

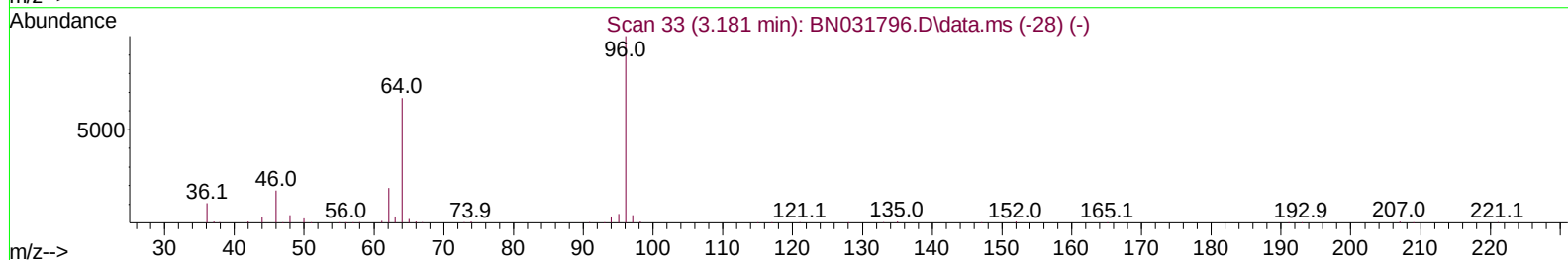
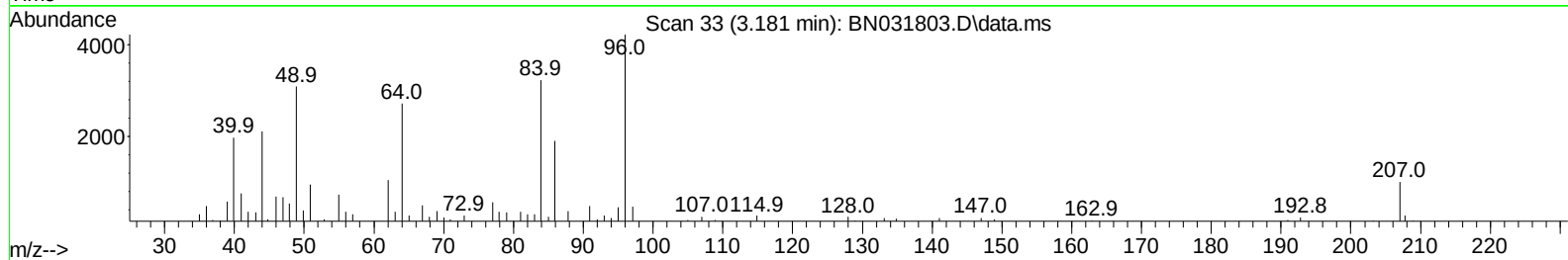
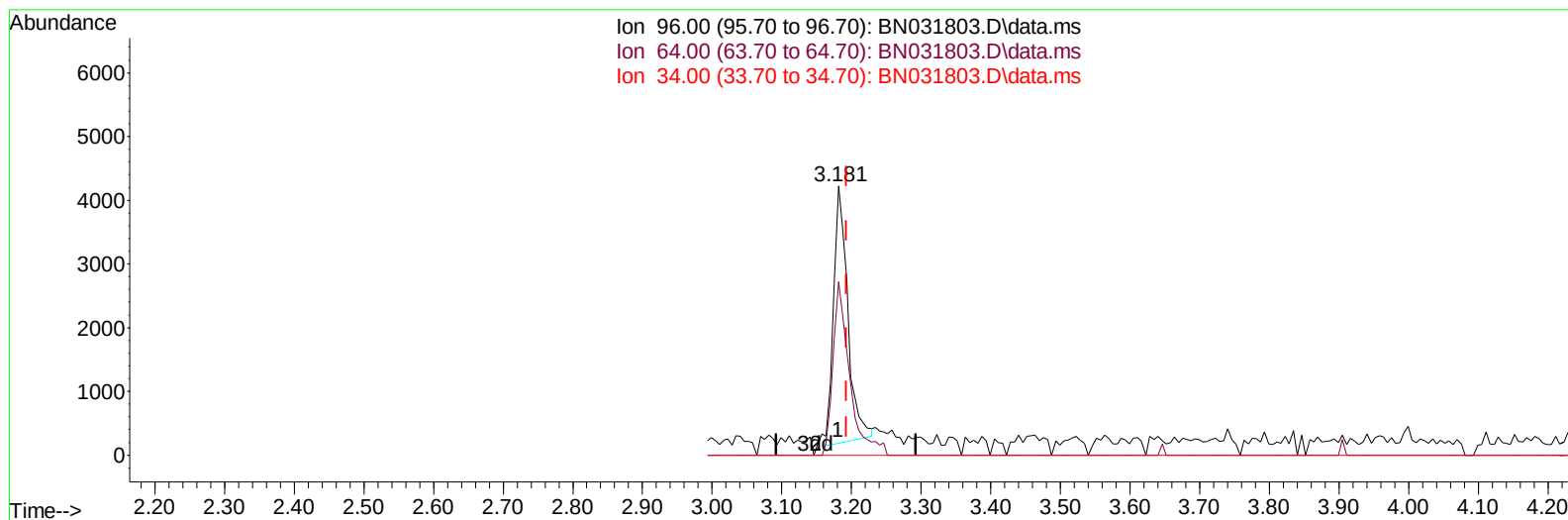
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060524\
Data File : BN031803.D
Acq On : 05 Jun 2024 17:06
Operator : MA/JU
Sample : P2591-12DL 5X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BHBQ3DL

Manual IntegrationsAPPROVED

Quant Time: Jun 05 17:41:03 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed May 29 15:13:07 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/05/2024
Supervised By :mohammad ahmed 06/07/2024



TIC: BN031803.D\data.ms

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

3.181min (-0.012) 0.80 ng/uL

response 5636

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	70.10	64.44
34.00	0.00	0.00
0.00	0.00	0.00

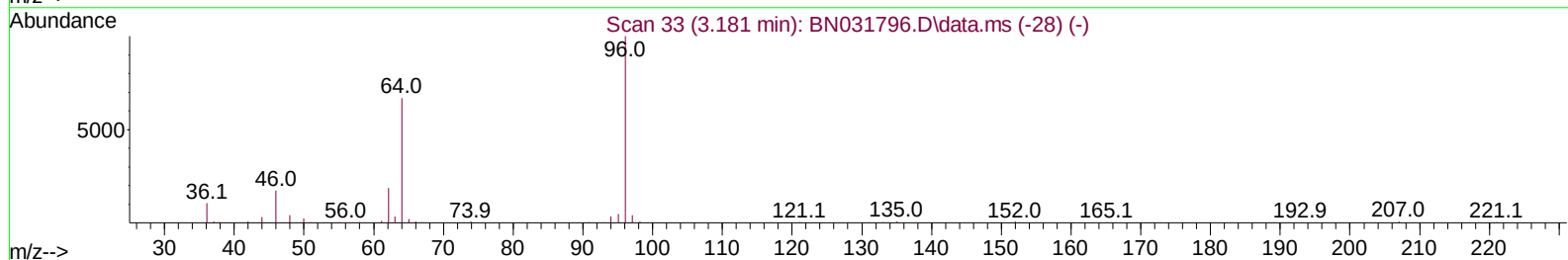
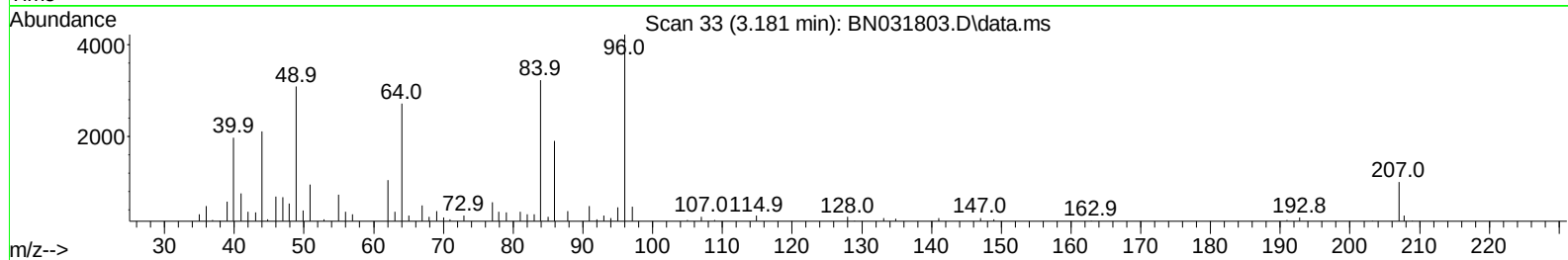
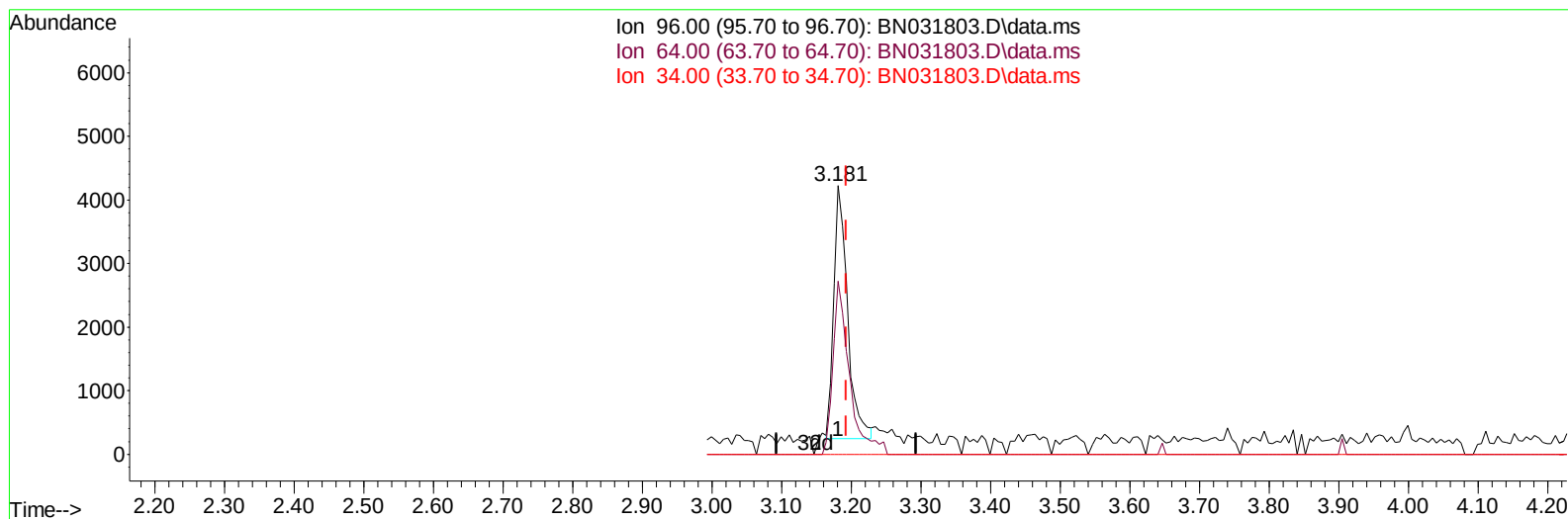
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060524\
Data File : BN031803.D
Acq On : 05 Jun 2024 17:06
Operator : MA/JU
Sample : P2591-12DL 5X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Manual IntegrationsAPPROVED

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TIC: BN031803.D\data.ms

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

3.181min (-0.012) 0.79 ng/uL m

response 5566

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	70.10	64.44
34.00	0.00	0.00
0.00	0.00	0.00

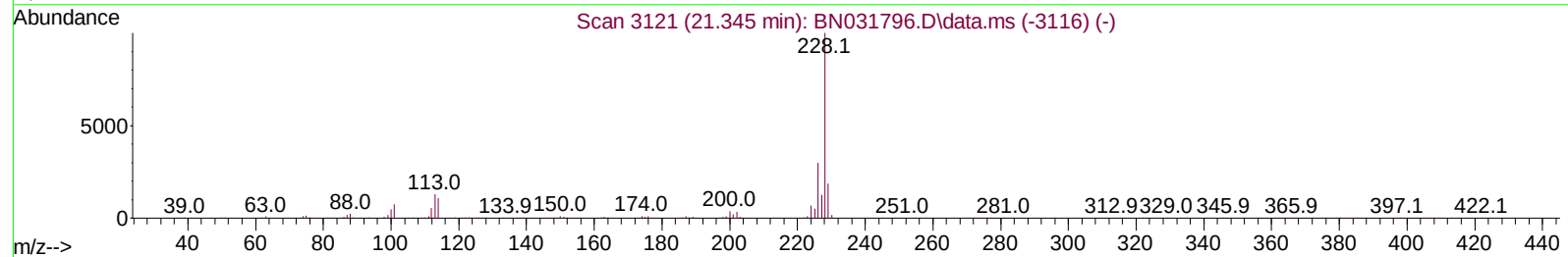
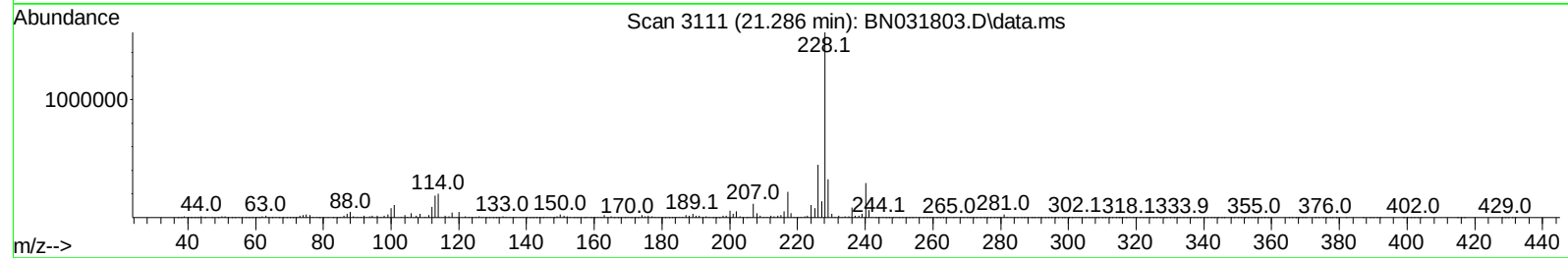
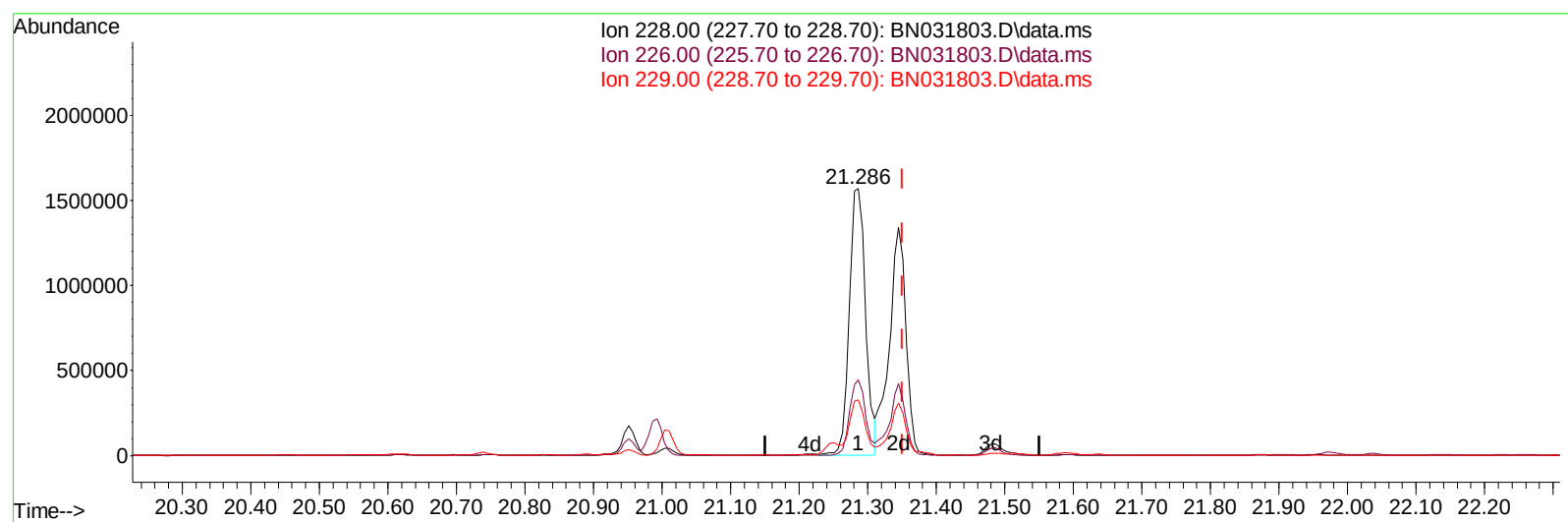
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060524\
 Data File : BN031803.D
 Acq On : 05 Jun 2024 17:06
 Operator : MA/JU
 Sample : P2591-12DL 5X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BHBQ3DL

Manual IntegrationsAPPROVED

Quant Time: Jun 05 17:41:03 2024
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TIC: BN031803.D\data.ms

(87) Chrysene

21.286min (-0.065) 19.65 ng/u1

response 2561529

Ion	Exp%	Act%
228.00	100.00	100.00
226.00	29.60	28.47
229.00	18.90	20.79
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060524\
 Data File : BN031803.D
 Acq On : 05 Jun 2024 17:06
 Operator : MA/JU
 Sample : P2591-12DL 5X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :

BNA_N

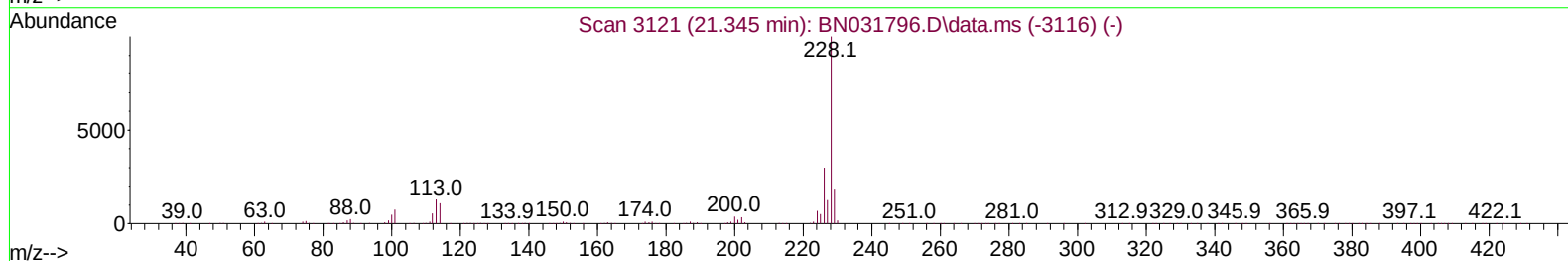
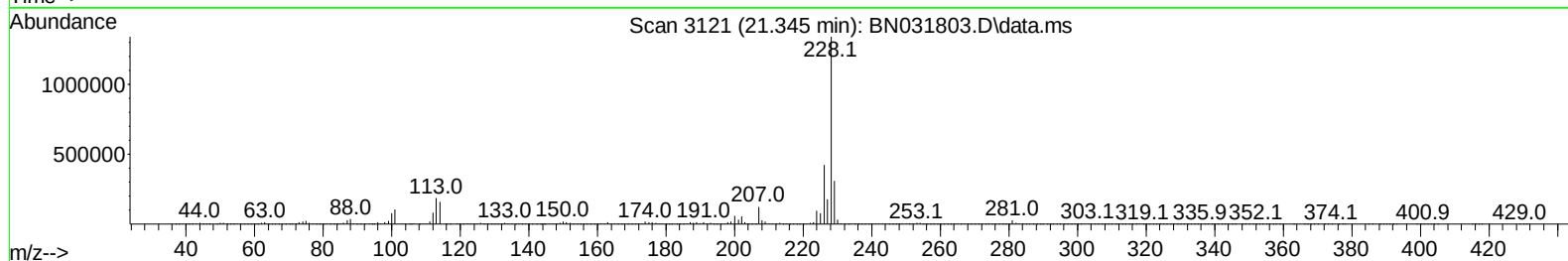
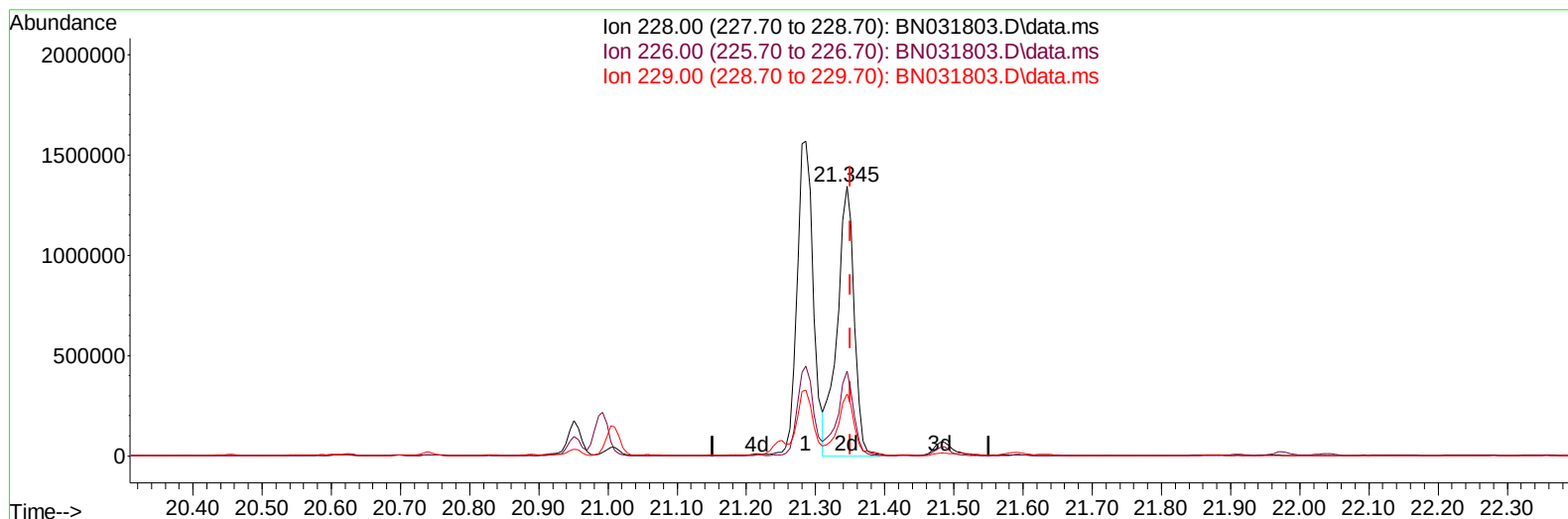
ClientSampleId :

BHBQ3DL

Manual IntegrationsAPPROVED

Quant Time: Jun 05 17:41:03 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
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TIC: BN031803.D\data.ms

(87) Chrysene

21.345min (-0.006) 17.61 ng/ul m

response 2295561

Ion	Exp%	Act%
228.00	100.00	100.00
226.00	29.60	31.62
229.00	18.90	23.01#
0.00	0.00	0.00

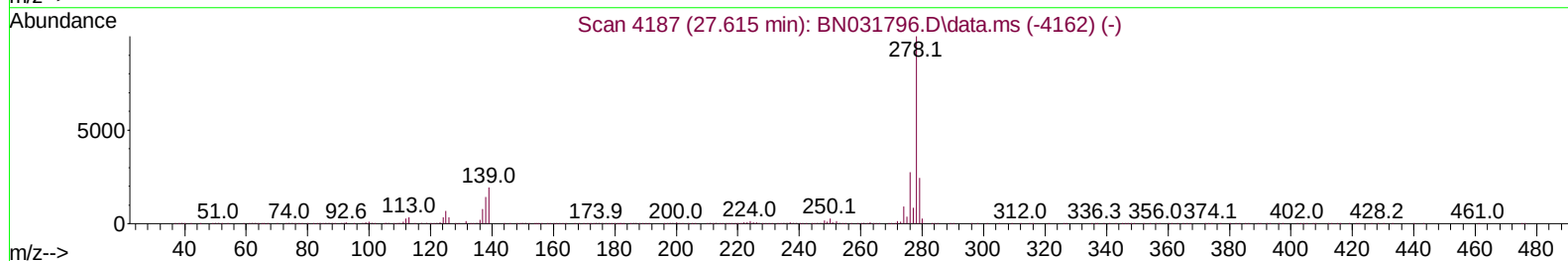
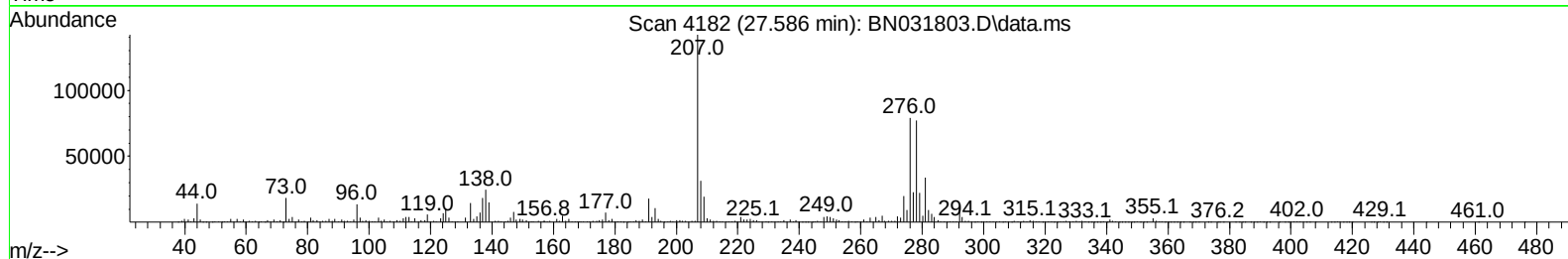
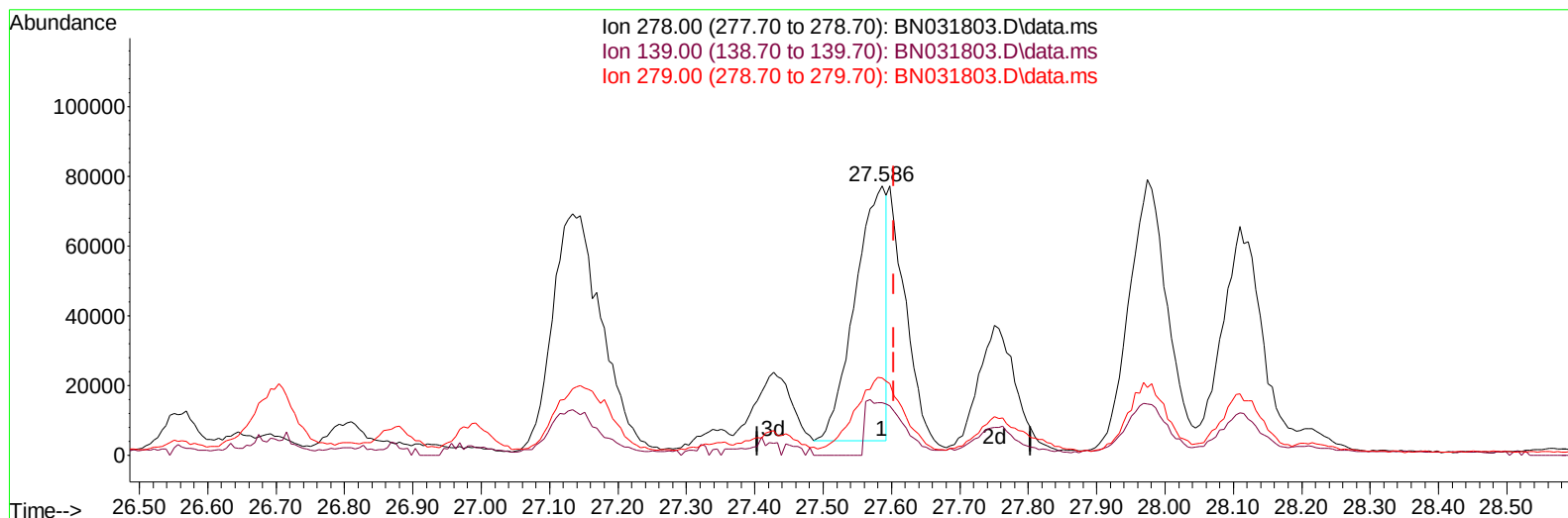
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060524\
Data File : BN031803.D
Acq On : 05 Jun 2024 17:06
Operator : MA/JU
Sample : P2591-12DL 5X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BHBQ3DL

Manual IntegrationsAPPROVED

Quant Time: Jun 05 17:41:03 2024
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TIC: BN031803.D\data.ms

(95) Dibenzo(a,h)anthracene

27.586min (-0.018) 1.70 ng/u1

response 232425

Ion	Exp%	Act%
278.00	100.00	100.00
139.00	18.80	19.47
279.00	21.70	28.66#
0.00	0.00	0.00

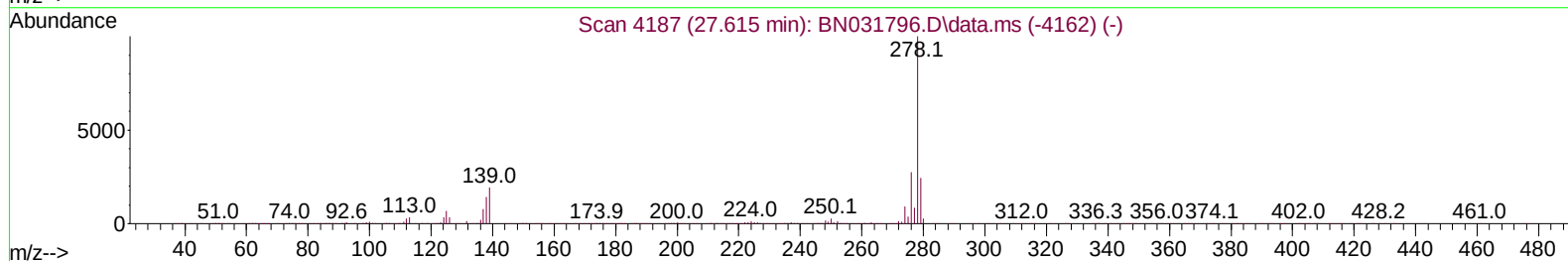
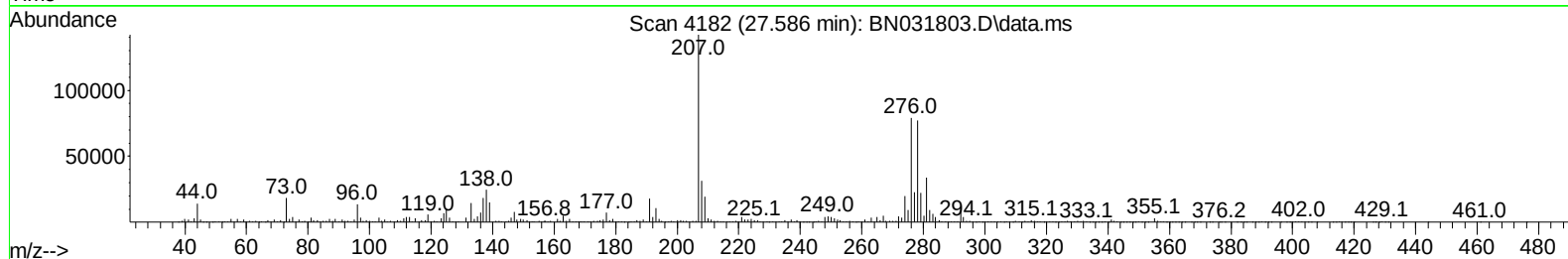
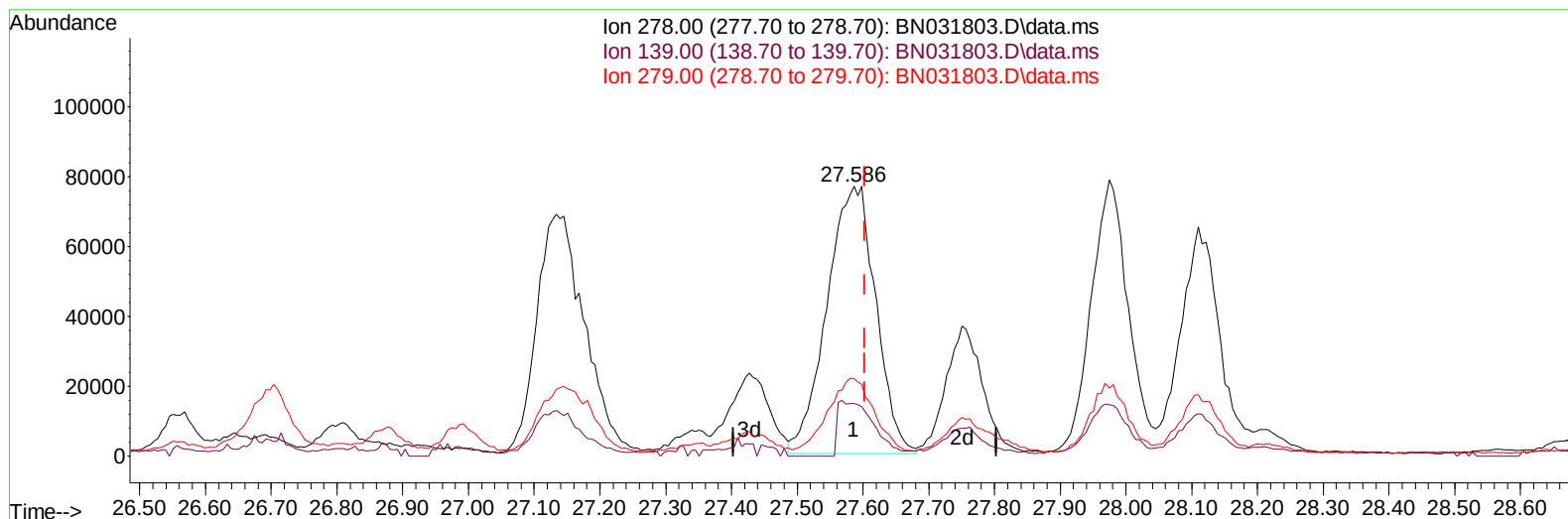
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060524\
Data File : BN031803.D
Acq On : 05 Jun 2024 17:06
Operator : MA/JU
Sample : P2591-12DL 5X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BHBQ3DL

Manual IntegrationsAPPROVED

Quant Time: Jun 05 17:41:03 2024
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Supervised By :mohammad ahmed 06/07/2024



TIC: BN031803.D\data.ms

(95) Dibenzo(a,h)anthracene

27.586min (-0.018) 2.90 ng/ul m

response 395551

Ion	Exp%	Act%
278.00	100.00	100.00
139.00	18.80	19.47
279.00	21.70	28.66#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060524\
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 Operator : MA/JU
 Sample : P2591-12DL 5X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BHBQ3DL

Manual IntegrationsAPPROVED

Quant Time: Jun 06 01:03:09 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed May 29 15:13:07 2024
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/05/2024
 Supervised By :mohammad ahmed 06/07/2024

Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.675	152		293098	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.458	136		1407136	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.316	164		962875	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.063	188		2054414	20.000	ng/ul	0.00
79) Chrysene-d12	21.304	240		2085434	20.000	ng/ul	0.00
88) Perylene-d12	24.239	264		2588374	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.181	96		5566m	0.786	ng/uL	-0.01
4) Pyridine-d5	3.611	84		16590	0.832	ng/ul	0.01
7) Phenol-d5	6.846	99		152752	6.079	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.011	67		85307	5.804	ng/ul	-0.02
11) 2-Chlorophenol-d4	7.205	132		116062	6.054	ng/ul	-0.01
15) 4-Methylphenol-d8	8.375	113		124917	6.099	ng/ul	-0.01
21) Nitrobenzene-d5	8.828	128		62160	5.777	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.546	143		63111	5.679	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.081	165		121925	5.876	ng/ul	0.00
31) 4-Chloroaniline-d4	10.599	131		130536	4.237	ng/ul	-0.01
46) Dimethylphthalate-d6	13.728	166		427964	6.380	ng/ul	-0.02
49) Acenaphthylene-d8	14.010	160		480657	6.232	ng/ul	0.00
54) 4-Nitrophenol-d4	14.528	143		54509	4.058	ng/ul	0.00
60) Fluorene-d10	15.310	176		381023	6.713	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	0.000	200		0d	0.000	ng/ul	
73) Anthracene-d10	17.163	188		608454	6.918	ng/ul	0.00
81) Pyrene-d10	19.463	212		709022	6.353	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.039	264		636887	5.153	ng/ul	0.00
Target Compounds							
						Qvalue	
52) Acenaphthene	14.381	153		76192	1.321	ng/ul	98
56) Dibenzofuran	14.716	168		114007	1.455	ng/ul	97
61) Fluorene	15.369	166		102698	1.627	ng/ul	98
72) Phenanthrene	17.104	178		2548576	24.352	ng/ul	97
74) Anthracene	17.198	178		548089	5.169	ng/ul	99
77) Carbazole	17.469	167		327602	3.623	ng/ul	99
80) Fluoranthene	19.128	202		4840138	36.292	ng/ul	99
82) Pyrene	19.492	202		3619826	26.454	ng/ul	99
85) Benzo(a)anthracene	21.286	228		2565927	18.382	ng/ul	100
87) Chrysene	21.345	228		2295561m	17.606	ng/ul	
90) Benzo(b)fluoranthene	23.316	252		3430496	22.173	ng/ul	97
91) Benzo(k)fluoranthene	23.374	252		1094096	7.207	ng/ul	99
93) Benzo(a)pyrene	24.104	252		1549442	10.693	ng/ul	96
94) Indeno(1,2,3-cd)pyrene	27.545	276		1125610	6.741	ng/ul	96
95) Dibenzo(a,h)anthracene	27.586	278		395551m	2.902	ng/ul	
96) Benzo(g,h,i)perylene	28.574	276		166717	1.256	ng/ul	99

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

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