

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112519\
 Data File : BF117981.D
 Acq On : 25 Nov 2019 22:23
 Operator : JU
 Sample : K5999-05MS
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 IDLEWILD-BH-05MS

Manual Integrations
 APPROVED

mohammad
 11/26/2019 12:49:12 PM

Quant Time: Nov 26 05:56:36 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 26 05:38:20 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	167172	20.00	ng	0.00
21) Naphthalene-d8	8.18	136	680615	20.00	ng	0.00
39) Acenaphthene-d10	9.95	164	383470	20.00	ng	0.00
64) Phenanthrene-d10	11.43	188	723978	20.00	ng	0.00
76) Chrysene-d12	14.09	240	473889	20.00	ng	0.00
87) Perylene-d12	15.58	264	421241	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.52	112	819099	86.73	ng	0.02
7) Phenol-d6	6.52	99	1085047	95.25	ng	0.00
23) Nitrobenzene-d5	7.46	82	673493	58.82	ng	0.00
42) 2,4,6-Tribromophenol	10.74	330	368724	90.82	ng	0.00
45) 2-Fluorobiphenyl	9.26	172	1122735	52.53	ng	0.00
79) Terphenyl-d14	13.02	244	1156660	44.61	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.77	88	145372	32.930	ng	96
3) Pyridine	3.52	79	371848	32.111	ng	97
4) n-Nitrosodimethylamine	3.47	42	190906	37.766	ng	97
6) Aniline	6.55	93	417707	27.749	ng	100
8) 2-Chlorophenol	6.68	128	458157	44.708	ng	96
9) Benzaldehyde	6.44	77	208079	25.125	ng	97
10) Phenol	6.53	94	541483m	45.745	ng	
11) bis(2-Chloroethyl)ether	6.63	93	399202	41.996	ng	96
12) 1,3-Dichlorobenzene	6.83	146	496751	41.174	ng	99
13) 1,4-Dichlorobenzene	6.91	146	507791	41.639	ng	99
14) 1,2-Dichlorobenzene	7.06	146	484239	41.519	ng	99
15) Benzyl Alcohol	7.03	79	406162	46.183	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.16	45	505424	34.535	ng	96
17) 2-Methylphenol	7.15	107	384812	44.342	ng	99
18) Hexachloroethane	7.40	117	184115	41.098	ng	92
19) n-Nitroso-di-n-propylamine	7.31	70	308577	43.910	ng	95
20) 3+4-Methylphenols	7.30	107	476502	44.232	ng	# 73
22) Acetophenone	7.30	105	631548	41.301	ng	# 98
24) Nitrobenzene	7.47	77	484396	41.011	ng	99
25) Isophorone	7.72	82	884089	42.130	ng	97
26) 2-Nitrophenol	7.79	139	263995	45.920	ng	97
27) 2,4-Dimethylphenol	7.83	122	443809	49.680	ng	96
28) bis(2-Chloroethoxy)methane	7.92	93	529173	39.738	ng	99
29) 2,4-Dichlorophenol	8.03	162	417482	43.084	ng	99
30) 1,2,4-Trichlorobenzene	8.12	180	449741	40.013	ng	99
31) Naphthalene	8.20	128	1397453	44.043	ng	98
32) Benzoic acid	7.96	122	286799	38.343	ng	99
33) 4-Chloroaniline	8.25	127	192081	13.522	ng	98
34) Hexachlorobutadiene	8.32	225	261851	39.846	ng	99
35) Caprolactam	8.63	113	122023m	39.622	ng	
36) 4-Chloro-3-methylphenol	8.73	107	447111	45.465	ng	94
37) 2-Methylnaphthalene	8.90	142	967171	45.113	ng	98
38) 1-Methylnaphthalene	9.00	142	917496	45.268	ng	100
40) 1,2,4,5-Tetrachlorobenzene	9.06	216	419900	37.705	ng	98

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112519\
 Data File : BF117981.D
 Acq On : 25 Nov 2019 22:23
 Operator : JU
 Sample : K5999-05MS
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 IDLEWILD-BH-05MS

Manual Integrations
 APPROVED

mohammad
 11/26/2019 12:49:12 PM

Quant Time: Nov 26 05:56:36 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 26 05:38:20 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	9.05	237	542901	86.992	nq	99
43) 2,4,6-Trichlorophenol	9.17	196	320424	40.171	nq	99
44) 2,4,5-Trichlorophenol	9.22	196	325410	39.292	nq	99
46) 1,1'-Biphenyl	9.36	154	1182280	41.557	nq	98
47) 2-Chloronaphthalene	9.39	162	898829	38.404	nq	97
48) 2-Nitroaniline	9.48	65	267409	37.845	nq	97
49) Acenaphthylene	9.80	152	1459895	42.784	nq	100
50) Dimethylphthalate	9.67	163	1308596	48.664	nq	99
51) 2,6-Dinitrotoluene	9.73	165	248028	42.203	nq	92
52) Acenaphthene	9.98	154	859850	39.337	nq	99
53) 3-Nitroaniline	9.89	138	134482	19.123	nq	100
54) 2,4-Dinitrophenol	10.00	184	239580	79.398	nq	# 87
55) Dibenzofuran	10.15	168	1278119	42.225	nq	98
56) 4-Nitrophenol	10.06	139	416633	79.372	nq	94
57) 2,4-Dinitrotoluene	10.13	165	331185	44.300	nq	98
58) Fluorene	10.50	166	983809	41.942	nq	100
59) 2,3,4,6-Tetrachlorophenol	10.27	232	275807	40.531	nq	96
60) Diethylphthalate	10.37	149	1122761	42.826	nq	100
61) 4-Chlorophenyl-phenylether	10.49	204	491914	41.324	nq	98
62) 4-Nitroaniline	10.52	138	255372	36.269	nq	95
63) Azobenzene	10.64	77	989921	41.504	nq	98
65) 4,6-Dinitro-2-methylphenol	10.54	198	166975	49.404	nq	95
66) n-Nitrosodiphenylamine	10.60	169	897964	42.618	nq	100
67) 4-Bromophenyl-phenylether	10.97	248	322522	41.287	nq	92
68) Hexachlorobenzene	11.04	284	363997	39.961	nq	98
69) Atrazine	11.13	200	318274	45.921	nq	98
70) Pentachlorophenol	11.24	266	404828	76.072	nq	98
71) Phenanthrene	11.46	178	1476713	40.570	nq	99
72) Anthracene	11.52	178	1522247	42.180	nq	99
73) Carbazole	11.67	167	1335410	42.345	nq	100
74) Di-n-butylphthalate	11.99	149	1732277	44.117	nq	100
75) Fluoranthene	12.65	202	1506337	45.535	nq	99
77) Benzidine	12.77	184	704367	45.377	nq	99
78) Pyrene	12.89	202	1506287	39.331	nq	99
80) Butylbenzylphthalate	13.50	149	710191	43.176	nq	96
81) Benzo(a)anthracene	14.07	228	1275335	40.726	nq	99
82) 3,3'-Dichlorobenzidine	14.03	252	299781	27.368	nq	99
83) Chrysene	14.12	228	1214128	40.011	nq	99
84) Bis(2-ethylhexyl)phthalate	14.06	149	991004	49.695	nq	98
85) Di-n-octyl phthalate	14.67	149	1583533	54.052	nq	98
86) Indeno(1,2,3-cd)pyrene	17.10	276	977580	37.852	nq	98
88) Benzo(b)fluoranthene	15.14	252	1076537	39.335	nq	99
89) Benzo(k)fluoranthene	15.17	252	1071156	41.539	nq	99
90) Benzo(a)pyrene	15.52	252	932583	38.751	nq	99
91) Dibenzo(a,h)anthracene	17.12	278	816628	39.424	nq	97
92) Benzo(g,h,i)perylene	17.56	276	789076	38.246	nq	98

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

