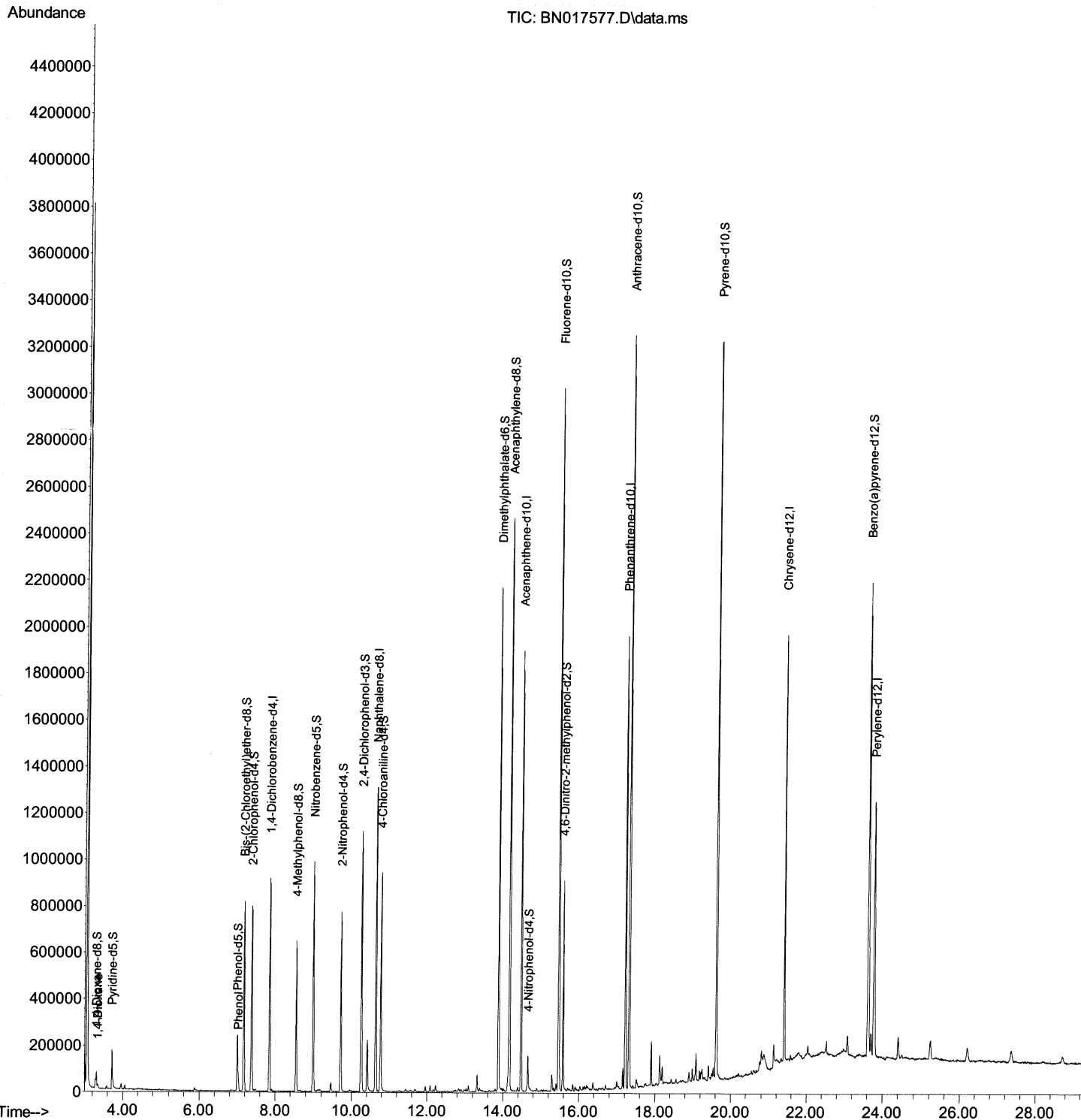Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112421\
Data File : BN017577.D
Acq On   : 23 Nov 2021 21:33
Operator : CG/JU
Sample   : M4753-02
Misc   :
ALS Vial : 18   Sample Multiplier: 1

Quant Time: Nov 24 01:00:02 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 :
A0012

Manual IntegrationsAPPROVED

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

TIC: BN017577.D\data.ms

```
                          Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112421\
Data File : BN017577.D
Acq On    : 23 Nov 2021  21:33
Operator  : CG/JU
Sample    : M4753-02
Misc      :
ALS Vial  : 18    Sample Multiplier: 1

Quant Time: Nov 24 01:00:02 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
```
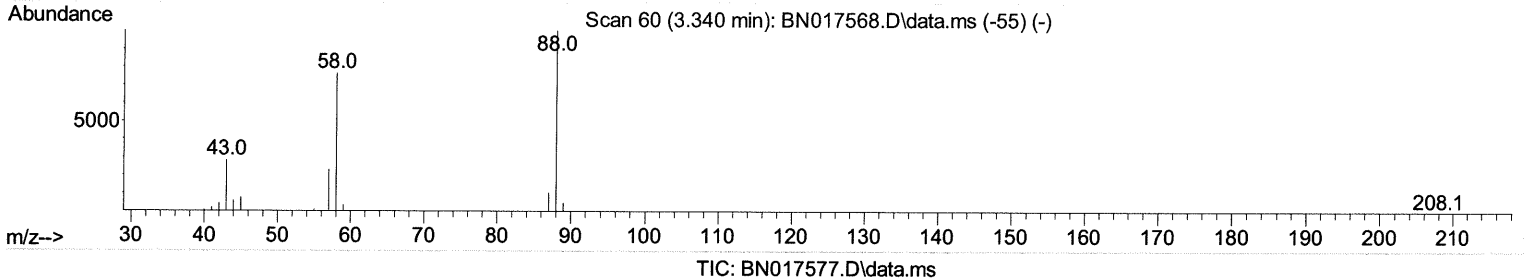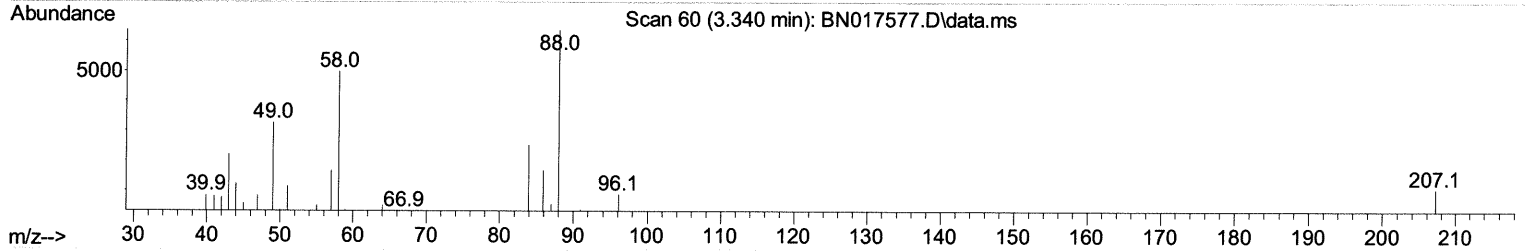
```
(2)  1,4-Dioxane

3.340min (+ 0.000)  1.32 ng/uL

response        8584

   Ion        Exp%        Act%

   88.00      100.00      100.00

   43.00       30.50       34.69

   58.00       77.40       78.00

    0.00        0.00        0.00
```
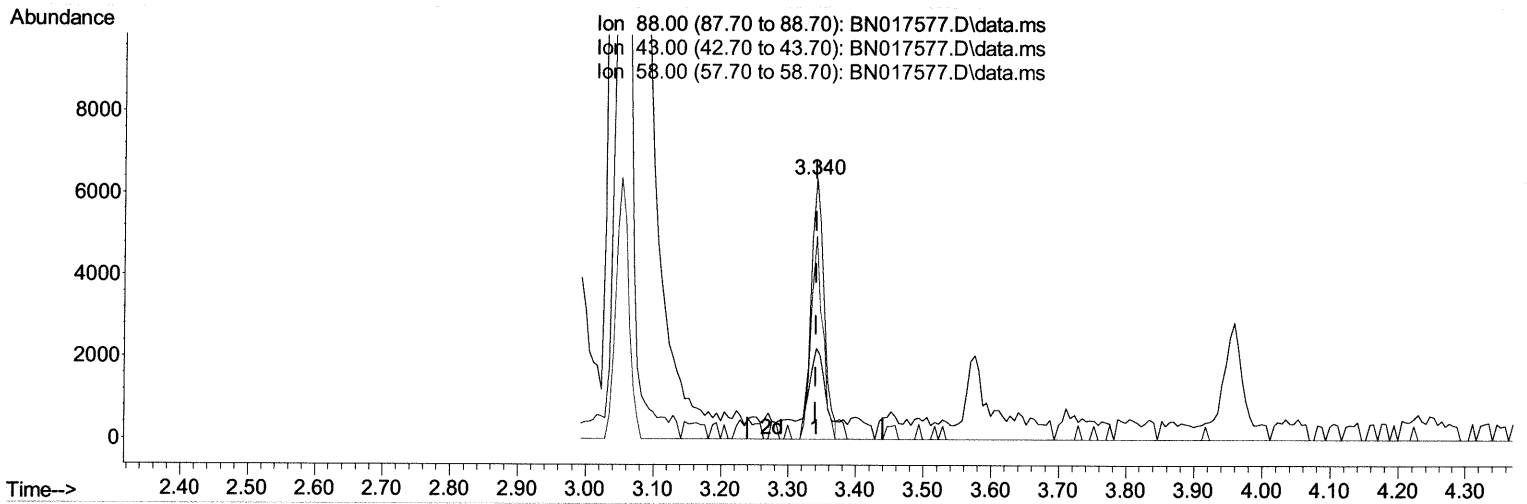
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112421\
Data File : BN017577.D
Acq On   : 23 Nov 2021 21:33
Operator : CG/JU
Sample   : M4753-02
Misc   :
ALS Vial : 18   Sample Multiplier: 1

Quant Time: Nov 24 01:00:02 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 :
A0012

Manual IntegrationsAPPROVED

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

Ion 88.00 (87.70 to 88.70): BN017577.D\data.ms
Ion 43.00 (42.70 to 43.70): BN017577.D\data.ms
Ion 58.00 (57.70 to 58.70): BN017577.D\data.ms

Scan 60 (3.340 min): BN017577.D\data.ms

Scan 60 (3.340 min): BN017568.D\data.ms (-55) (-)

TIC: BN017577.D\data.ms

(2) 1,4-Dioxane

3.340min (+ 0.000) 1.37 ng/uL m   11/30/21 JU

response   8895

Ion	Exp%	Act%
88.00	100.00	100.00
43.00	30.50	34.69
58.00	77.40	78.00
0.00	0.00	0.00

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```
                    Quantitation Report      (QT Reviewed)
```
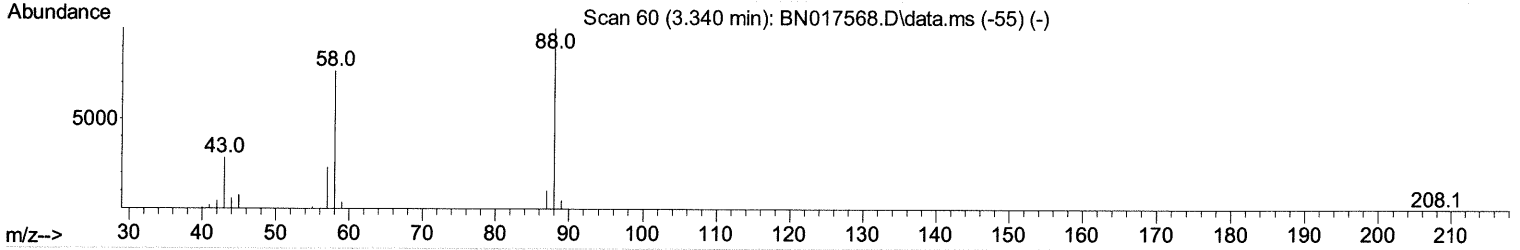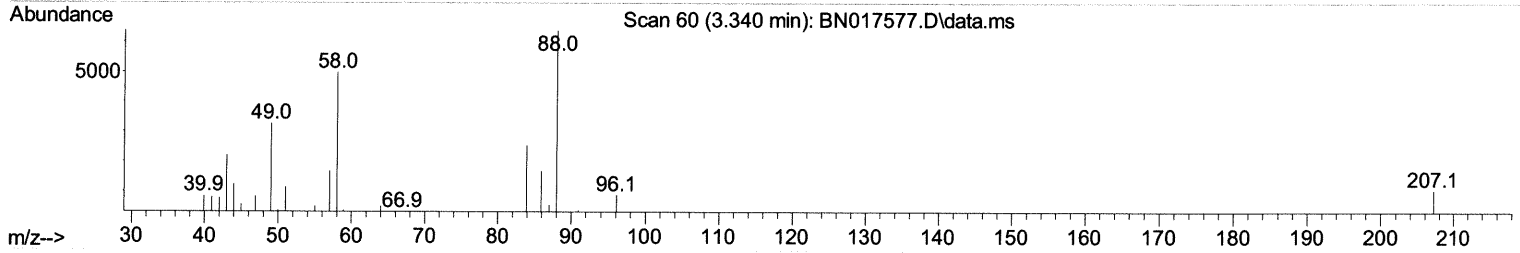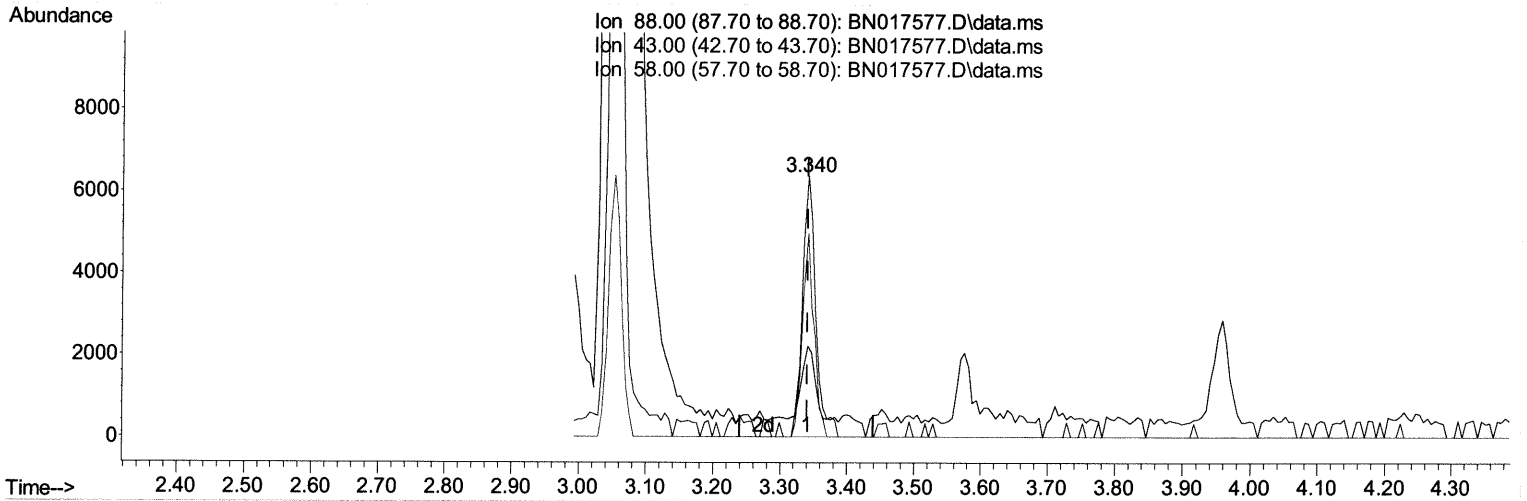
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112421\
Data File : BN017577.D
Acq On   : 23 Nov 2021 21:33
Operator : CG/JU
Sample   : M4753-02
Misc   :
ALS Vial : 18   Sample Multiplier: 1

Quant Time: Nov 24 01:00:02 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

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.840	152	235787	20.000	ng/ul	0.00
20) Naphthalene-d8	10.640	136	1044305	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.475	164	638180	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.216	188	1144329	20.000	ng/ul	0.00
79) Chrysene-d12	21.404	240	803892	20.000	ng/ul	0.00
88) Perylene-d12	23.768	264	771262	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.305	96	31687	4.803	ng/uL	0.00
4) Pyridine-d5	3.711	84	90232	5.212	ng/ul	0.00
7) Phenol-d5	7.005	99	131708	6.136	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.169	67	363378	27.015	ng/ul	0.00
11) 2-Chlorophenol-d4	7.375	132	358814	22.077	ng/ul	0.00
15) 4-Methylphenol-d8	8.546	113	243775	14.438	ng/ul	0.00
21) Nitrobenzene-d5	8.993	128	228145	28.836	ng/ul	0.00
24) 2-Nitrophenol-d4	9.716	143	234618	29.973	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.257	165	397831	25.816	ng/ul	0.00
31) 4-Chloroaniline-d4	10.769	131	498204	22.514	ng/ul	0.00
46) Dimethylphthalate-d6	13.887	166	1398498	30.140	ng/ul	0.00
49) Acenaphthylene-d8	14.169	160	1714685	28.782	ng/ul	0.00
54) 4-Nitrophenol-d4	14.657	143	39782	5.067	ng/ul	0.00
60) Fluorene-d10	15.469	176	1132920	28.948	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.581	200	169462	28.512	ng/ul	0.00
73) Anthracene-d10	17.316	188	1758844	33.068	ng/ul	0.00
81) Pyrene-d10	19.616	212	1717525	33.656	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.616	264	1382967	34.107	ng/ul	-0.01

Target Compounds						Qvalue
2) 1,4-Dioxane	3.340	88	8895m	1.365	ng/uL	
8) Phenol	7.028	94	27559	1.268	ng/ul	92

(handwritten: 11/30/21 JU)

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

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