

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030221\
 Data File : BF123350.D
 Acq On : 2 Mar 2021 22:34
 Operator : JU/CG
 Sample : M1527-01MSD
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ETGI-371MSD

Manual Integrations
 APPROVED

mohammad
 3/3/2021 11:01:48 AM

Quant Time: Mar 02 23:53:35 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF021921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 19 16:08:53 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.851	152	135562	20.00	ng	0.00
21) Naphthalene-d8	8.139	136	558022	20.00	ng	# 0.00
39) Acenaphthene-d10	9.898	164	287378	20.00	ng	0.00
64) Phenanthrene-d10	11.392	188	524127	20.00	ng	0.00
76) Chrysene-d12	14.045	240	334516	20.00	ng	0.00
86) Perylene-d12	15.515	264	294887	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.487	112	903259	95.91	ng	0.01
7) Phenol-d6	6.498	99	1186461	92.73	ng	0.00
23) Nitrobenzene-d5	7.422	82	810742	66.51	ng	0.00
42) 2,4,6-Tribromophenol	10.692	330	282466	96.53	ng	0.00
45) 2-Fluorobiphenyl	9.216	172	1221472	61.05	ng	0.00
79) Terphenyl-d14	12.980	244	1126036	64.73	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.669	88	162356	29.87	ng	89
3) Pyridine	3.422	79	466318	35.21	ng	90
4) n-Nitrosodimethylamine	3.381	42	239530	39.11	ng	85
6) Aniline	6.516	93	291818	19.31	ng	# 59
8) 2-Chlorophenol	6.640	128	451379	44.61	ng	89
9) Benzaldehyde	6.398	77	156041	20.13	ng	87
10) Phenol	6.510	94	602290	43.19	ng	90
11) bis(2-Chloroethyl)ether	6.587	93	453056	45.30	ng	93
12) 1,3-Dichlorobenzene	6.793	146	461597	43.86	ng	# 92
13) 1,4-Dichlorobenzene	6.869	146	465678	43.41	ng	# 94
14) 1,2-Dichlorobenzene	7.022	146	451287	45.23	ng	94
15) Benzyl Alcohol	6.998	79	456035	45.71	ng	92
16) 2,2'-oxybis(1-Chloropr...	7.128	45	617979	45.05	ng	86
17) 2-Methylphenol	7.116	107	385461	45.29	ng	# 91
18) Hexachloroethane	7.363	117	181982	46.97	ng	90
19) n-Nitroso-di-n-propyla...	7.275	70	357503	46.97	ng	97
20) 3+4-Methylphenols	7.269	107	475264	42.53	ng	92
22) Acetophenone	7.263	105	647989	44.77	ng	# 92
24) Nitrobenzene	7.440	77	549328	42.74	ng	91
25) Isophorone	7.681	82	968444	42.48	ng	97
26) 2-Nitrophenol	7.757	139	239077	44.40	ng	90
27) 2,4-Dimethylphenol	7.792	122	383904	46.06	ng	92
28) bis(2-Chloroethoxy)met...	7.887	93	577698	48.42	ng	99
29) 2,4-Dichlorophenol	7.998	162	370556	43.06	ng	93
30) 1,2,4-Trichlorobenzene	8.081	180	399498	42.70	ng	99
31) Naphthalene	8.163	128	1285288	42.03	ng	99
32) Benzoic acid	7.945	122	240791	39.17	ng	89
33) 4-Chloroaniline	8.210	127	96182	7.73	ng	# 93
34) Hexachlorobutadiene	8.275	225	222704	42.20	ng	99
35) Caprolactam	8.598	113	132844m	46.01	ng	
36) 4-Chloro-3-methylphenol	8.704	107	457435	44.72	ng	90
37) 2-Methylnaphthalene	8.851	142	879969	43.20	ng	100
38) 1-Methylnaphthalene	8.951	142	810374	42.53	ng	100
40) 1,2,4,5-Tetrachloroben...	9.022	216	384634	44.79	ng	# 97
41) Hexachlorocyclopentadiene	9.004	237	213840	53.15	ng	97
43) 2,4,6-Trichlorophenol	9.134	196	259955	42.68	ng	93

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.181	196	276775	42.47	ng	# 87
46) 1,1'-Biphenyl	9.322	154	1060466	44.48	ng	99
47) 2-Chloronaphthalene	9.345	162	808823	43.00	ng	98
48) 2-Nitroaniline	9.445	65	319722	47.44	ng	# 81
49) Acenaphthylene	9.763	152	1231899	43.18	ng	99
50) Dimethylphthalate	9.622	163	1037007	48.25	ng	99
51) 2,6-Dinitrotoluene	9.686	165	218115	44.04	ng	# 77
52) Acenaphthene	9.933	154	886931m	45.08	ng	
53) 3-Nitroaniline	9.857	138	110302	20.07	ng	83
54) 2,4-Dinitrophenol	9.969	184	158181	60.80	ng	88
55) Dibenzofuran	10.104	168	1119880	42.05	ng	93
56) 4-Nitrophenol	10.033	139	346225	79.04	ng	# 61
57) 2,4-Dinitrotoluene	10.092	165	295160	44.32	ng	# 78
58) Fluorene	10.451	166	895124	42.70	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.228	232	227931	43.83	ng	# 84
60) Diethylphthalate	10.322	149	996128	46.81	ng	98
61) 4-Chlorophenyl-phenyle...	10.439	204	435962	43.69	ng	93
62) 4-Nitroaniline	10.475	138	211948	38.30	ng	84
63) Azobenzene	10.604	77	1061192	43.28	ng	98
65) 4,6-Dinitro-2-methylph...	10.504	198	121975	34.57	ng	85
66) n-Nitrosodiphenylamine	10.563	169	794986	43.86	ng	99
67) 4-Bromophenyl-phenylether	10.933	248	257308	45.03	ng	96
68) Hexachlorobenzene	11.004	284	268727	44.65	ng	96
69) Atrazine	11.092	200	235623	41.14	ng	97
70) Pentachlorophenol	11.198	266	298623	78.57	ng	99
71) Phenanthrene	11.416	178	1340150	43.24	ng	99
72) Anthracene	11.469	178	1360380	43.99	ng	100
73) Carbazole	11.627	167	1194657	41.02	ng	100
74) Di-n-butylphthalate	11.951	149	1560505	47.92	ng	99
75) Fluoranthene	12.610	202	1389875	42.31	ng	99
77) Benzidine	12.727	184	385414	35.67	ng	100
78) Pyrene	12.839	202	1350105	51.70	ng	100
80) Butylbenzylphthalate	13.457	149	642937	58.24	ng	93
81) Benzo(a)anthracene	14.033	228	995439	43.79	ng	98
82) 3,3'-Dichlorobenzidine	13.992	252	255818	33.85	ng	# 96
83) Chrysene	14.074	228	968186	43.36	ng	99
84) Bis(2-ethylhexyl)phtha...	14.016	149	1226175	80.68	ng	99
85) Di-n-octyl phthalate	14.633	149	1380273	53.16	ng	96
87) Indeno(1,2,3-cd)pyrene	17.015	276	986874	48.42	ng	97
88) Benzo(b)fluoranthene	15.092	252	929559	44.88	ng	99
89) Benzo(k)fluoranthene	15.121	252	833104	44.37	ng	99
90) Benzo(a)pyrene	15.457	252	806604	43.68	ng	99
91) Dibenzo(a,h)anthracene	17.027	278	817128	46.73	ng	98
92) Benzo(g,h,i)perylene	17.462	276	794331	49.33	ng	98

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

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