

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG041823\
 Data File : BG057133.D
 Acq On : 18 Apr 2023 20:28
 Operator : CG/JU
 Sample : SSTDICC080
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICC080

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 04/19/2023
 Supervised By :Jagrut Upadhyay 04/19/2023

Quant Time: Apr 19 04:14:10 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG041823.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 19 03:30:08 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.396	152	10278	20.000	ng	0.00
21) Naphthalene-d8	11.239	136	63851	20.000	ng	0.00
39) Acenaphthene-d10	15.010	164	42566	20.000	ng	0.00
64) Phenanthrene-d10	17.748	188	98663	20.000	ng	# 0.00
76) Chrysene-d12	22.071	240	83511	20.000	ng	0.00
86) Perylene-d12	25.596	264	88857	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.905	112	87954	145.744	ng	0.00
7) Phenol-d6	7.538	99	132816	148.882	ng	0.00
23) Nitrobenzene-d5	0.000	82	0d	0.000	ng	
42) 2,4,6-Tribromophenol	16.497	330	108073	149.656	ng	0.00
45) 2-Fluorobiphenyl	13.642	172	501018	155.072	ng	0.00
79) Terphenyl-d14	20.321	244	827740	149.952	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.714	88	29636	80.644	ng	98
3) Pyridine	4.137	79	53045m	70.123	ng	
4) n-Nitrosodimethylamine	4.043	42	28851	76.247	ng	# 81
6) Aniline	7.708	93	84571	75.430	ng	99
8) 2-Chlorophenol	7.955	128	47472	77.752	ng	98
10) Phenol	7.562	94	67018	77.514	ng	95
11) bis(2-Chloroethyl)ether	7.797	93	46332	77.484	ng	90
12) 1,3-Dichlorobenzene	8.284	146	51537	73.006	ng	98
13) 1,4-Dichlorobenzene	8.431	146	53109	73.036	ng	99
14) 1,2-Dichlorobenzene	8.760	146	52004	70.261	ng	99
15) Benzyl Alcohol	8.637	79	56779	73.041	ng	98
16) 2,2'-oxybis(1-Chloropr...	8.913	45	52806m	59.112	ng	
17) 2-Methylphenol	8.830	107	49887	65.752	ng	# 94
18) Hexachloroethane	9.500	117	19288	64.572	ng	86
19) n-Nitroso-di-n-propyla...	9.206	70	47388	59.196	ng	# 89
20) 3+4-Methylphenols	9.165	107	69173	63.712	ng	98
29) 2,4-Dichlorophenol	10.875	162	73640	69.042	ng	97
30) 1,2,4-Trichlorobenzene	11.098	180	91244	76.738	ng	98
31) Naphthalene	11.292	128	261738	77.061	ng	99
32) Benzoic acid	10.552	122	34895m	71.414	ng	
33) 4-Chloroaniline	11.386	127	119064	77.298	ng	97
34) Hexachlorobutadiene	11.562	225	64731	74.840	ng	98
35) Caprolactam	12.179	113	27980	76.679	ng	98
36) 4-Chloro-3-methylphenol	12.484	107	91768	74.721	ng	99
37) 2-Methylnaphthalene	12.866	142	198926	76.891	ng	98
38) 1-Methylnaphthalene	13.084	142	185150	74.947	ng	99
40) 1,2,4,5-Tetrachloroben...	13.224	216	118640	82.808	ng	98
43) 2,4,6-Trichlorophenol	13.459	196	74802	79.260	ng	88
44) 2,4,5-Trichlorophenol	13.536	196	88974	80.696	ng	97
46) 1,1'-Biphenyl	13.853	154	256780	77.133	ng	99
47) 2-Chloronaphthalene	13.906	162	192858	80.602	ng	97
48) 2-Nitroaniline	14.100	65	67087	83.005	ng	98
49) Acenaphthylene	14.740	152	310531	80.470	ng	99
50) Dimethylphthalate	14.452	163	262053	76.139	ng	98
51) 2,6-Dinitrotoluene	14.581	165	57210	78.307	ng	96
52) Acenaphthene	15.081	154	187931	76.599	ng	99

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG041823\
 Data File : BG057133.D
 Acq On : 18 Apr 2023 20:28
 Operator : CG/JU
 Sample : SSTDICC080
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICC080

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 04/19/2023
 Supervised By :Jagrut Upadhyay 04/19/2023

Quant Time: Apr 19 04:14:10 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG041823.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 19 03:30:08 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
53) 3-Nitroaniline	14.916	138	58824	78.423	ng	90
54) 2,4-Dinitrophenol	15.116	184	37075	90.640	ng	97
55) Dibenzofuran	15.410	168	307775	78.000	ng	97
56) 4-Nitrophenol	15.204	139	42845	85.775	ng	98
57) 2,4-Dinitrotoluene	15.363	165	81487	79.123	ng	# 99
58) Fluorene	16.056	166	251615	73.218	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.627	232	74484	81.605	ng	# 95
60) Diethylphthalate	15.797	149	263180	75.204	ng	98
61) 4-Chlorophenyl-phenyle...	16.032	204	145773	73.384	ng	97
62) 4-Nitroaniline	16.074	138	59823	79.233	ng	95
63) Azobenzene	16.326	77	259727	80.339	ng	97
65) 4,6-Dinitro-2-methylph...	16.126	198	54723	81.883	ng	97
66) n-Nitrosodiphenylamine	16.244	169	218756	74.657	ng	98
67) 4-Bromophenyl-phenylether	16.925	248	94040	75.241	ng	98
68) Hexachlorobenzene	17.060	284	99310	74.485	ng	97
69) Atrazine	17.178	200	78905	71.659	ng	100
70) Pentachlorophenol	17.395	266	61705	87.418	ng	93
71) Phenanthrene	17.795	178	393911	76.984	ng	100
72) Anthracene	17.889	178	402724	76.206	ng	99
73) Carbazole	18.147	167	347698	78.608	ng	99
74) Di-n-butylphthalate	18.670	149	376344	72.608	ng	98
75) Fluoranthene	19.780	202	472700	78.685	ng	97
77) Benzidine	19.945	184	136978	59.748	ng	98
78) Pyrene	20.145	202	470128	75.194	ng	99
80) Butylbenzylphthalate	20.996	149	182795	79.559	ng	95
81) Benzo(a)anthracene	22.048	228	459825	77.041	ng	99
82) 3,3'-Dichlorobenzidine	21.942	252	151657	74.845	ng	99
83) Chrysene	22.124	228	432269	78.409	ng	99
84) Bis(2-ethylhexyl)phtha...	21.895	149	255657	81.108	ng	99
85) Di-n-octyl phthalate	23.199	149	422136	79.377	ng	100
87) Indeno(1,2,3-cd)pyrene	29.702	276	526501	79.553	ng	99
88) Benzo(b)fluoranthene	24.468	252	443340	78.019	ng	99
89) Benzo(k)fluoranthene	24.544	252	452856	78.767	ng	99
90) Benzo(a)pyrene	25.437	252	434077	79.125	ng	95
91) Dibenzo(a,h)anthracene	29.755	278	440048	79.232	ng	100
92) Benzo(g,h,i)perylene	30.977	276	421687	79.536	ng	99

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

