Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN110221\

Data File : BN017242.D

Acq On : 02 Nov 2021 16:07

Operator : CG/JU Sample : SSTDICV020

Misc

ALS Vial : 9 Sample Multiplier: 1

Quant Time: Nov 02 16:43:09 2021

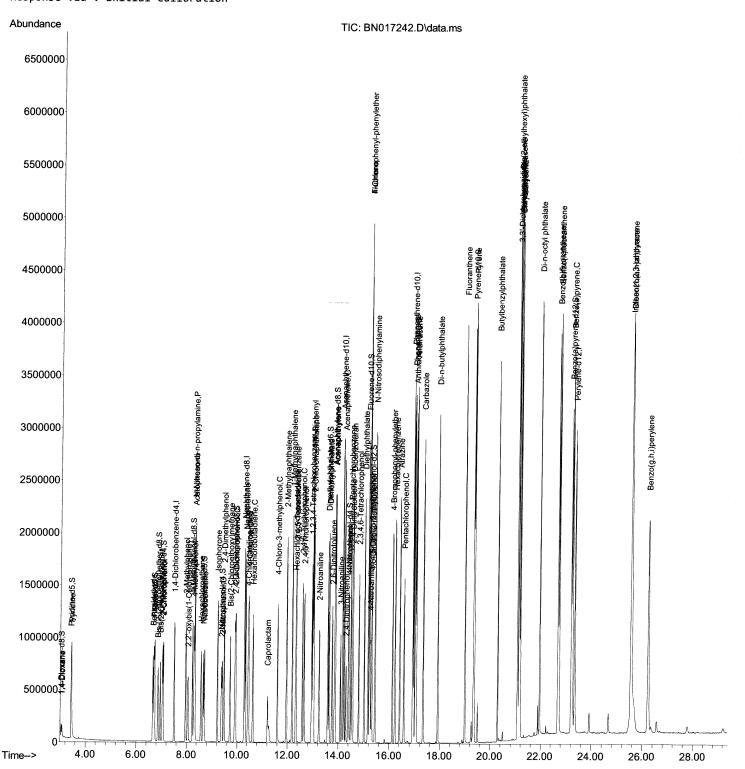
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN110221.M

Quant Title : SVOA CALIBRATION
QLast Update : Tue Nov 02 15:59:34 2021
Response via : Initial Calibration



## **Manual IntegrationsAPPROVED**

Reviewed By :Jagrut Upadhyay 11/02/2021 Supervised By :mohammad ahmed 11/08/2021



#### Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN110221\

Data File : BN017242.D

Acq On : 02 Nov 2021 16:07

Operator : CG/JU Sample : SSTDICV020

Misc

ALS Vial : 9 Sample Multiplier: 1

Quant Time: Nov 02 16:43:09 2021

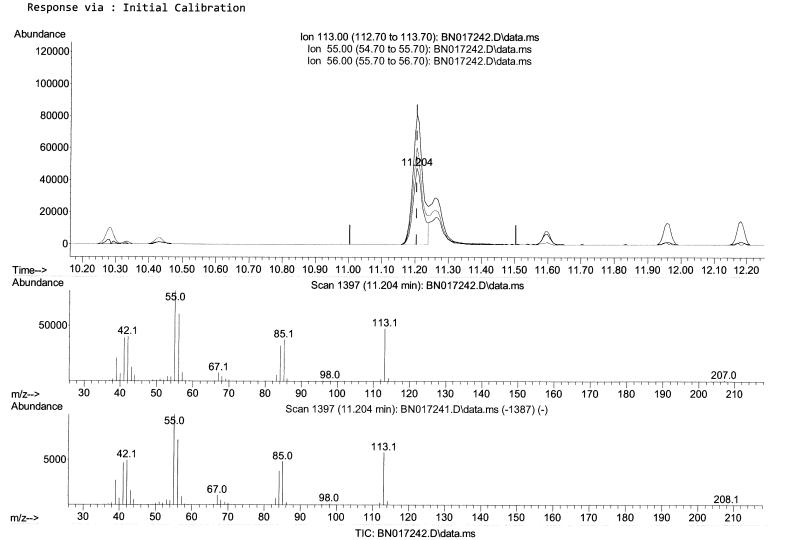
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN110221.M

Quant Title : SVOA CALIBRATION
QLast Update : Tue Nov 02 15:59:34 2021

Instrument : BNA\_N ClientSampleId : SICV243

## **Manual IntegrationsAPPROVED**

Reviewed By :Jagrut Upadhyay 11/02/2021 Supervised By :mohammad ahmed 11/08/2021



#### (34) Caprolactam

11.204min (-0.000) 11.67 ng/ul

response	97627	
Ion	Ежр%	Act%
113.00	100.00	100.00
55.00	172.30	170.57
56.00	123.70	126.98
0.00	0.00	0.00

#### Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN110221\

Data File : BN017242.D

Acq On : 02 Nov 2021 16:07

Operator : CG/JU Sample : SSTDICV020

Misc :

ALS Vial : 9 Sample Multiplier: 1

Quant Time: Nov 02 16:43:09 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN110221.M

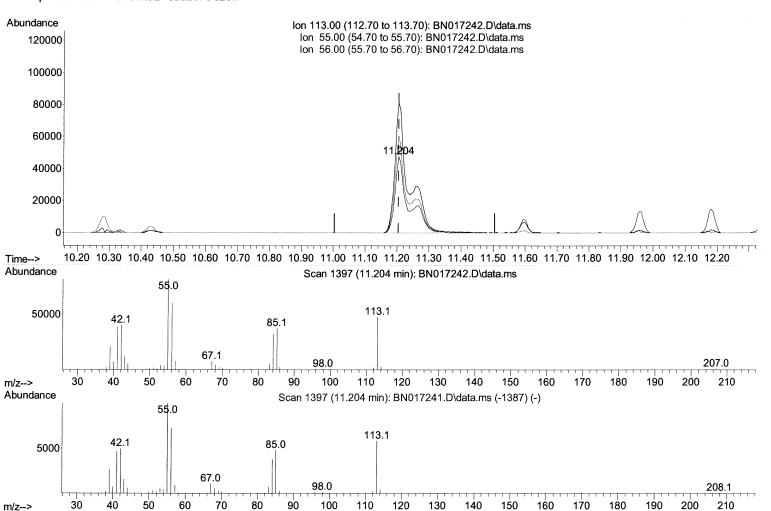
Quant Title : SVOA CALIBRATION

QLast Update : Tue Nov 02 15:59:34 2021 Response via : Initial Calibration



## **Manual IntegrationsAPPROVED**

Reviewed By :Jagrut Upadhyay 11/02/2021 Supervised By :mohammad ahmed 11/08/2021



TIC: BN017242.D\data.ms

# (34) Caprolactam

11.204min (-0.000) 16.93 ng/ul m 11/04/21 JA

response	141597	
Ion	Ехр%	Act%
113.00	100.00	100.00
55.00	172.30	170.57
56.00	123.70	126.98
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN110221\

Data File : BN017242.D

Acq On : 02 Nov 2021 16:07

Operator : CG/JU Sample : SSTDICV020

Misc :

ALS Vial : 9 Sample Multiplier: 1

Quant Time: Nov 02 16:43:09 2021

Quant Title : SVOA CALIBRATION

QLast Update : Tue Nov 02 15:59:34 2021 Response via : Initial Calibration Instrument : BNA\_N

ClientSampleId:

SICV243

#### Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/02/2021 Supervised By :mohammad ahmed 11/08/2021

Compound				Conc Units Dev	
Internal Standards					
1) 1,4-Dichlorobenzene-d4	7.504	152	303577	20 000 ng/u]	0.00
20) Naphthalene-d8	10.281		1475558	20.000 ng/ul 20.000 ng/ul	0.00 0.00
38) Acenaphthene-d10	14.169		992519	20.000 ng/ul	0.00
64) Phenanthrene-d10	16.927			20.000 ng/ul	0.00
79) Chrysene-d12	21.145			20.000 ng/ul	0.00
88) Perylene-d12	23.345	264	1853469	20.000 ng/ul	0.00
30, . c. yzee uzz	23.343	204	1000400	20.000 lig/u1	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.034	96	55326	7.203 ng/uL	0.00
4) Pyridine-d5	3.422	84	385464	17.986 ng/ul	0.00
7) Phenol-d5	6.693	99	495695	17.690 ng/ul	0.00
<pre>9) Bis-(2-Chloroethyl)eth</pre>	6.857	67	300321	17.955 ng/ul	0.00
<pre>11) 2-Chlorophenol-d4</pre>	7.040	132	400553	17.988 ng/ul	0.00
<pre>15) 4-Methylphenol-d8</pre>	8.228	113	410498	17.768 ng/ul	0.00
21) Nitrobenzene-d5	8.657	128	206273	18.017 ng/ul	0.00
24) 2-Nitrophenol-d4	9.375	143	231373	18.119 ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.910	165	410466	17.878 ng/ul	0.00
<pre>31) 4-Chloroaniline-d4</pre>	10.428	131	610937	17.755 ng/ul	0.00
46) Dimethylphthalate-d6	13.592	166	1249912	16.985 ng/ul	0.00
49) Acenaphthylene-d8	13.857	160	1601211	17.299 ng/ul	0.00
54) 4-Nitrophenol-d4	14.398	143	244423	16.922 ng/ul	0.00
60) Fluorene-d10	15.169	176	1071225	17.058 ng/ul	0.00
65) 4,6-Dinitro-2-methylph	15.304	200	222912	16.681 ng/ul	0.00
73) Anthracene-d10	17.022	188	1694533	17.188 ng/ul	0.00
81) Pyrene-d10	19.333	212	1967381	17.659 ng/ul	0.00
92) Benzo(a)pyrene-d12	23.204	264	2079625	20.768 ng/ul	0.00
Target Compounds				0	1
2) 1,4-Dioxane	3.069	88	E4602	Qva	
5) Pyridine	3.446	79	54603 395291	7.263 ng/uL	96 100
6) Benzaldehyde	6.657	7 <i>5</i> 77		18.105 ng/ul	100
8) Phenol	6.722	94	263363 500780	18.519 ng/ul	99
10) Bis(2-Chloroethyl)ether	6.951	93	400655	17.638 ng/ul	98
12) 2-Chlorophenol	7.075	128	414301	17.807 ng/ul	97
13) 2-Methylphenol	7.963	108	388942	18.200 ng/ul	98
14) 2,2'-oxybis(1-Chloropr	8.051	45	594965	17.829 ng/ul	99
16) Acetophenone	8.328	105	640290	17.857 ng/ul	99
17) N-Nitroso-di-n-propyla	8.322	70	320798	18.425 ng/ul 18.123 ng/ul	99
18) 4-Methylphenol	8.293	108	434783		98
19) Hexachloroethane	8.569	117	159259	17.997 ng/ul 17.910 ng/ul	94
22) Nitrobenzene	8.704	77	457091		93
23) Isophorone	9.222	82	901106	17.761 ng/ul 17.614 ng/ul	100
25) 2-Nitrophenol	9.410	139	248235	•	99
26) 2,4-Dimethylphenol	9.481	107	477159	18.034 ng/ul 17.664 ng/ul	96
27) Bis(2-Chloroethoxy)met	9.722	93	570839	٠,	99 100
29) 2,4-Dichlorophenol	9.940	162	407033	17.478 ng/ul	100
30) Naphthalene	10.334	128	1387107	17.920 ng/ul	95 100
32) 4-Chloroaniline	10.334	127	610899	17.493 ng/ul	100
33) Hexachlorobutadiene	10.431	225	244443	17.683 ng/ul 17.620 ng/ul	99 97
34) Caprolactam	11.204	113	141597m >		1004/2120
35) 4-Chloro-3-methylphenol	11.598	107	450514	17.686 ng/ul	98
==,	11.770	107	770714	17.000 Hg/ul	20

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Misc

ALS Vial : 9 Sample Multiplier: 1

Quant Time: Nov 02 16:43:09 2021

Quant Title : SVOA CALIBRATION QLast Update : Tue Nov 02 15:59:34 2021 Response via : Initial Calibration

### Instrument: BNA\_N ClientSampleId: SICV243

# **Manual IntegrationsAPPROVED**

Reviewed By: Jagrut Upadhyay 11/02/2021 Supervised By: mohammad ahmed 11/08/2021

Compound	R.T.	QIon	Response	Conc Units Dev	(Min)
36) 2-Methylnaphthalene	11.957	142	952259	17.308 ng/ul	99
37) 1-Methylnaphthalene	12.181	142	988779	17.493 ng/ul	96
39) 1,2,4,5-Tetrachloroben	12.339	216	483353	17.004 ng/ul	97
40) Hexachlorocyclopentadiene	12.316	237	297836	16.444 ng/ul	91
41) 2,4,6-Trichlorophenol	12.586	196	329314	17.443 ng/ul	97
42) 2,4,5-Trichlorophenol	12.657		361660	17.469 ng/ul	90
43) 1,1'-Biphenyl	12.992	154	1298016	17.390 ng/ul	100
44) 2-Chloronaphthalene	13.028	162	989618	17.191 ng/ul	97
45) 2-Nitroaniline	13.245	65	286908	17.910 ng/ul	96
47) Dimethylphthalate	13.639	163	1240074	16.921 ng/ul	99
48) 2,6-Dinitrotoluene	13.763	165	256622	17.847 ng/ul	99
50) Acenaphthylene	13.886	152	1652822	17.369 ng/ul	99
51) 3-Nitroaniline	14.086	138	254722	16.857 ng/ul	97
52) Acenaphthene	14.233	153	1057856	17.153 ng/ul	99
53) 2,4-Dinitrophenol	14.298	184	149990	16.011 ng/ul	94
55) 4-Nitrophenol	14.410	109	173828	17.327 ng/ul	94
56) Dibenzofuran	14.575	168	1507224	17.223 ng/ul	98
57) 2,4-Dinitrotoluene	14.551	165	378059	17.770 ng/ul	99
58) 2,3,4,6-Tetrachlorophenol	14.804	232	300380	17.084 ng/ul	99
59) Diethylphthalate	15.022	149	1275150	17.159 ng/ul	98
61) Fluorene	15.227	166	1194708	17.091 ng/ul	99
62) 4-Chlorophenyl-phenyle	15.227	204	589882	16.962 ng/ul	99
63) 4-Nitroaniline	15.257	138	259744	17.276 ng/ul	97
66) 4,6-Dinitro-2-methylph	15.316	198	226580	17.079 ng/ul	97
67) N-Nitrosodiphenylamine	15.445	169	1060930	17.159 ng/ul	97
68) 4-Bromophenyl-phenylether	16.127	248	371419	17.190 ng/ul	98
69) Hexachlorobenzene	16.227	284	424589	16.902 ng/ul	98
70) Atrazine	16.410	200	398339	17.550 ng/ul	99
71) Pentachlorophenol	16.580	266	254555	16.477 ng/ul	92
72) Phenanthrene	16.969	178	1946581	17.062 ng/ul	99
74) Anthracene	17.057	178	1983829	17.274 ng/ul	98
75) 1,2,3,4-Tetrachloroben	12.951	216	508615	17.378 ng/uL	99
76) Pentachlorobenzene	14.492	250	520083	17.098 ng/uL	97
77) Carbazole	17.333	167	1812915	17.616 ng/ul	99
78) Di-n-butylphthalate	17.921	149	2184221	17.619 ng/ul	99
80) Fluoranthene	18.992	202	2314958	17.646 ng/ul	96
82) Pyrene	19.363	202	2349981	17.489 ng/ul	100
83) Butylbenzylphthalate	20.292	149	1011069	18.091 ng/ul	98
84) 3,3'-Dichlorobenzidine	21.074	252	833521	17.590 ng/ul	98
85) Benzo(a)anthracene	21.127	228	2342152	17.349 ng/ul	99
86) Bis(2-ethylhexyl)phtha	21.086	149	1548964	18.185 ng/ul	99
87) Chrysene 89) Di-n-octyl phthalate	21.180	228	2300670	17.545 ng/ul	97
90) Benzo(b)fluoranthene	21.957	149	2641263	21.626 ng/ul	100
91) Benzo(k)fluoranthene	22.686	252	2491050	20.169 ng/ul	98
93) Benzo(a)pyrene	22.727	252	2480180	21.315 ng/ul	99
94) Indeno(1,2,3-cd)pyrene	23.245	252 276	2466096	20.551 ng/ul	98 07
95) Dibenzo(a,h)anthracene	25.556 25.574	276 278	2821141	19.945 ng/ul 20.006 ng/ul	97 100
96) Benzo(g,h,i)perylene	26.233	276 276	2398677 2316105	19.351 ng/ul	100 97
(B)::/pc://ciic				IIB/UI	<i>э,</i>

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