

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082321\
 Data File : BG049941.D
 Acq On : 23 Aug 2021 20:15
 Operator : CG/JU
 Sample : M3463-11MSD
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SWCS-20-0306MSD

Manual Integrations
 APPROVED

mohammad
 8/24/2021 11:19:12 AM

Quant Time: Aug 24 02:28:17 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG080321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 23 16:13:24 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.980	152	29842	20.000	ng	0.00
21) Naphthalene-d8	10.800	136	122950	20.000	ng	-0.01
39) Acenaphthene-d10	14.636	164	83666	20.000	ng	-0.01
64) Phenanthrene-d10	17.391	188	175146	20.000	ng	#-0.01
76) Chrysene-d12	21.668	240	148312	20.000	ng	-0.01
86) Perylene-d12	24.905	264	147373	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.537	112	149685	89.916	ng	0.00
7) Phenol-d6	7.152	99	204228	81.044	ng	0.00
23) Nitrobenzene-d5	9.155	82	197453	71.066	ng	0.00
42) 2,4,6-Tribromophenol	16.134	330	91306	74.246	ng	-0.01
45) 2-Fluorobiphenyl	13.256	172	345847	60.148	ng	0.00
79) Terphenyl-d14	19.999	244	417589	51.137	ng	-0.01
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.369	88	24248	33.332	ng	96
3) Pyridine	3.780	79	59610	29.893	ng	88
4) n-Nitrosodimethylamine	3.698	42	44519	41.519	ng	# 82
6) Aniline	7.299	93	110928	41.535	ng	94
8) 2-Chlorophenol	7.546	128	93141	51.317	ng	94
9) Benzaldehyde	7.111	77	65760	39.109	ng	# 78
10) Phenol	7.182	94	120797	49.472	ng	92
11) bis(2-Chloroethyl)ether	7.393	93	76994	46.702	ng	96
12) 1,3-Dichlorobenzene	7.869	146	97024	46.675	ng	93
13) 1,4-Dichlorobenzene	8.016	146	97767	46.464	ng	95
14) 1,2-Dichlorobenzene	8.333	146	94031	47.173	ng	94
15) Benzyl Alcohol	8.227	79	102312	48.345	ng	96
16) 2,2'-oxybis(1-Chloropr...	8.503	45	86287	46.062	ng	95
17) 2-Methylphenol	8.439	107	83033	51.641	ng	96
18) Hexachloroethane	9.061	117	38618	47.744	ng	93
19) n-Nitroso-di-n-propyla...	8.791	70	70929	42.061	ng	94
20) 3+4-Methylphenols	8.768	107	112255	49.711	ng	98
22) Acetophenone	8.809	105	149138	44.648	ng	# 93
24) Nitrobenzene	9.196	77	117681	40.139	ng	92
25) Isophorone	9.719	82	195607	39.453	ng	97
26) 2-Nitrophenol	9.907	139	53569	47.877	ng	99
27) 2,4-Dimethylphenol	9.978	122	90565	54.765	ng	92
28) bis(2-Chloroethoxy)met...	10.201	93	106054	48.899	ng	91
29) 2,4-Dichlorophenol	10.454	162	96905	47.752	ng	95
30) 1,2,4-Trichlorobenzene	10.659	180	101364	45.364	ng	97
31) Naphthalene	10.853	128	279309	43.377	ng	98
32) Benzoic acid	10.154	122	47497	41.428	ng	# 74
33) 4-Chloroaniline	10.965	127	72911	28.681	ng	94
34) Hexachlorobutadiene	11.129	225	70714	43.550	ng	96
35) Caprolactam	11.752	113	28952m	46.174	ng	
36) 4-Chloro-3-methylphenol	12.104	107	110577	47.172	ng	# 92
37) 2-Methylnaphthalene	12.463	142	216680	44.670	ng	97
38) 1-Methylnaphthalene	12.680	142	195295	42.839	ng	97
40) 1,2,4,5-Tetrachloroben...	12.833	216	117720	47.753	ng	98
41) Hexachlorocyclopentadiene	12.803	237	89863	68.262	ng	94
43) 2,4,6-Trichlorophenol	13.079	196	80427	48.393	ng	91

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082321\
 Data File : BG049941.D
 Acq On : 23 Aug 2021 20:15
 Operator : CG/JU
 Sample : M3463-11MSD
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SWCS-20-0306MSD

Manual Integrations
 APPROVED

mohammad
 8/24/2021 11:19:12 AM

Quant Time: Aug 24 02:28:17 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG080321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 23 16:13:24 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.156	196	89251	49.501	ng	# 86
46) 1,1'-Biphenyl	13.467	154	268671	45.087	ng	98
47) 2-Chloronaphthalene	13.514	162	219091	46.032	ng	98
48) 2-Nitroaniline	13.720	65	78815	45.098	ng	92
49) Acenaphthylene	14.360	152	344592	45.681	ng	96
50) Dimethylphthalate	14.090	163	285555	45.237	ng	97
51) 2,6-Dinitrotoluene	14.213	165	62434	44.922	ng	90
52) Acenaphthene	14.707	154	206333	45.406	ng	95
53) 3-Nitroaniline	14.548	138	46217	34.099	ng	# 94
54) 2,4-Dinitrophenol	14.771	184	75481	95.663	ng	94
55) Dibenzofuran	15.041	168	323178	43.839	ng	99
56) 4-Nitrophenol	14.883	139	85071	85.913	ng	93
57) 2,4-Dinitrotoluene	15.012	165	88894	45.538	ng	92
58) Fluorene	15.688	166	268956	44.887	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.276	232	72532	43.313	ng	96
60) Diethylphthalate	15.453	149	286585	45.140	ng	98
61) 4-Chlorophenyl-phenyle...	15.676	204	145182	45.397	ng	96
62) 4-Nitroaniline	15.717	138	61542	44.141	ng	89
63) Azobenzene	15.970	77	274490	40.848	ng	87
65) 4,6-Dinitro-2-methylph...	15.782	198	49136	43.645	ng	91
66) n-Nitrosodiphenylamine	15.893	169	233418	47.166	ng	96
67) 4-Bromophenyl-phenylether	16.575	248	100147	48.550	ng	95
68) Hexachlorobenzene	16.698	284	109350	47.715	ng	96
69) Atrazine	16.845	200	99748	50.388	ng	99
70) Pentachlorophenol	17.051	266	91142	74.168	ng	94
71) Phenanthrene	17.438	178	419939	45.573	ng	96
72) Anthracene	17.526	178	421207	46.544	ng	99
73) Carbazole	17.797	167	366366	42.745	ng	97
74) Di-n-butylphthalate	18.343	149	450794	45.118	ng	98
75) Fluoranthene	19.447	202	472426	41.663	ng	99
77) Benzidine	19.624	184	265720	69.925	ng	98
78) Pyrene	19.811	202	468124	46.298	ng	97
80) Butylbenzylphthalate	20.687	149	177520	46.100	ng	97
81) Benzo(a)anthracene	21.650	228	419507	43.433	ng	98
82) 3,3'-Dichlorobenzidine	21.562	252	117948	41.824	ng	97
83) Chrysene	21.715	228	402860	43.634	ng	98
84) Bis(2-ethylhexyl)phtha...	21.539	149	240016	44.032	ng	94
85) Di-n-octyl phthalate	22.749	149	393830	42.769	ng	98
87) Indeno(1,2,3-cd)pyrene	28.617	276	493164	46.401	ng	98
88) Benzo(b)fluoranthene	23.877	252	422741m	43.737	ng	
89) Benzo(k)fluoranthene	23.941	252	413773	45.895	ng	98
90) Benzo(a)pyrene	24.752	252	407448	46.581	ng	# 98
91) Dibenzo(a,h)anthracene	28.670	278	421750	47.091	ng	# 97
92) Benzo(g,h,i)perylene	29.786	276	414955	47.292	ng	97

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082321\
Data File : BG049941.D
Acq On : 23 Aug 2021 20:15
Operator : CG/JU
Sample : M3463-11MSD
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SWCS-20-0306MSD

Manual Integrations
APPROVED

mohammad
8/24/2021 11:19:12 AM

Quant Time: Aug 24 02:28:17 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG080321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Aug 23 16:13:24 2021
Response via : Initial Calibration

