

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082621\
 Data File : BF125223.D
 Acq On : 26 Aug 2021 12:55
 Operator : JU/CG
 Sample : M3426-13MS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SWCS-10-0510MS

Manual Integrations
 APPROVED

mohammad
 8/27/2021 5:07:00 PM

Quant Time: Aug 26 13:20:29 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF081821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 24 12:39:21 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.916	152	236119	20.000	ng	0.00
21) Naphthalene-d8	8.198	136	935032	20.000	ng	# 0.00
39) Acenaphthene-d10	9.957	164	497503	20.000	ng	0.00
64) Phenanthrene-d10	11.439	188	902473	20.000	ng	# 0.00
76) Chrysene-d12	14.080	240	464778	20.000	ng	0.00
86) Perylene-d12	15.574	264	474004	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.545	112	1694186	108.603	ng	0.02
7) Phenol-d6	6.545	99	2159899	107.055	ng	0.00
23) Nitrobenzene-d5	7.481	82	1449005	78.017	ng	0.00
42) 2,4,6-Tribromophenol	10.745	330	662601	133.423	ng	0.00
45) 2-Fluorobiphenyl	9.275	172	2371160	66.427	ng	0.00
79) Terphenyl-d14	13.027	244	2376604	81.461	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.769	88	309693	40.232	ng	# 100
3) Pyridine	3.534	79	708992	34.890	ng	# 90
4) n-Nitrosodimethylamine	3.487	42	402400	39.573	ng	# 26
6) Aniline	6.581	93	1030796	43.495	ng	96
8) 2-Chlorophenol	6.698	128	687984	43.232	ng	# 81
9) Benzaldehyde	6.469	77	536574	37.896	ng	94
10) Phenol	6.563	94	952540	45.083	ng	100
11) bis(2-Chloroethyl)ether	6.651	93	735921	44.226	ng	87
12) 1,3-Dichlorobenzene	6.857	146	756715	41.034	ng	98
13) 1,4-Dichlorobenzene	6.934	146	769289	41.866	ng	97
14) 1,2-Dichlorobenzene	7.087	146	734689	42.499	ng	98
15) Benzyl Alcohol	7.057	79	670767	43.180	ng	97
16) 2,2'-oxybis(1-Chloropr...	7.192	45	1363685	40.864	ng	95
17) 2-Methylphenol	7.169	107	618996	46.304	ng	95
18) Hexachloroethane	7.428	117	276106	42.911	ng	# 88
19) n-Nitroso-di-n-propyla...	7.334	70	512721	42.250	ng	# 74
20) 3+4-Methylphenols	7.322	107	698469	41.623	ng	# 77
22) Acetophenone	7.328	105	954384	39.660	ng	96
24) Nitrobenzene	7.498	77	819158	40.477	ng	# 87
25) Isophorone	7.739	82	1488339	41.540	ng	# 89
26) 2-Nitrophenol	7.816	139	391673	48.469	ng	# 83
27) 2,4-Dimethylphenol	7.851	122	630853	44.593	ng	94
28) bis(2-Chloroethoxy)met...	7.945	93	895396	45.135	ng	# 97
29) 2,4-Dichlorophenol	8.057	162	601864	41.716	ng	96
30) 1,2,4-Trichlorobenzene	8.139	180	633657	40.276	ng	95
31) Naphthalene	8.222	128	1853456	38.705	ng	100
32) Benzoic acid	7.975	122	435310	45.336	ng	89
33) 4-Chloroaniline	8.269	127	696034	36.870	ng	# 93
34) Hexachlorobutadiene	8.334	225	392037	39.079	ng	99
35) Caprolactam	8.645	113	209539m	56.072	ng	
36) 4-Chloro-3-methylphenol	8.745	107	665023	45.151	ng	91
37) 2-Methylnaphthalene	8.910	142	1302337	41.157	ng	98
38) 1-Methylnaphthalene	9.010	142	1210168	40.954	ng	96
40) 1,2,4,5-Tetrachloroben...	9.075	216	630515	38.791	ng	98
41) Hexachlorocyclopentadiene	9.063	237	703867	82.422	ng	99
43) 2,4,6-Trichlorophenol	9.186	196	464832	40.615	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.228	196	512121	42.532	ng	92
46) 1,1'-Biphenyl	9.375	154	1497719	37.341	ng	97
47) 2-Chloronaphthalene	9.404	162	1264768	38.804	ng	96
48) 2-Nitroaniline	9.498	65	484512	47.955	ng	88
49) Acenaphthylene	9.816	152	1896059	39.921	ng	97
50) Dimethylphthalate	9.675	163	1501258	43.544	ng	# 97
51) 2,6-Dinitrotoluene	9.739	165	339835	45.458	ng	93
52) Acenaphthene	9.992	154	1300250	44.160	ng	98
53) 3-Nitroaniline	9.910	138	327530	40.627	ng	86
54) 2,4-Dinitrophenol	10.016	184	355845	120.506	ng	# 79
55) Dibenzofuran	10.163	168	1663051	39.247	ng	97
56) 4-Nitrophenol	10.069	139	570698	89.392	ng	# 75
57) 2,4-Dinitrotoluene	10.145	165	454261	50.207	ng	# 90
58) Fluorene	10.504	166	1282695	40.944	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.275	232	393054	43.987	ng	# 86
60) Diethylphthalate	10.375	149	1460088	43.441	ng	97
61) 4-Chlorophenyl-phenyle...	10.492	204	658535	41.575	ng	88
62) 4-Nitroaniline	10.527	138	366460	50.550	ng	# 78
63) Azobenzene	10.657	77	1542401	40.913	ng	91
65) 4,6-Dinitro-2-methylph...	10.551	198	241545	50.570	ng	82
66) n-Nitrosodiphenylamine	10.616	169	1261067	39.025	ng	98
67) 4-Bromophenyl-phenylether	10.986	248	451457	41.658	ng	# 88
68) Hexachlorobenzene	11.051	284	478431	42.914	ng	# 86
69) Atrazine	11.139	200	423847	45.709	ng	92
70) Pentachlorophenol	11.245	266	503821	77.397	ng	90
71) Phenanthrene	11.469	178	1946347	39.141	ng	96
72) Anthracene	11.522	178	1987860	40.557	ng	98
73) Carbazole	11.674	167	1737404	38.790	ng	99
74) Di-n-butylphthalate	11.998	149	2227032	41.900	ng	99
75) Fluoranthene	12.657	202	1932129	39.159	ng	96
77) Benzidine	12.774	184	809644	49.879	ng	98
78) Pyrene	12.886	202	1913042	47.984	ng	95
80) Butylbenzylphthalate	13.498	149	778622	49.702	ng	90
81) Benzo(a)anthracene	14.074	228	1404032	41.927	ng	99
82) 3,3'-Dichlorobenzidine	14.033	252	480588	46.297	ng	98
83) Chrysene	14.110	228	1362152	41.131	ng	97
84) Bis(2-ethylhexyl)phtha...	14.051	149	989617	45.962	ng	99
85) Di-n-octyl phthalate	14.668	149	1519132	42.421	ng	# 98
87) Indeno(1,2,3-cd)pyrene	17.098	276	1717528	50.296	ng	92
88) Benzo(b)fluoranthene	15.139	252	1382434	39.920	ng	# 95
89) Benzo(k)fluoranthene	15.168	252	1297112	40.140	ng	99
90) Benzo(a)pyrene	15.515	252	1363447	44.572	ng	97
91) Dibenzo(a,h)anthracene	17.115	278	1433188	48.553	ng	95
92) Benzo(g,h,i)perylene	17.562	276	1481268	52.045	ng	# 91

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

