

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062321\
 Data File : BG049066.D
 Acq On : 23 Jun 2021 18:09
 Operator : CG/JU
 Sample : M2800-01MSD
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BORING-ADDITIONMSD

Manual Integrations
 APPROVED

mohammad
 6/24/2021 9:50:18 AM

Quant Time: Jun 24 02:18:10 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG061121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 16 12:04:50 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.111	152	28674	20.000	ng	0.00
21) Naphthalene-d8	10.931	136	112731	20.000	ng	# 0.00
39) Acenaphthene-d10	14.750	164	86840	20.000	ng	0.00
64) Phenanthrene-d10	17.500	188	224113	20.000	ng	# 0.00
76) Chrysene-d12	21.789	240	226148	20.000	ng	0.00
86) Perylene-d12	25.109	264	240143	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.661	112	168104	91.262	ng	0.00
7) Phenol-d6	7.259	99	222181	80.477	ng	0.00
23) Nitrobenzene-d5	9.274	82	155640	59.533	ng	0.00
42) 2,4,6-Tribromophenol	16.237	330	144690	112.460	ng	0.00
45) 2-Fluorobiphenyl	13.375	172	333235	54.209	ng	0.00
79) Terphenyl-d14	20.103	244	719918	58.301	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.528	88	31139	34.628	ng	# 76
3) Pyridine	3.933	79	79178	32.457	ng	91
4) n-Nitrosodimethylamine	3.845	42	60477	51.769	ng	# 81
6) Aniline	7.424	93	120272	34.593	ng	93
8) 2-Chlorophenol	7.670	128	98194	48.365	ng	85
9) Benzaldehyde	7.241	77	75491	52.844	ng	90
10) Phenol	7.288	94	131993	46.278	ng	93
11) bis(2-Chloroethyl)ether	7.518	93	90080	42.567	ng	86
12) 1,3-Dichlorobenzene	7.999	146	104771	46.195	ng	98
13) 1,4-Dichlorobenzene	8.146	146	104969	46.423	ng	99
14) 1,2-Dichlorobenzene	8.463	146	100914	46.392	ng	91
15) Benzyl Alcohol	8.346	79	115106	45.525	ng	95
16) 2,2'-oxybis(1-Chloropr...	8.640	45	153831	38.418	ng	83
17) 2-Methylphenol	8.552	107	88865	47.511	ng	98
18) Hexachloroethane	9.198	117	43538	47.996	ng	95
19) n-Nitroso-di-n-propyla...	8.916	70	84438	39.128	ng	92
20) 3+4-Methylphenols	8.881	107	122502	47.903	ng	96
22) Acetophenone	8.934	105	165169	48.158	ng	# 98
24) Nitrobenzene	9.321	77	131266	44.574	ng	94
25) Isophorone	9.844	82	225824	39.649	ng	98
26) 2-Nitrophenol	10.032	139	56571	56.631	ng	97
27) 2,4-Dimethylphenol	10.091	122	97851	54.027	ng	95
28) bis(2-Chloroethoxy)met...	10.326	93	125036	47.410	ng	# 97
29) 2,4-Dichlorophenol	10.573	162	102556	50.472	ng	93
30) 1,2,4-Trichlorobenzene	10.790	180	115754	49.688	ng	94
31) Naphthalene	10.984	128	307859	45.840	ng	99
32) Benzoic acid	10.232	122	75235	63.597	ng	# 77
33) 4-Chloroaniline	11.084	127	53908	18.905	ng	98
34) Hexachlorobutadiene	11.266	225	80925	48.739	ng	97
35) Caprolactam	11.865	113	38379m	65.341	ng	
36) 4-Chloro-3-methylphenol	12.206	107	124577	50.160	ng	94
37) 2-Methylnaphthalene	12.582	142	234137	46.156	ng	96
38) 1-Methylnaphthalene	12.800	142	217252	45.618	ng	97
40) 1,2,4,5-Tetrachloroben...	12.946	216	136817	49.616	ng	98
41) Hexachlorocyclopentadiene	12.929	237	194742	109.858	ng	90
43) 2,4,6-Trichlorophenol	13.187	196	102382	55.491	ng	96

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.264	196	112955	59.206	ng	94
46) 1,1'-Biphenyl	13.587	154	298849	46.678	ng	98
47) 2-Chloronaphthalene	13.628	162	245265	47.972	ng	97
48) 2-Nitroaniline	13.828	65	98886	51.758	ng	96
49) Acenaphthylene	14.474	152	400571	47.017	ng	94
50) Dimethylphthalate	14.204	163	349124	51.612	ng	100
51) 2,6-Dinitrotoluene	14.321	165	77074	55.949	ng	91
52) Acenaphthene	14.815	154	308997m	56.270	ng	
53) 3-Nitroaniline	14.650	138	46430	33.291	ng	# 81
54) 2,4-Dinitrophenol	14.856	184	98070	153.690	ng	96
55) Dibenzofuran	15.150	168	393749	46.129	ng	97
56) 4-Nitrophenol	14.962	139	139409	122.608	ng	91
57) 2,4-Dinitrotoluene	15.109	165	117514	69.449	ng	91
58) Fluorene	15.796	166	337822	48.399	ng	96
59) 2,3,4,6-Tetrachlorophenol	15.373	232	110828	57.390	ng	# 96
60) Diethylphthalate	15.561	149	382588	52.221	ng	97
61) 4-Chlorophenyl-phenyle...	15.790	204	191518	52.474	ng	95
62) 4-Nitroaniline	15.814	138	79801	56.578	ng	90
63) Azobenzene	16.078	77	362082	43.194	ng	89
65) 4,6-Dinitro-2-methylph...	15.867	198	67820	63.815	ng	91
66) n-Nitrosodiphenylamine	16.002	169	306783	43.925	ng	96
67) 4-Bromophenyl-phenylether	16.683	248	135048	48.260	ng	93
68) Hexachlorobenzene	16.807	284	151547	49.967	ng	95
69) Atrazine	16.948	200	138395	61.734	ng	99
70) Pentachlorophenol	17.147	266	200416	109.953	ng	98
71) Phenanthrene	17.541	178	588044	46.632	ng	98
72) Anthracene	17.635	178	599441	47.306	ng	97
73) Carbazole	17.899	167	560692	46.161	ng	98
74) Di-n-butylphthalate	18.458	149	686301	50.377	ng	98
75) Fluoranthene	19.550	202	764657	50.285	ng	98
77) Benzidine	19.721	184	369310	81.065	ng	98
78) Pyrene	19.909	202	764112	46.551	ng	98
80) Butylbenzylphthalate	20.790	149	299182	48.310	ng	96
81) Benzo(a)anthracene	21.771	228	770046	47.424	ng	99
82) 3,3'-Dichlorobenzidine	21.677	252	196174	37.780	ng	98
83) Chrysene	21.836	228	739060	47.472	ng	97
84) Bis(2-ethylhexyl)phtha...	21.666	149	428376	48.527	ng	95
85) Di-n-octyl phthalate	22.917	149	733978	49.024	ng	98
87) Indeno(1,2,3-cd)pyrene	28.916	276	947893	52.449	ng	# 97
88) Benzo(b)fluoranthene	24.057	252	824600	48.169	ng	# 95
89) Benzo(k)fluoranthene	24.127	252	795517	48.978	ng	98
90) Benzo(a)pyrene	24.956	252	808625	51.747	ng	99
91) Dibenzo(a,h)anthracene	28.992	278	792282	52.454	ng	# 97
92) Benzo(g,h,i)perylene	30.115	276	786858	51.662	ng	97

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

