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

Data File : BN017579.D

Acq On : 23 Nov 2021 23:20

Operator : CG/JU Sample : M4753-03

Misc

ALS Vial : 21 Sample Multiplier: 1

Quant Time: Nov 24 01:01:00 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

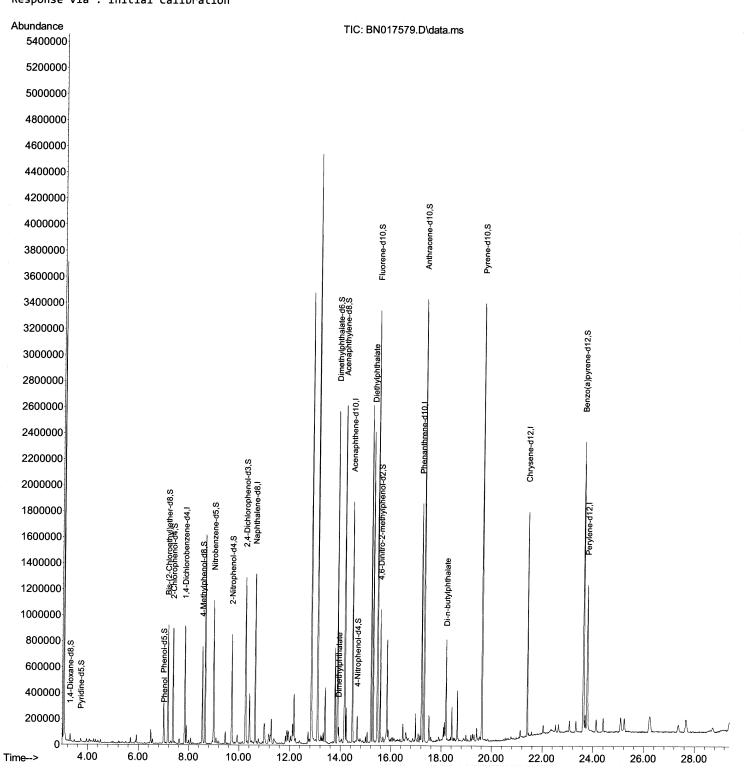
Instrument : BNA\_N

ClientSampleId :

A0016

# **Manual IntegrationsAPPROVED**

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



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

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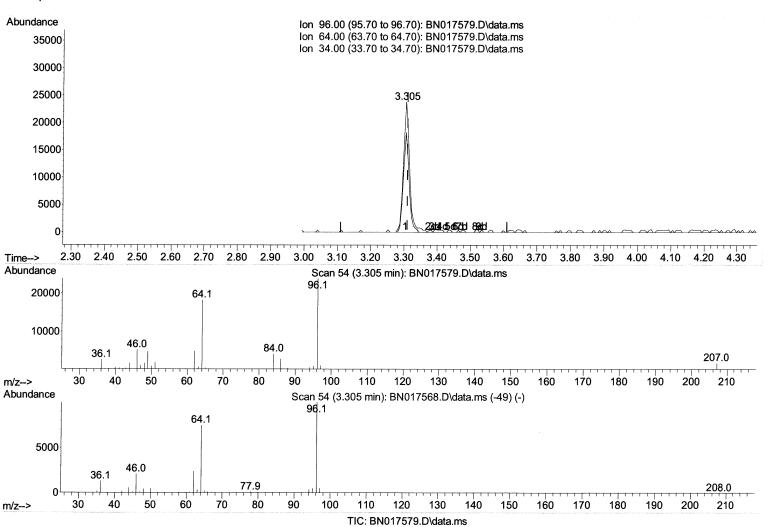
Quant Title : SVOA CALIBRATION

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



#### (3) 1,4-Dioxane-d8 (S)

3.305min (-0.005) 4.61 ng/uL

response	30401	
Ion	Ежр%	Act%
96.00	100.00	100.00
64.00	74.70	76.61
34.00	0.00	0.00
0.00	0.00	0.00

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

Data File : BN017579.D

Acq On : 23 Nov 2021 23:20

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Misc :

ALS Vial : 21 Sample Multiplier: 1

Quant Time: Nov 24 01:01:00 2021

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

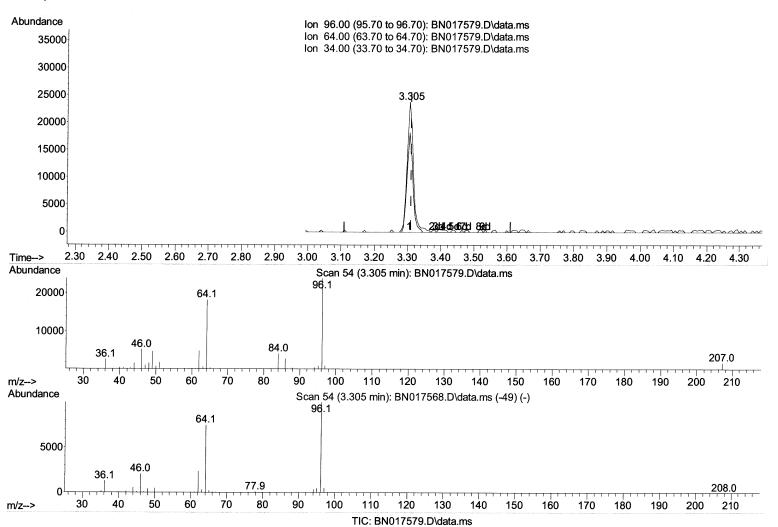
Quant Title : SVOA CALIBRATION

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



# Manual IntegrationsAPPROVED

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### (3) 1,4-Dioxane-d8 (S)

3.305min (-0.005) 4.73 ng/uL m (//30/21)y

response	31176	
Ion	Ежр%	Act%
96.00	100.00	100.00
64.00	74.70	76.61
34.00	0.00	0.00
0.00	0.00	0.00

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

Data File : BN017579.D

Acq On : 23 Nov 2021 23:20

Operator : CG/JU Sample : M4753-03

Misc :

ALS Vial : 21 Sample Multiplier: 1

Quant Time: Nov 24 01:01:00 2021

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

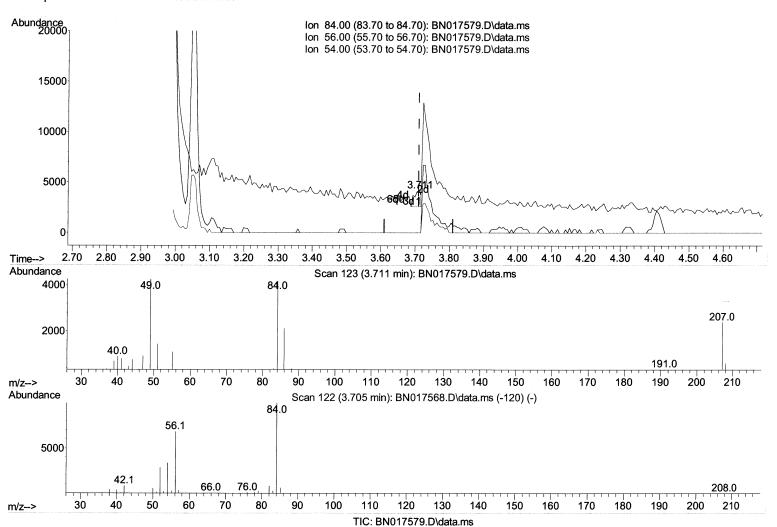
Quant Title : SVOA CALIBRATION

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



# **Manual IntegrationsAPPROVED**

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#### (4) Pyridine-d5 (S)

3.711min (+ 0.000) 0.13 ng/ul

response	2246			
Ion	Ехр%	Act%		
84.00	100.00	100.00		
56.00	68.80	0.00#		
54.00	32.90	0.00#		
0.00	0.00	0.00		

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

Data File : BN017579.D

Acq On : 23 Nov 2021 23:20

Operator : CG/JU Sample : M4753-03

Misc :

ALS Vial : 21 Sample Multiplier: 1

Quant Time: Nov 24 01:01:00 2021

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

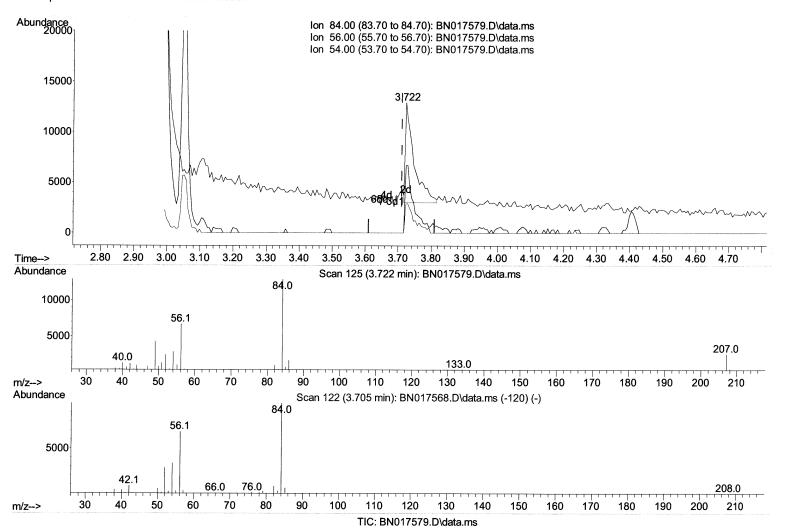
Quant Title : SVOA CALIBRATION

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



#### (4) Pyridine-d5 (S)

3.722min (+ 0.012) 1.05 ng/ul m [[/36/2[JU

response	18144	
Ion	Ехр%	Act%
84.00	100.00	100.00
56.00	68.80	51.75#
54.00	32.90	22.15#
0.00	0.00	0.00

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

Data File : BN017579.D

Acq On : 23 Nov 2021 23:20

Operator : CG/JU Sample : M4753-03

Misc

ALS Vial : 21 Sample Multiplier: 1

Quant Time: Nov 24 01:01:00 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 : A0016

### **Manual IntegrationsAPPROVED**

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

	Compound	R.T.	QIon	Response	Conc Un	its Dev	(Min)
Inte	ernal Standards						
1)	1,4-Dichlorobenzene-d4	7.846	152	235767	20.000	ng/ul	0.00
20)	Naphthalene-d8	10.640	136	1040441		ng/ul	0.00
	Acenaphthene-d10	14.475	164	619695	20.000	ng/ul	0.00
64)	Phenanthrene-d10	17.216	188	1081493	20.000	ng/ul	0.00
79)	Chrysene-d12	21.404	240	757042	20.000	ng/ul	0.00
88)	Perylene-d12	23.768	264	755375	20.000	ng/ul	0.00
Syst	em Monitoring Compounds						
3)	1,4-Dioxane-d8	3.305	96	3 <b>11</b> 76m \$		ng/uL、	0.00/1/20/21
4)	Pyridine-d5	3.722	84	18144m <sup>&lt;</sup>	1.048	ng/ul	0.00 11/30/21
	Phenol-d5	7.005	99	148311		ng/ul	0.00
9)	Bis-(2-Chloroethyl)eth	7.169	67	401781	29.872	ng/ul	0.00
11)	2-Chlorophenol-d4	7.375	132	397403	24.453	ng/ul	0.00
15)	4-Methylphenol-d8	8.546	113	272993	16.170	ng/ul	0.00
21)	Nitrobenzene-d5	8.993	128	250814	31.819	ng/ul	0.00
24)	2-Nitrophenol-d4	9.716	143	253659	32.526	ng/ul	0.00
28)	2,4-Dichlorophenol-d3	10.257	165	458465	29.861	ng/ul	0.00
31)	4-Chloroaniline-d4	10.769	131	20427		ng/ul	0.00
46)	Dimethylphthalate-d6	13.887	166	1627158	36.114	ng/ul	0.00
49)	Acenaphthylene-d8	14.169	160	1834961	31.720	ng/ul	0.00
54)	4-Nitrophenol-d4	14.669	143	46567	6.108	-	0.00
60)	Fluorene-d10	15.469	176	1266854	33.336	_	0.00
65)	4,6-Dinitro-2-methylph	15.581	200	189322	33.704	ng/ul	0.00
73)	Anthracene-d10	17.316	188	1911710	38.031	ng/ul	0.00
81)	Pyrene-d10	19.610	212	1859402	38.691	ng/ul	0.00
92)	Benzo(a)pyrene-d12	23.615	264	1525404	38.411	ng/ul	-0.01
arg	et Compounds					Qva	lue
8)	Phenol	7.028	94	30668	1.411	ng/ul#	92
47)	Dimethylphthalate	13.934	163	60124		ng/ul	97
	Diethylphthalate	15.298	149	1194900	26.715		99
	Di-n-butylphthalate	18.198	149	362990	5.948		98

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