

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG043024\
 Data File : BG061059.D
 Acq On : 30 Apr 2024 11:32
 Operator : MA/JU
 Sample : SSTDICC005
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

SSTDICC005

Manual Integrations

APPROVED

Reviewed By :Jagrut Upadhyay 05/01/2024

Supervised By :mohammad ahmed 05/03/2024

Quant Time: May 01 04:48:28 2024

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG043024.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Wed May 01 04:45:52 2024

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.959	152	17399	20.000	ng	# 0.00
21) Naphthalene-d8	10.756	136	70277	20.000	ng	# 0.00
39) Acenaphthene-d10	14.586	164	65131	20.000	ng	0.00
64) Phenanthrene-d10	17.330	188	182901	20.000	ng	0.00
76) Chrysene-d12	21.584	240	207160	20.000	ng	0.00
86) Perylene-d12	24.722	264	235870	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.550	112	8629	10.086	ng	0.00
7) Phenol-d6	7.130	99	11962	9.452	ng	0.00
23) Nitrobenzene-d5	9.116	82	14392	10.152	ng	0.00
42) 2,4,6-Tribromophenol	16.073	330	13319	10.231	ng	0.00
45) 2-Fluorobiphenyl	13.206	172	43653	9.911	ng	0.00
79) Terphenyl-d14	19.945	244	126280	10.138	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.464	88	1987	5.744	ng	92
3) Pyridine	3.858	79	4970	5.151	ng	86
4) n-Nitrosodimethylamine	3.776	42	4670	4.775	ng	# 82
6) Aniline	7.283	93	5855	4.667	ng	# 94
8) 2-Chlorophenol	7.536	128	4878	4.648	ng	85
9) Benzaldehyde	7.101	77	4275m	4.205	ng	
10) Phenol	7.160	94	6038	4.691	ng	87
11) bis(2-Chloroethyl)ether	7.383	93	4229	4.756	ng	93
12) 1,3-Dichlorobenzene	7.853	146	6073m	4.908	ng	
13) 1,4-Dichlorobenzene	7.994	146	6487	5.073	ng	90
14) 1,2-Dichlorobenzene	8.311	146	6173m	5.003	ng	
15) Benzyl Alcohol	8.194	79	5631	4.932	ng	90
16) 2,2'-oxybis(1-Chloropr...	8.488	45	9304	4.822	ng	94
17) 2-Methylphenol	8.405	107	4168	4.544	ng	94
18) Hexachloroethane	9.046	117	2206	4.834	ng	98
19) n-Nitroso-di-n-propyla...	8.758	70	5218	5.509	ng	# 74
20) 3+4-Methylphenols	8.723	107	5892	4.453	ng	96
22) Acetophenone	8.776	105	8769	5.010	ng	# 95
24) Nitrobenzene	9.157	77	6748	4.866	ng	# 89
25) Isophorone	9.680	82	13900	5.304	ng	97
26) 2-Nitrophenol	9.863	139	3168	4.821	ng	# 81
27) 2,4-Dimethylphenol	9.933	122	3599	4.751	ng	83
28) bis(2-Chloroethoxy)met...	10.162	93	6727	5.230	ng	# 89
29) 2,4-Dichlorophenol	10.403	162	5859	4.936	ng	82
30) 1,2,4-Trichlorobenzene	10.626	180	7299	5.104	ng	# 93
31) Naphthalene	10.808	128	18967	5.193	ng	97
32) Benzoic acid	10.004	122	3426	4.097	ng	# 60
33) 4-Chloroaniline	10.914	127	7028	4.825	ng	# 86
34) Hexachlorobutadiene	11.096	225	6243	5.087	ng	85
35) Caprolactam	11.655	113	2357	5.514	ng	# 82
36) 4-Chloro-3-methylphenol	12.042	107	6883	4.809	ng	96
37) 2-Methylnaphthalene	12.412	142	13499	4.932	ng	97
38) 1-Methylnaphthalene	12.630	142	14657	5.319	ng	86
40) 1,2,4,5-Tetrachloroben...	12.777	216	11077	5.096	ng	98
41) Hexachlorocyclopentadiene	12.759	237	3312	3.980	ng	81
43) 2,4,6-Trichlorophenol	13.018	196	6749m	4.836	ng	

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44) 2,4,5-Trichlorophenol	13.094	196	7360	4.660	ng	96
46) 1,1'-Biphenyl	13.417	154	21269	5.093	ng	94
47) 2-Chloronaphthalene	13.464	162	16581	4.958	ng	91
48) 2-Nitroaniline	13.658	65	6191	4.989	ng	# 80
49) Acenaphthylene	14.304	152	25583	4.981	ng	98
50) Dimethylphthalate	14.040	163	26812	5.330	ng	98
51) 2,6-Dinitrotoluene	14.152	165	5768	5.360	ng	# 84
52) Acenaphthene	14.651	154	16017	4.999	ng	98
53) 3-Nitroaniline	14.481	138	5162	5.243	ng	99
54) 2,4-Dinitrophenol	14.692	184	3018	6.055	ng	# 73
55) Dibenzofuran	14.980	168	27596	5.099	ng	98
56) 4-Nitrophenol	14.798	139	4250	5.341	ng	# 79
57) 2,4-Dinitrotoluene	14.945	165	7760	4.874	ng	# 96
58) Fluorene	15.632	166	23593	5.038	ng	94
59) 2,3,4,6-Tetrachlorophenol	15.215	232	8124	4.997	ng	# 96
60) Diethylphthalate	15.403	149	27853	5.490	ng	# 96
61) 4-Chlorophenyl-phenyle...	15.626	204	14213	5.055	ng	87
62) 4-Nitroaniline	15.644	138	5221	4.889	ng	93
63) Azobenzene	15.920	77	20521	5.196	ng	96
65) 4,6-Dinitro-2-methylph...	15.703	198	5164	4.551	ng	96
66) n-Nitrosodiphenylamine	15.838	169	22345	4.927	ng	# 88
67) 4-Bromophenyl-phenylether	16.519	248	10008	4.645	ng	93
68) Hexachlorobenzene	16.637	284	13567	5.059	ng	# 88
69) Atrazine	16.790	200	10613	5.674	ng	96
70) Pentachlorophenol	16.984	266	7075	4.294	ng	98
71) Phenanthrene	17.371	178	42831	5.121	ng	97
72) Anthracene	17.465	178	41357	4.956	ng	96
73) Carbazole	17.730	167	39083	5.327	ng	96
74) Di-n-butylphthalate	18.294	149	46739	5.357	ng	98
75) Fluoranthene	19.381	202	62170	5.426	ng	98
77) Benzidine	19.563	184	9259	5.012	ng	95
78) Pyrene	19.745	202	65595	4.980	ng	99
80) Butylbenzylphthalate	20.638	149	20988	5.034	ng	98
81) Benzo(a)anthracene	21.566	228	71601	5.282	ng	96
82) 3,3'-Dichlorobenzidine	21.484	252	24985	5.121	ng	96
83) Chrysene	21.631	228	67381	5.305	ng	96
84) Bis(2-ethylhexyl)phtha...	21.496	149	27141	4.985	ng	# 98
85) Di-n-octyl phthalate	22.677	149	48122	5.010	ng	# 93
87) Indeno(1,2,3-cd)pyrene	28.270	276	81384	5.206	ng	# 89
88) Benzo(b)fluoranthene	23.723	252	68629	5.207	ng	97
89) Benzo(k)fluoranthene	23.793	252	68379m	5.401	ng	
90) Benzo(a)pyrene	24.569	252	61478	5.354	ng	96
91) Dibenzo(a,h)anthracene	28.329	278	66990	5.175	ng	# 94
92) Benzo(g,h,i)perylene	29.375	276	65458	5.088	ng	99

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

