

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102220\
 Data File : BF122083.D
 Acq On : 22 Oct 2020 14:57
 Operator : JU/CG
 Sample : L4436-08MSD
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 36AMSD

Manual Integrations
 APPROVED

mohammad
 10/23/2020 1:32:00 PM

Quant Time: Oct 22 15:29:50 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 21 12:04:42 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.728	152	116887	20.00	ng	0.00
21) Naphthalene-d8	8.016	136	465500	20.00	ng	0.00
39) Acenaphthene-d10	9.763	164	242117	20.00	ng	0.00
64) Phenanthrene-d10	11.251	188	452889	20.00	ng	0.00
76) Chrysene-d12	13.892	240	278549	20.00	ng	0.00
86) Perylene-d12	15.321	264	264137	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.357	112	813281	102.69	ng	0.01
7) Phenol-d6	6.387	99	1003326	98.42	ng	0.00
23) Nitrobenzene-d5	7.304	82	718377	79.14	ng	0.00
42) 2,4,6-Tribromophenol	10.563	330	333948	111.50	ng	0.00
45) 2-Fluorobiphenyl	9.086	172	1189285	72.27	ng	0.00
79) Terphenyl-d14	12.833	244	1266624	86.66	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.440	88	116619	33.48	ng	99
3) Pyridine	3.169	79	345630	37.74	ng	99
4) n-Nitrosodimethylamine	3.128	42	188252	42.53	ng	96
6) Aniline	6.398	93	392415	31.26	ng	# 20
8) 2-Chlorophenol	6.522	128	352341	44.02	ng	98
9) Benzaldehyde	6.281	77	150143	23.08	ng	99
10) Phenol	6.398	94	442982	42.10	ng	81
11) bis(2-Chloroethyl)ether	6.469	93	360841	44.36	ng	99
12) 1,3-Dichlorobenzene	6.669	146	373302	41.46	ng	99
13) 1,4-Dichlorobenzene	6.745	146	378108	41.83	ng	99
14) 1,2-Dichlorobenzene	6.898	146	351600	42.55	ng	98
15) Benzyl Alcohol	6.887	79	313486	47.54	ng	97
16) 2,2'-oxybis(1-Chloropr...	7.010	45	523149	41.80	ng	# 44
17) 2-Methylphenol	7.004	107	281780	41.17	ng	# 89
18) Hexachloroethane	7.234	117	141689	43.39	ng	98
19) n-Nitroso-di-n-propyla...	7.157	70	258832	44.59	ng	98
20) 3+4-Methylphenols	7.157	107	365705	44.31	ng	# 73
22) Acetophenone	7.145	105	488378	40.78	ng	# 89
24) Nitrobenzene	7.322	77	394464	40.66	ng	99
25) Isophorone	7.563	82	718581	41.87	ng	99
26) 2-Nitrophenol	7.639	139	191633	42.89	ng	97
27) 2,4-Dimethylphenol	7.681	122	311236	40.86	ng	98
28) bis(2-Chloroethoxy)met...	7.769	93	442052	41.17	ng	100
29) 2,4-Dichlorophenol	7.887	162	298122	41.80	ng	98
30) 1,2,4-Trichlorobenzene	7.957	180	306690	40.44	ng	98
31) Naphthalene	8.034	128	998453	40.04	ng	99
32) Benzoic acid	7.851	122	177138	41.86	ng	96
33) 4-Chloroaniline	8.092	127	247131	23.27	ng	99
34) Hexachlorobutadiene	8.145	225	188216	41.85	ng	99
35) Caprolactam	8.486	113	99901m	44.11	ng	
36) 4-Chloro-3-methylphenol	8.592	107	331189	43.23	ng	98
37) 2-Methylnaphthalene	8.728	142	649363	40.21	ng	98
38) 1-Methylnaphthalene	8.822	142	601985	40.25	ng	99
40) 1,2,4,5-Tetrachloroben...	8.892	216	314183	40.20	ng	100
41) Hexachlorocyclopentadiene	8.875	237	133643	62.25	ng	99
43) 2,4,6-Trichlorophenol	9.016	196	204207	42.35	ng	100

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.063	196	213598	41.02	ng	100
46) 1,1'-Biphenyl	9.186	154	786690	39.96	ng	99
47) 2-Chloronaphthalene	9.216	162	589411	38.50	ng	100
48) 2-Nitroaniline	9.322	65	220790	43.72	ng	97
49) Acenaphthylene	9.628	152	928228	39.48	ng	100
50) Dimethylphthalate	9.492	163	961609	54.31	ng	99
51) 2,6-Dinitrotoluene	9.569	165	167709	43.18	ng	# 83
52) Acenaphthene	9.798	154	569687	40.73	ng	99
53) 3-Nitroaniline	9.739	138	129352	29.29	ng	94
54) 2,4-Dinitrophenol	9.857	184	151516	76.84	ng	# 83
55) Dibenzofuran	9.975	168	820236	40.10	ng	99
56) 4-Nitrophenol	9.928	139	172282	71.70	ng	88
57) 2,4-Dinitrotoluene	9.975	165	207376	43.38	ng	# 82
58) Fluorene	10.316	166	679275	41.22	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.098	232	178676	44.35	ng	99
60) Diethylphthalate	10.192	149	764883	44.80	ng	99
61) 4-Chlorophenyl-phenyle...	10.304	204	333715	42.23	ng	97
62) 4-Nitroaniline	10.357	138	164231	37.22	ng	91
63) Azobenzene	10.469	77	779207	43.16	ng	97
65) 4,6-Dinitro-2-methylph...	10.392	198	123405	41.34	ng	98
66) n-Nitrosodiphenylamine	10.428	169	644660	41.32	ng	99
67) 4-Bromophenyl-phenylether	10.792	248	225198	43.04	ng	97
68) Hexachlorobenzene	10.863	284	251810	42.52	ng	98
69) Atrazine	10.957	200	184841	48.46	ng	99
70) Pentachlorophenol	11.069	266	183605	78.96	ng	99
71) Phenanthrene	11.275	178	1046052	42.12	ng	99
72) Anthracene	11.327	178	1052128	40.85	ng	100
73) Carbazole	11.486	167	919503	40.15	ng	99
74) Di-n-butylphthalate	11.810	149	1215043	46.12	ng	99
75) Fluoranthene	12.463	202	1094654	41.11	ng	99
77) Benzidine	12.586	184	436332	50.81	ng	99
78) Pyrene	12.692	202	1060245	49.95	ng	99
80) Butylbenzylphthalate	13.304	149	460817	54.74	ng	97
81) Benzo(a)anthracene	13.880	228	806358	44.18	ng	99
82) 3,3'-Dichlorobenzidine	13.839	252	223159	32.96	ng	98
83) Chrysene	13.921	228	779488	43.50	ng	99
84) Bis(2-ethylhexyl)phtha...	13.857	149	641841	56.16	ng	99
85) Di-n-octyl phthalate	14.468	149	972049	52.58	ng	98
87) Indeno(1,2,3-cd)pyrene	16.751	276	904637	52.75	ng	98
88) Benzo(b)fluoranthene	14.915	252	775648	43.44	ng	99
89) Benzo(k)fluoranthene	14.945	252	720965	42.51	ng	100
90) Benzo(a)pyrene	15.268	252	687684	44.59	ng	100
91) Dibenzo(a,h)anthracene	16.745	278	736638	52.17	ng	99
92) Benzo(g,h,i)perylene	17.162	276	779775	55.18	ng	98

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

