

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG051822\
 Data File : BG053615.D
 Acq On : 18 May 2022 22:42
 Operator : CG/JU
 Sample : N2907-02MSD
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 GC-2MSD

Quant Time: May 19 05:13:23 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG050522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 11 06:00:46 2022
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Christian Giraldo 05/19/2022
 Supervised By :Jagrut Upadhyay 05/19/2022

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.076	152	22685	20.000	ng	-0.01
21) Naphthalene-d8	10.890	136	95806	20.000	ng	0.00
39) Acenaphthene-d10	14.708	164	77286	20.000	ng	0.00
64) Phenanthrene-d10	17.452	188	193620	20.000	ng	0.00
76) Chrysene-d12	21.734	240	190777	20.000	ng	0.00
86) Perylene-d12	25.006	264	199231	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.644	112	118277	88.487	ng	0.00
7) Phenol-d6	7.242	99	176397	86.955	ng	0.00
23) Nitrobenzene-d5	9.239	82	123018	54.659	ng	-0.01
42) 2,4,6-Tribromophenol	16.195	330	98892	86.757	ng	0.00
45) 2-Fluorobiphenyl	13.328	172	269457	50.947	ng	0.00
79) Terphenyl-d14	20.054	244	553000	53.987	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.518	88	24369	34.747	ng	# 90
3) Pyridine	3.923	79	60735	35.460	ng	95
4) n-Nitrosodimethylamine	3.835	42	40470	48.506	ng	# 94
6) Aniline	7.401	93	105149	46.799	ng	98
8) 2-Chlorophenol	7.647	128	70519	50.466	ng	95
9) Benzaldehyde	7.213	77	60244m	46.746	ng	
10) Phenol	7.271	94	103080	51.955	ng	96
11) bis(2-Chloroethyl)ether	7.489	93	69451	46.936	ng	98
12) 1,3-Dichlorobenzene	7.971	146	75474	45.784	ng	95
13) 1,4-Dichlorobenzene	8.112	146	77808	46.774	ng	96
14) 1,2-Dichlorobenzene	8.435	146	73904	47.311	ng	98
15) Benzyl Alcohol	8.317	79	75747m	44.663	ng	
16) 2,2'-oxybis(1-Chloropr...	8.599	45	109802	45.593	ng	97
17) 2-Methylphenol	8.523	107	66784	49.802	ng	94
18) Hexachloroethane	9.163	117	28371	44.912	ng	95
19) n-Nitroso-di-n-propyla...	8.881	70	66909	45.970	ng	96
20) 3+4-Methylphenols	8.852	107	94246	49.763	ng	97
22) Acetophenone	8.899	105	116211	44.320	ng	# 97
24) Nitrobenzene	9.281	77	100861	46.302	ng	97
25) Isophorone	9.803	82	189864	50.126	ng	99
26) 2-Nitrophenol	9.997	139	41091	46.574	ng	95
27) 2,4-Dimethylphenol	10.056	122	74035	55.760	ng	94
28) bis(2-Chloroethoxy)met...	10.285	93	98805	45.356	ng	99
29) 2,4-Dichlorophenol	10.544	162	81225	50.687	ng	94
30) 1,2,4-Trichlorobenzene	10.749	180	81749	44.900	ng	98
31) Naphthalene	10.937	128	219787	45.099	ng	99
32) Benzoic acid	10.232	122	32668m	40.888	ng	
33) 4-Chloroaniline	11.049	127	43082	21.109	ng	96
34) Hexachlorobutadiene	11.225	225	60874	45.089	ng	96
35) Caprolactam	11.824	113	29437m	50.550	ng	
36) 4-Chloro-3-methylphenol	12.177	107	96183	50.178	ng	100
37) 2-Methylnaphthalene	12.541	142	164585	45.254	ng	98
38) 1-Methylnaphthalene	12.758	142	159440m	44.933	ng	
40) 1,2,4,5-Tetrachloroben...	12.911	216	109202	44.826	ng	99
41) Hexachlorocyclopentadiene	12.887	237	95285	94.254	ng	98
43) 2,4,6-Trichlorophenol	13.146	196	76413	48.052	ng	95

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG051822\
 Data File : BG053615.D
 Acq On : 18 May 2022 22:42
 Operator : CG/JU
 Sample : N2907-02MSD
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 GC-2MSD

Quant Time: May 19 05:13:23 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG050522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 11 06:00:46 2022
 Response via : Initial Calibration

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 05/19/2022
 Supervised By : Jagrut Upadhyay 05/19/2022

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.234	196	84281	49.798	ng	98
46) 1,1'-Biphenyl	13.539	154	236634	44.918	ng	98
47) 2-Chloronaphthalene	13.586	162	182011	44.098	ng	99
48) 2-Nitroaniline	13.786	65	76704	49.874	ng	97
49) Acenaphthylene	14.432	152	320866	48.717	ng	98
50) Dimethylphthalate	14.156	163	268755	45.761	ng	99
51) 2,6-Dinitrotoluene	14.280	165	59225	49.498	ng	98
52) Acenaphthene	14.773	154	176844	45.055	ng	97
53) 3-Nitroaniline	14.609	138	41635	32.205	ng	95
54) 2,4-Dinitrophenol	14.826	184	51910	69.256	ng	96
55) Dibenzofuran	15.102	168	315124	44.697	ng	99
56) 4-Nitrophenol	14.932	139	88475	95.814	ng	87
57) 2,4-Dinitrotoluene	15.073	165	85456	49.272	ng	# 93
58) Fluorene	15.754	166	259306	45.832	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.337	232	78919	45.959	ng	98
60) Diethylphthalate	15.513	149	281391	46.383	ng	99
61) 4-Chlorophenyl-phenyle...	15.736	204	144427	45.747	ng	98
62) 4-Nitroaniline	15.778	138	62486	46.599	ng	98
63) Azobenzene	16.030	77	284836	45.851	ng	99
65) 4,6-Dinitro-2-methylph...	15.836	198	48868	42.357	ng	96
66) n-Nitrosodiphenylamine	15.954	169	237522	46.724	ng	99
67) 4-Bromophenyl-phenylether	16.635	248	104719	46.310	ng	92
68) Hexachlorobenzene	16.765	284	114410	46.198	ng	97
69) Atrazine	16.900	200	102369	47.834	ng	96
70) Pentachlorophenol	17.111	266	119142	88.791	ng	91
71) Phenanthrene	17.493	178	455755	46.292	ng	97
72) Anthracene	17.587	178	462300	47.773	ng	98
73) Carbazole	17.857	167	424412	44.750	ng	98
74) Di-n-butylphthalate	18.403	149	495535	46.458	ng	99
75) Fluoranthene	19.502	202	582222	45.673	ng	99
77) Benzidine	19.678	184	351426	85.465	ng	99
78) Pyrene	19.866	202	581344	47.073	ng	99
80) Butylbenzylphthalate	20.736	149	213839	47.219	ng	96
81) Benzo(a)anthracene	21.717	228	566241	46.535	ng	99
82) 3,3'-Dichlorobenzidine	21.623	252	134713	43.657	ng	# 95
83) Chrysene	21.781	228	517002	45.492	ng	97
84) Bis(2-ethylhexyl)phtha...	21.605	149	304726	48.751	ng	99
85) Di-n-octyl phthalate	22.827	149	510767	48.324	ng	100
87) Indeno(1,2,3-cd)pyrene	28.772	276	636310	45.857	ng	# 95
88) Benzo(b)fluoranthene	23.972	252	599004	50.794	ng	100
89) Benzo(k)fluoranthene	24.037	252	544344	46.973	ng	99
90) Benzo(a)pyrene	24.859	252	554586	55.252	ng	100
91) Dibenzo(a,h)anthracene	28.825	278	548073	47.935	ng	98
92) Benzo(g,h,i)perylene	29.947	276	543425	47.948	ng	97

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

