

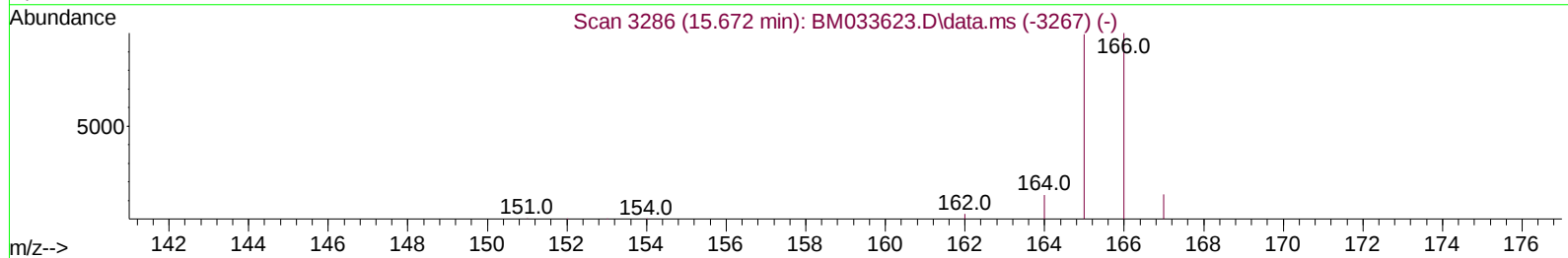
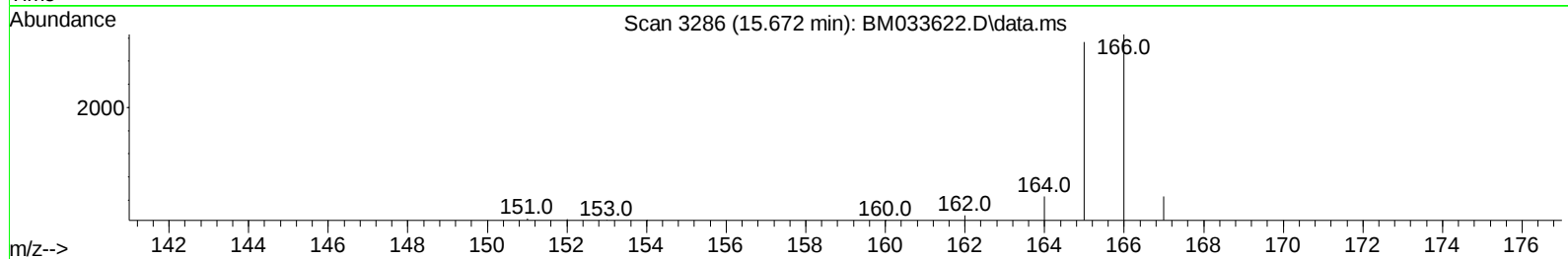
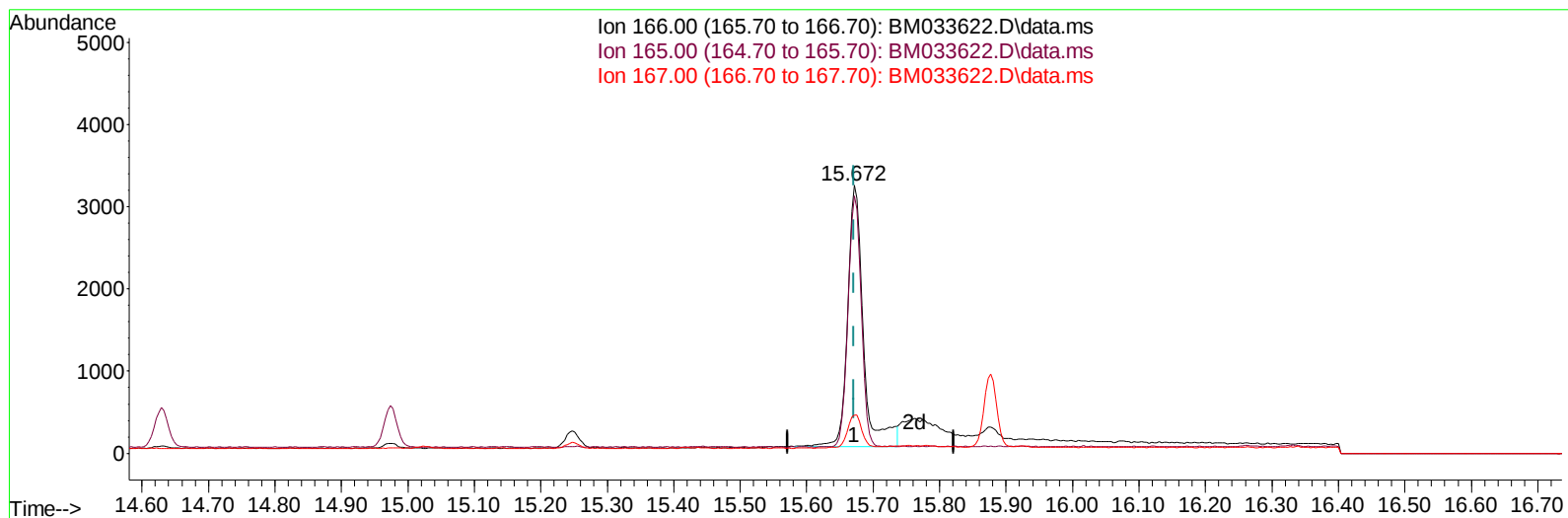
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010422\
Data File : BM033622.D
Acq On : 04 Jan 2022 12:59
Operator : CG/JU
Sample : SST00.237
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SST00.2037

Manual IntegrationsAPPROVED

Quant Time: Jan 04 14:19:03 2022
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM010422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jan 04 14:17:24 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 01/04/2022
Supervised By :Yogesh Patel 01/06/2022



TIC: BM033622.D\data.ms

(12) Fluorene

15.672min (+ 0.000) 0.22 ng/u1

response 5241

Ion	Exp%	Act%
166.00	100.00	100.00
165.00	101.40	96.07
167.00	18.60	14.33#
0.00	0.00	0.00

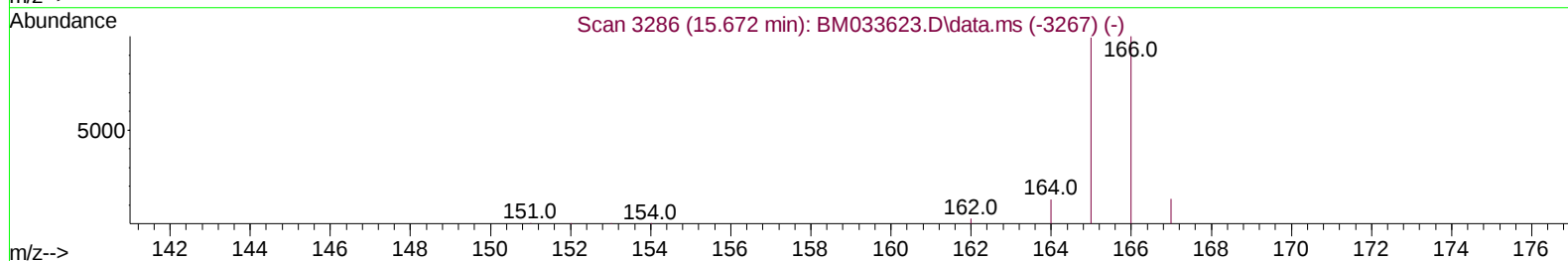
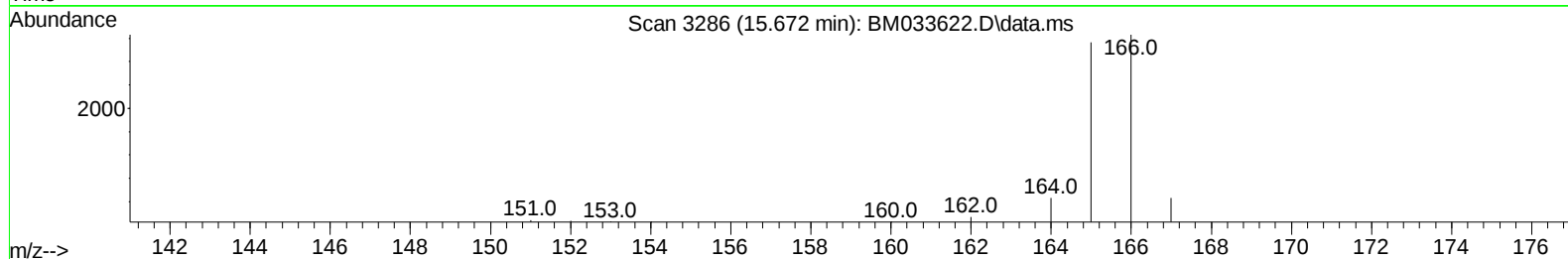
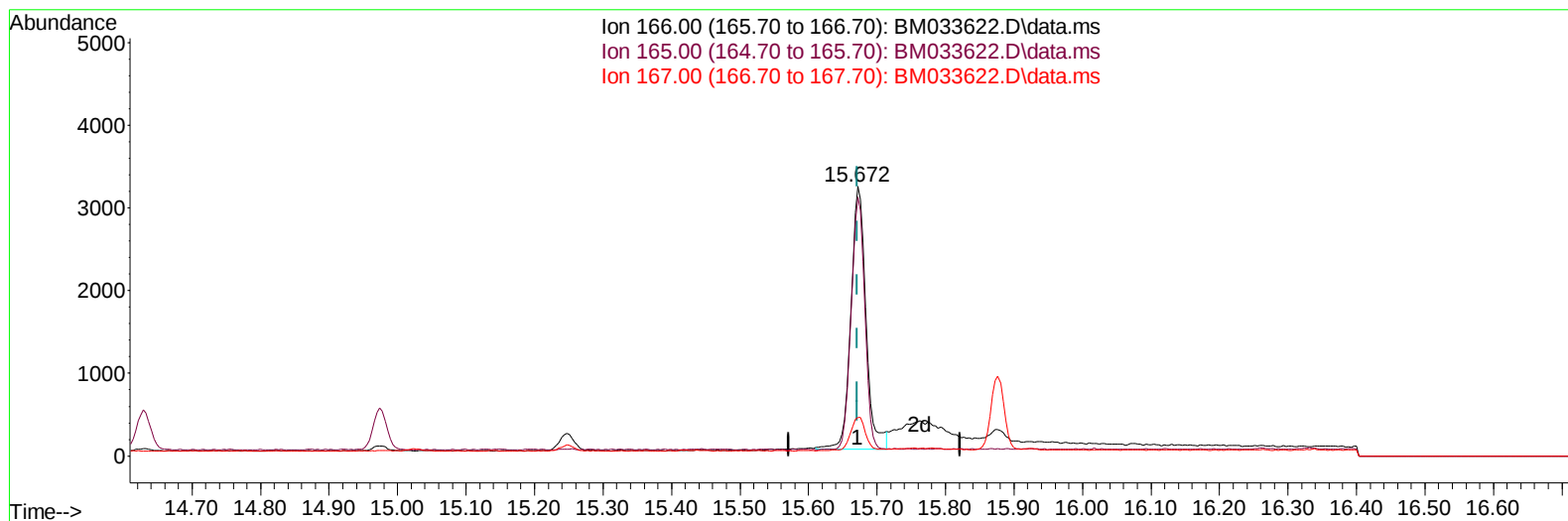
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010422\
 Data File : BM033622.D
 Acq On : 04 Jan 2022 12:59
 Operator : CG/JU
 Sample : SST0.237
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST0.2037

Manual IntegrationsAPPROVED

Quant Time: Jan 04 14:19:03 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM010422.M
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 Supervised By :Yogesh Patel 01/06/2022



TIC: BM033622.D\data.ms

(12) Fluorene

15.672min (+ 0.000) 0.20 ng/ul m

response 4884

Ion	Exp%	Act%
166.00	100.00	100.00
165.00	101.40	96.07
167.00	18.60	14.33#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010422\
 Data File : BM033622.D
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 Operator : CG/JU
 Sample : SST00.237
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST00.2037

Manual IntegrationsAPPROVED

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.991	152	3332	0.400	ng/ul	0.00
4) Naphthalene-d8	10.810	136	10812	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.629	164	5309	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.365	188	9439	0.400	ng/ul	0.00
17) Chrysene-d12	21.528	240	6780	0.400	ng/ul	# 0.00
23) Perylene-d12	23.962	264	7114	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.331	96	750	0.264	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.398	152	3002	0.208	ng/ul	0.00
18) Fluoranthene-d10	19.384	212	4320	0.276	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.369	88	823	0.223	ng/ul	98
5) Naphthalene	10.860	128	6435	0.180	ng/ul#	89
7) 2-Methylnaphthalene	12.470	142	4098	0.168	ng/ul	99
8) 1-Methylnaphthalene	12.686	142	3986	0.175	ng/ul	99
10) Acenaphthylene	14.348	152	5000	0.176	ng/ul#	92
11) Acenaphthene	14.690	153	4009	0.185	ng/ul	97
12) Fluorene	15.672	166	4884m	0.203	ng/ul	
14) Pentachlorophenol	17.011	266	765	0.162	ng/ul	99
15) Phenanthrene	17.406	178	6178	0.202	ng/ul	96
16) Anthracene	17.497	178	5620	0.196	ng/ul	94
19) Fluoranthene	19.414	202	6485	0.312	ng/ul	95
20) Pyrene	19.772	202	6573	0.302	ng/ul#	91
21) Benzo(a)anthracene	21.514	228	5776	0.310	ng/ul	99
22) Chrysene	21.564	228	5990	0.285	ng/ul	98
24) Benzo(b)fluoranthene	23.216	252	6388	0.284	ng/ul	92
25) Benzo(k)fluoranthene	23.267	252	6201	0.263	ng/ul#	90
26) Benzo(a)pyrene	23.849	252	5477	0.263	ng/ul#	87
27) Indeno(1,2,3-cd)pyrene	26.506	276	7269	0.252	ng/ul#	77
28) Dibenzo(a,h)anthracene	26.521	278	5709	0.247	ng/ul	95
29) Benzo(g,h,i)perylene	27.282	276	6247	0.245	ng/ul#	92

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

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