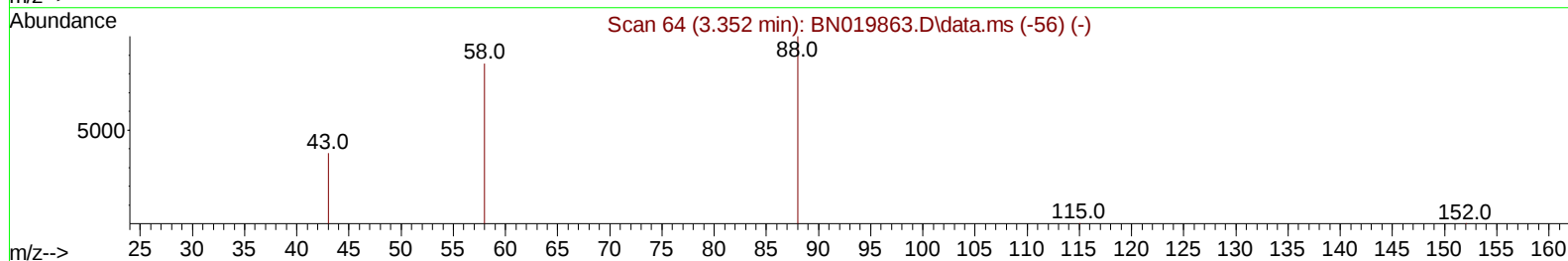
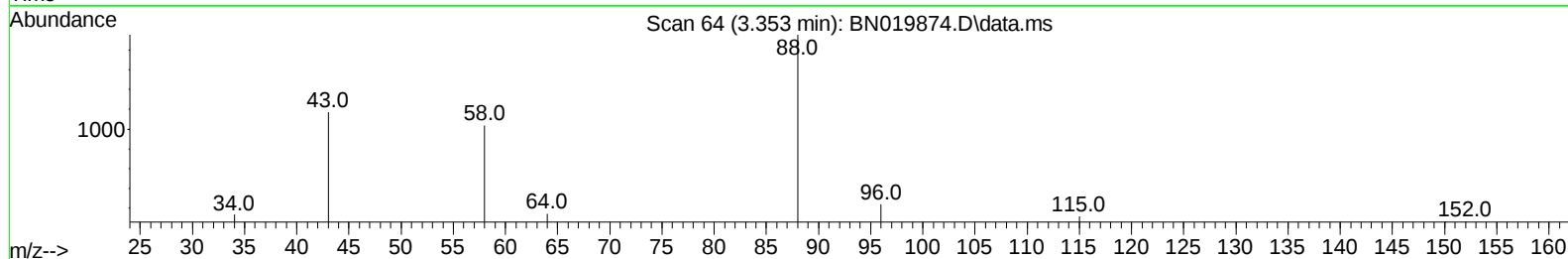
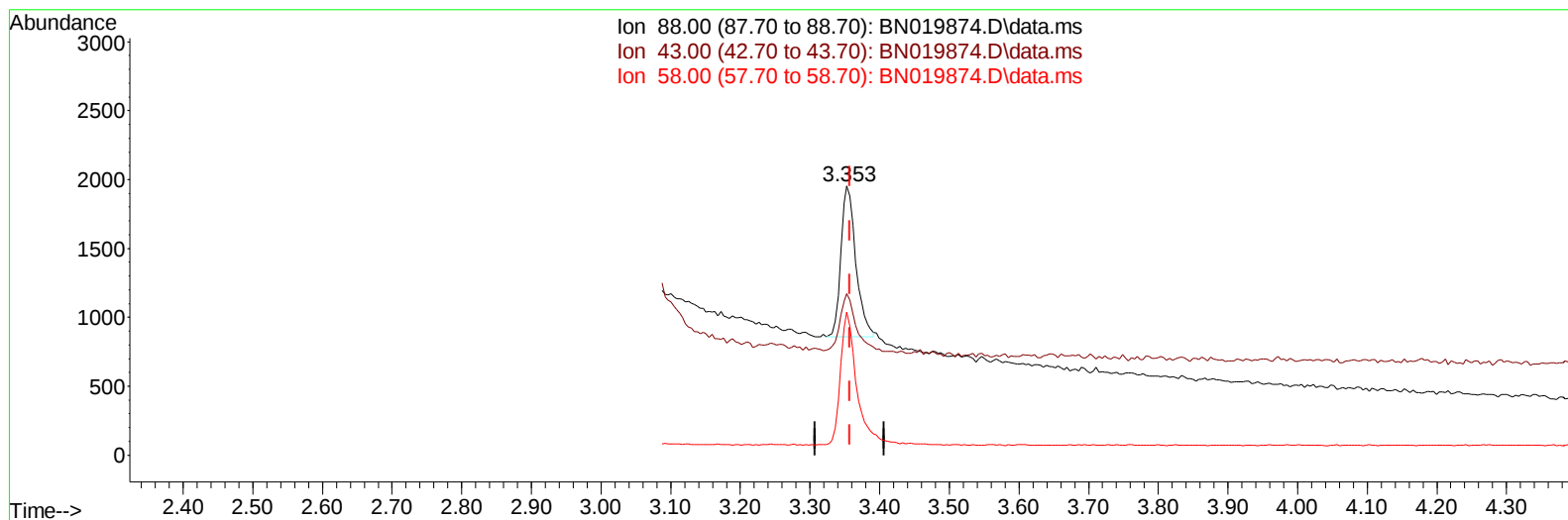


Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN052022\
Data File : BN019874.D
Acq On : 20 May 2022 18:16
Operator : CG/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 13 Sample Multiplier: 1

Quant Time: May 21 01:53:52 2022
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN042922.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue May 17 01:41:05 2022
Response via : Initial Calibration



TIC: BN019874.D\data.ms

(2) 1,4-Dioxane

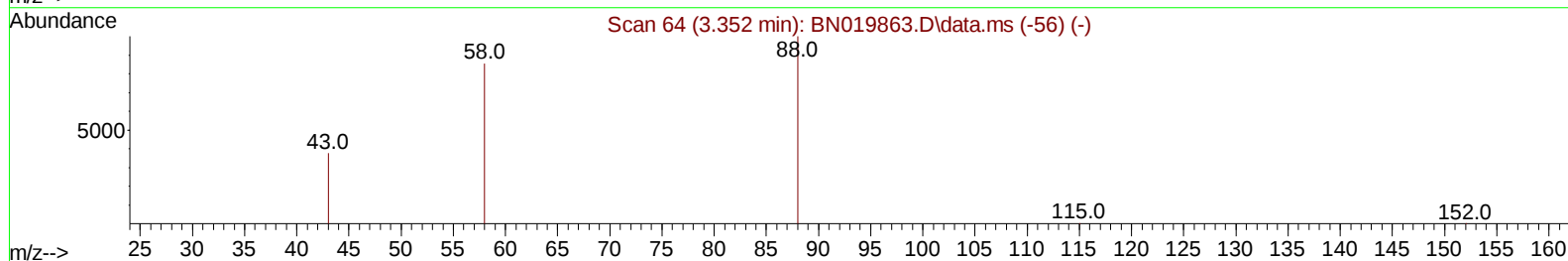
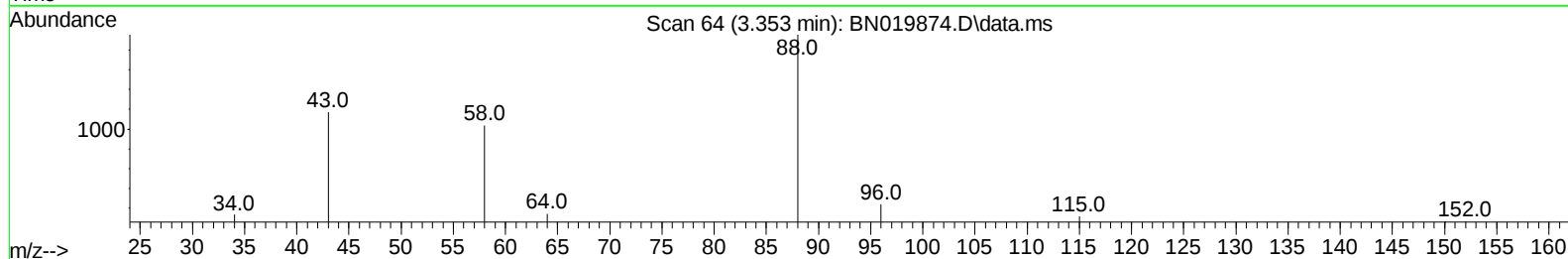
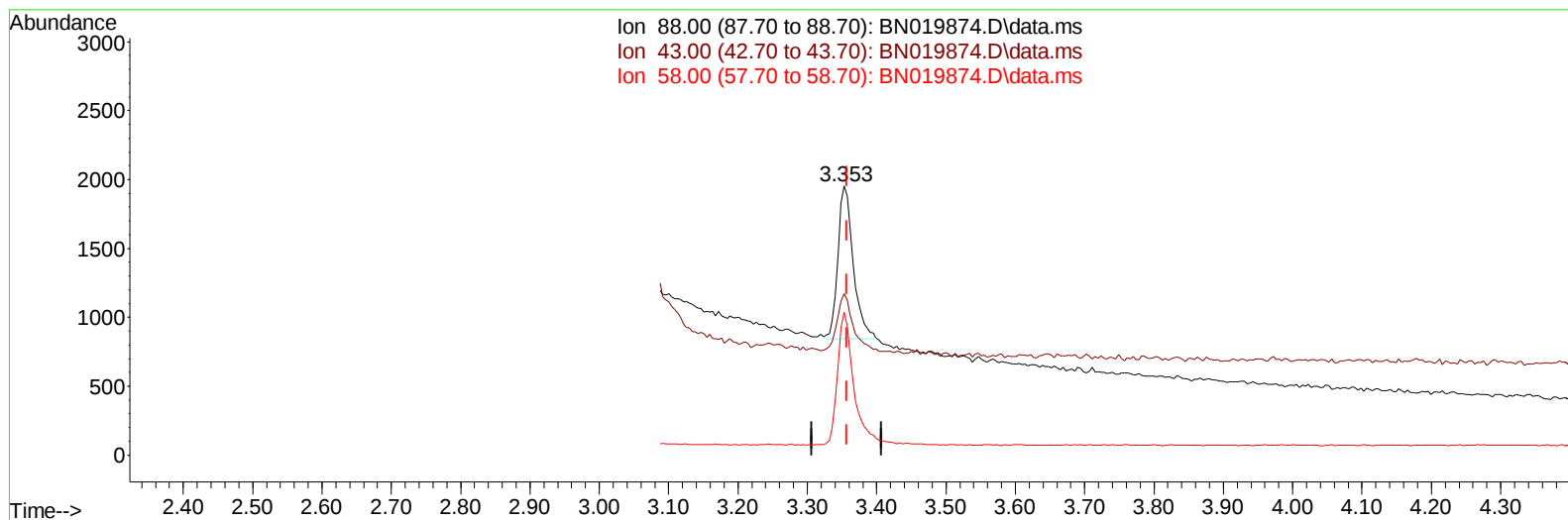
3.353min (-0.004) 0.36 ng/u1

response 1631

Ion	Exp%	Act%
88.00	100.00	100.00
43.00	48.70	60.04#
58.00	49.80	53.23
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN052022\
Data File : BN019874.D
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Misc :
ALS Vial : 13 Sample Multiplier: 1

Quant Time: May 21 01:53:52 2022
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN042922.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
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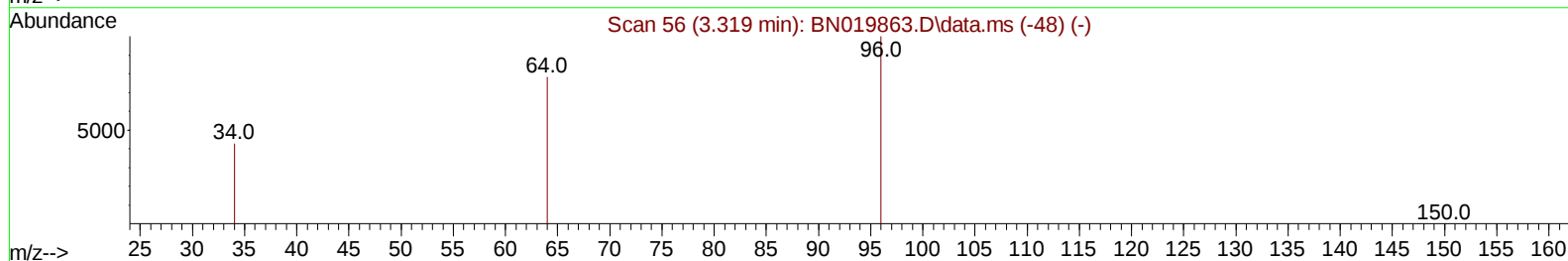
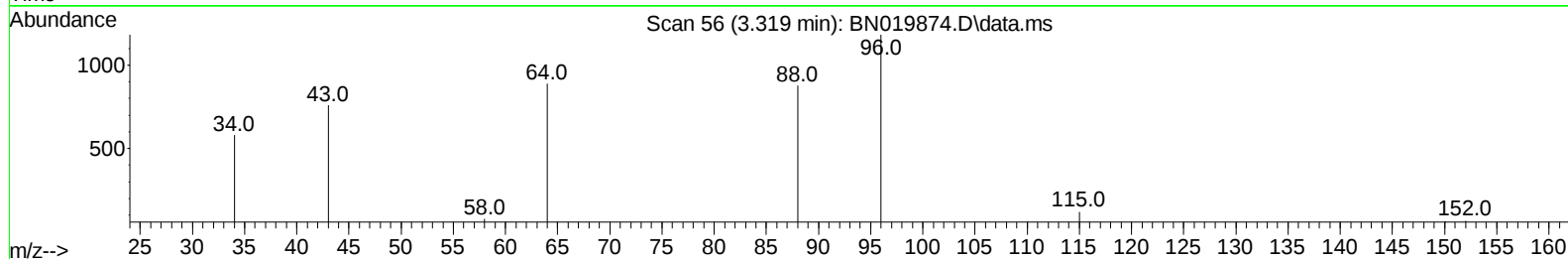
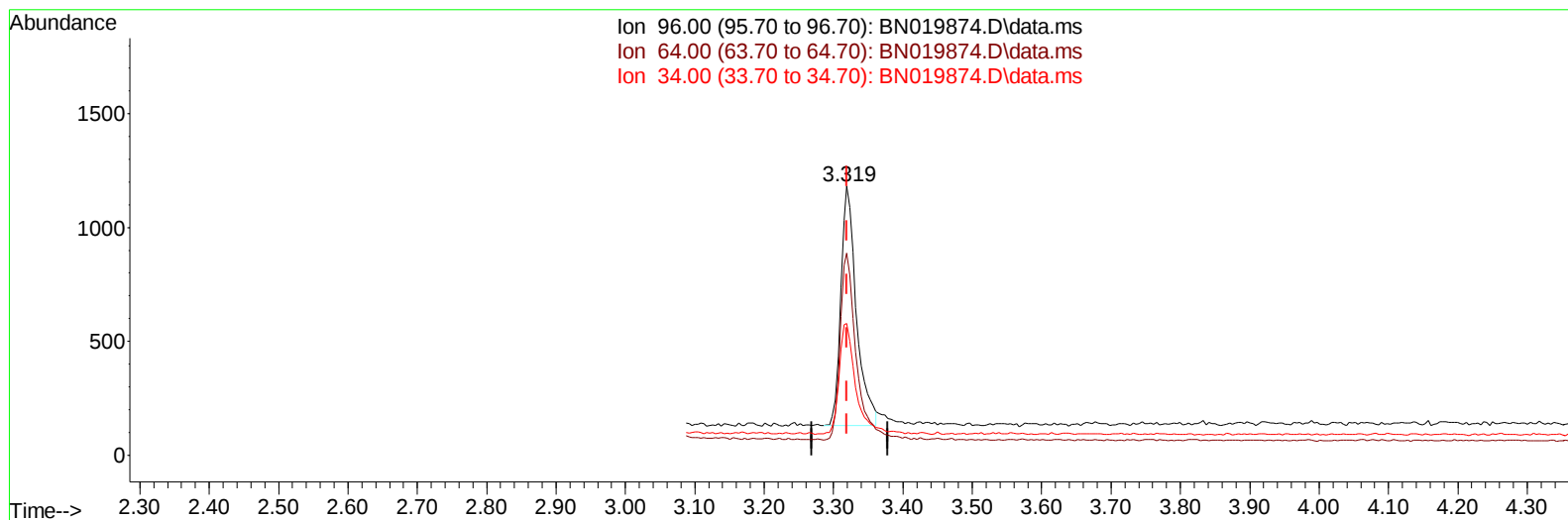
TIC: BN019874.D\data.ms

(2) 1,4-Dioxane**3.353min (-0.004) 0.37 ng/ul m****response 1686**

Ion	Exp%	Act%
88.00	100.00	100.00
43.00	48.70	60.04#
58.00	49.80	53.23
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN052022\
Data File : BN019874.D
Acq On : 20 May 2022 18:16
Operator : CG/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 13 Sample Multiplier: 1

Quant Time: May 21 01:53:52 2022
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN042922.M
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QLast Update : Tue May 17 01:41:05 2022
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TIC: BN019874.D\data.ms

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

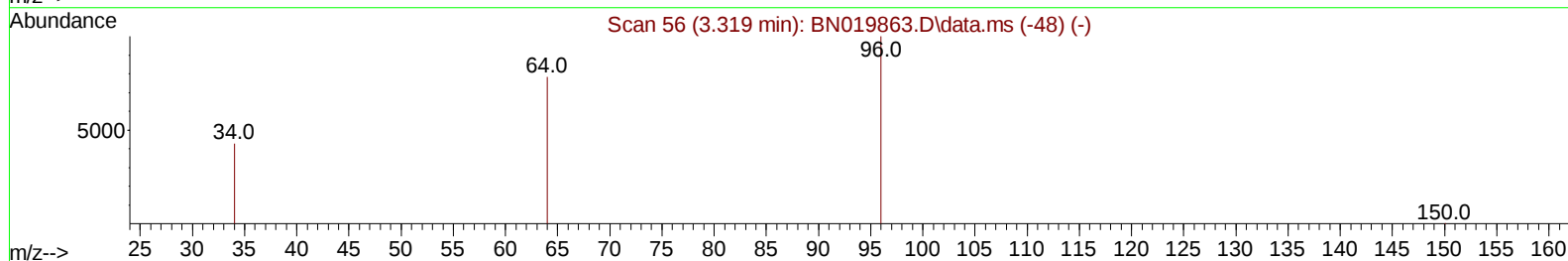
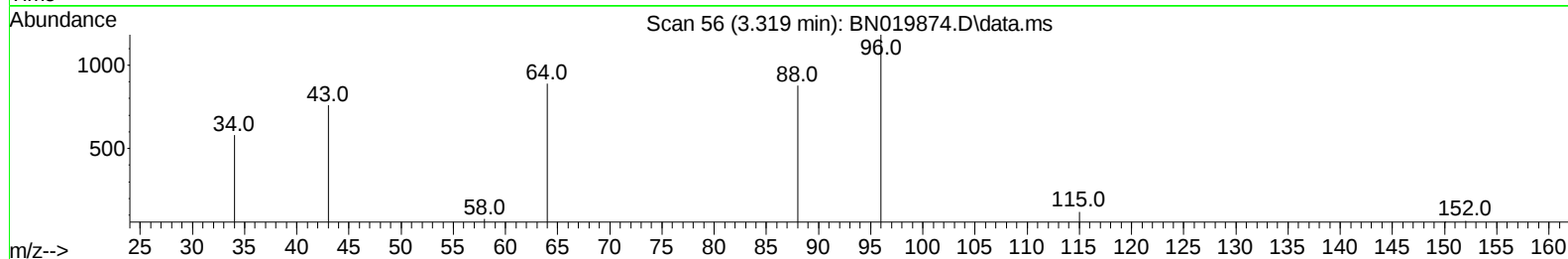
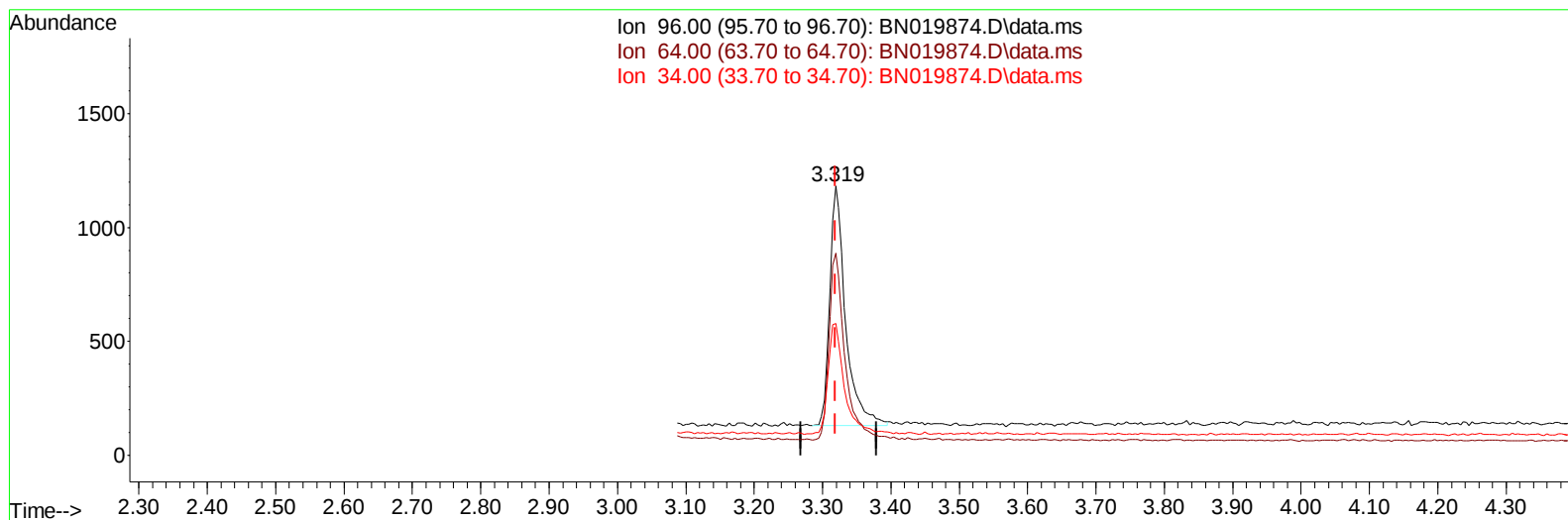
3.319min (+ 0.000) 0.38 ng/u1

response 1629

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	40.10	75.13#
34.00	22.70	48.98#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN052022\
Data File : BN019874.D
Acq On : 20 May 2022 18:16
Operator : CG/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 13 Sample Multiplier: 1

Quant Time: May 21 01:53:52 2022
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN042922.M
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TIC: BN019874.D\data.ms

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

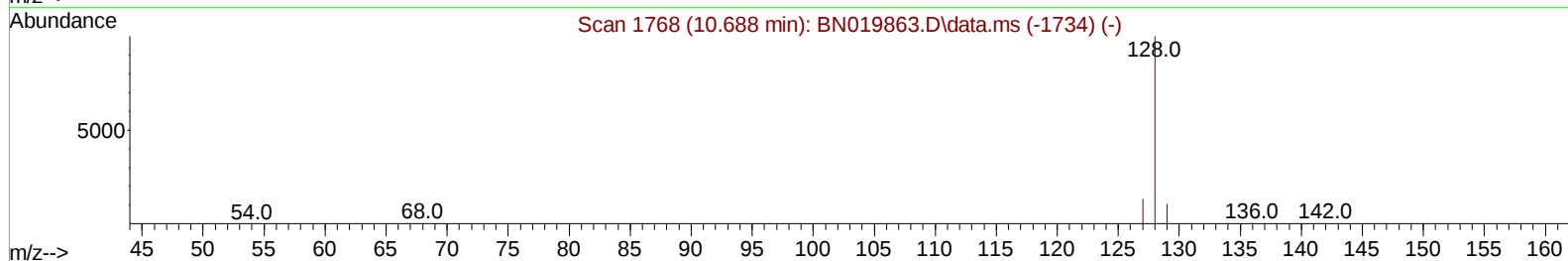
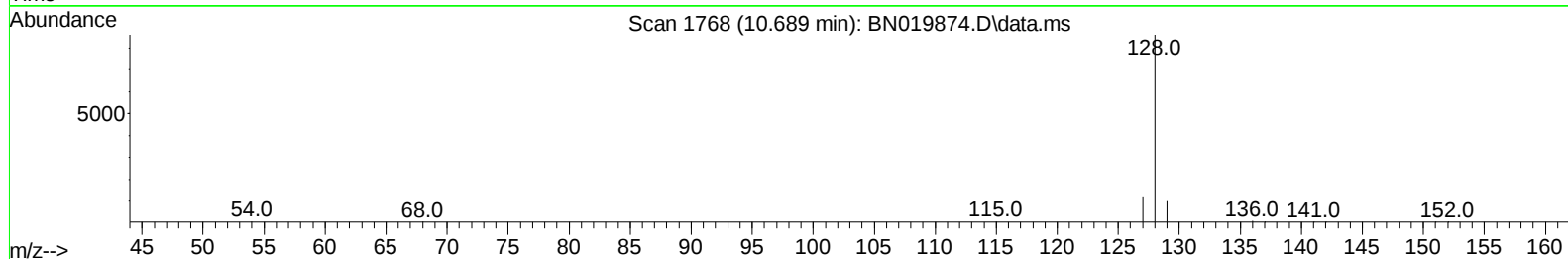
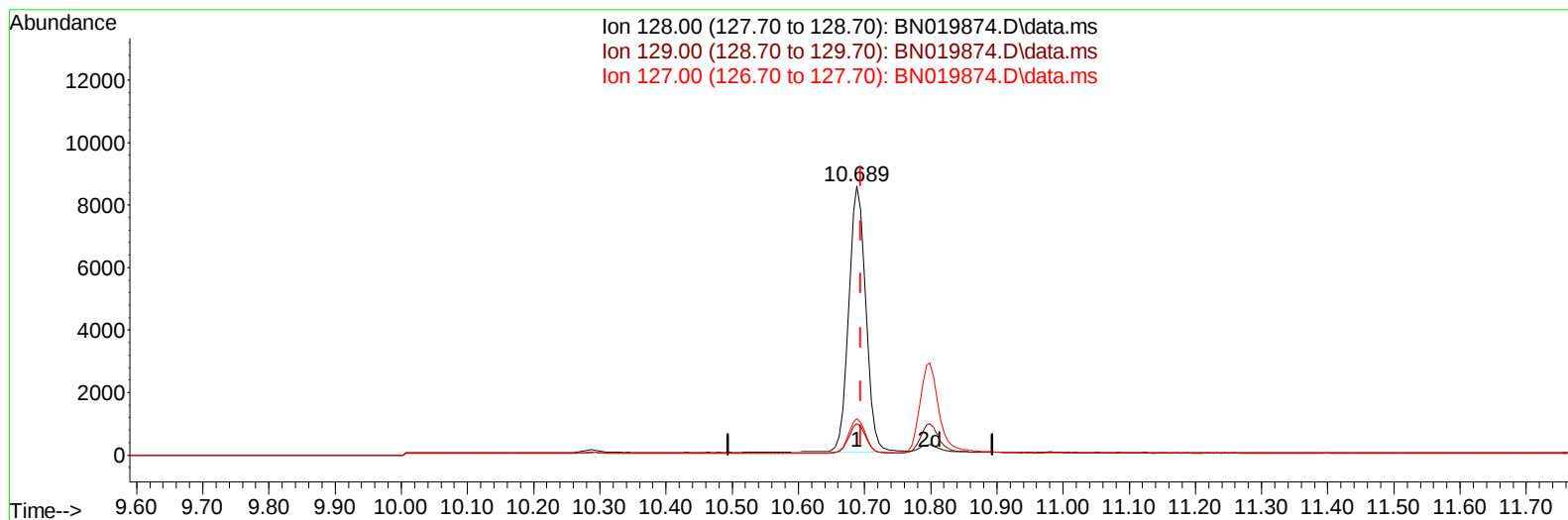
3.319min (+ 0.000) 0.40 ng/ul m

response 1706

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	40.10	75.13#
34.00	22.70	48.98#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN052022\
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Operator : CG/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 13 Sample Multiplier: 1

Quant Time: May 21 01:53:52 2022
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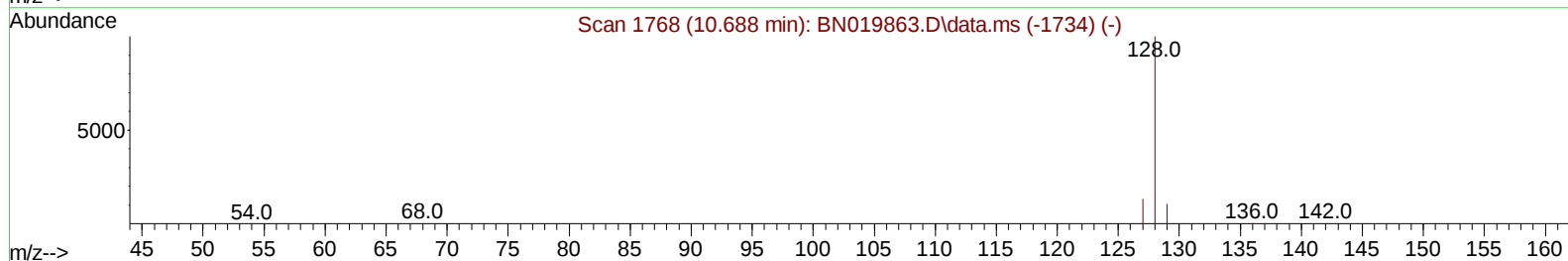
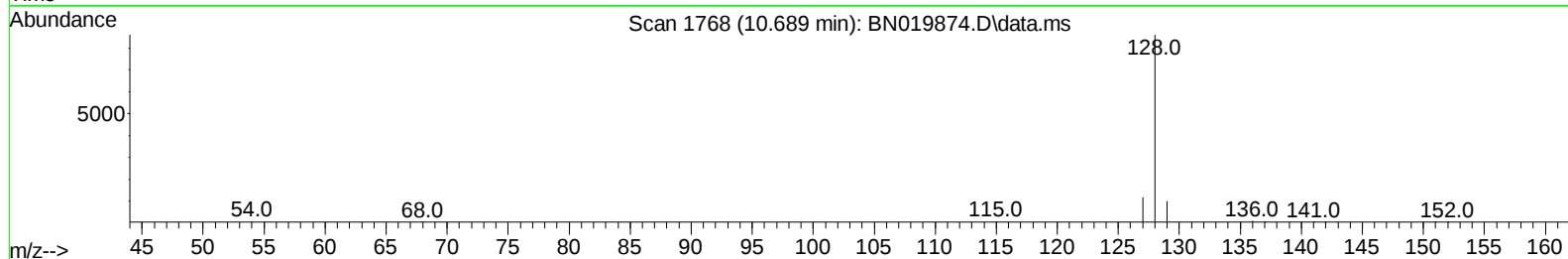
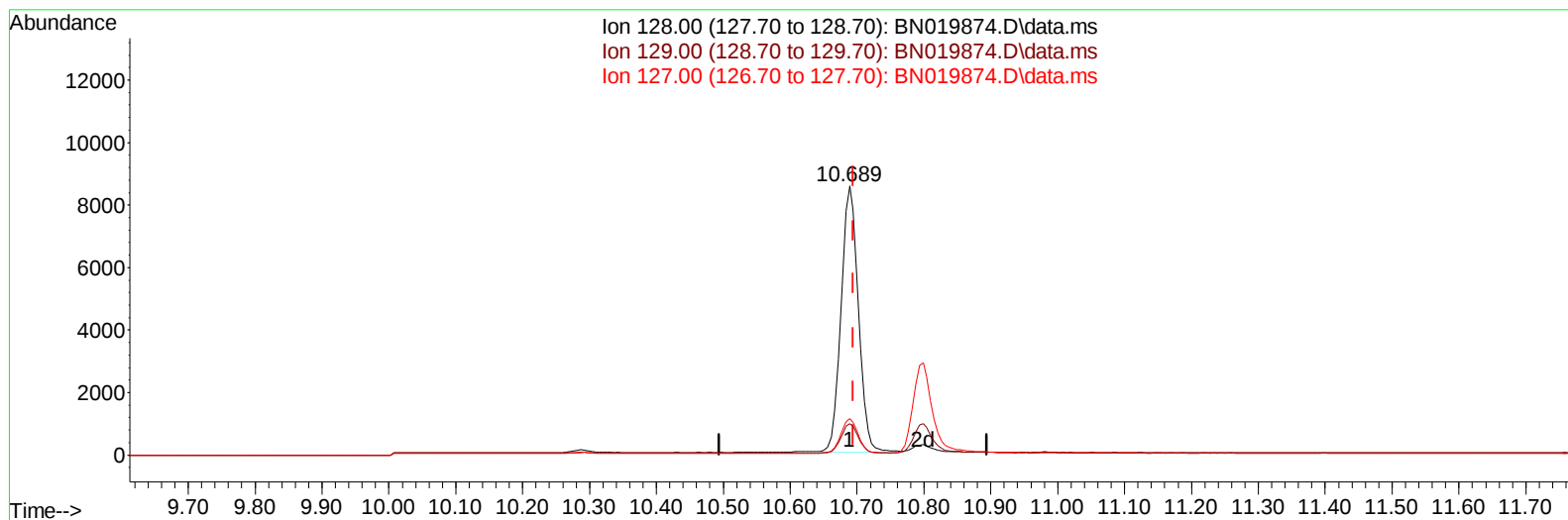
TIC: BN019874.D\data.ms

(5) Naphthalene**10.689min (-0.005) 0.39 ng/ul****response 15299**

Ion	Exp%	Act%
128.00	100.00	100.00
129.00	11.90	11.67
127.00	13.70	13.66
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN052022\
Data File : BN019874.D
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Operator : CG/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 13 Sample Multiplier: 1

Quant Time: May 21 01:53:52 2022
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN042922.M
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TIC: BN019874.D\data.ms

(5) Naphthalene**10.689min (-0.005) 0.39 ng/ul m****response 15246**

Ion	Exp%	Act%
128.00	100.00	100.00
129.00	11.90	11.67
127.00	13.70	13.66
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN052022\
 Data File : BN019874.D
 Acq On : 20 May 2022 18:16
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Quant Time: May 21 01:53:52 2022
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 17 01:41:05 2022
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.855	152	3583	0.400	ng/ul	0.00
4) Naphthalene-d8	10.639	136	12107	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.474	164	7691	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.213	188	16313	0.400	ng/ul	0.00
17) Chrysene-d12	21.400	240	15619	0.400	ng/ul	0.00
23) Perylene-d12	23.738	264	12943	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.319	96	1706m	0.402	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.229	152	8652	0.472	ng/ul	0.00
18) Fluoranthene-d10	19.239	212	19980	0.404	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.353	88	1686m	0.374	ng/ul	
5) Naphthalene	10.689	128	15246m	0.389	ng/ul	
7) 2-Methylnaphthalene	12.300	142	10262	0.430	ng/ul#	100
8) 1-Methylnaphthalene	12.520	142	10344	0.468	ng/ul#	100
10) Acenaphthylene	14.192	152	16174	0.425	ng/ul	99
11) Acenaphthene	14.534	153	11504	0.376	ng/ul	99
12) Fluorene	15.520	166	13279	0.396	ng/ul	99
14) Pentachlorophenol	16.867	266	1272	0.362	ng/ul	91
15) Phenanthrene	17.255	178	22438	0.374	ng/ul	99
16) Anthracene	17.344	178	20436	0.417	ng/ul	99
19) Fluoranthene	19.271	202	26008	0.364	ng/ul	98
20) Pyrene	19.634	202	26618	0.373	ng/ul	98
21) Benzo(a)anthracene	21.382	228	23697	0.406	ng/ul	100
22) Chrysene	21.435	228	24198	0.352	ng/ul	99
24) Benzo(b)fluoranthene	23.028	252	22287	0.339	ng/ul	92
25) Benzo(k)fluoranthene	23.074	252	21063	0.333	ng/ul	98
26) Benzo(a)pyrene	23.633	252	20433	0.365	ng/ul#	78
27) Indeno(1,2,3-cd)pyrene	26.141	276	20418	0.293	ng/ul#	96
28) Dibenzo(a,h)anthracene	26.158	278	16743	0.287	ng/ul#	82
29) Benzo(g,h,i)perylene	26.875	276	16912	0.265	ng/ul	97

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

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