

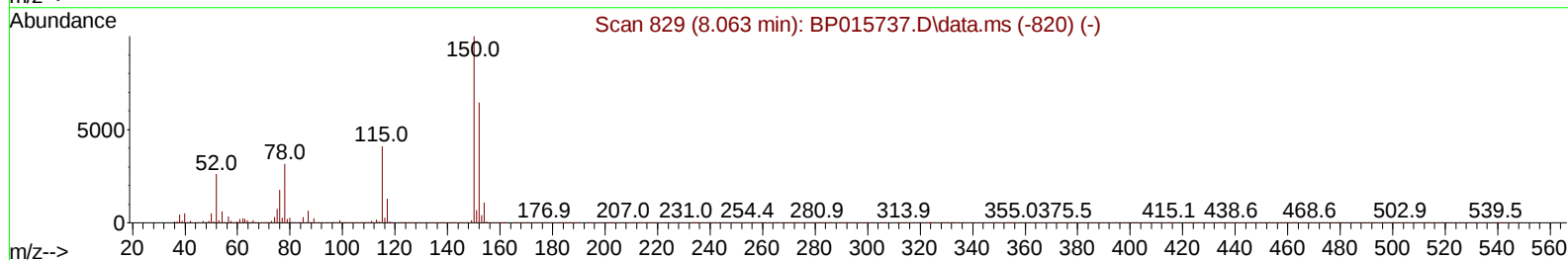
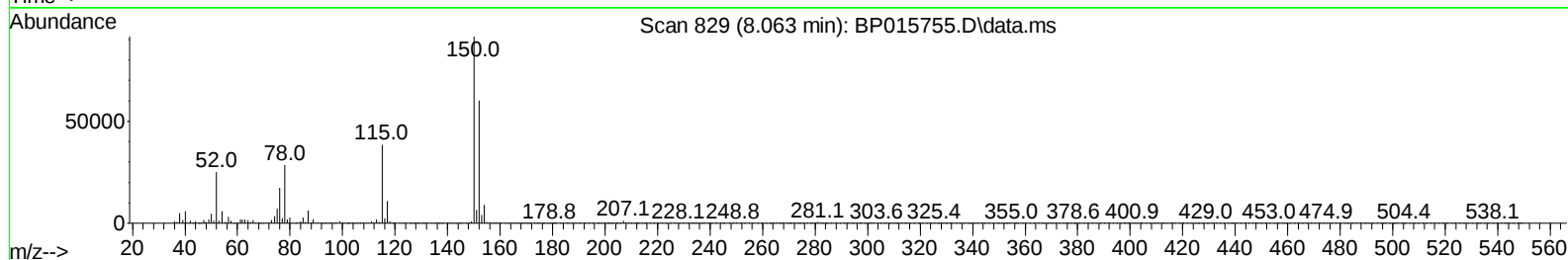
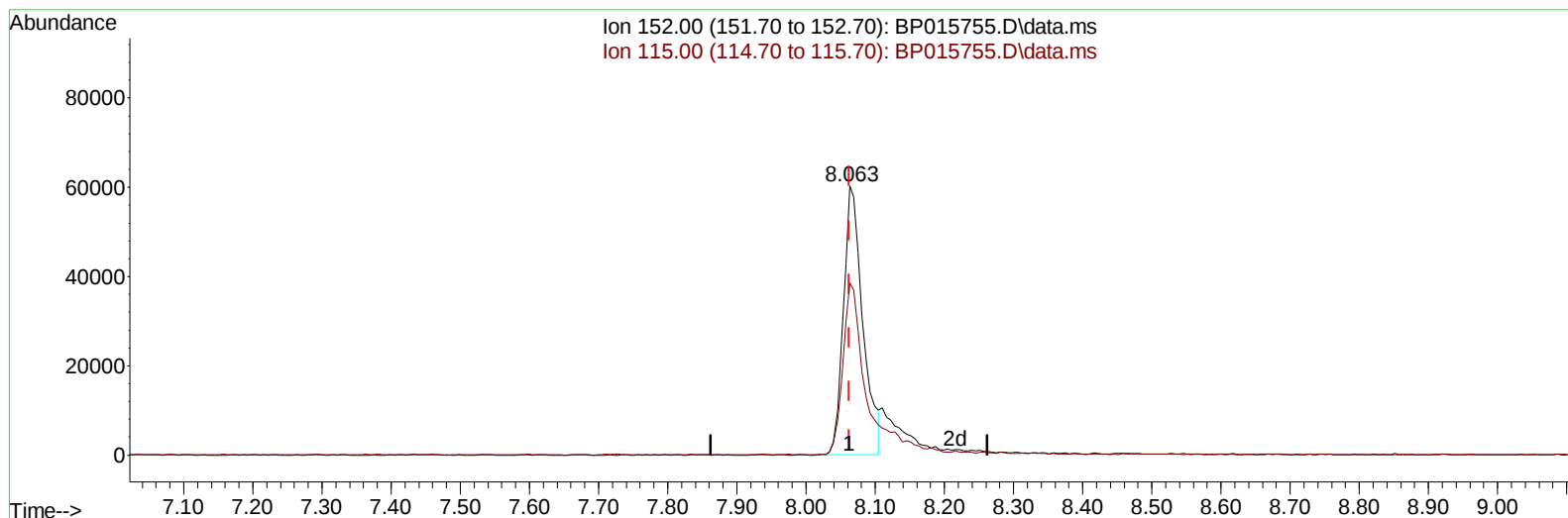
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062723\
 Data File : BP015755.D
 Acq On : 27 Jun 2023 19:44
 Operator : MA/JU
 Sample : 03411-10
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SVOC-GPC2-BLANK

Manual IntegrationsAPPROVED

Quant Time: Jun 28 03:19:53 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jun 27 00:56:58 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 06/28/2023
 Supervised By :mohammad ahmed 06/28/2023



TIC: BP015755.D\data.ms

(1) 1,4-Dichlorobenzene-d4 (I)

8.063min (+ 0.000) 20.00 ng/u1

response 117279

Ion	Exp%	Act%
152.00	100.00	100.00
115.00	65.50	63.91
0.00	0.00	0.00
0.00	0.00	0.00

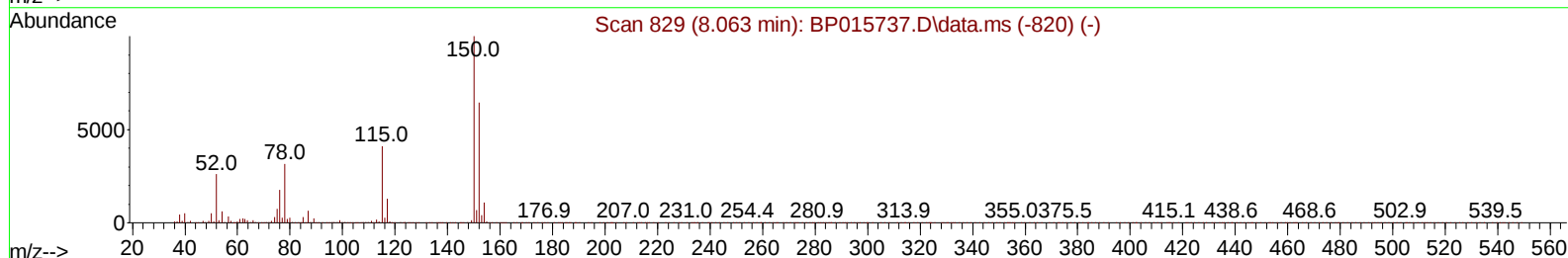
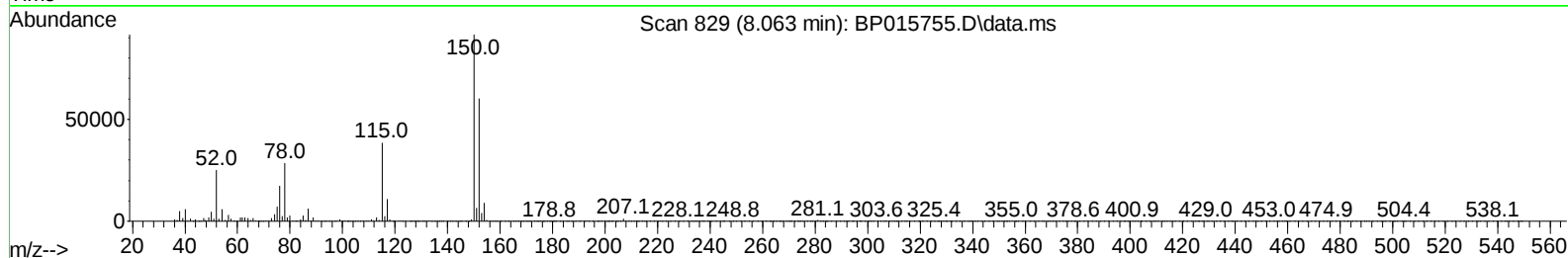
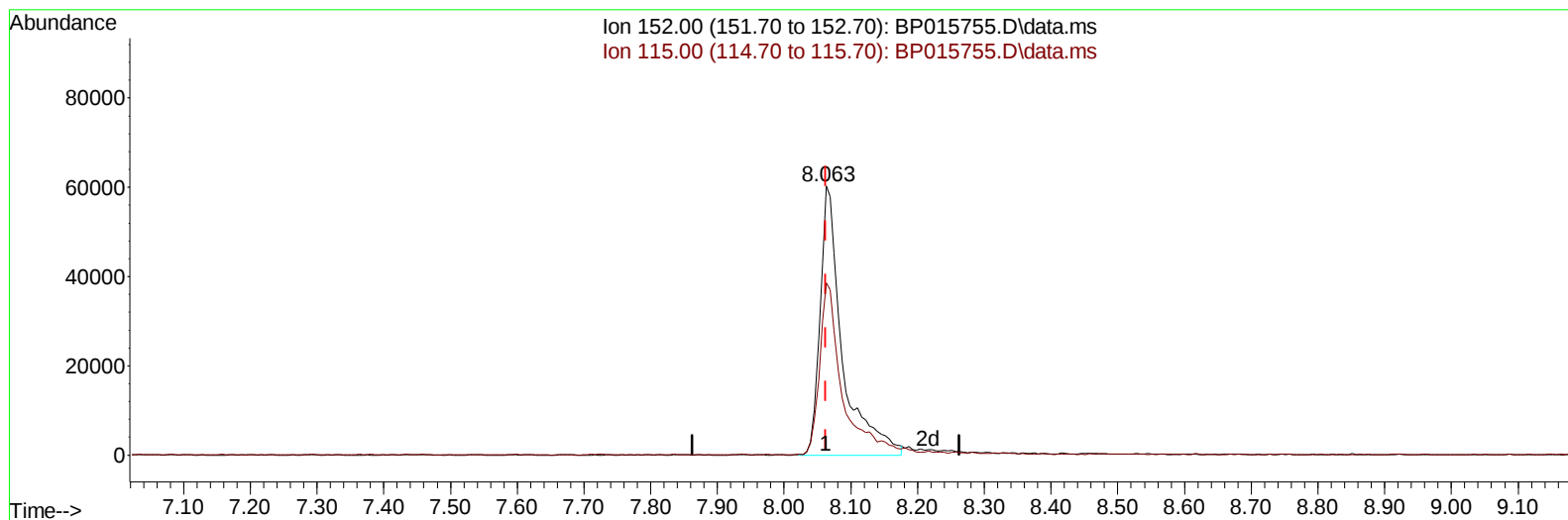
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(1) 1,4-Dichlorobenzene-d4 (I)

8.063min (+ 0.000) 20.00 ng/ul m

response 140361

Ion	Exp%	Act%
152.00	100.00	100.00
115.00	65.50	63.91
0.00	0.00	0.00
0.00	0.00	0.00

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Quant Time: Jun 28 03:38:57 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.063	152	140361m	20.000	ng/ul	0.00
20) Naphthalene-d8	10.898	136	581115	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.716	164	350416	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.486	188	675925	20.000	ng/ul	0.01
79) Chrysene-d12	21.568	240	541905	20.000	ng/ul	0.00
88) Perylene-d12	24.115	264	693028	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	0.000	96	0d	0.000	ng/uL	
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	0.000	99	0d	0.000	ng/ul	
9) Bis-(2-Chloroethyl)eth...	0.000	67	0d	0.000	ng/ul	
11) 2-Chlorophenol-d4	0.000	132	0d	0.000	ng/ul	
15) 4-Methylphenol-d8	0.000	113	0d	0.000	ng/ul	
21) Nitrobenzene-d5	0.000	128	0d	0.000	ng/ul	
24) 2-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
28) 2,4-Dichlorophenol-d3	0.000	165	0d	0.000	ng/ul	
31) 4-Chloroaniline-d4	0.000	131	0d	0.000	ng/ul	
46) Dimethylphthalate-d6	0.000	166	0d	0.000	ng/ul	
49) Acenaphthylene-d8	0.000	160	0d	0.000	ng/ul	
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
60) Fluorene-d10	0.000	176	0d	0.000	ng/ul	
65) 4,6-Dinitro-2-methylph...	0.000	200	0d	0.000	ng/ul	
73) Anthracene-d10	0.000	188	0d	0.000	ng/ul	
81) Pyrene-d10	0.000	212	0d	0.000	ng/ul	
92) Benzo(a)pyrene-d12	0.000	264	0d	0.000	ng/ul	

Target Compounds Qvalue

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

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