

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG011224\
 Data File : BG060321.D
 Acq On : 12 Jan 2024 18:30
 Operator : MA/JU
 Sample : P1104-04MS
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TP-1MS

Quant Time: Jan 12 23:43:26 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG010824.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 09 04:07:12 2024
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 01/15/2024
 Supervised By :mohammad ahmed 01/16/2024

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.931	152	301426	20.000	ng	0.00
21) Naphthalene-d8	10.734	136	1036275	20.000	ng	0.00
39) Acenaphthene-d10	14.565	164	754748	20.000	ng	0.00
64) Phenanthrene-d10	17.314	188	1895619	20.000	ng	0.00
76) Chrysene-d12	21.580	240	1775265	20.000	ng	0.00
86) Perylene-d12	24.741	264	1960057	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.505	112	1958937	106.893	ng	0.00
7) Phenol-d6	7.097	99	2908568	107.182	ng	0.00
23) Nitrobenzene-d5	9.095	82	2003456	91.292	ng	0.00
42) 2,4,6-Tribromophenol	16.057	330	1543378	138.986	ng	0.00
45) 2-Fluorobiphenyl	13.190	172	4504686	82.981	ng	0.00
79) Terphenyl-d14	19.935	244	5970591	84.706	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.419	88	228252	27.266	ng	# 69
3) Pyridine	3.813	79	537799	22.770	ng	97
4) n-Nitrosodimethylamine	3.719	42	225416	30.495	ng	96
6) Aniline	7.256	93	569106	20.981	ng	99
8) 2-Chlorophenol	7.496	128	734353	40.433	ng	95
9) Benzaldehyde	7.068	77	139420m	8.951	ng	
10) Phenol	7.126	94	1056290	38.823	ng	99
11) bis(2-Chloroethyl)ether	7.350	93	840115	37.893	ng	96
12) 1,3-Dichlorobenzene	7.820	146	804862	35.306	ng	99
13) 1,4-Dichlorobenzene	7.966	146	924776	39.982	ng	98
14) 1,2-Dichlorobenzene	8.284	146	807656	33.958	ng	99
15) Benzyl Alcohol	8.166	79	732517m	41.700	ng	
16) 2,2'-oxybis(1-Chloropr...	8.448	45	1062152m	39.472	ng	
17) 2-Methylphenol	8.372	107	802169	42.925	ng	99
18) Hexachloroethane	9.012	117	275803	34.009	ng	97
19) n-Nitroso-di-n-propyla...	8.730	70	726413	38.677	ng	# 96
20) 3+4-Methylphenols	8.701	107	1093736	43.408	ng	96
22) Acetophenone	8.754	105	1435786	50.404	ng	# 99
24) Nitrobenzene	9.136	77	977453	42.965	ng	98
25) Isophorone	9.653	82	1914458	43.439	ng	98
26) 2-Nitrophenol	9.847	139	411740	49.224	ng	95
27) 2,4-Dimethylphenol	9.899	122	641051	57.999	ng	99
28) bis(2-Chloroethoxy)met...	10.140	93	1188978	44.154	ng	100
29) 2,4-Dichlorophenol	10.381	162	879017	49.077	ng	97
30) 1,2,4-Trichlorobenzene	10.593	180	934473	41.937	ng	93
31) Naphthalene	10.781	128	2280535	41.983	ng	99
32) Benzoic acid	10.035	122	316904	35.768	ng	90
33) 4-Chloroaniline	10.898	127	102427	4.894	ng	# 93
34) Hexachlorobutadiene	11.063	225	582870	40.154	ng	97
35) Caprolactam	11.668	113	253435m	43.932	ng	
36) 4-Chloro-3-methylphenol	12.015	107	881275	48.433	ng	99
37) 2-Methylnaphthalene	12.391	142	1711622	42.485	ng	98
38) 1-Methylnaphthalene	12.608	142	1699834	42.018	ng	99
40) 1,2,4,5-Tetrachloroben...	12.761	216	1217660	45.426	ng	99
41) Hexachlorocyclopentadiene	12.737	237	1193568	134.117	ng	99
43) 2,4,6-Trichlorophenol	12.996	196	817334	47.913	ng	97

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG011224\
 Data File : BG060321.D
 Acq On : 12 Jan 2024 18:30
 Operator : MA/JU
 Sample : P1104-04MS
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TP-1MS

Quant Time: Jan 12 23:43:26 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG010824.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 09 04:07:12 2024
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 01/15/2024
 Supervised By :mohammad ahmed 01/16/2024

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.072	196	905764	48.140	ng	96
46) 1,1'-Biphenyl	13.401	154	2577113	45.235	ng	99
47) 2-Chloronaphthalene	13.442	162	2138664	43.819	ng	99
48) 2-Nitroaniline	13.648	65	618543	50.889	ng	99
49) Acenaphthylene	14.288	152	3294470	49.018	ng	98
50) Dimethylphthalate	14.024	163	2584557	44.811	ng	100
51) 2,6-Dinitrotoluene	14.142	165	591070	48.027	ng	92
52) Acenaphthene	14.629	154	1873456	44.766	ng	99
53) 3-Nitroaniline	14.471	138	251077	22.389	ng	98
54) 2,4-Dinitrophenol	14.676	184	491053	67.331	ng	96
55) Dibenzofuran	14.964	168	3195269	45.737	ng	97
56) 4-Nitrophenol	14.788	139	831958	100.020	ng	99
57) 2,4-Dinitrotoluene	14.929	165	821676	48.688	ng	97
58) Fluorene	15.616	166	2687431	46.452	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.193	232	838499	52.481	ng	99
60) Diethylphthalate	15.387	149	2489044	44.952	ng	100
61) 4-Chlorophenyl-phenyle...	15.605	204	1422567	44.636	ng	97
62) 4-Nitroaniline	15.640	138	571949	47.921	ng	97
63) Azobenzene	15.898	77	2576857	46.011	ng	99
65) 4,6-Dinitro-2-methylph...	15.693	198	431219	41.120	ng	96
66) n-Nitrosodiphenylamine	15.822	169	2376305	47.211	ng	98
67) 4-Bromophenyl-phenylether	16.504	248	931620	44.727	ng	98
68) Hexachlorobenzene	16.621	284	1091165	44.256	ng	99
69) Atrazine	16.768	200	951852	50.152	ng	97
70) Pentachlorophenol	16.962	266	1240752	93.361	ng	94
71) Phenanthrene	17.355	178	4095088	44.694	ng	96
72) Anthracene	17.449	178	4183935	45.737	ng	97
73) Carbazole	17.720	167	3829390	44.403	ng	99
74) Di-n-butylphthalate	18.278	149	3923509	44.215	ng	99
75) Fluoranthene	19.371	202	4643832	42.197	ng	97
77) Benzidine	19.553	184	653736	16.950	ng	98
78) Pyrene	19.735	202	4784893	42.207	ng	98
80) Butylbenzylphthalate	20.622	149	1922349	49.287	ng	97
81) Benzo(a)anthracene	21.556	228	4899874	44.584	ng	96
82) 3,3'-Dichlorobenzidine	21.480	252	1107323	26.309	ng	99
83) Chrysene	21.627	228	4699591	43.479	ng	98
84) Bis(2-ethylhexyl)phtha...	21.462	149	2815053	47.763	ng	96
85) Di-n-octyl phthalate	22.655	149	4659304	48.784	ng	100
87) Indeno(1,2,3-cd)pyrene	28.354	276	6419875	47.106	ng	# 92
88) Benzo(b)fluoranthene	23.742	252	5295739	45.735	ng	99
89) Benzo(k)fluoranthene	23.807	252	5277206	46.184	ng	100
90) Benzo(a)pyrene	24.594	252	4914716	46.985	ng	98
91) Dibenzo(a,h)anthracene	28.413	278	5178787	46.530	ng	99
92) Benzo(g,h,i)perylene	29.482	276	4949295	43.388	ng	99

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG011224\
Data File : BG060321.D
Acq On : 12 Jan 2024 18:30
Operator : MA/JU
Sample : P1104-04MS
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
TP-1MS

Quant Time: Jan 12 23:43:26 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG010824.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jan 09 04:07:12 2024
Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 01/15/2024
Supervised By :mohammad ahmed 01/16/2024

