

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101320\
 Data File : BF121939.D
 Acq On : 13 Oct 2020 15:39
 Operator : JU/CG
 Sample : L4369-08MS
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HD-4MS

Manual Integrations
 APPROVED

mohammad
 10/14/2020 12:53:26 PM

Quant Time: Oct 13 16:01:39 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 12 16:22:33 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.745	152	114917	20.00	ng	0.00
21) Naphthalene-d8	8.028	136	439974	20.00	ng	0.00
39) Acenaphthene-d10	9.781	164	218647	20.00	ng	0.00
64) Phenanthrene-d10	11.263	188	385085	20.00	ng	-0.01
76) Chrysene-d12	13.904	240	252378	20.00	ng	-0.01
86) Perylene-d12	15.339	264	264904	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.369	112	603496	77.51	ng	0.00
7) Phenol-d6	6.393	99	760861	75.91	ng	0.00
23) Nitrobenzene-d5	7.310	82	517866	60.36	ng	-0.01
42) 2,4,6-Tribromophenol	10.569	330	207710	76.80	ng	-0.01
45) 2-Fluorobiphenyl	9.104	172	876974	59.01	ng	0.00
79) Terphenyl-d14	12.845	244	722541	54.56	ng	-0.02
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.505	88	126133	36.83	ng	100
3) Pyridine	3.228	79	366635	40.72	ng	99
4) n-Nitrosodimethylamine	3.187	42	175033	40.22	ng	98
6) Aniline	6.410	93	255014	20.66	ng	# 43
8) 2-Chlorophenol	6.534	128	318044	40.41	ng	96
9) Benzaldehyde	6.293	77	176151	29.03	ng	99
10) Phenol	6.404	94	405028	39.16	ng	98
11) bis(2-Chloroethyl)ether	6.481	93	318646	39.85	ng	99
12) 1,3-Dichlorobenzene	6.687	146	344040	38.87	ng	97
13) 1,4-Dichlorobenzene	6.763	146	346239	38.96	ng	97
14) 1,2-Dichlorobenzene	6.910	146	325717	40.09	ng	97
15) Benzyl Alcohol	6.898	79	279828	43.16	ng	100
16) 2,2'-oxybis(1-Chloropr...	7.022	45	478502	38.88	ng	76
17) 2-Methylphenol	7.010	107	255999	38.05	ng	95
18) Hexachloroethane	7.251	117	125655	39.14	ng	97
19) n-Nitroso-di-n-propyla...	7.163	70	223152	39.10	ng	100
20) 3+4-Methylphenols	7.169	107	323023	39.81	ng	# 68
22) Acetophenone	7.157	105	427489	37.76	ng	93
24) Nitrobenzene	7.328	77	347629	37.91	ng	99
25) Isophorone	7.569	82	609135	37.55	ng	100
26) 2-Nitrophenol	7.645	139	164506	38.96	ng	92
27) 2,4-Dimethylphenol	7.692	122	273656	38.01	ng	100
28) bis(2-Chloroethoxy)met...	7.781	93	389637	38.39	ng	99
29) 2,4-Dichlorophenol	7.892	162	259548	38.50	ng	99
30) 1,2,4-Trichlorobenzene	7.969	180	269844	37.65	ng	98
31) Naphthalene	8.051	128	890912	37.80	ng	99
32) Benzoic acid	7.845	122	155615	39.45	ng	98
33) 4-Chloroaniline	8.104	127	144240	14.37	ng	99
34) Hexachlorobutadiene	8.163	225	162508	38.23	ng	99
35) Caprolactam	8.481	113	79877m	37.31	ng	
36) 4-Chloro-3-methylphenol	8.598	107	275172	38.01	ng	99
37) 2-Methylnaphthalene	8.739	142	579726	37.98	ng	100
38) 1-Methylnaphthalene	8.839	142	532372	37.66	ng	99
40) 1,2,4,5-Tetrachloroben...	8.910	216	268790	38.09	ng	99
41) Hexachlorocyclopentadiene	8.892	237	101351	56.11	ng	99
43) 2,4,6-Trichlorophenol	9.022	196	178250	40.94	ng	97

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44) 2,4,5-Trichlorophenol	9.069	196	184795	39.30	ng	97
46) 1,1'-Biphenyl	9.204	154	682837	38.41	ng	100
47) 2-Chloronaphthalene	9.228	162	537238	38.86	ng	99
48) 2-Nitroaniline	9.334	65	181082	39.71	ng	96
49) Acenaphthylene	9.645	152	834048	39.28	ng	100
50) Dimethylphthalate	9.504	163	702168	43.92	ng	99
51) 2,6-Dinitrotoluene	9.575	165	137683	39.26	ng	90
52) Acenaphthene	9.816	154	485593	38.45	ng	99
53) 3-Nitroaniline	9.739	138	81412	20.41	ng	98
54) 2,4-Dinitrophenol	9.863	184	57660	37.41	ng	99
55) Dibenzofuran	9.986	168	695757	37.67	ng	99
56) 4-Nitrophenol	9.922	139	167067	77.00	ng	# 81
57) 2,4-Dinitrotoluene	9.981	165	169742	39.32	ng	# 88
58) Fluorene	10.328	166	564743	37.95	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.110	232	146922	40.38	ng	98
60) Diethylphthalate	10.204	149	606152	39.32	ng	100
61) 4-Chlorophenyl-phenyle...	10.322	204	271235	38.00	ng	99
62) 4-Nitroaniline	10.357	138	123055	30.88	ng	98
63) Azobenzene	10.480	77	623255	38.23	ng	98
65) 4,6-Dinitro-2-methylph...	10.392	198	50774	22.29	ng	96
66) n-Nitrosodiphenylamine	10.439	169	524881	39.57	ng	99
67) 4-Bromophenyl-phenylether	10.810	248	173708	39.04	ng	97
68) Hexachlorobenzene	10.875	284	193329	38.39	ng	97
69) Atrazine	10.969	200	137125	42.28	ng	99
70) Pentachlorophenol	11.080	266	143991	73.45	ng	99
71) Phenanthrene	11.292	178	827073	39.17	ng	100
72) Anthracene	11.339	178	835774	38.16	ng	100
73) Carbazole	11.498	167	719509	36.95	ng	100
74) Di-n-butylphthalate	11.827	149	902151	40.27	ng	99
75) Fluoranthene	12.474	202	810881	35.82	ng	100
77) Benzidine	12.604	184	330257	42.45	ng	99
78) Pyrene	12.704	202	807579	41.99	ng	100
80) Butylbenzylphthalate	13.321	149	319484	41.89	ng	99
81) Benzo(a)anthracene	13.898	228	653743	39.53	ng	99
82) 3,3'-Dichlorobenzidine	13.857	252	196881	32.09	ng	99
83) Chrysene	13.933	228	641816	39.53	ng	100
84) Bis(2-ethylhexyl)phtha...	13.880	149	436196	42.13	ng	99
85) Di-n-octyl phthalate	14.492	149	706318	42.17	ng	100
87) Indeno(1,2,3-cd)pyrene	16.762	276	790983	45.99	ng	99
88) Benzo(b)fluoranthene	14.927	252	718228	40.11	ng	99
89) Benzo(k)fluoranthene	14.957	252	597557	35.13	ng	99
90) Benzo(a)pyrene	15.280	252	623707	40.33	ng	98
91) Dibenzo(a,h)anthracene	16.768	278	640403	45.22	ng	99
92) Benzo(g,h,i)perylene	17.186	276	665325	46.94	ng	98

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

