

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF123020\
 Data File : BF122882.D
 Acq On : 30 Dec 2020 16:58
 Operator : JU/CG
 Sample : L5242-03MSD
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SAMPLE-2(484+26)MSD

Manual Integrations
 APPROVED

mohammad
 12/31/2020 10:45:09 AM

Quant Time: Dec 30 17:44:16 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF120520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 15 18:14:23 2020
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.822	152	117243	20.00	ng	0.00
21) Naphthalene-d8	8.104	136	438792	20.00	ng	0.00
39) Acenaphthene-d10	9.863	164	221427	20.00	ng	0.00
64) Phenanthrene-d10	11.351	188	370595	20.00	ng	0.00
76) Chrysene-d12	13.998	240	294442	20.00	ng	0.01
86) Perylene-d12	15.462	264	341965	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.445	112	783353	105.78	ng	0.01
7) Phenol-d6	6.463	99	944184	96.13	ng	0.00
23) Nitrobenzene-d5	7.386	82	614193	70.13	ng	0.00
42) 2,4,6-Tribromophenol	10.657	330	283809	97.92	ng	0.00
45) 2-Fluorobiphenyl	9.181	172	1065722	65.10	ng	0.00
79) Terphenyl-d14	12.933	244	990088	58.86	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.604	88	134496	38.64	ng	100
3) Pyridine	3.357	79	377911	41.08	ng	98
4) n-Nitrosodimethylamine	3.304	42	156486	42.41	ng	95
6) Aniline	6.487	93	255080	22.06	ng	# 87
8) 2-Chlorophenol	6.610	128	395466	48.45	ng	96
9) Benzaldehyde	6.369	77	93556	15.74	ng	97
10) Phenol	6.481	94	460412	45.24	ng	94
11) bis(2-Chloroethyl)ether	6.557	93	370231	47.62	ng	98
12) 1,3-Dichlorobenzene	6.763	146	446236	46.20	ng	99
13) 1,4-Dichlorobenzene	6.839	146	450319	46.24	ng	99
14) 1,2-Dichlorobenzene	6.992	146	422414	44.93	ng	99
15) Benzyl Alcohol	6.969	79	319349	47.22	ng	100
16) 2,2'-oxybis(1-Chloropr...	7.098	45	453449	40.90	ng	86
17) 2-Methylphenol	7.087	107	305703	47.11	ng	95
18) Hexachloroethane	7.334	117	153432	44.21	ng	97
19) n-Nitroso-di-n-propyla...	7.239	70	256859	45.66	ng	99
20) 3+4-Methylphenols	7.239	107	377372	46.04	ng	97
22) Acetophenone	7.234	105	497694	45.62	ng	# 98
24) Nitrobenzene	7.410	77	391957	44.99	ng	97
25) Isophorone	7.651	82	702199	46.55	ng	98
26) 2-Nitrophenol	7.728	139	213769	50.31	ng	89
27) 2,4-Dimethylphenol	7.769	122	334349	53.68	ng	98
28) bis(2-Chloroethoxy)met...	7.857	93	456505	44.19	ng	99
29) 2,4-Dichlorophenol	7.969	162	331978	45.64	ng	98
30) 1,2,4-Trichlorobenzene	8.051	180	366003	44.41	ng	99
31) Naphthalene	8.128	128	1096179	45.27	ng	100
32) Benzoic acid	7.910	122	220521	44.44	ng	100
33) 4-Chloroaniline	8.181	127	106785	10.55	ng	99
34) Hexachlorobutadiene	8.245	225	219957	43.37	ng	99
35) Caprolactam	8.563	113	93498m	42.68	ng	
36) 4-Chloro-3-methylphenol	8.675	107	326327	45.55	ng	99
37) 2-Methylnaphthalene	8.822	142	729054	45.52	ng	98
38) 1-Methylnaphthalene	8.922	142	670291	44.24	ng	99
40) 1,2,4,5-Tetrachloroben...	8.986	216	350453	47.78	ng	99
41) Hexachlorocyclopentadiene	8.975	237	128336	39.58	ng	97
43) 2,4,6-Trichlorophenol	9.104	196	225501	44.52	ng	97

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF123020\
 Data File : BF122882.D
 Acq On : 30 Dec 2020 16:58
 Operator : JU/CG
 Sample : L5242-03MSD
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SAMPLE-2(484+26)MSD

Manual Integrations
 APPROVED

mohammad
 12/31/2020 10:45:09 AM

Quant Time: Dec 30 17:44:16 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF120520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 15 18:14:23 2020
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.151	196	238158	45.49	ng	97
46) 1,1'-Biphenyl	9.286	154	863769	47.00	ng	99
47) 2-Chloronaphthalene	9.310	162	682094	44.80	ng	98
48) 2-Nitroaniline	9.410	65	181142	40.81	ng	96
49) Acenaphthylene	9.728	152	1059648	47.12	ng	99
50) Dimethylphthalate	9.586	163	883654	49.97	ng	100
51) 2,6-Dinitrotoluene	9.651	165	175602	47.02	ng	97
52) Acenaphthene	9.898	154	616448	42.37	ng	98
53) 3-Nitroaniline	9.816	138	85581	19.83	ng	93
54) 2,4-Dinitrophenol	9.939	184	134327	73.56	ng	# 87
55) Dibenzofuran	10.069	168	922586	45.08	ng	100
56) 4-Nitrophenol	9.998	139	219717	71.40	ng	98
57) 2,4-Dinitrotoluene	10.057	165	222695	44.76	ng	99
58) Fluorene	10.410	166	717642	44.08	ng	97
59) 2,3,4,6-Tetrachlorophenol	10.192	232	200431	43.57	ng	98
60) Diethylphthalate	10.286	149	793398	46.13	ng	99
61) 4-Chlorophenyl-phenyle...	10.404	204	356191	44.44	ng	95
62) 4-Nitroaniline	10.433	138	148878	33.99	ng	97
63) Azobenzene	10.563	77	689210	43.59	ng	96
65) 4,6-Dinitro-2-methylph...	10.469	198	102241	45.24	ng	97
66) n-Nitrosodiphenylamine	10.522	169	631012	48.19	ng	99
67) 4-Bromophenyl-phenylether	10.892	248	235777	46.95	ng	93
68) Hexachlorobenzene	10.963	284	246224	44.35	ng	99
69) Atrazine	11.051	200	175628	41.30	ng	98
70) Pentachlorophenol	11.163	266	246101	83.67	ng	100
71) Phenanthrene	11.374	178	1045906	47.49	ng	98
72) Anthracene	11.427	178	1095079	50.14	ng	99
73) Carbazole	11.580	167	901531	45.52	ng	99
74) Di-n-butylphthalate	11.910	149	1194184	47.98	ng	99
75) Fluoranthene	12.569	202	1513386	65.53	ng	99
77) Benzidine	12.686	184	374728	58.84	ng	99
78) Pyrene	12.798	202	1503101	66.03	ng	100
80) Butylbenzylphthalate	13.410	149	461388	45.00	ng	95
81) Benzo(a)anthracene	13.986	228	1232326	62.63	ng	99
82) 3,3'-Dichlorobenzidine	13.945	252	259897	36.88	ng	97
83) Chrysene	14.027	228	1300631	66.42	ng	99
84) Bis(2-ethylhexyl)phtha...	13.968	149	707450	50.38	ng	99
85) Di-n-octyl phthalate	14.580	149	1198814	51.47	ng	97
87) Indeno(1,2,3-cd)pyrene	16.939	276	1467196	65.36	ng	99
88) Benzo(b)fluoranthene	15.045	252	1726884	71.97	ng	100
89) Benzo(k)fluoranthene	15.074	252	1194086	54.43	ng	98
90) Benzo(a)pyrene	15.404	252	1237086	58.93	ng	99
91) Dibenzo(a,h)anthracene	16.951	278	1059683	57.68	ng	99
92) Benzo(g,h,i)perylene	17.380	276	1248291	70.21	ng	99

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

