#### Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG120221\

Data File : BG051329.D

Acq On : 3 Dec 2021 14:38

Operator : CG/JU Sample : SSTDCCC020

Misc

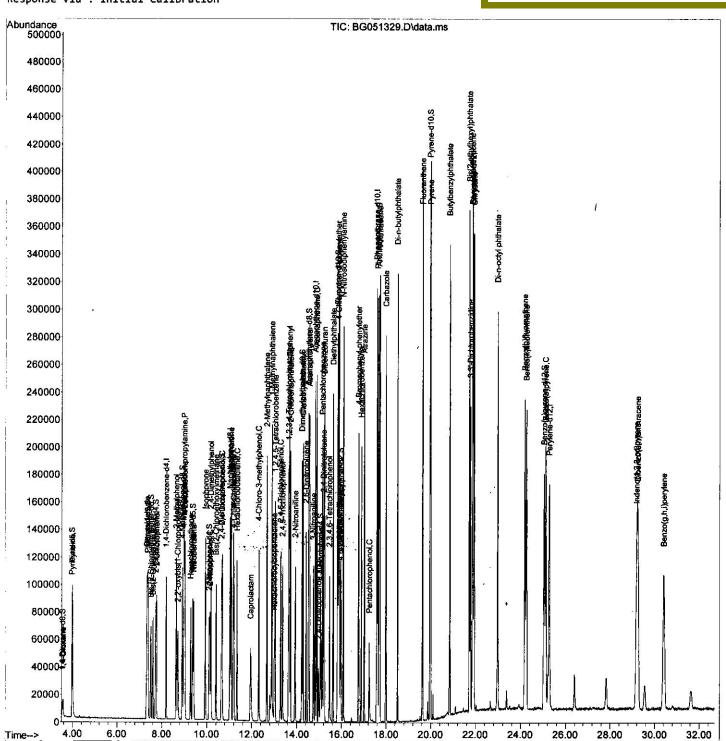
ALS Vial : 34 Sample Multiplier: 1

Quant Time: Dec 03 15:24:27 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG112321.M

Quant Title : SVOA CALIBRATION QLast Update : Fri Dec 03 15:23:09 2021 Response via : Initial Calibration Instrument :
BNA\_G
LabSampleId :
SSTDCCC020

# **Manual IntegrationsAPPROVED**



SFAM-EPA-BG112321.M Fri Dec 03 15:28:23 2021

#### Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG120221\

Data File: BG051329.D

Acq On : 3 Dec 2021 14:38

Operator : CG/JU Sample : SSTDCCC020

Misc

ALS Vial : 34 Sample Multiplier: 1

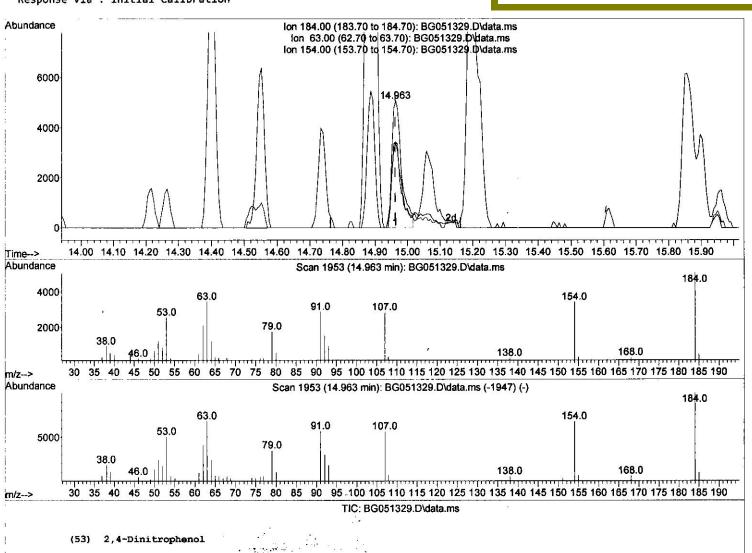
Quant Time: Dec 03 15:24:27 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG112321.M

Quant Title : SVOA CALIBRATION QLast Update : Fri Dec 03 15:23:09 2021 Response via : Initial Calibration Instrument :
BNA\_G
LabSampleId :
SSTDCCC020

# **Manual IntegrationsAPPROVED**

Reviewed By :Jagrut Upadhyay 12/03/2021 Supervised By :mohammad ahmed 12/05/2021



14.963min ( 0.000) 13.67 ng/ul

response	10554	
Ion	Ежрв	Act*
184.00	100.00	100.00
63.00	82.70	68.00
154.00	67.00	67.47
0.00	0.00	0.00

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Operator : CG/JU Sample : SSTDCCC020

Misc

ALS Vial : 34 Sample Multiplier: 1

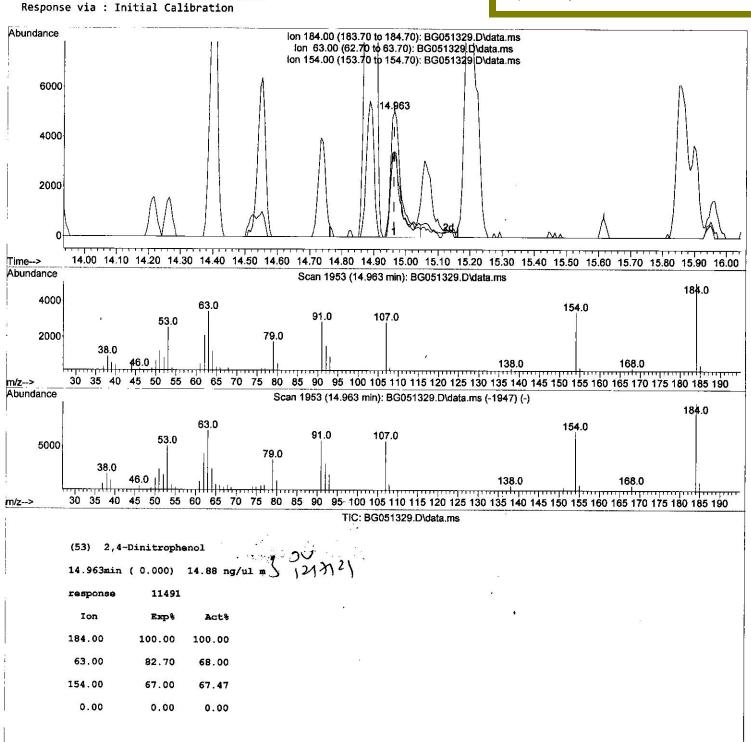
Quant Time: Dec 03 15:24:27 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG112321.M

Quant Title : SVOA CALIBRATION

QLast Update : Fri Dec 03 15:23:09 2021 Response via : Initial Calibration Instrument :
BNA\_G
LabSampleId :
SSTDCCC020

# **Manual IntegrationsAPPROVED**



# Quantitation Report (QT Reviewed)

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Acq On : 3 Dec 2021 14:38 Operator : CG/JU

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Misc

ALS Vial : 34 Sample Multiplier: 1

Quant Time: Dec 03 15:24:27 2021

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Quant Title : SVOA CALIBRATION

QLast Update : Fri Dec 03 15:23:09 2021 Response via : Initial Calibration

nstrument :	
BNA_G	
abSampleId	
STDCCC020	

# **Manual Integrations APPROVED**

Compound	R.T. (	QIon	Response	Conc Units Dev(	Min)
Internal Standards					
1) 1,4-Dichlorobenzene-d4	8.189	152	28595	20.000 ng/ul	0.00
20) Naphthalene-d8	11.015	136	127984	20.000 ng/ul	0.00
38) Acenaphthene-d10	14.822	164	85403	20.000 ng/ul	0.00
64) Phenanthrene-d10	17.572	188	192236	20.000 ng/ul	0.00
79) Chrysene-d12	21.873	240	177305	20.000 ng/ul	0.00
88) Perylene-d12	25.269	264	177897	20.000 ng/ul	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.530	96	6360	7.729 ng/uL	0.00
4) Pyridine-d5	3.959	84	45954	19.032 ng/ul	0.00
7) Phenol-d5	7.355	99	53208	18.827 ng/ul	0.00
<ol><li>Bis-(2-Chloroethyl)eth</li></ol>	7.507	67	33901	19.100 ng/ul	0.00
11) 2-Chlorophenol-d4	7.725	132	38972	19.150 ng/ul	0.00
<pre>15) 4-Methylphenol-d8</pre>	8.906	113	43436	19.046 ng/ul	0.00
21) Nitrobenzene-d5	9.370	128	20887	19.333 ng/ul	0.00 0.00
24) 2-Nitrophenol-d4	10.093	143	23671	19.423 ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.645	165	40877	19.769 ng/ul 19.327 ng/ul	0.00
31) 4-Chloroaniline-d4	11.156	131	58476	18.976 ng/ul	0.00
46) Dimethylphthalate-d6	14.217	166	124698	19.590 ng/ul	0.00
49) Acenaphthylene-d8	14.523	160	162327 16882	15.871 ng/ul	0.00
54) 4-Nitrophenol-d4	15.046	143	113822	19.235 ng/ul	0.00
60) Fluorene-d10	15.815	176 200	18117	15.273 ng/ul	0.00
65) 4,6-Dinitro-2-methylph	15.944 17.672	188	179905	19.568 ng/ul	0.00
73) Anthracene-d10	19.952	212	210118	19.585 ng/ul	0.00
81) Pyrene-d10 92) Benzo(a)pyrene-d12	25.040	264	182861	19.247 ng/ul	0.00
				Ove	alue
Target Compounds		-00	CAC7	6.969 ng/uL	96
<ol><li>1,4-Dioxane</li></ol>	3.565	88	6467	19.636 ng/ul	99
5) Pyridine	3.982	79	49336	22.177 ng/ul	95
6) Benzaldehyde	7.325	77	39914	19.191 ng/ul	99
8) Phenol	7.378	94	56185 42629	19.246 ng/ul	92
10) Bis(2-Chloroethyl)ether	7.595	93 128	39915	19.247 ng/ul	94
12) 2-Chlorophenol	7.754 8.641	108	41568		95
13) 2-Methylphenol	8.706		62927	19.688 ng/ul	98
14) 2,2'-oxybis(1-Chloropr	9.017		68162	19.323 ng/ul	98
16) Acetophenone	8.994	70	39484	19.478 ng/ul	95
17) N-Nitroso-di-n-propyla	8.970		44361	19.023 ng/ul	96
<pre>18) 4-Methylphenol 19) Hexachloroethane</pre>	9.276	117	16847	19.233 ng/ul	96.
	9.411	77	56346	19.890 ng/ul	95
22) Nitrobenzene	9.928	82	107394	19.513 ng/ul	100
<pre>23) Isophorone 25) 2-Nitrophenol</pre>	10.128	139	23963	18.983 ng/ul	97
26) 2,4-Dimethylphenol	10.175	107	52064	20.173 ng/ul	98
27) Bis(2-Chloroethoxy)met	10.404	93	59278	19.510 ng/ul	98
29) 2,4-Dichlorophenol	10.668	162	40224	19.762 ng/ul	98
30) Naphthalene	11.068	128	133808	19.215 ng/ul	98
32) 4-Chloroaniline	11.179			19.121 ng/ul	99
33) Hexachlorobutadiene	11.332	225		18.427 ng/ul	95
34) Caprolactam	11.937	113		19.112 ng/ul	90
35) 4-Chloro-3-methylphenol	12.302	107		19.877 ng/ul	97

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ALS Vial : 34 Sample Multiplier: 1

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Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG112321.M

Quant Title : SVOA CALIBRATION QLast Update : Fri Dec 03 15:23:09 2021 Response via : Initial Calibration Instrument:
BNA\_G
LabSampleId:
SSTDCCC020

# **Manual IntegrationsAPPROVED**

Compound	R.T.	QIon	Response	Conc Unit	s Dev(N	in)	
26) 2 Mothylpanhthalana	12.660	142	90289	19.061 n	g/ul	100	
36) 2-Methylnaphthalene 37) 1-Methylnaphthalene	12.878		92898	19.063 n		94	
39) 1,2,4,5-Tetrachloroben	13.024		51891	19.354 n	- 10 TO 10 T	97	
40) Hexachlorocyclopentadiene	12.989		15307	14.125 n	7 C	97	
41) 2,4,6-Trichlorophenol	13.271		32210	19.144 n	- 10 MONE	99	
42) 2,4,5-Trichlorophenol	13.353		33253	18.873 n	- S	96	
43) 1,1'-Biphenyl	13.659		124728	19.554 n		98	
44) 2-Chloronaphthalene	13.706		97982	19.310 n	-	99	
45) 2-Nitroaniline	13.917		36296	20.668 n	-	94	
47) Dimethylphthalate	14.264		127120	19.112 n		99	
48) 2,6-Dinitrotoluene	14.399		26650	19.074 n		97	200
50) Acenaphthylene	14.552		159515	19.485 n	g/ul	98	
51) 3-Nitroaniline	14.734		29086	21.061 n		96	20
52) Acenaphthene	14.887		105623 4	19.563 n	g/ul	97	12/2/2/
53) 2,4-Dinitrophenol	14.963		11491m	14.879 n			12/11
55) 4-Nitrophenol	15.063	0.0000000000000000000000000000000000000	15328		220	93	
56) Dibenzofuran	15.222		150369	19.309 n		100	
57) 2,4-Dinitrotoluene	15.192		39532	19.810 n	<del></del>	97	
58) 2,3,4,6-Tetrachlorophenol	15.451		23946	17.307 n	and the same of th	95	
59) Diethylphthalate	15.615		133464	19.116 n	<del></del> 3	99	
61) Fluorene	15.868		121528	19.482 m		98	
62) 4-Chlorophenyl-phenyle	15.850		63421	18.866 r	0.00	96	
63) 4-Nitroaniline	15.903		30072	22.376 r		99	
66) 4,6-Dinitro-2-methylph	15.962		18061	15.787 r	g/ul#	99	
67) N-Nitrosodiphenylamine	16.068		107849	19.597 r		99	
68) 4-Bromophenyl-phenylether	16.749		38611	18.740 r	ig/ul	93	
69) Hexachlorobenzene	16.873		40486	19.271 r	ng/ul	99	
70) Atrazine	17.008		44961	19.439 r	ng/ul	99	
71) Pentachlorophenol	17.231		13518	14.521 r	ng/ul	98	
72) Phenanthrene	17.619		207790	19.577 r	ig/ul	99	
74) Anthracene	17.707		209524	19.876 r	ng/ul	98	
75) 1,2,3,4-Tetrachloroben	13.630	216	54797	19.543 r	ng/uL	98	
76) Pentachlorobenzene	15.140	250	50472	19.318 r	ng/uL	97	
77) Carbazole	17.983	167	187054	20.216 r	ng/ul	99	. T
78) Di-n-butylphthalate	18.506	149	237531	19.909 r	ng/ul	99	*
80) Fluoranthene	19.617	202	258129	19.590 r	ng/ul	98	
82) Pyrene	19.981	202	254735	19.763 r	ng/ul	98	
83) Butylbenzylphthalate	20,839	149	104890	19.574 r	ng/ul	95	
84) 3,3'-Dichlorobenzidine	21.761	252	82718	20.038	44.00	98	
85) Benzo(a)anthracene	21.855	228	234844	19.528 r	ng/ul	99	
86) Bis(2-ethylhexyl)phtha	21.714	149	150929	19.573	ng/ul	99.	
87) Chrysene	21.926		224623	19.443		100	
89) Di-n-octyl phthalate	22.972	149	257022	19.943	ng/ul	100	•
90) Benzo(b)fluoranthene	24.182	252	230446	19.195 ו		98	
91) Benzo(k)fluoranthene	24.252			19.403		99	
93) Benzo(a)pyrene	25.104	252	221681	19.355		98	
94) Indeno(1,2,3-cd)pyrene	29.182			19.267		99	
95) Dibenzo(a,h)anthracene	29.235			19.183		98	
96) Benzo(g,h,i)perylene	30.416	276	205713	19.077	ng/ul	97	

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