

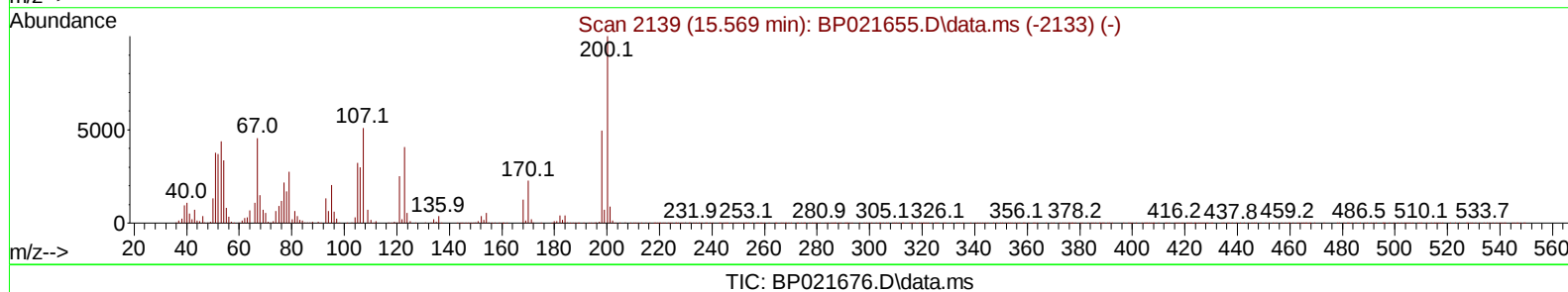
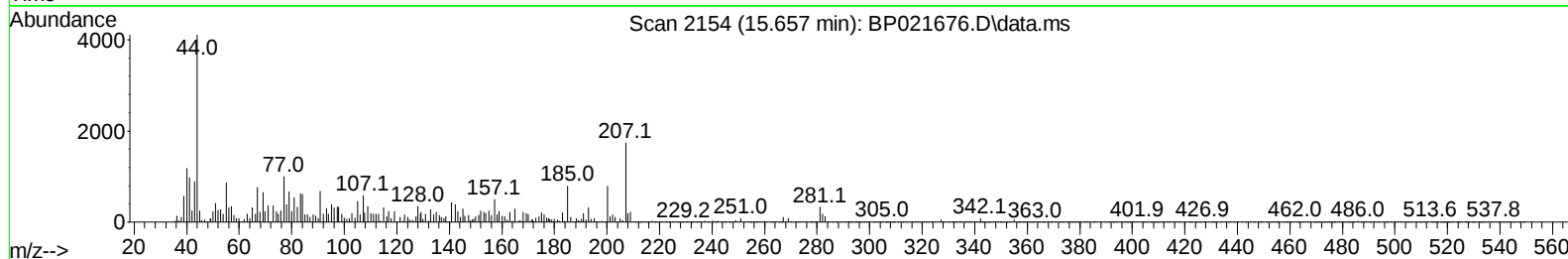
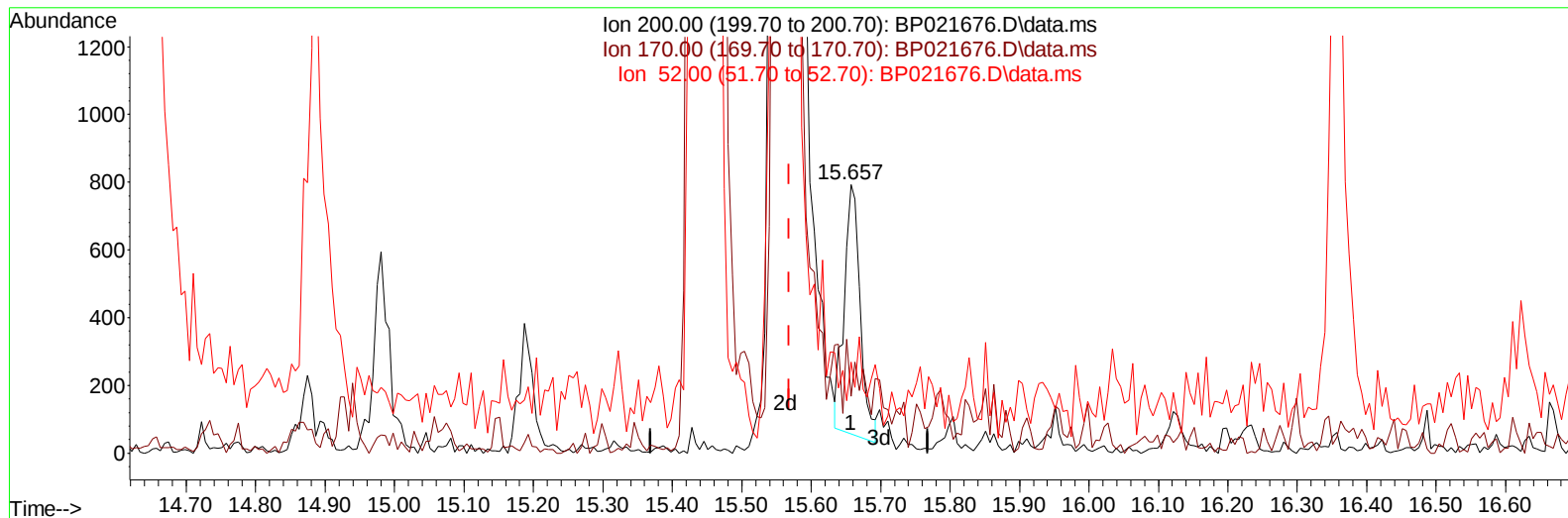
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
Data File : BP021676.D
Acq On : 27 Aug 2024 18:15
Operator : RC/JU
Sample : P3675-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCXY0

Manual IntegrationsAPPROVED

Quant Time: Aug 27 22:48:19 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Aug 27 05:11:38 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 08/28/2024
Supervised By :mohammad ahmed 08/29/2024



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

15.657min (+ 0.088) 0.11 ng/u1

response 1177

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	23.80	23.68
52.00	41.90	33.88
0.00	0.00	0.00

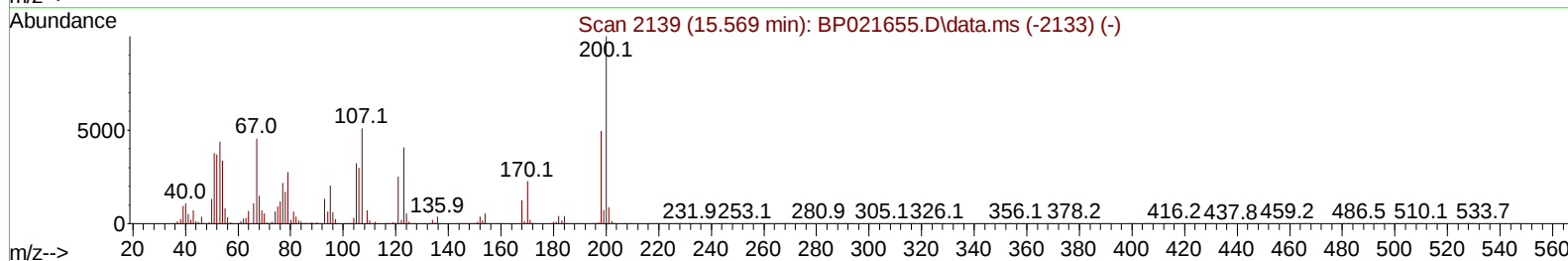
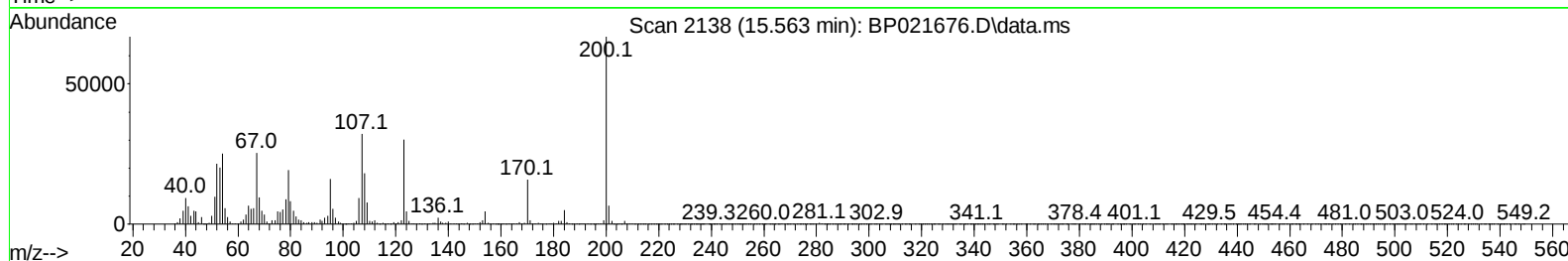
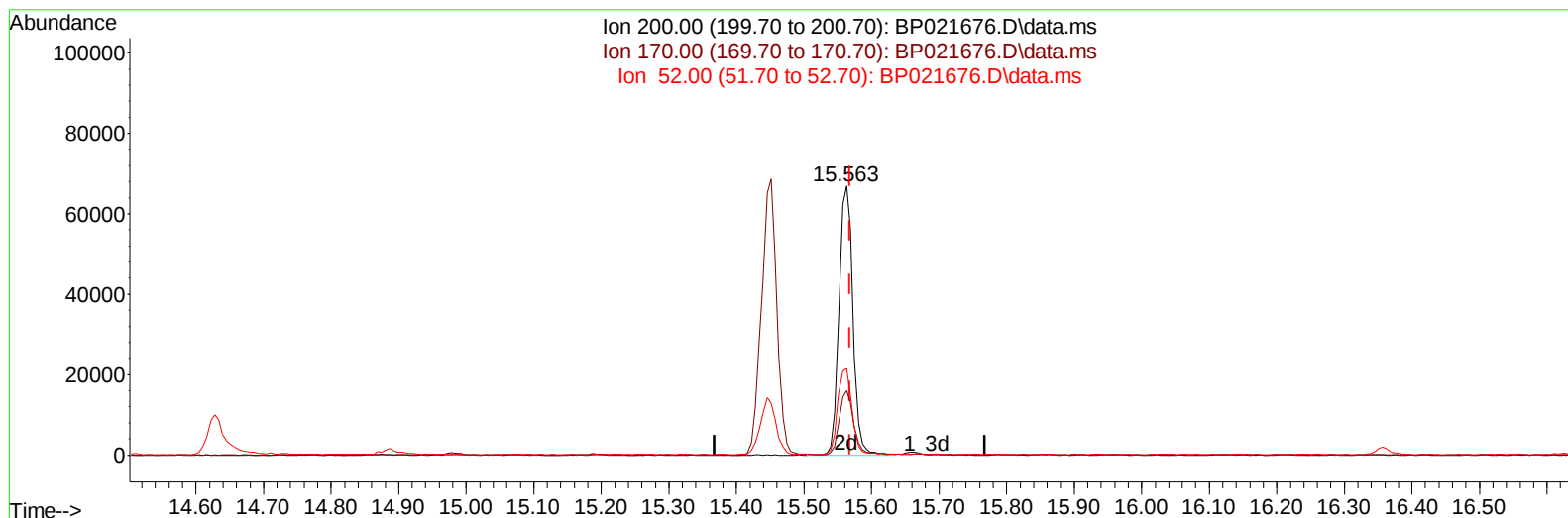
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TIC: BP021676.D\data.ms

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

15.563min (-0.006) 9.16 ng/ul m

response 95609

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	23.80	23.91
52.00	41.90	32.22#
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.793	152	323630	20.000	ng/ul	0.00
20) Naphthalene-d8	10.581	136	1334117	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.440	164	863365	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.239	188	1897910	20.000	ng/ul	-0.01
79) Chrysene-d12	21.692	240	1861901	20.000	ng/ul	0.00
88) Perylene-d12	25.115	264	2134082	20.000	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	23867	2.448	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	6.958	99	515849	14.936	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.122	67	291903	15.530	ng/ul	0.00
11) 2-Chlorophenol-d4	7.328	132	344142	15.434	ng/ul	0.00
15) 4-Methylphenol-d8	8.493	113	397992	14.912	ng/ul	0.00
21) Nitrobenzene-d5	8.934	128	161137	15.013	ng/ul	0.00
24) 2-Nitrophenol-d4	9.657	143	159760	14.663	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.199	165	316200	14.727	ng/ul	0.00
31) 4-Chloroaniline-d4	10.710	131	23244	0.744	ng/ul	0.00
46) Dimethylphthalate-d6	13.840	166	1009539	15.282	ng/ul	0.00
49) Acenaphthylene-d8	14.128	160	1176139	15.857	ng/ul	0.00
54) 4-Nitrophenol-d4	14.628	143	140192	11.154	ng/ul	0.00
60) Fluorene-d10	15.451	176	915353	16.481	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.563	200	95609m	9.159	ng/ul	0.00
73) Anthracene-d10	17.345	188	1376717	15.285	ng/ul	0.00
81) Pyrene-d10	19.710	212	1765212	16.585	ng/ul	-0.01
92) Benzo(a)pyrene-d12	24.874	264	1831411	16.094	ng/ul	0.00

Target Compounds Qvalue

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

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