

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091922\
 Data File : BG054969.D
 Acq On : 19 Sep 2022 12:10
 Operator : CG/JU
 Sample : SSTDICC005
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

SSTDICC005

Manual Integrations

APPROVED

Reviewed By :Christian Giraldo 09/20/2022
 Supervised By :Jagrut Upadhyay 09/20/2022

Quant Time: Sep 19 16:03:14 2022

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG091922.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Mon Sep 19 16:00:32 2022

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.061	152	31845	20.000	ng	0.00
21) Naphthalene-d8	10.887	136	123207	20.000	ng	# 0.00
39) Acenaphthene-d10	14.706	164	86214	20.000	ng	0.00
64) Phenanthrene-d10	17.456	188	217033	20.000	ng	0.00
76) Chrysene-d12	21.733	240	236263	20.000	ng	0.00
86) Perylene-d12	25.035	264	250601	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.599	112	16432	8.985	ng	0.00
7) Phenol-d6	7.221	99	20800	8.317	ng	0.00
23) Nitrobenzene-d5	9.236	82	20669	8.807	ng	0.00
42) 2,4,6-Tribromophenol	16.198	330	8939	8.298	ng	0.00
45) 2-Fluorobiphenyl	13.325	172	58381	10.178	ng	0.00
79) Terphenyl-d14	20.053	244	121958	10.228	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.407	88	4491m	5.143	ng	
3) Pyridine	3.836	79	10596m	4.350	ng	
4) n-Nitrosodimethylamine	3.748	42	3874	3.679	ng	# 84
6) Aniline	7.379	93	13396	4.508	ng	93
8) 2-Chlorophenol	7.626	128	9613	4.811	ng	97
9) Benzaldehyde	7.197	77	8112	4.949	ng	97
10) Phenol	7.250	94	11454	4.453	ng	89
11) bis(2-Chloroethyl)ether	7.473	93	9935	4.803	ng	88
12) 1,3-Dichlorobenzene	7.949	146	11851	5.125	ng	93
13) 1,4-Dichlorobenzene	8.096	146	12031	5.159	ng	93
14) 1,2-Dichlorobenzene	8.419	146	11182	5.056	ng	96
15) Benzyl Alcohol	8.308	79	7215	3.993	ng	95
16) 2,2'-oxybis(1-Chloropr...	8.584	45	11455	3.539	ng	93
17) 2-Methylphenol	8.513	107	7728	4.364	ng	97
18) Hexachloroethane	9.154	117	3993	4.721	ng	92
19) n-Nitroso-di-n-propyla...	8.866	70	7169	4.168	ng	96
20) 3+4-Methylphenols	8.848	107	10997	4.478	ng	94
22) Acetophenone	8.895	105	14664	4.757	ng	# 91
24) Nitrobenzene	9.283	77	10266	4.340	ng	93
25) Isophorone	9.800	82	20130	4.600	ng	98
26) 2-Nitrophenol	10.006	139	5170	4.948	ng	98
27) 2,4-Dimethylphenol	10.053	122	7446	4.497	ng	98
28) bis(2-Chloroethoxy)met...	10.282	93	11225	4.502	ng	# 97
29) 2,4-Dichlorophenol	10.546	162	8352	4.435	ng	97
30) 1,2,4-Trichlorobenzene	10.746	180	10784	5.011	ng	95
31) Naphthalene	10.940	128	32827	4.966	ng	98
33) 4-Chloroaniline	11.051	127	12306	4.566	ng	96
34) Hexachlorobutadiene	11.210	225	7515	5.450	ng	97
35) Caprolactam	11.809	113	3100	4.666	ng	# 74
36) 4-Chloro-3-methylphenol	12.174	107	9123	4.430	ng	98
37) 2-Methylnaphthalene	12.544	142	22241	4.800	ng	94
38) 1-Methylnaphthalene	12.761	142	22195	4.924	ng	97
40) 1,2,4,5-Tetrachloroben...	12.908	216	12704	4.888	ng	# 97
43) 2,4,6-Trichlorophenol	13.155	196	6384	3.701	ng	95
44) 2,4,5-Trichlorophenol	13.231	196	7350	3.794	ng	98
46) 1,1'-Biphenyl	13.543	154	30009	4.687	ng	97

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091922\
 Data File : BG054969.D
 Acq On : 19 Sep 2022 12:10
 Operator : CG/JU
 Sample : SSTDICC005
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICC005

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 09/20/2022
 Supervised By : Jagrut Upadhyay 09/20/2022

Quant Time: Sep 19 16:03:14 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG091922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 19 16:00:32 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
47) 2-Chloronaphthalene	13.590	162	23146	4.760	ng	98
48) 2-Nitroaniline	13.795	65	6113	4.012	ng	97
49) Acenaphthylene	14.436	152	39726	4.951	ng	99
50) Dimethylphthalate	14.154	163	34275	5.160	ng	# 99
51) 2,6-Dinitrotoluene	14.283	165	6775	4.987	ng	86
52) Acenaphthene	14.771	154	23527	4.565	ng	99
53) 3-Nitroaniline	14.618	138	7154	4.704	ng	91
55) Dibenzofuran	15.105	168	39038	5.155	ng	98
57) 2,4-Dinitrotoluene	15.076	165	9231	4.912	ng	# 81
58) Fluorene	15.752	166	32968	5.167	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.346	232	5714	3.272	ng	# 86
60) Diethylphthalate	15.505	149	34550	5.209	ng	98
61) 4-Chlorophenyl-phenyle...	15.740	204	16466	5.232	ng	94
62) 4-Nitroaniline	15.781	138	7231	4.657	ng	90
63) Azobenzene	16.028	77	26568	4.215	ng	94
66) n-Nitrosodiphenylamine	15.952	169	29018	4.636	ng	94
67) 4-Bromophenyl-phenylether	16.633	248	11795	4.950	ng	95
68) Hexachlorobenzene	16.756	284	13894	5.187	ng	# 93
69) Atrazine	16.898	200	11978	5.425	ng	97
71) Phenanthrene	17.497	178	57925	5.010	ng	97
72) Anthracene	17.585	178	55883	4.972	ng	97
73) Carbazole	17.861	167	51373	4.961	ng	97
74) Di-n-butylphthalate	18.396	149	65915	5.030	ng	98
75) Fluoranthene	19.506	202	77437	5.506	ng	94
77) Benzidine	19.683	184	36768	6.283	ng	99
78) Pyrene	19.865	202	78752	4.780	ng	99
80) Butylbenzylphthalate	20.734	149	29705	4.323	ng	93
81) Benzo(a)anthracene	21.710	228	78862	4.923	ng	98
82) 3,3'-Dichlorobenzidine	21.621	252	26103	4.648	ng	# 96
83) Chrysene	21.780	228	75878	4.954	ng	98
84) Bis(2-ethylhexyl)phtha...	21.586	149	42723	4.315	ng	98
85) Di-n-octyl phthalate	22.820	149	75859	4.518	ng	# 91
87) Indeno(1,2,3-cd)pyrene	28.854	276	91827	4.827	ng	# 94
88) Benzo(b)fluoranthene	23.977	252	73347	4.666	ng	98
89) Benzo(k)fluoranthene	24.048	252	76232	4.818	ng	97
90) Benzo(a)pyrene	24.876	252	63871	4.756	ng	96
91) Dibenzo(a,h)anthracene	28.907	278	76189	4.916	ng	99
92) Benzo(g,h,i)perylene	30.047	276	76810	4.905	ng	# 92

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

Manual Integrations

Quant Time: Sep 19 16:03:14 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG091922.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Sep 19 16:00:32 2022
Response via : Initial Calibration

