

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101821\
 Data File : BG050620.D
 Acq On : 18 Oct 2021 16:49
 Operator : CG/JU
 Sample : M4249-04MSD
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TP1-(D)MSD

Manual Integrations
 APPROVED

mohammad
 10/19/2021 12:09:06 PM

Quant Time: Oct 18 17:28:45 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 14 10:56:22 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.253	152	43965	20.00	ng	0.00
21) Naphthalene-d8	11.079	136	195247	20.00	ng	# 0.00
39) Acenaphthene-d10	14.875	164	133691	20.00	ng	0.00
64) Phenanthrene-d10	17.619	188	293810	20.00	ng	# 0.00
76) Chrysene-d12	21.920	240	250040	20.00	ng	# 0.00
86) Perylene-d12	25.333	264	250667	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.792	112	349976	119.05	ng	0.00
7) Phenol-d6	7.401	99	472064	110.91	ng	0.00
23) Nitrobenzene-d5	9.428	82	326242	69.17	ng	0.00
42) 2,4,6-Tribromophenol	16.356	330	132173	103.65	ng	0.00
45) 2-Fluorobiphenyl	13.500	172	593215	66.78	ng	0.00
79) Terphenyl-d14	20.204	244	911016	71.01	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.653	88	55712	36.59	ng	91
3) Pyridine	4.058	79	126503	30.41	ng	# 92
4) n-Nitrosodimethylamine	3.976	42	84224	42.97	ng	# 52
6) Aniline	7.572	93	159084	32.17	ng	# 90
8) 2-Chlorophenol	7.813	128	142323	50.28	ng	89
9) Benzaldehyde	7.384	77	106320	39.57	ng	93
10) Phenol	7.425	94	205023	49.54	ng	86
11) bis(2-Chloroethyl)ether	7.666	93	144856	44.85	ng	86
12) 1,3-Dichlorobenzene	8.142	146	144778	44.31	ng	98
13) 1,4-Dichlorobenzene	8.289	146	147860	44.89	ng	94
14) 1,2-Dichlorobenzene	8.612	146	141563	44.99	ng	97
15) Benzyl Alcohol	8.488	79	161128	47.90	ng	91
16) 2,2'-oxybis(1-Chloropr...	8.776	45	237105	45.01	ng	87
17) 2-Methylphenol	8.688	107	135113	49.62	ng	92
18) Hexachloroethane	9.346	117	57477	46.10	ng	95
19) n-Nitroso-di-n-propyla...	9.058	70	137090	44.19	ng	# 88
20) 3+4-Methylphenols	9.017	107	186256	48.60	ng	94
22) Acetophenone	9.082	105	226611	40.82	ng	# 96
24) Nitrobenzene	9.470	77	201710	43.01	ng	96
25) Isophorone	9.992	82	347381	41.52	ng	# 95
26) 2-Nitrophenol	10.180	139	79367	46.09	ng	# 92
27) 2,4-Dimethylphenol	10.227	122	146551	49.26	ng	95
28) bis(2-Chloroethoxy)met...	10.468	93	197840	41.36	ng	94
29) 2,4-Dichlorophenol	10.721	162	141009	46.50	ng	97
30) 1,2,4-Trichlorobenzene	10.938	180	140525	40.80	ng	98
31) Naphthalene	11.132	128	441206	42.22	ng	99
32) Benzoic acid	10.368	122	85062	38.52	ng	# 73
33) 4-Chloroaniline	11.232	127	65934	15.04	ng	97
34) Hexachlorobutadiene	11.403	225	86411	39.96	ng	93
35) Caprolactam	12.008	113	53992m	46.52	ng	
36) 4-Chloro-3-methylphenol	12.337	107	172782	45.59	ng	# 89
37) 2-Methylnaphthalene	12.719	142	313016	43.41	ng	94
38) 1-Methylnaphthalene	12.936	142	292660	41.85	ng	94
40) 1,2,4,5-Tetrachloroben...	13.077	216	157620	39.92	ng	97
41) Hexachlorocyclopentadiene	13.054	237	199217	87.31	ng	92
43) 2,4,6-Trichlorophenol	13.312	196	117163	42.11	ng	91

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101821\
 Data File : BG050620.D
 Acq On : 18 Oct 2021 16:49
 Operator : CG/JU
 Sample : M4249-04MSD
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TP1-(D)MSD

Manual Integrations
 APPROVED

mohammad
 10/19/2021 12:09:06 PM

Quant Time: Oct 18 17:28:45 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 14 10:56:22 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.383	196	127627	43.88	ng	92
46) 1,1'-Biphenyl	13.712	154	389471	39.94	ng	93
47) 2-Chloronaphthalene	13.759	162	311314	41.58	ng	99
48) 2-Nitroaniline	13.958	65	130430	43.79	ng	94
49) Acenaphthylene	14.599	152	511335	41.68	ng	97
50) Dimethylphthalate	14.317	163	417873	41.35	ng	99
51) 2,6-Dinitrotoluene	14.446	165	91091	45.00	ng	89
52) Acenaphthene	14.940	154	360650m	45.75	ng	
53) 3-Nitroaniline	14.775	138	54438	22.74	ng	87
54) 2,4-Dinitrophenol	14.981	184	102336	81.50	ng	87
55) Dibenzofuran	15.269	168	485341	39.80	ng	98
56) 4-Nitrophenol	15.075	139	162457	84.36	ng	# 80
57) 2,4-Dinitrotoluene	15.227	165	126950	44.50	ng	93
58) Fluorene	15.921	166	387316	41.35	ng	96
59) 2,3,4,6-Tetrachlorophenol	15.492	232	102843	40.59	ng	94
60) Diethylphthalate	15.668	149	436568	41.12	ng	96
61) 4-Chlorophenyl-phenyle...	15.903	204	206498	40.30	ng	98
62) 4-Nitroaniline	15.938	138	99080	39.78	ng	# 68
63) Azobenzene	16.197	77	491901	40.02	ng	82
65) 4,6-Dinitro-2-methylph...	15.985	198	69406	39.33	ng	82
66) n-Nitrosodiphenylamine	16.115	169	346888	41.55	ng	96
67) 4-Bromophenyl-phenylether	16.796	248	122986	41.54	ng	97
68) Hexachlorobenzene	16.920	284	127155	43.21	ng	# 93
69) Atrazine	17.055	200	141669	43.08	ng	98
70) Pentachlorophenol	17.260	266	154922	77.34	ng	97
71) Phenanthrene	17.660	178	656027	42.37	ng	97
72) Anthracene	17.754	178	659724	42.88	ng	98
73) Carbazole	18.018	167	606648	39.56	ng	98
74) Di-n-butylphthalate	18.553	149	757061	42.07	ng	# 97
75) Fluoranthene	19.658	202	778259	40.89	ng	96
77) Benzidine	19.828	184	243549	30.05	ng	97
78) Pyrene	20.016	202	784638	44.23	ng	98
80) Butylbenzylphthalate	20.886	149	331897	44.23	ng	95
81) Benzo(a)anthracene	21.896	228	700548	42.34	ng	99
82) 3,3'-Dichlorobenzidine	21.802	252	151552	27.64	ng	# 98
83) Chrysene	21.967	228	672929	43.24	ng	96
84) Bis(2-ethylhexyl)phtha...	21.773	149	477959	44.83	ng	# 96
85) Di-n-octyl phthalate	23.048	149	808848	45.48	ng	97
87) Indeno(1,2,3-cd)pyrene	29.258	276	773123	44.28	ng	# 89
88) Benzo(b)fluoranthene	24.240	252	704426	45.89	ng	# 93
89) Benzo(k)fluoranthene	24.311	252	671665	44.47	ng	# 98
90) Benzo(a)pyrene	25.175	252	686412	52.59	ng	# 94
91) Dibenzo(a,h)anthracene	29.329	278	646015	44.74	ng	# 91
92) Benzo(g,h,i)perylene	30.486	276	639804	45.24	ng	# 91

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

