

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102621\
 Data File : BF125939.D
 Acq On : 26 Oct 2021 17:55
 Operator : JU/CG
 Sample : M4340-04MS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TH-EDMS

Manual Integrations
 APPROVED

mohammad
 10/27/2021 11:48:57 AM

Quant Time: Oct 27 09:03:02 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF102521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 25 18:49:16 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.869	152	174656	20.000	ng	0.00
21) Naphthalene-d8	8.151	136	721750	20.000	ng	# 0.00
39) Acenaphthene-d10	9.910	164	403298	20.000	ng	0.00
64) Phenanthrene-d10	11.392	188	710391	20.000	ng	0.00
76) Chrysene-d12	14.027	240	344459	20.000	ng	# 0.00
86) Perylene-d12	15.498	264	348783	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.487	112	1133360	96.657	ng	0.00
7) Phenol-d6	6.492	99	1434649	89.205	ng	0.00
23) Nitrobenzene-d5	7.428	82	995212	71.224	ng	0.00
42) 2,4,6-Tribromophenol	10.698	330	358727	102.035	ng	0.00
45) 2-Fluorobiphenyl	9.228	172	1494119	65.516	ng	0.00
79) Terphenyl-d14	12.974	244	1446665	68.698	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.722	88	229943	40.133	ng	# 100
3) Pyridine	3.475	79	524632	34.582	ng	# 83
4) n-Nitrosodimethylamine	3.410	42	419795	43.143	ng	# 72
6) Aniline	6.528	93	770438	40.670	ng	# 90
8) 2-Chlorophenol	6.651	128	566365	47.854	ng	# 76
9) Benzaldehyde	6.416	77	351429	36.518	ng	94
10) Phenol	6.510	94	748317	48.257	ng	91
11) bis(2-Chloroethyl)ether	6.604	93	578172	46.299	ng	# 80
12) 1,3-Dichlorobenzene	6.810	146	568837	43.509	ng	96
13) 1,4-Dichlorobenzene	6.887	146	571377	43.688	ng	95
14) 1,2-Dichlorobenzene	7.039	146	552284	43.565	ng	97
15) Benzyl Alcohol	7.010	79	563778	46.306	ng	96
16) 2,2'-oxybis(1-Chloropr...	7.145	45	1345394	43.992	ng	97
17) 2-Methylphenol	7.116	107	478854	46.842	ng	# 92
18) Hexachloroethane	7.381	117	230127	46.410	ng	97
19) n-Nitroso-di-n-propyla...	7.287	70	421027	45.821	ng	# 91
20) 3+4-Methylphenols	7.269	107	544370	41.657	ng	87
22) Acetophenone	7.281	105	770177	41.331	ng	# 89
24) Nitrobenzene	7.451	77	667164	46.244	ng	88
25) Isophorone	7.692	82	1201860	45.167	ng	# 86
26) 2-Nitrophenol	7.763	139	277913	59.413	ng	# 68
27) 2,4-Dimethylphenol	7.804	122	515291	49.610	ng	90
28) bis(2-Chloroethoxy)met...	7.898	93	720835	43.178	ng	# 96
29) 2,4-Dichlorophenol	8.004	162	459711	45.150	ng	91
30) 1,2,4-Trichlorobenzene	8.092	180	498786	41.767	ng	97
31) Naphthalene	8.175	128	1546545	42.141	ng	99
32) Benzoic acid	7.928	122	389914	51.135	ng	89
33) 4-Chloroaniline	8.216	127	415622	27.482	ng	# 88
34) Hexachlorobutadiene	8.292	225	316740	42.202	ng	100
35) Caprolactam	8.592	113	160688m	48.209	ng	
36) 4-Chloro-3-methylphenol	8.698	107	539816	45.302	ng	88
37) 2-Methylnaphthalene	8.863	142	1065956	45.125	ng	90
38) 1-Methylnaphthalene	8.963	142	992184	43.519	ng	90
40) 1,2,4,5-Tetrachloroben...	9.033	216	453281	40.800	ng	97
41) Hexachlorocyclopentadiene	9.022	237	600443	100.067	ng	99
43) 2,4,6-Trichlorophenol	9.139	196	341955	45.099	ng	92

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44) 2,4,5-Trichlorophenol	9.175	196	351034	43.882	ng	95
46) 1,1'-Biphenyl	9.328	154	1237828	41.905	ng	98
47) 2-Chloronaphthalene	9.351	162	941649	42.281	ng	94
48) 2-Nitroaniline	9.445	65	415965	52.291	ng #	71
49) Acenaphthylene	9.769	152	1451553	43.147	ng	97
50) Dimethylphthalate	9.633	163	1099212	44.000	ng	96
51) 2,6-Dinitrotoluene	9.686	165	262048	53.181	ng #	60
52) Acenaphthene	9.939	154	1046664	47.873	ng	98
53) 3-Nitroaniline	9.857	138	193797	34.308	ng #	78
54) 2,4-Dinitrophenol	9.963	184	236128	111.196	ng	87
55) Dibenzofuran	10.116	168	1356373	41.536	ng	96
56) 4-Nitrophenol	10.016	139	412764	94.033	ng #	63
57) 2,4-Dinitrotoluene	10.092	165	348588	57.595	ng #	90
58) Fluorene	10.457	166	968484	47.310	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.227	232	296646	45.431	ng #	89
60) Diethylphthalate	10.333	149	1176499	46.169	ng	98
61) 4-Chlorophenyl-phenyle...	10.445	204	467133	38.873	ng	94
62) 4-Nitroaniline	10.475	138	255021	49.204	ng	88
63) Azobenzene	10.610	77	1289194	43.864	ng	95
65) 4,6-Dinitro-2-methylph...	10.498	198	166808	52.780	ng	94
66) n-Nitrosodiphenylamine	10.569	169	919632	43.220	ng	96
67) 4-Bromophenyl-phenylether	10.939	248	327792	43.194	ng	92
68) Hexachlorobenzene	11.004	284	356027	44.501	ng #	89
69) Atrazine	11.098	200	315907	46.878	ng	89
70) Pentachlorophenol	11.192	266	391287	86.632	ng	95
71) Phenanthrene	11.416	178	1551795	42.098	ng	97
72) Anthracene	11.469	178	1614039	43.855	ng	99
73) Carbazole	11.622	167	1381967	39.500	ng	99
74) Di-n-butylphthalate	11.957	149	1683173	45.157	ng	99
75) Fluoranthene	12.604	202	1504114	40.477	ng	97
77) Benzidine	12.721	184	657396	61.184	ng	98
78) Pyrene	12.833	202	1516132	47.297	ng	96
80) Butylbenzylphthalate	13.457	149	582375	55.850	ng #	86
81) Benzo(a)anthracene	14.015	228	996113	42.966	ng	99
82) 3,3'-Dichlorobenzidine	13.980	252	296633	40.352	ng	100
83) Chrysene	14.057	228	1065079	45.407	ng	98
84) Bis(2-ethylhexyl)phtha...	14.021	149	672740	59.551	ng #	99
85) Di-n-octyl phthalate	14.633	149	1184377	67.674	ng #	94
87) Indeno(1,2,3-cd)pyrene	16.968	276	1190693	47.387	ng #	91
88) Benzo(b)fluoranthene	15.068	252	1098831	47.730	ng #	96
89) Benzo(k)fluoranthene	15.098	252	922922	41.736	ng	100
90) Benzo(a)pyrene	15.439	252	1005663	54.110	ng #	96
91) Dibenzo(a,h)anthracene	16.986	278	1005249	49.384	ng #	95
92) Benzo(g,h,i)perylene	17.409	276	1024706	48.182	ng #	89

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

