

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG043024\
 Data File : BG061060.D
 Acq On : 30 Apr 2024 12:13
 Operator : MA/JU
 Sample : SSTDICC010
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

SSTDICC010

Manual Integrations

APPROVED

Reviewed By :Jagrut Upadhyay 05/01/2024
 Supervised By :mohammad ahmed 05/03/2024

Quant Time: May 01 04:49:08 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	7.964	152	17677	20.000	ng	0.00
21) Naphthalene-d8	10.755	136	76628	20.000	ng	# 0.00
39) Acenaphthene-d10	14.585	164	72556	20.000	ng	0.00
64) Phenanthrene-d10	17.329	188	202493	20.000	ng	0.00
76) Chrysene-d12	21.589	240	218523	20.000	ng	0.00
86) Perylene-d12	24.721	264	247795	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.549	112	17482	20.113	ng	0.00
7) Phenol-d6	7.129	99	24436	19.004	ng	0.00
23) Nitrobenzene-d5	9.109	82	29902	19.344	ng	0.00
42) 2,4,6-Tribromophenol	16.072	330	28641	19.749	ng	0.00
45) 2-Fluorobiphenyl	13.210	172	95158	19.394	ng	0.00
79) Terphenyl-d14	19.944	244	256067	19.489	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.463	88	4098	11.661	ng	84
3) Pyridine	3.857	79	10043	10.245	ng	# 81
4) n-Nitrosodimethylamine	3.769	42	10620	10.689	ng	# 89
6) Aniline	7.288	93	12543	9.840	ng	97
8) 2-Chlorophenol	7.529	128	10703	10.037	ng	90
9) Benzaldehyde	7.100	77	8806	10.119	ng	94
10) Phenol	7.159	94	12508	9.566	ng	90
11) bis(2-Chloroethyl)ether	7.382	93	8945	9.902	ng	96
12) 1,3-Dichlorobenzene	7.852	146	12598	10.021	ng	95
13) 1,4-Dichlorobenzene	7.999	146	12968	9.981	ng	93
14) 1,2-Dichlorobenzene	8.316	146	12316	9.825	ng	93
15) Benzyl Alcohol	8.199	79	10234	8.823	ng	# 84
16) 2,2'-oxybis(1-Chloropr...	8.492	45	19518	9.957	ng	96
17) 2-Methylphenol	8.404	107	9198	9.871	ng	86
18) Hexachloroethane	9.039	117	5025	10.839	ng	98
19) n-Nitroso-di-n-propyla...	8.757	70	9345	9.712	ng	94
20) 3+4-Methylphenols	8.728	107	13507	10.047	ng	98
22) Acetophenone	8.780	105	18591	9.742	ng	# 90
24) Nitrobenzene	9.156	77	14445	9.552	ng	91
25) Isophorone	9.679	82	29286	10.249	ng	94
26) 2-Nitrophenol	9.867	139	7255	10.126	ng	96
27) 2,4-Dimethylphenol	9.932	122	8146	9.862	ng	93
28) bis(2-Chloroethoxy)met...	10.161	93	14033	10.007	ng	92
29) 2,4-Dichlorophenol	10.408	162	12393	9.576	ng	96
30) 1,2,4-Trichlorobenzene	10.619	180	14929	9.573	ng	# 78
31) Naphthalene	10.813	128	39052	9.806	ng	96
32) Benzoic acid	10.014	122	8595m	9.425	ng	
33) 4-Chloroaniline	10.913	127	15262	9.609	ng	96
34) Hexachlorobutadiene	11.095	225	13714	10.249	ng	# 90
35) Caprolactam	11.665	113	4896	10.505	ng	# 84
36) 4-Chloro-3-methylphenol	12.035	107	16020	10.265	ng	98
37) 2-Methylnaphthalene	12.411	142	30543	10.234	ng	90
38) 1-Methylnaphthalene	12.629	142	29611	9.856	ng	97
40) 1,2,4,5-Tetrachloroben...	12.782	216	22192	9.165	ng	98
41) Hexachlorocyclopentadiene	12.764	237	7687	8.293	ng	85
43) 2,4,6-Trichlorophenol	13.017	196	14634	9.414	ng	92

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.093	196	17050	9.691	ng	96
46) 1,1'-Biphenyl	13.416	154	45129	9.701	ng	96
47) 2-Chloronaphthalene	13.463	162	35655	9.571	ng	95
48) 2-Nitroaniline	13.663	65	13437	9.720	ng	97
49) Acenaphthylene	14.303	152	54657	9.553	ng	97
50) Dimethylphthalate	14.039	163	55662	9.932	ng	96
51) 2,6-Dinitrotoluene	14.156	165	11822	9.862	ng	91
52) Acenaphthene	14.650	154	34006	9.527	ng	94
53) 3-Nitroaniline	14.485	138	11056	10.080	ng	# 72
54) 2,4-Dinitrophenol	14.697	184	6517	9.665	ng	# 87
55) Dibenzofuran	14.985	168	59896	9.935	ng	97
56) 4-Nitrophenol	14.803	139	8621	9.726	ng	90
57) 2,4-Dinitrotoluene	14.944	165	17814	10.045	ng	# 88
58) Fluorene	15.637	166	51742	9.918	ng	96
59) 2,3,4,6-Tetrachlorophenol	15.208	232	17571	9.702	ng	98
60) Diethylphthalate	15.402	149	56613	10.018	ng	99
61) 4-Chlorophenyl-phenyle...	15.625	204	31971	10.206	ng	96
62) 4-Nitroaniline	15.643	138	11868	9.977	ng	99
63) Azobenzene	15.919	77	44163	10.039	ng	95
65) 4,6-Dinitro-2-methylph...	15.708	198	11322	9.012	ng	99
66) n-Nitrosodiphenylamine	15.843	169	48600	9.678	ng	97
67) 4-Bromophenyl-phenylether	16.518	248	22589	9.470	ng	91
68) Hexachlorobenzene	16.642	284	29053	9.785	ng	98
69) Atrazine	16.789	200	21134	10.205	ng	90
70) Pentachlorophenol	16.983	266	16519	9.056	ng	# 87
71) Phenanthrene	17.370	178	90760	9.803	ng	98
72) Anthracene	17.464	178	90072	9.749	ng	98
73) Carbazole	17.729	167	84263	10.373	ng	96
74) Di-n-butylphthalate	18.299	149	97220	10.065	ng	98
75) Fluoranthene	19.380	202	129943	10.244	ng	99
77) Benzidine	19.562	184	22714	7.944	ng	98
78) Pyrene	19.744	202	135143	9.728	ng	99
80) Butylbenzylphthalate	20.637	149	41395	9.413	ng	96
81) Benzo(a)anthracene	21.565	228	138157	9.662	ng	98
82) 3,3'-Dichlorobenzidine	21.483	252	49295	9.579	ng	99
83) Chrysene	21.630	228	131589	9.822	ng	98
84) Bis(2-ethylhexyl)phtha...	21.495	149	54286	9.453	ng	98
85) Di-n-octyl phthalate	22.676	149	93732	9.251	ng	97
87) Indeno(1,2,3-cd)pyrene	28.263	276	155614	9.476	ng	# 99
88) Benzo(b)fluoranthene	23.728	252	131563	9.501	ng	99
89) Benzo(k)fluoranthene	23.792	252	130486	9.810	ng	100
90) Benzo(a)pyrene	24.568	252	114201	9.467	ng	# 98
91) Dibenzo(a,h)anthracene	28.328	278	130212	9.575	ng	97
92) Benzo(g,h,i)perylene	29.374	276	131016	9.694	ng	# 98

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

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