.80	146	841707	75.762 ng	98
13) 1,4-Dichlorobenzene	6.87	146	847749	75.779 ng	98
14) 1,2-Dichlorobenzene	7.03	146	799705	76.120 ng	99
15) Benzyl Alcohol	7.00	79	617644	77.561 ng	99
16) 2,2'-oxybis(1-Chloropropan	7.13	45	688619	68.065 ng	97
17) 2-Methylphenol	7.12	107	604001	79.851 ng	98
18) Hexachloroethane	7.37	117	320029	77.636 ng	95
19) n-Nitroso-di-n-propylamine	7.28	70	501546	80.907 ng	97
20) 3+4-Methylphenols	7.27	107	741950	78.091 ng	93
22) Acetophenone	7.27	105	1041786	77.841 ng	# 92
24) Nitrobenzene	7.45	77	757208	77.832 ng	95
25) Isophorone	7.69	82	1328302	79.131 ng	98
26) 2-Nitrophenol	7.76	139	417687	80.326 ng	98
27) 2,4-Dimethylphenol	7.80	122	550422	79.356 ng	100
28) bis(2-Chloroethoxy)methane	7.89	93	853781	77.838 ng	100
29) 2,4-Dichlorophenol	8.00	162	633273	78.318 ng	99
30) 1,2,4-Trichlorobenzene	8.09	180	728991	77.145 ng	99
31) Naphthalene	8.17	128	2138516	76.437 ng	100
32) Benzoic acid	7.97	122	507897	81.100 ng	99
33) 4-Chloroaniline	8.22	127	931377	77.099 ng	99
34) Hexachlorobutadiene	8.28	225	436076	76.957 ng	99
35) Caprolactam	8.62	113	198329m	79.535 ng	
36) 4-Chloro-3-methylphenol	8.71	107	673734	79.177 ng	100
37) 2-Methylnaphthalene	8.86	142	1471384	77.683 ng	98
38) 1-Methylnaphthalene	8.96	142	1390531	78.188 ng	99
40) 1,2,4,5-Tetrachlorobenzene	9.03	216	687484	76.774 ng	100
41) Hexachlorocyclopentadiene	9.01	237	304725	75.952 ng	99

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122419\
 Data File : BF118305.D
 Acq On : 24 Dec 2019 13:20
 Operator : JU
 Sample : SSTDICC080
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC080

Manual Integrations
 APPROVED

mohammad
 12/26/2019 4:03:22 PM

Quant Time: Dec 24 15:59:48 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF122419.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 24 14:40:48 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) 2,4,6-Trichlorophenol	9.14	196	472266	78.990	nq	100
44) 2,4,5-Trichlorophenol	9.19	196	483190	78.006	nq	97
46) 1,1'-Biphenyl	9.33	154	1781765	76.428	nq	99
47) 2-Chloronaphthalene	9.35	162	1446379	77.127	nq	98
48) 2-Nitroaniline	9.45	65	422924	79.082	nq	98
49) Acenaphthylene	9.77	152	2107271	76.186	nq	98
50) Dimethylphthalate	9.63	163	1721646	78.853	nq	99
51) 2,6-Dinitrotoluene	9.70	165	371465	78.242	nq	95
52) Acenaphthene	9.95	154	1304380	77.262	nq	99
53) 3-Nitroaniline	9.87	138	439923	78.737	nq	90
54) 2,4-Dinitrophenol	9.98	184	198840	84.183	nq	94
55) Dibenzofuran	10.12	168	1857511	75.088	nq	96
56) 4-Nitrophenol	10.03	139	290266	74.860	nq	97
57) 2,4-Dinitrotoluene	10.10	165	496461	78.346	nq	# 84
58) Fluorene	10.46	166	1539807	78.326	nq	99
59) 2,3,4,6-Tetrachlorophenol	10.23	232	404681	79.188	nq	97
60) Diethylphthalate	10.33	149	1674855	77.858	nq	99
61) 4-Chlorophenyl-phenylether	10.45	204	768856	77.729	nq	99
62) 4-Nitroaniline	10.49	138	453905	77.894	nq	99
63) Azobenzene	10.61	77	1418905	77.266	nq	99
65) 4,6-Dinitro-2-methylphenol	10.52	198	253440	84.639	nq	98
66) n-Nitrosodiphenylamine	10.57	169	1351238	79.803	nq	98
67) 4-Bromophenyl-phenylether	10.94	248	493623	79.753	nq	97
68) Hexachlorobenzene	11.01	284	577324	79.248	nq	99
69) Atrazine	11.10	200	394535	91.887	nq	97
70) Pentachlorophenol	11.20	266	275881	80.527	nq	97
71) Phenanthrene	11.43	178	2214862	76.843	nq	98
72) Anthracene	11.48	178	2220350	77.221	nq	99
73) Carbazole	11.63	167	1960669	76.001	nq	100
74) Di-n-butylphthalate	11.96	149	2489173	77.097	nq	99
75) Fluoranthene	12.62	202	2193602	74.438	nq	99
78) Pyrene	12.85	202	2185687	82.227	nq	98
80) Butylbenzylphthalate	13.46	149	1005665	83.540	nq	95
81) Benzo(a)anthracene	14.04	228	1821366	78.092	nq	99
82) 3,3'-Dichlorobenzidine	14.00	252	570021	74.416	nq	99
83) Chrysene	14.08	228	1718864	77.151	nq	99
84) Bis(2-ethylhexyl)phthalate	14.02	149	1277632	80.488	nq	100
85) Di-n-octyl phthalate	14.63	149	1849683	76.924	nq	99
86) Indeno(1,2,3-cd)pyrene	17.05	276	1643302	87.662	nq	100
88) Benzo(b)fluoranthene	15.10	252	1567852	74.545	nq	99
89) Benzo(k)fluoranthene	15.13	252	1539147	76.176	nq	99
90) Benzo(a)pyrene	15.48	252	1432874	77.137	nq	98
91) Dibenzo(a,h)anthracene	17.06	278	1361375	85.447	nq	98
92) Benzo(g,h,i)perylene	17.50	276	1338056	87.397	nq	99

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

