

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG041024\
 Data File : BG060849.D
 Acq On : 10 Apr 2024 14:00
 Operator : MA/JU
 Sample : PB160074BS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

PB160028BS

Manual Integrations

APPROVED

Reviewed By :Yogesh Patel 04/12/2024

Supervised By :mohammad ahmed 04/12/2024

Quant Time: Apr 10 14:33:37 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG032924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Mar 30 03:12:26 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.942	152	58076	20.000	ng	-0.02
21) Naphthalene-d8	10.745	136	239507	20.000	ng	-0.01
39) Acenaphthene-d10	14.576	164	194386	20.000	ng	-0.01
64) Phenanthrene-d10	17.325	188	460638	20.000	ng	-0.01
76) Chrysene-d12	21.591	240	432488	20.000	ng	-0.01
86) Perylene-d12	24.717	264	495602	20.000	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.527	112	461078	132.259	ng	0.00
7) Phenol-d6	7.114	99	650029	125.613	ng	0.00
23) Nitrobenzene-d5	9.100	82	419337	99.905	ng	-0.01
42) 2,4,6-Tribromophenol	16.068	330	369666	167.528	ng	0.00
45) 2-Fluorobiphenyl	13.207	172	1088391	82.174	ng	-0.01
79) Terphenyl-d14	19.952	244	2015397	82.055	ng	-0.01
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.448	88	42983	30.073	ng	94
3) Pyridine	3.835	79	124901	28.970	ng	88
4) n-Nitrosodimethylamine	3.753	42	99013	38.420	ng	97
6) Aniline	7.267	93	151599	23.199	ng	100
8) 2-Chlorophenol	7.513	128	177118	44.814	ng	98
9) Benzaldehyde	7.085	77	29053m	10.181	ng	
10) Phenol	7.137	94	239386	45.328	ng	99
11) bis(2-Chloroethyl)ether	7.367	93	160194	37.570	ng	97
12) 1,3-Dichlorobenzene	7.837	146	177078	43.084	ng	97
13) 1,4-Dichlorobenzene	7.983	146	182997	43.642	ng	95
14) 1,2-Dichlorobenzene	8.301	146	178196	43.364	ng	98
15) Benzyl Alcohol	8.177	79	176219	42.039	ng	95
16) 2,2'-oxybis(1-Chloropr...	8.465	45	398424	41.405	ng	100
17) 2-Methylphenol	8.389	107	164756	41.324	ng	98
18) Hexachloroethane	9.029	117	65183	43.769	ng	88
19) n-Nitroso-di-n-propyla...	8.753	70	149808	37.962	ng	# 95
20) 3+4-Methylphenols	8.712	107	229455	42.028	ng	96
22) Acetophenone	8.765	105	283190	42.881	ng	# 96
24) Nitrobenzene	9.141	77	213449	45.819	ng	97
25) Isophorone	9.670	82	391753	40.522	ng	99
26) 2-Nitrophenol	9.852	139	92199	54.757	ng	91
27) 2,4-Dimethylphenol	9.916	122	140563	36.195	ng	97
28) bis(2-Chloroethoxy)met...	10.152	93	235533	41.267	ng	99
29) 2,4-Dichlorophenol	10.392	162	184173	51.645	ng	99
30) 1,2,4-Trichlorobenzene	10.610	180	192522	47.676	ng	97
31) Naphthalene	10.792	128	545974	42.141	ng	98
32) Benzoic acid	10.057	122	127016	48.993	ng	96
33) 4-Chloroaniline	10.898	127	89027	14.794	ng	96
34) Hexachlorobutadiene	11.086	225	129796	48.229	ng	98
35) Caprolactam	11.667	113	62673m	44.373	ng	
36) 4-Chloro-3-methylphenol	12.026	107	219239	48.125	ng	96
37) 2-Methylnaphthalene	12.402	142	392429	41.199	ng	99
38) 1-Methylnaphthalene	12.619	142	372528	41.400	ng	98
40) 1,2,4,5-Tetrachloroben...	12.772	216	257626	46.100	ng	99
41) Hexachlorocyclopentadiene	12.754	237	297064	104.542	ng	100
43) 2,4,6-Trichlorophenol	13.007	196	183349	50.953	ng	95

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG041024\
 Data File : BG060849.D
 Acq On : 10 Apr 2024 14:00
 Operator : MA/JU
 Sample : PB160074BS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB160028BS

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 04/12/2024
 Supervised By :mohammad ahmed 04/12/2024

Quant Time: Apr 10 14:33:37 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG032924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Mar 30 03:12:26 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.077	196	204535	47.629	ng	98
46) 1,1'-Biphenyl	13.412	154	583017	42.462	ng	98
47) 2-Chloronaphthalene	13.454	162	457028	43.811	ng	99
48) 2-Nitroaniline	13.647	65	173349	50.747	ng	93
49) Acenaphthylene	14.300	152	779677	44.482	ng	99
50) Dimethylphthalate	14.041	163	641943	41.781	ng	99
51) 2,6-Dinitrotoluene	14.147	165	132790	48.251	ng	97
52) Acenaphthene	14.640	154	526368m	47.829	ng	
53) 3-Nitroaniline	14.476	138	83980	29.026	ng	99
54) 2,4-Dinitrophenol	14.681	184	158370	148.085	ng	89
55) Dibenzofuran	14.981	168	722979	41.649	ng	100
56) 4-Nitrophenol	14.793	139	233951	104.158	ng	97
57) 2,4-Dinitrotoluene	14.940	165	194829	53.723	ng	94
58) Fluorene	15.627	166	597560	42.891	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.204	232	194262	51.908	ng	# 96
60) Diethylphthalate	15.410	149	635221	41.094	ng	98
61) 4-Chlorophenyl-phenyle...	15.627	204	322631	43.773	ng	95
62) 4-Nitroaniline	15.645	138	139279	49.581	ng	95
63) Azobenzene	15.921	77	600156	39.843	ng	95
65) 4,6-Dinitro-2-methylph...	15.704	198	114360	58.376	ng	97
66) n-Nitrosodiphenylamine	15.839	169	566084	43.004	ng	97
67) 4-Bromophenyl-phenylether	16.520	248	227902	48.750	ng	97
68) Hexachlorobenzene	16.632	284	258781	49.036	ng	96
69) Atrazine	16.791	200	218390	47.443	ng	97
70) Pentachlorophenol	16.979	266	341814	99.265	ng	98
71) Phenanthrene	17.372	178	1028859	42.946	ng	99
72) Anthracene	17.461	178	1042601	42.853	ng	99
73) Carbazole	17.725	167	917031	42.750	ng	99
74) Di-n-butylphthalate	18.307	149	1069726	39.999	ng	99
75) Fluoranthene	19.382	202	1248217	42.938	ng	98
77) Benzidine	19.558	184	312596	27.449	ng	99
78) Pyrene	19.740	202	1290140	42.653	ng	99
80) Butylbenzylphthalate	20.645	149	465163	42.158	ng	99
81) Benzo(a)anthracene	21.568	228	1336662	43.845	ng	99
82) 3,3'-Dichlorobenzidine	21.485	252	310743	30.188	ng	# 96
83) Chrysene	21.632	228	1265475	43.067	ng	100
84) Bis(2-ethylhexyl)phtha...	21.509	149	683731	38.878	ng	# 95
85) Di-n-octyl phthalate	22.701	149	1177669	41.105	ng	99
87) Indeno(1,2,3-cd)pyrene	28.283	276	1623465	46.517	ng	98
88) Benzo(b)fluoranthene	23.736	252	1295947m	45.516	ng	
89) Benzo(k)fluoranthene	23.800	252	1315794	45.150	ng	# 98
90) Benzo(a)pyrene	24.576	252	1228228	44.389	ng	98
91) Dibenzo(a,h)anthracene	28.348	278	1325103	45.041	ng	95
92) Benzo(g,h,i)perylene	29.382	276	1235460	43.169	ng	97

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

Manual Integrations

Quant Time: Apr 10 14:33:37 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG032924.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Mar 30 03:12:26 2024
Response via : Initial Calibration

