Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG110821\

Data File : BG050953.D

Acq On : 10 Nov 2021 21:45

Operator : CG/JU Sample : SSTDCCC020

Misc

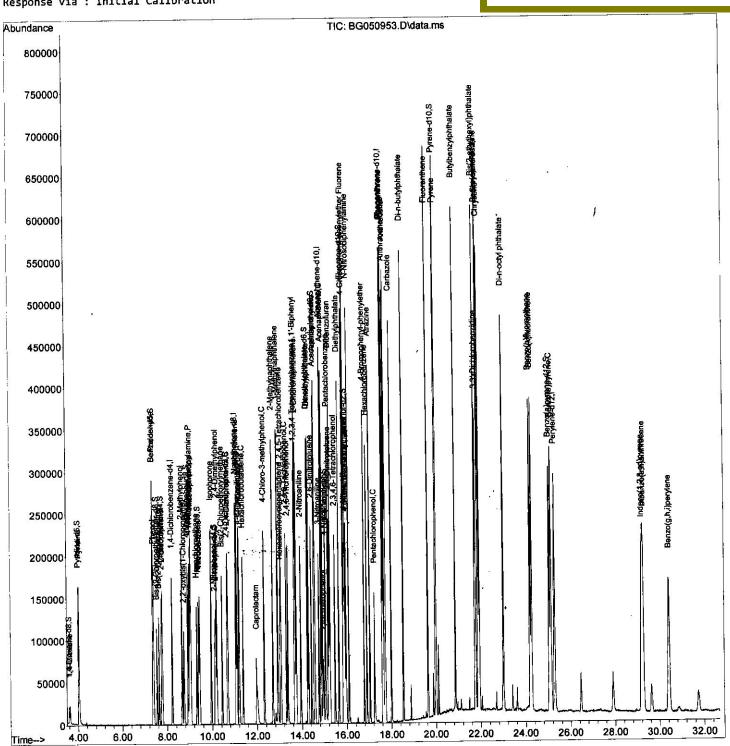
ALS Vial : 31 Sample Multiplier: 1

Quant Time: Nov 11 01:45:02 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG110321.M

Quant Title : SVOA CALIBRATION QLast Update : Tue Nov 02 14:49:05 2021 Response via : Initial Calibration Instrument :
BNA\_G
LabSampleId :
SSTDCCC020

#### **Manual IntegrationsAPPROVED**



SFAM-EPA-BG110321.M Thu Nov 11 01:47:25 2021

# Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG110821\

Data File: BG050953.D

Acq On : 10 Nov 2021 21:45

Operator : CG/JU Sample : SSTDCCC020

Misc

ALS Vial : 31 Sample Multiplier: 1

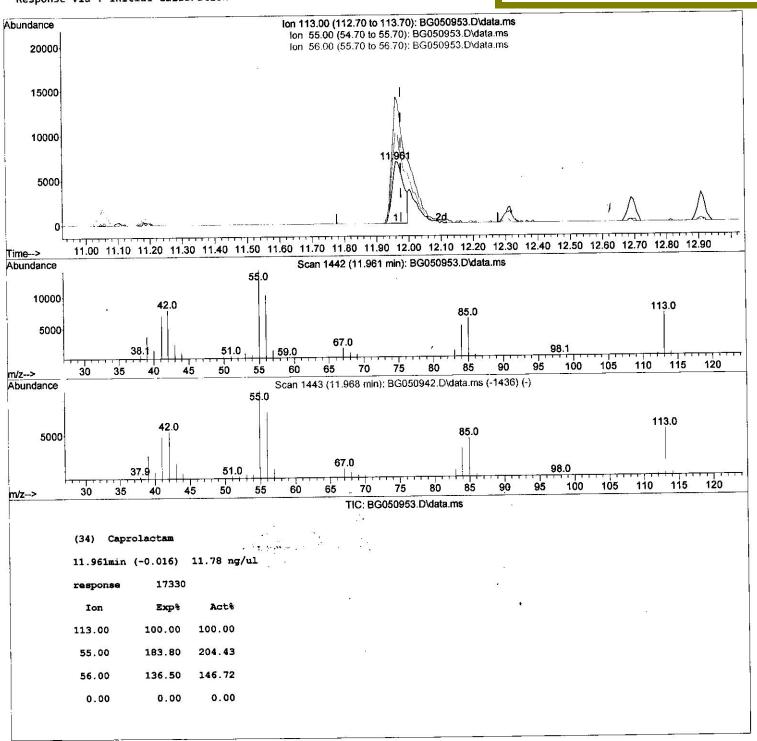
Quant Time: Nov 11 01:45:02 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG110321.M

Quant Title : SVOA CALIBRATION
QLast Update : Tue Nov 02 14:49:05 2021
Response via : Initial Calibration

Instrument :
BNA\_G
LabSampleId :
SSTDCCC020

# **Manual IntegrationsAPPROVED**



## Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG110821\

Data File : BG050953.D

Acq On : 10 Nov 2021 21:45

Operator : CG/JU Sample : SSTDCCC020

Misc

ALS Vial : 31 Sample Multiplier: 1

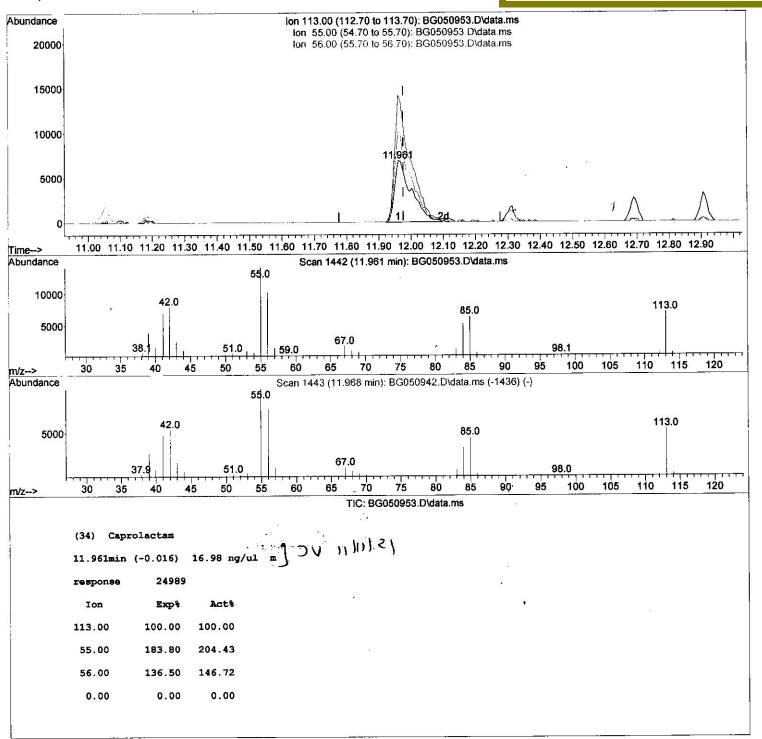
Quant Time: Nov 11 01:45:02 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG110321.M

Quant Title : SVOA CALIBRATION
QLast Update : Tue Nov 02 14:49:05 2021
Response via : Initial Calibration

Instrument :
BNA\_G
LabSampleId :
SSTDCCC020

#### **Manual IntegrationsAPPROVED**



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG110821\

Data File : BG050953.D

Acq On : 10 Nov 2021 21:45

Operator : CG/JU : SSTDCCC020 Sample

Misc

ALS Vial : 31 Sample Multiplier: 1

Quant Time: Nov 11 01:45:02 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG110321.M

Quant Title : SVOA CALIBRATION QLast Update : Tue Nov 02 14:49:05 2021 Response via : Initial Calibration

Compound	R.T.	QIon	Response (	Conc Units Dev	(Min)
Internal Standards		453	40676	20.00 ng/ul	0.00
<ol> <li>1,4-Dichlorobenzene-d4</li> </ol>	8.224	152	48676	20.00 ng/ul	-0.01
20) Naphthalene-d8	11.050	136	223282	20.00 ng/ul	-0.01
38) Acenaphthene-d10	14.852	164	157395	20.00 ng/ul	-0.01
64) Phenanthrene-d10	17.596	188	347701	20.00 ng/ul	-0.02
79) Chrysene-d12	21.890	240	291630		-0.02
88) Perylene-d12	25.287	264	280467	20.00 ng/ul	-0.02
System Monitoring Compounds			40040	6.66 ng/uL	0.00
3) 1,4-Dioxane-d8	3.588	96	10048		0.00
4) Pyridine-d5	4.011	84	77782	17.24 ng/ul	-0.01
7) Phenol-d5	7.366	99	91208	17.56 ng/ul	0.00
<ol><li>Bis-(2-Chloroethyl)eth</li></ol>	7.543	67	58773	17.52 ng/ul	
11) 2-Chlorophenol-d4	7.754	132	65077	18.08 ng/ul	-0.01
<pre>15) 4-Methylphenol-d8</pre>	8.923	113	72172	17.65 ng/ul	-0.01
21) Nitrobenzene-d5	9.399	128	34366	18.11 ng/ul	-0.01
24) 2-Nitrophenol-d4	10.122	143	38713	18.35 ng/ul	-0.02
28) 2,4-Dichlorophenol-d3	10.662	165	66873	18.82 ng/ul	-0.01
31) 4-Chloroaniline-d4	11.180	131	99099	18.41 ng/ul	-0.01
46) Dimethylphthalate-d6	14.241	166	217820	18.09 ng/ul	-0.01
49) Acenaphthylene-d8	14.546	160	276955	18.46 ng/ul	-0.01
54) 4-Nitrophenol-d4	15.034	143	35769	16.38 ng/ul	-0.01
60) Fluorene-d10	15.839	176	196551	18.43 ng/ul	-0.01
65) 4,6-Dinitro-2-methylph	15.956	200	34161	16.21 ng/ul	0.00
73) Anthracene-d10	17.695	188	304525	18.52 ng/ul	-0.01
81) Pyrene-d10	19.969	212	342746	18.20 ng/ul	-0.01
92) Benzo(a)pyrene-d12	25.051	264	295727	19.07 ng/ul	-0.02
32) Belled(d)p). and all					
Target Compounds				Ç	value
2) 1,4-Dioxane	3.624	88	10870	6.56 ng/uL	92
5) Pyridine	4.029	79	80824	17.31 ng/ul	97
6) Benzaldehyde	7.360		62045	18.94 ng/ul	93
8) Phenol	7.396		95202	17.72 ng/ul	99
10) Bis(2-Chloroethyl)ether	7.637		72047.	17.92 ng/ul	96
12) 2-Chlorophenol	7.784		67869	18.57 ng/ul	99
13) 2-Methylphenol	8.659		69811	17.58 ng/ul	100
14) 2,2'-oxybis(1-Chloropr	8.741		107960	17.05 ng/ul	99
16) Acetophenone	9.053	D vonasnovan	4	17.98 ng/ul	98
17) N-Nitroso-di-n-propyla	9.029	10.00	66635	17.39 ng/ul	98
18) 4-Methylphenol	8.988			17.86 ng/ul	95
19) Hexachloroethane	9.317			17.60 ng/ul	92 .
	9.440		95821	18.11 ng/ul	99
22) Nitrobenzene	9.963			18.04 ng/ul	99
23) Isophorone	10.157			18.58 ng/ul	99
25) 2-Nitrophenol	10.204			18.68 ng/ul	99
26) 2,4-Dimethylphenol	10.439		to the state of the state of	18.32 ng/ul	97
27) Bis(2-Chloroethoxy)met	10.433			19.24 ng/ul	97
29) 2,4-Dichlorophenol	11.103			18.56 ng/ul	98
30) Naphthalene				18.48 ng/ul	96
32) 4-Chloroaniline	11.209			19.17 ng/ul	939,330
33) Hexachlorobutadiene	11.373				
34) Caprolactam	11.961			18.67 ng/ul	
35) 4-Chloro-3-methylphenol	12.308	10/	02070	10.07 hg/u1	

Instrument: BNA\_G **LabSampleld**: SSTDCCC020

### **Manual IntegrationsAPPROVED**

Reviewed By :Jagrut Upadhyay 11/11/2021 Supervised By :mohammad ahmed 11/11/2021

) 11/11/5/

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG110821\

Data File : BG050953.D

: 10 Nov 2021 21:45 Acq On

Operator : CG/JU : SSTDCCC020 Sample

Misc

ALS Vial : 31 Sample Multiplier: 1

Quant Time: Nov 11 01:45:02 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG110321.M

Quant Title : SVOA CALIBRATION QLast Update : Tue Nov 02 14:49:05 2021

Response via : Initial Calibration

Compound	R.T. Ç	[]Ion	Response	Conc Units Dev(	Min)
36) 2-Methylnaphthalene	12.690	142	156301	18.79 ng/ul	98
37) 1-Methylnaphthalene	12.907	142	158879	18.85 ng/ul	99
39) 1,2,4,5-Tetrachloroben	13.054	216	86442	18.85 ng/ul	96
40) Hexachlorocyclopentadiene	13.024	237	40484	18.38 ng/ul#	98
41) 2,4,6-Trichlorophenol	13.289	196	55355	18.45 ng/ul	98
42) 2,4,5-Trichlorophenol	13.365	196	57479	17.84 ng/ul	99
43) 1,1'-Biphenyl	13.682	154	212880	18.50 ng/ul	98
44) 2-Chloronaphthalene	13.735	162	166180	18.43 ng/ul	99
45) 2-Nitroaniline	13.935	65	61592	17.20 ng/ul	90
47) Dimethylphthalate	14.288	163	218297	18.13 ng/ul	100
48) 2,6-Dinitrotoluene	14.417	165	45087	17.89 ng/ul	96
50) Acenaphthylene	14.576	152	275666	18.33 ng/ul	97
51) 3-Nitroaniline	14.752	138	46691	17.91 ng/ul	91
52) Acenaphthene	14.910	153	180510	18.26 ng/ul	98
53) 2,4-Dinitrophenol	14.963	184	17252	12.41 ng/ul	91
55) 4-Nitrophenol	15.051	109	32976	16.47 ng/ul	94
56) Dibenzofuran	15.245	168	261267	18.46 ng/ul	98
57) 2,4-Dinitrotoluene	15.204	165	65383	18.18 ng/ul	97
58) 2,3,4,6-Tetrachlorophenol	15.469	232	46064	18.19 ng/ul	99
59) Diethylphthalate	15.645	149	231215	17.94 ng/ul	100
61) Fluorene	15.892	166	202415	18.07 ng/ul	98 97
62) 4-Chlorophenyl-phenyle	15.874	204	109168	18.72 ng/ul	97
63) 4-Nitroaniline	15.915	138	46266	17.89 ng/ul	97 97
66) 4,6-Dinitro-2-methylph	15.968	198	33980	16.53 ng/ul	98
67) N-Nitrosodiphenylamine	16.091	169	181963	18.72 ng/ul	93
68) 4-Bromophenyl-phenylether	16.773	248	67018	19.38 ng/ul	98
69) Hexachlorobenzene	16.896	284	69061	19.42 ng/ul	96
70) Atrazine	17.031	200	75120	18.23 ng/ul 18.20 ng/ul	95
71) Pentachlorophenol	17.243	266	29708	18.51 ng/ul	99
72) Phenanthrene	17.637	178	343488	18.55 ng/ul	99
74) Anthracene	17.731	178	345420	19.60 ng/uL	99
75) 1,2,3,4-Tetrachloroben	13.653	216	92795	19.71 ng/uL	98
76) Pentachlorobenzene	15.163	250	86482 312391	18.72 ng/ul	99
77) Carbazole	17.995	167	395319	18.02 ng/ul	100
78) Di-n-butylphthalate	18.530	149 202	408584		98
80) Fluoranthene	19.634 19.999			18.30 ng/ul	97
82) Pyrene	20.862				98
83) Butylbenzylphthalate	10.0	252		17.84 ng/ul	96
84) 3,3'-Dichlorobenzidine	21.779 21.873			18.52 ng/ul	99
85) Benzo(a)anthracene	21.738			17.96 ng/ul	98 .
86) Bis(2-ethylhexyl)phtha	21.937			18.27 ng/ul	99
87) Chrysene	23.007			18.00 ng/ul	100
89) Di-n-octyl phthalate	24.200			18.93 ng/ul	99
90) Benzo(b)fluoranthene 91) Benzo(k)fluoranthene	24.270			18.42 ng/ul	99
93) Benzo(k) Tuoranchene 93) Benzo(a) pyrene	25.128				98
94) Indeno(1,2,3-cd)pyrene	29.188				97
95) Dibenzo(a,h)anthracene	29.252				97
96) Benzo(g,h,i)perylene	30.416				98
20/ PEHZO(83H) 1/bei 3 1 1 1 1 1			. <b></b> -		

<sup>(#) =</sup> qualifier out of range (m) = manual integration (+) = signals summed

Instrument: BNA\_G **LabSampleId**: SSTDCCC020

### **Manual IntegrationsAPPROVED**