

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060623\
 Data File : BG057660.D
 Acq On : 7 Jun 2023 7:19
 Operator : CG/JU
 Sample : PB153141BS
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB153141BS

Manual Integrations
 APPROVED

Reviewed By :Yogesh
 Patel

Quant Time: Jun 07 08:02:40 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG052723.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 01 03:51:58 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.336	152	23694	20.000	ng	0.00
21) Naphthalene-d8	11.174	136	104439	20.000	ng	0.00
39) Acenaphthene-d10	14.957	164	80117	20.000	ng	0.00
64) Phenanthrene-d10	17.694	188	201020	20.000	ng	0.00
76) Chrysene-d12	22.006	240	180641	20.000	ng	0.00
86) Perylene-d12	25.490	264	196685	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.857	112	228480	167.528	ng	0.00
7) Phenol-d6	7.490	99	333589	156.293	ng	0.00
23) Nitrobenzene-d5	9.517	82	204577	91.567	ng	0.00
42) 2,4,6-Tribromophenol	16.443	330	160318	124.985	ng	0.00
45) 2-Fluorobiphenyl	13.582	172	513292	92.073	ng	0.00
79) Terphenyl-d14	20.267	244	998811	95.779	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.643	88	20712	36.654	ng	93
3) Pyridine	4.066	79	60347	34.933	ng	99
4) n-Nitrosodimethylamine	3.977	42	29437	42.115	ng	80
6) Aniline	7.649	93	99272	38.447	ng	98
8) 2-Chlorophenol	7.902	128	81285	55.516	ng	93
9) Benzaldehyde	7.467	77	45828	32.408	ng	94
10) Phenol	7.514	94	109766	54.336	ng	95
11) bis(2-Chloroethyl)ether	7.737	93	75397	49.984	ng	97
12) 1,3-Dichlorobenzene	8.225	146	79670	51.074	ng	99
13) 1,4-Dichlorobenzene	8.371	146	82605	52.052	ng	99
14) 1,2-Dichlorobenzene	8.700	146	80985	52.319	ng	94
15) Benzyl Alcohol	8.577	79	81860	49.858	ng	95
16) 2,2'-oxybis(1-Chloropr...	8.859	45	98038	52.341	ng	95
17) 2-Methylphenol	8.783	107	75094	50.619	ng	97
18) Hexachloroethane	9.429	117	29464	54.222	ng	91
19) n-Nitroso-di-n-propyla...	9.141	70	69044	46.562	ng	97
20) 3+4-Methylphenols	9.112	107	103465	49.682	ng	96
22) Acetophenone	9.170	105	128262	45.928	ng	# 97
24) Nitrobenzene	9.564	77	100000	44.094	ng	100
25) Isophorone	10.081	82	187817	42.770	ng	98
26) 2-Nitrophenol	10.281	139	46569	47.892	ng	96
27) 2,4-Dimethylphenol	10.322	122	85094	50.617	ng	96
28) bis(2-Chloroethoxy)met...	10.551	93	109673	48.164	ng	99
29) 2,4-Dichlorophenol	10.827	162	87426	49.514	ng	97
30) 1,2,4-Trichlorobenzene	11.033	180	89579	46.550	ng	97
31) Naphthalene	11.226	128	252608	45.723	ng	97
32) Benzoic acid	10.480	122	31415m	34.255	ng	
33) 4-Chloroaniline	11.332	127	69869	27.501	ng	98
34) Hexachlorobutadiene	11.491	225	58968	44.303	ng	97
35) Caprolactam	12.108	113	30809m	42.090	ng	
36) 4-Chloro-3-methylphenol	12.437	107	97894	46.839	ng	98
37) 2-Methylnaphthalene	12.807	142	187390	43.216	ng	98
38) 1-Methylnaphthalene	13.024	142	176477	42.649	ng	95
40) 1,2,4,5-Tetrachloroben...	13.171	216	120900	45.975	ng	99
41) Hexachlorocyclopentadiene	13.136	237	63535	88.959	ng	98
43) 2,4,6-Trichlorophenol	13.406	196	75351	45.514	ng	96

06/09/2023
 Supervised By :mohammad
 Ahmed

06/13/2023

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060623\
 Data File : BG057660.D
 Acq On : 7 Jun 2023 7:19
 Operator : CG/JU
 Sample : PB153141BS
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

PB153141BS

Manual Integrations

APPROVED

Reviewed By :Yogesh
 Patel

Quant Time: Jun 07 08:02:40 2023

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG052723.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Thu Jun 01 03:51:58 2023

Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
44) 2,4,5-Trichlorophenol	13.494	196	80772	40.964	ng	95	06/09/2023
46) 1,1'-Biphenyl	13.794	154	268112	46.823	ng	97	Supervised By :mohammad
47) 2-Chloronaphthalene	13.846	162	202160	46.872	ng	98	ahmed
48) 2-Nitroaniline	14.046	65	67210	44.487	ng	97	
49) Acenaphthylene	14.686	152	327452	44.835	ng	100	
50) Dimethylphthalate	14.393	163	278167	42.126	ng	100	
51) 2,6-Dinitrotoluene	14.528	165	63443	44.884	ng	96	06/13/2023
52) Acenaphthene	15.021	154	202352	45.279	ng	98	
53) 3-Nitroaniline	14.869	138	45846	31.097	ng	95	
54) 2,4-Dinitrophenol	15.086	184	62404	73.850	ng	96	
55) Dibenzofuran	15.350	168	334104	43.753	ng	100	
56) 4-Nitrophenol	15.180	139	80589	82.041	ng	94	
57) 2,4-Dinitrotoluene	15.321	165	92400	43.659	ng	98	
58) Fluorene	15.996	166	279083	43.286	ng	99	
59) 2,3,4,6-Tetrachlorophenol	15.579	232	76505	41.416	ng	97	
60) Diethylphthalate	15.738	149	293142	42.534	ng	99	
61) 4-Chlorophenyl-phenyle...	15.973	204	155169	43.184	ng	97	
62) 4-Nitroaniline	16.026	138	66666	41.919	ng	89	
63) Azobenzene	16.267	77	272437	44.449	ng	97	
65) 4,6-Dinitro-2-methylph...	16.090	198	56382	46.065	ng	96	
66) n-Nitrosodiphenylamine	16.190	169	249457	47.745	ng	98	
67) 4-Bromophenyl-phenylether	16.866	248	104317	46.262	ng	99	
68) Hexachlorobenzene	17.007	284	113547	48.851	ng	98	
69) Atrazine	17.124	200	111599	49.750	ng	99	
70) Pentachlorophenol	17.353	266	82435	69.604	ng	97	
71) Phenanthrene	17.741	178	473110	45.991	ng	98	
72) Anthracene	17.835	178	483038	45.773	ng	98	
73) Carbazole	18.099	167	438417	46.440	ng	99	
74) Di-n-butylphthalate	18.611	149	513924	45.857	ng	99	
75) Fluoranthene	19.733	202	582426	43.613	ng	99	
77) Benzidine	19.897	184	277721	50.807	ng	99	
78) Pyrene	20.091	202	595472	46.297	ng	99	
80) Butylbenzylphthalate	20.937	149	225938	48.233	ng	99	
81) Benzo(a)anthracene	21.982	228	582127	46.450	ng	99	
82) 3,3'-Dichlorobenzidine	21.877	252	163761	35.435	ng	#	99
83) Chrysene	22.053	228	533200	45.022	ng	99	
84) Bis(2-ethylhexyl)phtha...	21.818	149	325961	48.783	ng	98	
85) Di-n-octyl phthalate	23.104	149	553241	49.952	ng	99	
87) Indeno(1,2,3-cd)pyrene	29.537	276	655720	47.817	ng	98	
88) Benzo(b)fluoranthene	24.373	252	586673	48.459	ng	98	
89) Benzo(k)fluoranthene	24.450	252	582880	47.706	ng	98	
90) Benzo(a)pyrene	25.331	252	520006	44.044	ng	98	
91) Dibenzo(a,h)anthracene	29.590	278	546468	47.883	ng	98	
92) Benzo(g,h,i)perylene	30.812	276	530569	47.669	ng	98	

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

Manual Integrations APPROVED

Quant Time: Jun 07 08:02:40 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG052723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 01 03:51:58 2023
Response via : Initial Calibration

