Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN112421\

Data File : BN017572.D Acq On : 23 Nov 2021 18:33

Operator : CG/JU Sample : PB140905BS

Misc

ALS Vial : 13 Sample Multiplier: 1

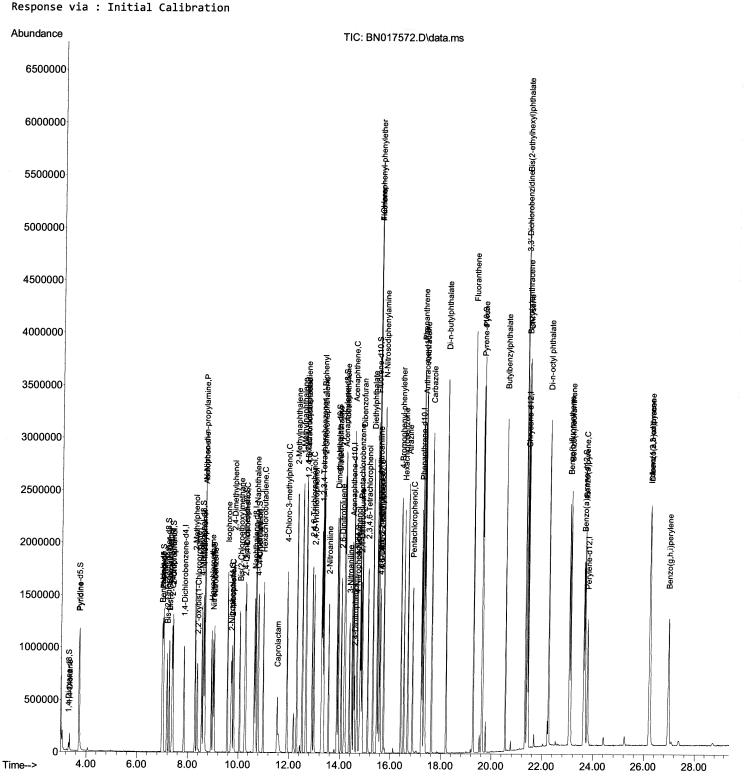
Quant Time: Nov 24 00:58:36 2021

 $\label{thm:lem1_BNA_N\ethods\SFAM-EPA-BN112221.M} Quant \ \mbox{Methods} : Z:\svoasrv\HPCHEM1\BNA_N\mbox{Methods} : Z$ 

Quant Title : SVOA CALIBRATION QLast Update : Mon Nov 22 16:16:36 2021 Instrument :
BNA\_N
ClientSampleId :
SLCS905

## **Manual IntegrationsAPPROVED**

Reviewed By :Jagrut Upadhyay 11/24/2021 Supervised By :mohammad ahmed 11/26/2021



SFAM-EPA-BN112221.M Wed Nov 24 01:19:00 2021

### Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN112421\

Data File : BN017572.D

Acq On : 23 Nov 2021 18:33

Operator : CG/JU Sample : PB140905BS

Misc

ALS Vial : 13 Sample Multiplier: 1

Quant Time: Nov 24 00:58:36 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN112221.M

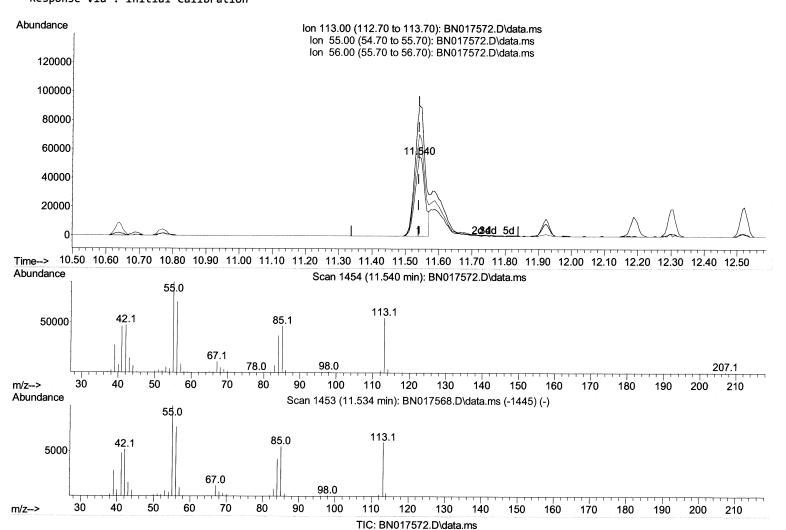
Quant Title : SVOA CALIBRATION

QLast Update : Mon Nov 22 16:16:36 2021 Response via : Initial Calibration

Instrument : BNA\_N ClientSampleId : SLCS905

# **Manual IntegrationsAPPROVED**

Reviewed By :Jagrut Upadhyay 11/24/2021 Supervised By :mohammad ahmed 11/26/2021



### (34) Caprolactam

11.540min (+ 0.000) 21.39 ng/ul

response	112438	
Ion	Ехр%	Act%
113.00	100.00	100.00
55.00	176.80	164.12
56.00	129.90	128.17
0.00	0.00	0.00

#### Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN112421\

Data File : BN017572.D

Acq On : 23 Nov 2021 18:33

Operator : CG/JU Sample : PB140905BS

Misc

ALS Vial : 13 Sample Multiplier: 1

Quant Time: Nov 24 00:58:36 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN112221.M

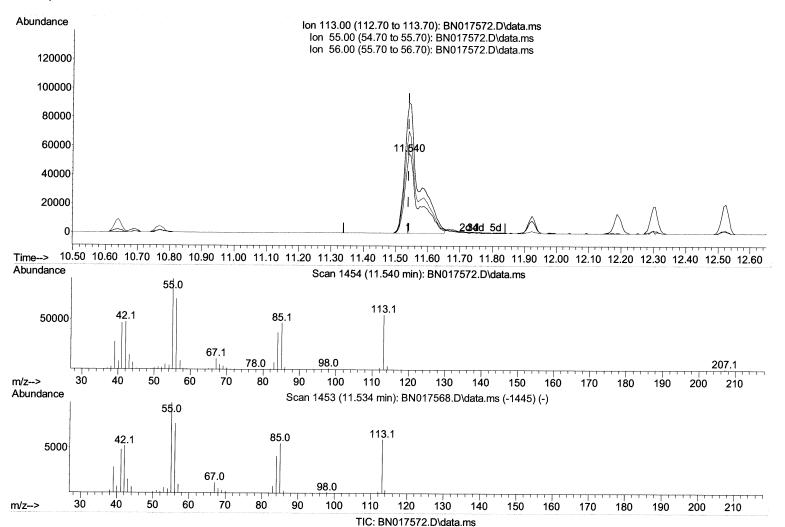
Quant Title : SVOA CALIBRATION

QLast Update : Mon Nov 22 16:16:36 2021 Response via : Initial Calibration



## **Manual IntegrationsAPPROVED**

Reviewed By :Jagrut Upadhyay 11/24/2021 Supervised By :mohammad ahmed 11/26/2021



#### (34) Caprolactam

11.540min (+ 0.000) 31.74 ng/ul m 1/(30/2) J

response	166820	
Ion	Ехр%	Act%
113.00	100.00	100.00
55.00	176.80	164.12
56.00	129.90	128.17
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN112421\

Data File : BN017572.D

Acq On : 23 Nov 2021 18:33

Operator : CG/JU Sample : PB140905BS

Misc

ALS Vial : 13 Sample Multiplier: 1

Quant Time: Nov 24 00:58:36 2021

Quant Title : SVOA CALIBRATION

QLast Update : Mon Nov 22 16:16:36 2021 Response via : Initial Calibration Instrument : BNA\_N ClientSampleId : SLCS905

# **Manual IntegrationsAPPROVED**

Reviewed By: Jagrut Upadhyay 11/24/2021 Supervised By: mohammad ahmed 11/26/2021

Compound	R.T.	QIon	Response	Conc Units Dev	(Min)
Internal Standards					
1) 1,4-Dichlorobenzene-d4	7.840	152	267417	20.000 ng/ul	0.00
20) Naphthalene-d8	10.640		1164599	20.000 ng/ul	0.00
38) Acenaphthene-d10	14.481		701606	20.000 ng/ul	0.00
64) Phenanthrene-d10	17.222	188	1339441	20.000 ng/ul	0.00
79) Chrysene-d12	21.410	240	886002	20.000 ng/ul	0.00
88) Perylene-d12	23.774	264	816877	20.000 ng/ul	0.00
System Monitoring Compounds					
<ol><li>3) 1,4-Dioxane-d8</li></ol>	3.305	96	36173	4.834 ng/uL	0.00
<ol><li>4) Pyridine-d5</li></ol>	3.705	84	482970	24.597 ng/ul	0.00
7) Phenol-d5	7.005	99	680061	27.935 ng/ul	0.00
<pre>9) Bis-(2-Chloroethyl)eth</pre>	7.175	67	409584	26.848 ng/ul	0.00
11) 2-Chlorophenol-d4	7.375	132	514768	27.926 ng/ul	0.00
<pre>15) 4-Methylphenol-d8</pre>	8.552	113	532663	27.816 ng/ul	0.00
21) Nitrobenzene-d5	8.999	128	255262	28.931 ng/ul	0.00
24) 2-Nitrophenol-d4	9.722	143	267268	30.617 ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.258	165	495619	28.839 ng/ul	0.00
31) 4-Chloroaniline-d4	10.769	131	635255	25.742 ng/ul	0.00
46) Dimethylphthalate-d6	13.887	166	1446832	28.363 ng/ul	0.00
49) Acenaphthylene-d8	14.169	160	1832384	27.977 ng/ul	0.00
54) 4-Nitrophenol-d4	14.663	143	259270	30.035 ng/ul	0.00
60) Fluorene-d10	15.469	176	1183089	27.497 ng/ul	0.00
65) 4,6-Dinitro-2-methylph	15.587	200	198140	28.481 ng/ul	0.00
73) Anthracene-d10	17.322	188	1740136	27.951 ng/ul	0.00
81) Pyrene-d10	19.616	212	1774900	31.557 ng/ul	0.00
92) Benzo(a)pyrene-d12	23.622	264	1215669	28.307 ng/ul	0.00
arget Compounds				Qva	lue
2) 1,4-Dioxane	3.340	88	76045	10.293 ng/uL	96
5) Pyridine	3.728	79	514751	26.051 ng/ul	91
<ol><li>Benzaldehyde</li></ol>	6.975	77	399337	30.922 ng/ul	89
8) Phenol	7.034	94	714299	28.970 ng/ul	95
<pre>10) Bis(2-Chloroethyl)ether</pre>	7.264	93	570418	28.215 ng/ul	96
12) 2-Chlorophenol	7.405	128	546661	28.780 ng/ul	94
13) 2-Methylphenol	8.281	108	544584	29.325 ng/ul	98
<pre>14) 2,2'-oxybis(1-Chloropr</pre>	8.375	45	759424	27.863 ng/ul	96
l6) Acetophenone	8.663	105	826916	28.876 ng/ul	96
17) N-Nitroso-di-n-propyla	8.652	70	438668	29.390 ng/ul	94
<pre>18) 4-Methylphenol</pre>	8.611	108	577795	29.044 ng/ul	99
l9) Hexachloroethane	8.928	117	227273	28.318 ng/ul	97
22) Nitrobenzene	9.040	77	643091	28.841 ng/ul	97
23) Isophorone	9.563	82	1252016	30.000 ng/ul	97
25) 2-Nitrophenol	9.752	139	299869	31.143 ng/ul#	91
26) 2,4-Dimethylphenol	9.816	107	636854	28.599 ng/ul	94
7) Bis(2-Chloroethoxy)met	10.052	93	758371	27.918 ng/ul	98
9) 2,4-Dichlorophenol	10.287	162	504417	29.108 ng/ul	97
0) Naphthalene	10.693	128	1742099	27.960 ng/ul	100
2) 4-Chloroaniline	10.793	127	655885	26.195 ng/ul	98
3) Hexachlorobutadiene	10.987	225	322910	28.329 ng/ul	98
4) Caprolactam	11.540	113		31.742 ng/ul>	11/30/21
5) 4-Chloro-3-methylphenol					

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN112421\

Data File : BN017572.D

Acq On : 23 Nov 2021 18:33

Operator : CG/JU Sample : PB140905BS

Misc :

ALS Vial : 13 Sample Multiplier: 1

Quant Time: Nov 24 00:58:36 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN112221.M

Quant Title : SVOA CALIBRATION

QLast Update : Mon Nov 22 16:16:36 2021 Response via : Initial Calibration Instrument : BNA\_N ClientSampleld : SLCS905

## Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/24/2021 Supervised By :mohammad ahmed 11/26/2021

Compound	R.T.	QIon	Response	Conc Units Dev	(Min)
36) 2-Methylnaphthalene	12.299	142	1169002	28.113 ng/ul	99
37) 1-Methylnaphthalene	12.522			27.889 ng/ul	97
39) 1,2,4,5-Tetrachloroben	12.675		578737	27.811 ng/ul	96
40) Hexachlorocyclopentadiene	12.663	237		25.836 ng/ul	96
41) 2,4,6-Trichlorophenol	12.910	196	387875	30.510 ng/ul	98
42) 2,4,5-Trichlorophenol	12.981	196	416202	30.023 ng/ul	99
43) 1,1'-Biphenyl	13.316	154	1562952	28.245 ng/ul	99
44) 2-Chloronaphthalene	13.351	162	1188676	28.251 ng/ul	97
45) 2-Nitroaniline	13.551	65	379988	32.629 ng/ul	93
47) Dimethylphthalate	13.934	163	1498948	29.645 ng/ul	99
48) 2,6-Dinitrotoluene	14.051	165	312504	32.932 ng/ul	92
50) Acenaphthylene	14.198	152	1917408	28.725 ng/ul	100
51) 3-Nitroaniline	14.375	138	293700	29.234 ng/ul	92
52) Acenaphthene	14.540	153	1233850	28.670 ng/ul	99
53) 2,4-Dinitrophenol	14.581	184	140390	29.430 ng/ul	95
55) 4-Nitrophenol	14.681	109	230016	30.551 ng/ul#	83
56) Dibenzofuran	14.875	168	1744457	28.236 ng/ul	94
57) 2,4-Dinitrotoluene	14.834	165	434191	32.290 ng/ul	95
58) 2,3,4,6-Tetrachlorophenol	15.098	232	315318	30.742 ng/ul#	93
59) Diethylphthalate	15.304	149	1534821	30.309 ng/ul	99
61) Fluorene	15.528	166	1301109	28.055 ng/ul	98
62) 4-Chlorophenyl-phenyle	15.522	204	636745	28.134 ng/ul	94
63) 4-Nitroaniline	15.540	138	282929	33.401 ng/ul#	86
66) 4,6-Dinitro-2-methylph	15.598	198	213973	30.635 ng/ul#	85
67) N-Nitrosodiphenylamine	15.734	169	1206175	29.874 ng/ul	98
68) 4-Bromophenyl-phenylether	16.416	248	382644	28.933 ng/ul	92
69) Hexachlorobenzene	16.534	284	425039	30.157 ng/ul	91
70) Atrazine	16.687	200	432363	30.322 ng/ul	95
71) Pentachlorophenol	16.875	266	240644	30.027 ng/ul	94
72) Phenanthrene	17.263	178	2140827	29.293 ng/ul	99
74) Anthracene	17.357	178	2099943	28.975 ng/ul	98
75) 1,2,3,4-Tetrachloroben	13.281	216	594110	28.289 ng/uL	98
76) Pentachlorobenzene	14.804	250	528702	27.686 ng/uL	96
77) Carbazole	17.622	167	1872238	30.100 ng/ul	98
78) Di-n-butylphthalate	18.198	149	2435350	32.219 ng/ul	99
80) Fluoranthene 82) Pyrene	19.281	202	2247356	33.217 ng/ul	99
83) Butylbenzylphthalate	19.645	202	2198905	32.464 ng/ul	97
84) 3,3'-Dichlorobenzidine	20.551	149	915178	37.011 ng/ul	93
85) Benzo(a)anthracene	21.328	252	505670	30.003 ng/ul	97
86) Bis(2-ethylhexyl)phtha	21.392	228	1739828	29.644 ng/ul	99
	21.333	149 228	1304694	36.456 ng/ul#	98
	21.445 22.251		1721673	29.854 ng/ul	99
	23.057	149	2061690	33.231 ng/ul	100
04) 5 (1) 65	23.104	252 252	1627508	28.814 ng/ul	99
	23.674	252	1551164 1537270	29.084 ng/ul	100
	26.221	276	1752115	28.665 ng/ul 31.465 ng/ul	99
	26.233	278	1484087	31.379 ng/ul	99 98
	26.968	276	1470785	30.903 ng/ul	96
				116/UI	

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