

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF123019\
 Data File : BF118394.D
 Acq On : 30 Dec 2019 21:24
 Operator : JU
 Sample : K6456-03MS
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-04MS

Manual Integrations
 APPROVED

mohammad
 12/31/2019 12:04:35 PM

Quant Time: Dec 31 07:35:57 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 15:58:52 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	142180	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	560312	20.00	ng	0.00
39) Acenaphthene-d10	9.90	164	285880	20.00	ng	0.00
64) Phenanthrene-d10	11.40	188	469483	20.00	ng	0.00
76) Chrysene-d12	14.06	240	286975	20.00	ng	0.00
87) Perylene-d12	15.55	264	274845	20.00	ng	0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.48	112	857366	102.93	ng	0.02
7) Phenol-d6	6.50	99	1100697	106.53	ng	0.01
23) Nitrobenzene-d5	7.42	82	660442	67.42	ng	0.00
42) 2,4,6-Tribromophenol	10.70	330	333391	94.43	ng	0.01
45) 2-Fluorobiphenyl	9.22	172	1374509	73.39	ng	0.00
79) Terphenyl-d14	12.99	244	1260091	72.85	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.65	88	121750	36.770	ng	98
3) Pyridine	3.42	79	289198	32.782	ng	99
4) n-Nitrosodimethylamine	3.37	42	146578	40.187	ng	98
6) Aniline	6.52	93	295644	22.533	ng	93
8) 2-Chlorophenol	6.65	128	408084	43.259	ng	98
9) Benzaldehyde	6.40	77	219256	36.193	ng	99
10) Phenol	6.51	94	484932	41.756	ng	92
11) bis(2-Chloroethyl)ether	6.59	93	352661	42.460	ng	97
12) 1,3-Dichlorobenzene	6.79	146	442793	40.756	ng	99
13) 1,4-Dichlorobenzene	6.87	146	451118	40.880	ng	98
14) 1,2-Dichlorobenzene	7.03	146	422120	40.516	ng	97
15) Benzyl Alcohol	7.00	79	325639	41.392	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.13	45	366280	41.328	ng	64
17) 2-Methylphenol	7.12	107	319769	42.263	ng	97
18) Hexachloroethane	7.36	117	151012	37.286	ng	97
19) n-Nitroso-di-n-propylamine	7.27	70	261664	41.108	ng	98
20) 3+4-Methylphenols	7.27	107	417574	43.121	ng	# 81
22) Acetophenone	7.27	105	552238	40.214	ng	98
24) Nitrobenzene	7.45	77	400272	40.788	ng	96
25) Isophorone	7.68	82	715152	41.887	ng	99
26) 2-Nitrophenol	7.76	139	218914	42.116	ng	95
27) 2,4-Dimethylphenol	7.80	122	343148	48.386	ng	99
28) bis(2-Chloroethoxy)methane	7.89	93	459165	41.122	ng	99
29) 2,4-Dichlorophenol	8.01	162	344818	42.243	ng	97
30) 1,2,4-Trichlorobenzene	8.09	180	376995	39.647	ng	99
31) Naphthalene	8.16	128	1175842	41.288	ng	99
32) Benzoic acid	7.95	122	195862	34.514	ng	95
33) 4-Chloroaniline	8.22	127	114584	9.385	ng	98
34) Hexachlorobutadiene	8.27	225	223364	39.462	ng	98
35) Caprolactam	8.60	113	107730m	42.273	ng	
36) 4-Chloro-3-methylphenol	8.72	107	359585	41.030	ng	97
37) 2-Methylnaphthalene	8.86	142	816603	41.881	ng	99
38) 1-Methylnaphthalene	8.96	142	758684	41.300	ng	100
40) 1,2,4,5-Tetrachlorobenzene	9.03	216	354636	41.287	ng	100

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF123019\
 Data File : BF118394.D
 Acq On : 30 Dec 2019 21:24
 Operator : JU
 Sample : K6456-03MS
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-04MS

Manual Integrations
 APPROVED

mohammad
 12/31/2019 12:04:35 PM

Quant Time: Dec 31 07:35:57 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 15:58:52 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	9.00	237	63180	24.534	nq	98
43) 2,4,6-Trichlorophenol	9.15	196	242095	42.189	nq	98
44) 2,4,5-Trichlorophenol	9.19	196	258392	43.772	nq	99
46) 1,1'-Biphenyl	9.32	154	957846	42.361	nq	99
47) 2-Chloronaphthalene	9.35	162	742992	40.827	nq	99
48) 2-Nitroaniline	9.45	65	203316	39.554	nq	99
49) Acenaphthylene	9.77	152	1137065	42.129	nq	100
50) Dimethylphthalate	9.62	163	1004848	46.826	nq	99
51) 2,6-Dinitrotoluene	9.69	165	194670	42.078	nq	100
52) Acenaphthene	9.94	154	668531	40.665	nq	99
53) 3-Nitroaniline	9.86	138	104717	19.285	nq	92
54) 2,4-Dinitrophenol	9.99	184	37811	20.343	nq	95
55) Dibenzofuran	10.11	168	1011462	41.638	nq	97
56) 4-Nitrophenol	10.05	139	254378	75.884	nq	94
57) 2,4-Dinitrotoluene	10.10	165	244851	39.608	nq	# 86
58) Fluorene	10.46	166	783845	39.973	nq	99
59) 2,3,4,6-Tetrachlorophenol	10.24	232	192441	38.368	nq	98
60) Diethylphthalate	10.33	149	910254	42.876	nq	100
61) 4-Chlorophenyl-phenylether	10.45	204	403233	41.192	nq	99
62) 4-Nitroaniline	10.49	138	167211	29.642	nq	93
63) Azobenzene	10.61	77	740588	40.727	nq	97
65) 4,6-Dinitro-2-methylphenol	10.52	198	35933	14.188	nq	98
66) n-Nitrosodiphenylamine	10.57	169	711754	47.303	nq	99
67) 4-Bromophenyl-phenylether	10.94	248	245340	44.653	nq	95
68) Hexachlorobenzene	11.01	284	267249	41.323	nq	98
69) Atrazine	11.10	200	237280	51.095	nq	97
70) Pentachlorophenol	11.21	266	210901	75.299	nq	97
71) Phenanthrene	11.43	178	1041257	40.684	nq	99
72) Anthracene	11.48	178	1071898	41.842	nq	99
73) Carbazole	11.63	167	898585	39.530	nq	99
74) Di-n-butylphthalate	11.96	149	1348908	46.800	nq	100
75) Fluoranthene	12.62	202	931027	36.045	nq	99
77) Benzidine	12.74	184	770081	97.934	nq	98
78) Pyrene	12.85	202	914534	39.025	nq	100
80) Butylbenzylphthalate	13.46	149	466016	44.165	nq	97
81) Benzo(a)anthracene	14.04	228	836381	41.637	nq	98
82) 3,3'-Dichlorobenzidine	14.00	252	330741	48.562	nq	100
83) Chrysene	14.09	228	786518	40.892	nq	100
84) Bis(2-ethylhexyl)phthalate	14.02	149	704516	50.777	nq	100
85) Di-n-octyl phthalate	14.63	149	1143916	56.049	nq	99
86) Indeno(1,2,3-cd)pyrene	17.07	276	626164	38.207	nq	100
88) Benzo(b)fluoranthene	15.11	252	792804	43.455	nq	99
89) Benzo(k)fluoranthene	15.14	252	718511	41.016	nq	98
90) Benzo(a)pyrene	15.49	252	640797	39.871	nq	99
91) Dibenzo(a,h)anthracene	17.08	278	535863	38.422	nq	97
92) Benzo(g,h,i)perylene	17.53	276	485092	36.006	nq	98

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

