

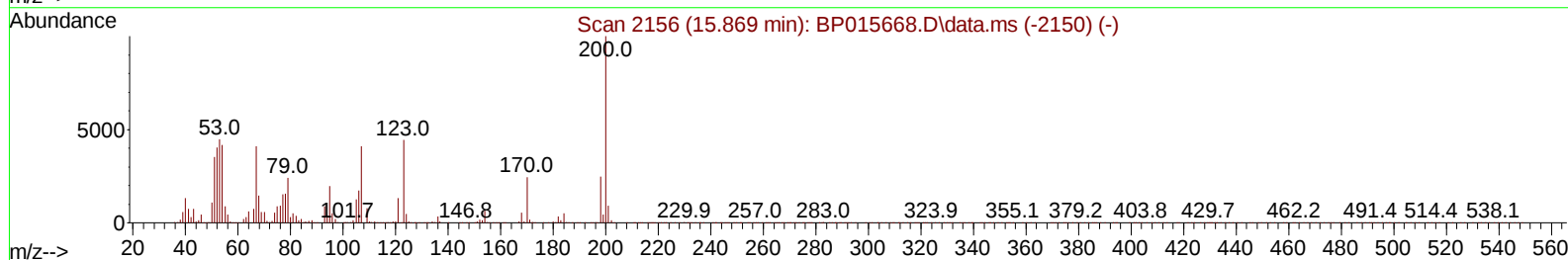
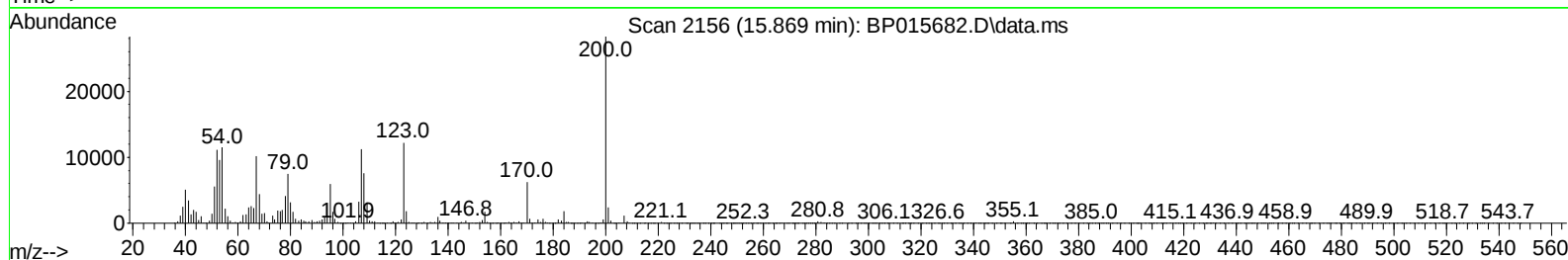
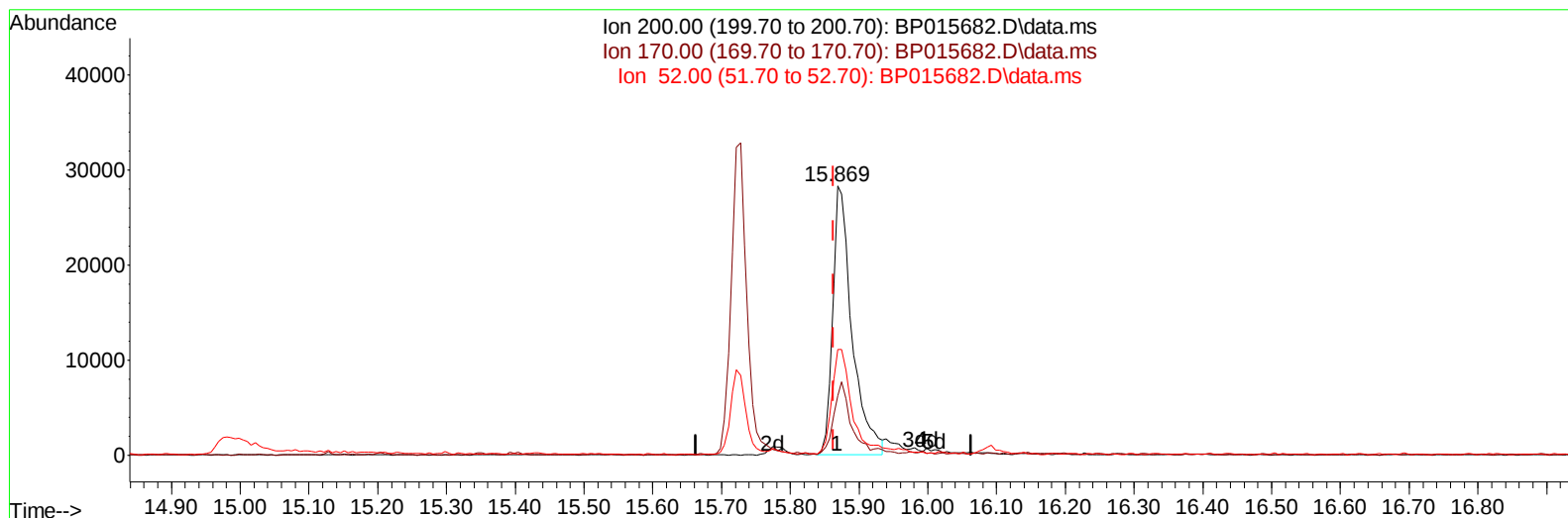
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015682.D
Acq On : 23 Jun 2023 21:08
Operator : MA/JU
Sample : 03084-05
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFM2

Manual IntegrationsAPPROVED

Quant Time: Jun 23 22:38:04 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jun 22 22:01:49 2023
Response via : Initial Calibration

Reviewed By :Yogesh Patel 06/26/2023
Supervised By :mohammad ahmed 06/27/2023



TIC: BP015682.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.869min (+ 0.006) 13.91 ng/u1

response 55298

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	22.30	22.16
52.00	46.20	39.26
0.00	0.00	0.00

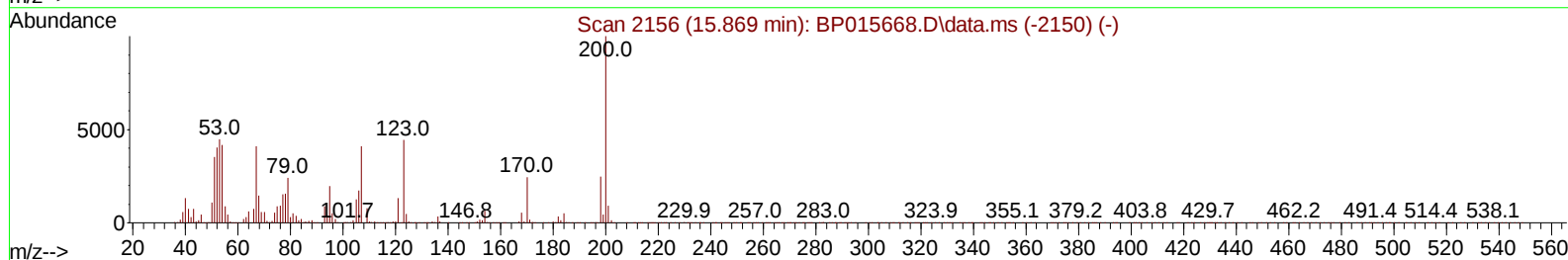
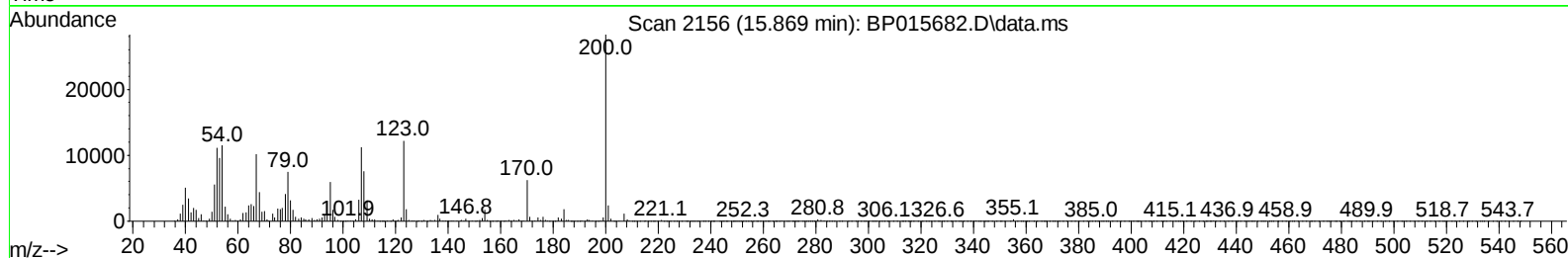
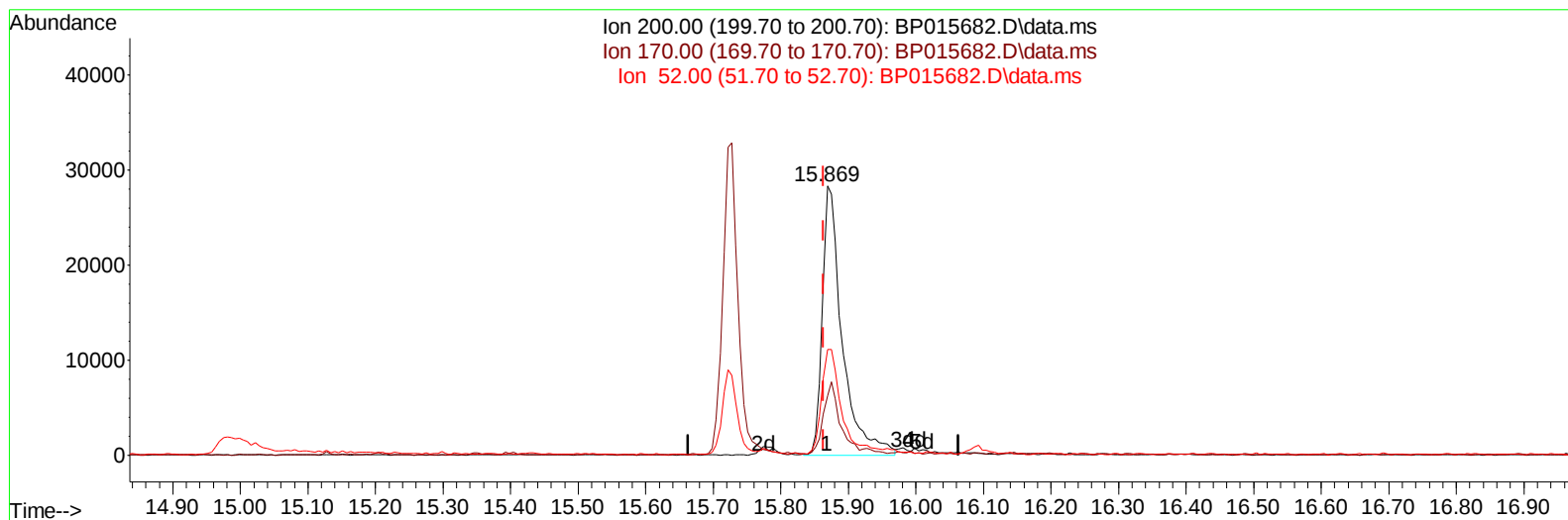
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TIC: BP015682.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)**15.869min (+ 0.006) 14.55 ng/ul m****response 57832**

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	22.30	22.16
52.00	46.20	39.26
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
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 Acq On : 23 Jun 2023 21:08
 Operator : MA/JU
 Sample : 03084-05
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCFM2

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 06/26/2023
 Supervised By :mohammad ahmed 06/27/2023

Quant Time: Jun 24 00:30:30 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jun 22 22:01:49 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.081	152	147305	20.000	ng/ul	0.00
20) Naphthalene-d8	10.904	136	558461	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.728	164	303410	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.492	188	627500	20.000	ng/ul	0.00
79) Chrysene-d12	21.580	240	608805	20.000	ng/ul	0.00
88) Perylene-d12	24.139	264	756931	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.399	96	14378	3.614	ng/uL	0.00
4) Pyridine-d5	3.852	84	51117	4.947	ng/ul	0.01
7) Phenol-d5	7.246	99	231835	17.084	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.399	67	162677	20.072	ng/ul	0.00
11) 2-Chlorophenol-d4	7.605	132	200787	20.407	ng/ul	0.00
15) 4-Methylphenol-d8	8.805	113	114129	10.247	ng/ul	0.00
21) Nitrobenzene-d5	9.263	128	96494	22.008	ng/ul	0.00
24) 2-Nitrophenol-d4	9.993	143	104580	20.887	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.546	165	161350	17.665	ng/ul	0.01
31) 4-Chloroaniline-d4	11.069	131	109183	8.328	ng/ul	0.02
46) Dimethylphthalate-d6	14.134	166	517364	21.624	ng/ul	0.00
49) Acenaphthylene-d8	14.422	160	595946	22.779	ng/ul	0.00
54) 4-Nitrophenol-d4	14.981	143	48670	11.359	ng/ul	0.03
60) Fluorene-d10	15.722	176	427949	21.211	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.869	200	57832m	14.548	ng/ul	0.00
73) Anthracene-d10	17.592	188	597875	20.941	ng/ul	0.00
81) Pyrene-d10	19.816	212	779779	23.931	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.974	264	832602	21.922	ng/ul	0.00

Target Compounds Qvalue

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

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Misc :
ALS Vial : 16 Sample Multiplier: 3

Instrument :
BNA_P
ClientSampleId :
DCFM2

Manual IntegrationsAPPROVED

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