

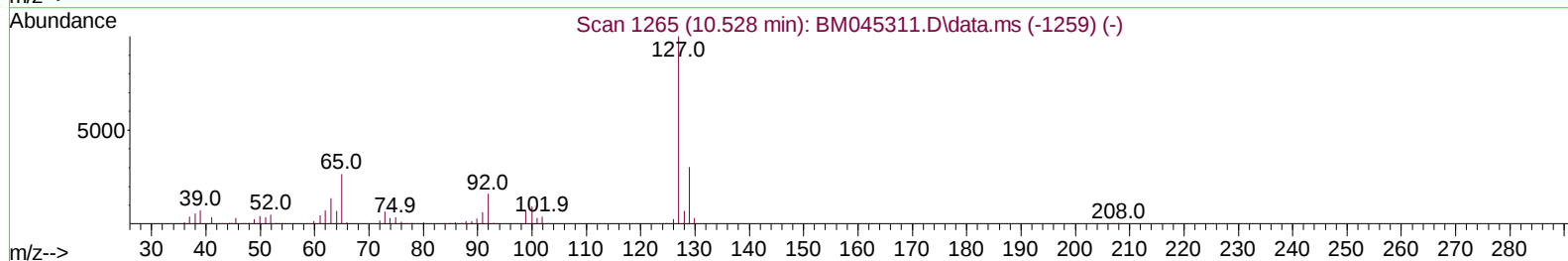
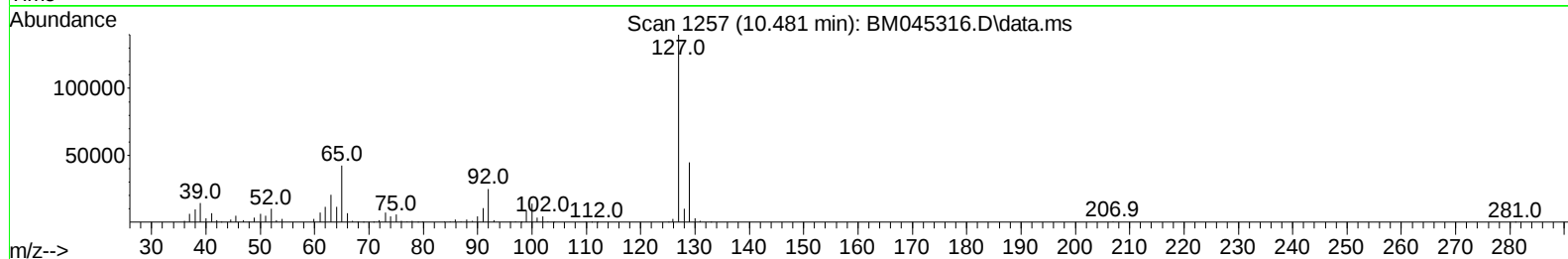
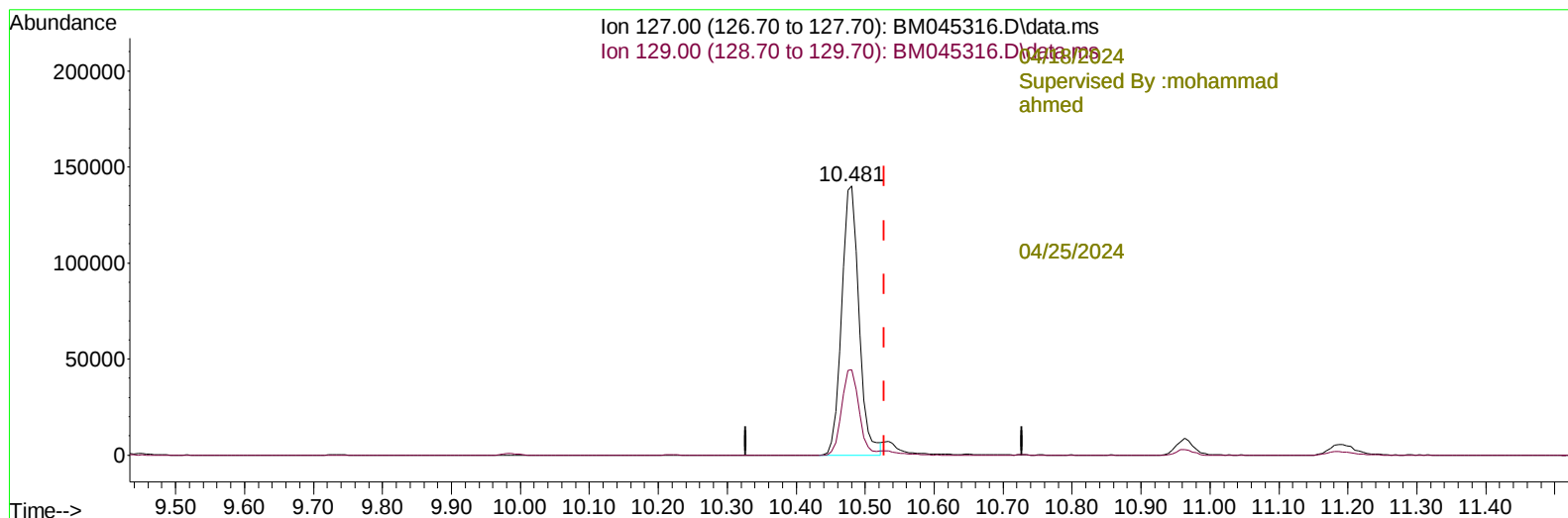
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041724\
Data File : BM045316.D
Acq On : 17 Apr 2024 14:36
Operator : MA/JU
Sample : P2051-12DL 5X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0RH9DL

Manual IntegrationsAPPROVED

Quant Time: Apr 17 17:04:30 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040524.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Apr 17 12:42:01 2024
Response via : Initial Calibration

Reviewed By :Yogesh
Patel



TIC: BM045316.D\data.ms

(32) 4-Chloroaniline

10.481min (-0.047) 59.95 ng/ul

response 243456

Ion	Exp%	Act%
127.00	100.00	100.00
129.00	32.70	31.79
0.00	0.00	0.00
0.00	0.00	0.00

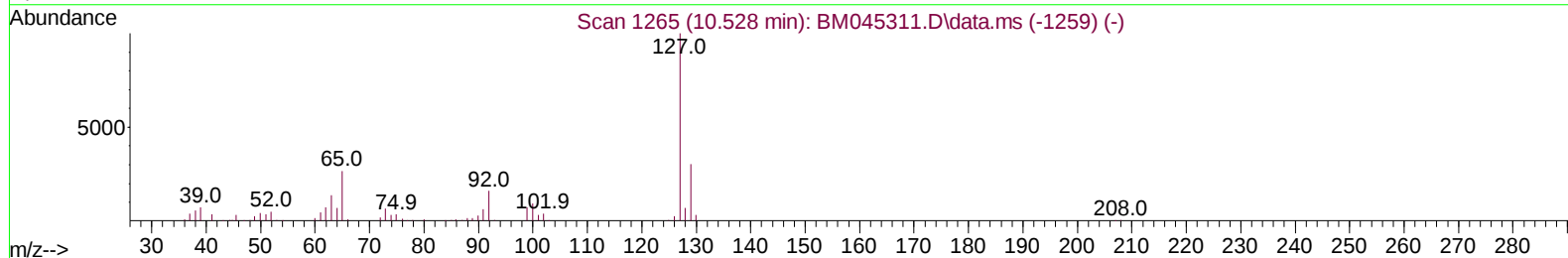
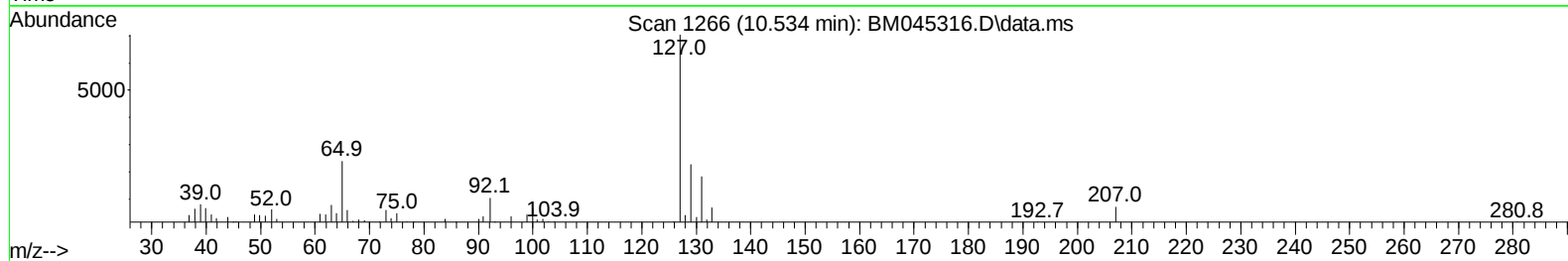
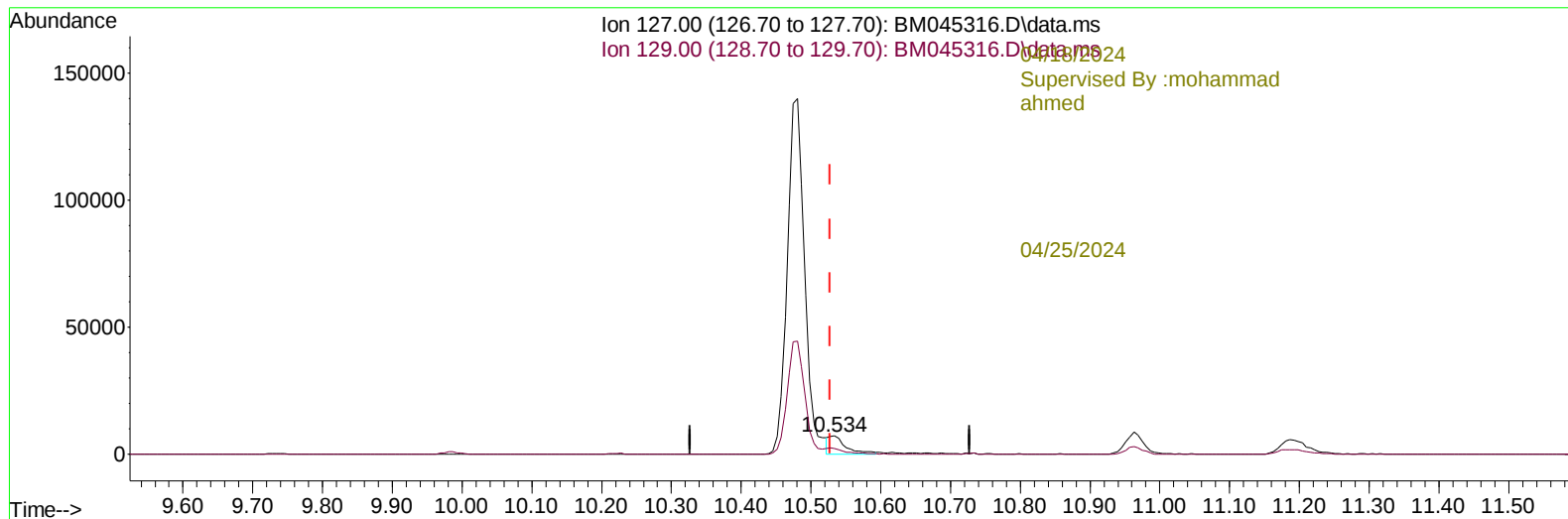
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 Patel



TIC: BM045316.D\data.ms

(32) 4-Chloroaniline

10.534min (+ 0.006) 2.92 ng/ul m

response 11842

Ion	Exp%	Act%
127.00	100.00	100.00
129.00	32.70	32.33
0.00	0.00	0.00
0.00	0.00	0.00

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 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :

BNA_M

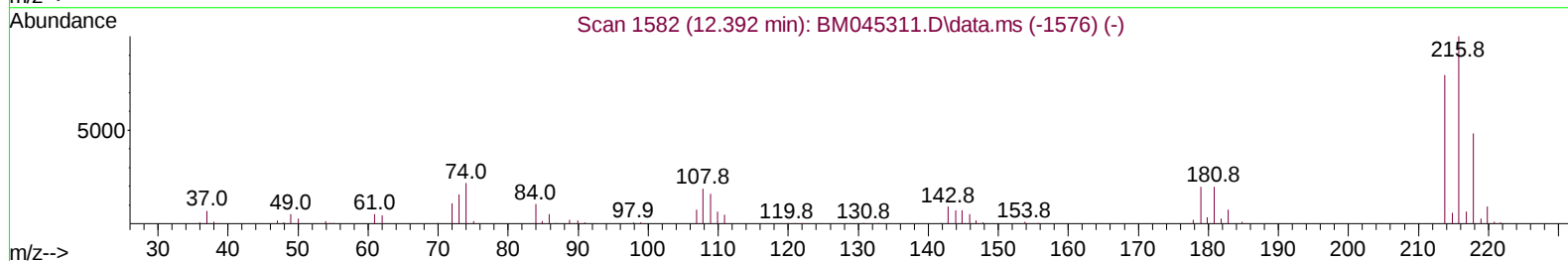
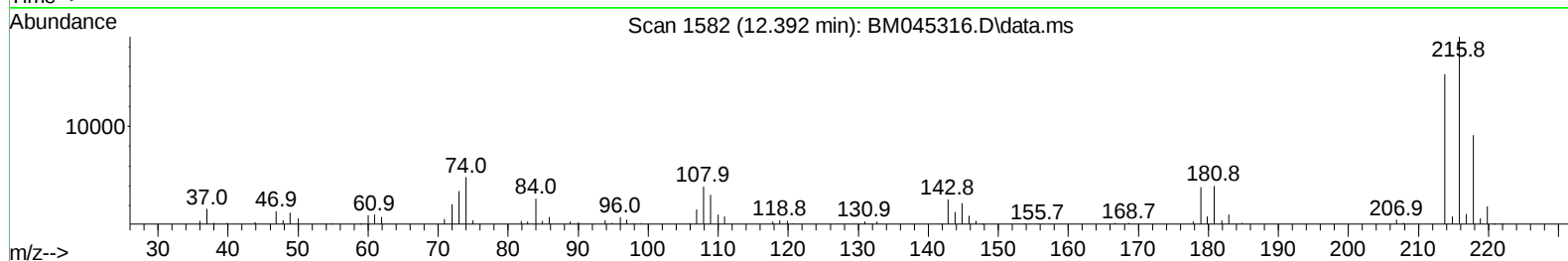
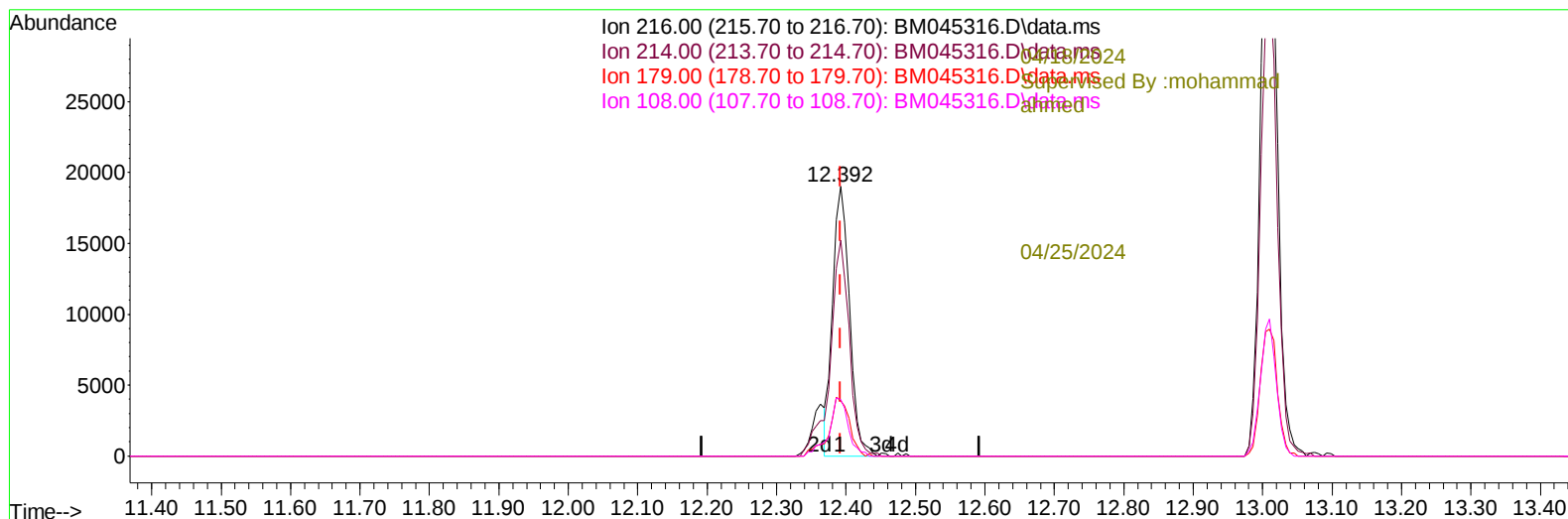
ClientSampleId :

C0RH9DL

Manual IntegrationsAPPROVED

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 Patel

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TIC: BM045316.D\data.ms

(39) 1,2,4,5-Tetrachlorobenzene

12.392min (+ 0.000) 10.95 ng/u1

response 32042

Ion	Exp%	Act%
216.00	100.00	100.00
214.00	79.40	80.03
179.00	20.80	20.42
108.00	18.70	20.78

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041724\
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 Acq On : 17 Apr 2024 14:36
 Operator : MA/JU
 Sample : P2051-12DL 5X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :

BNA_M

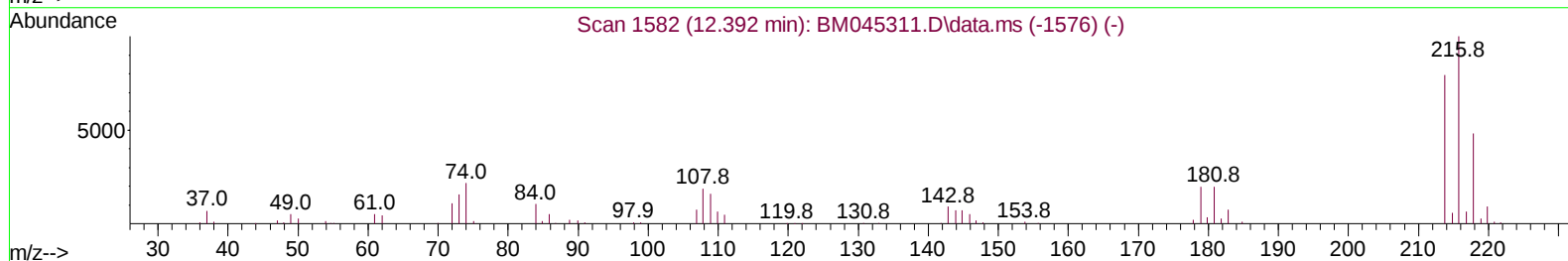
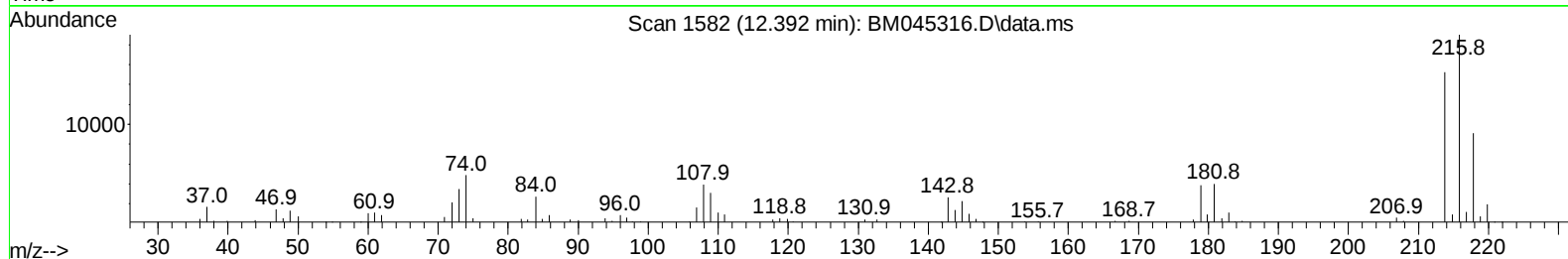
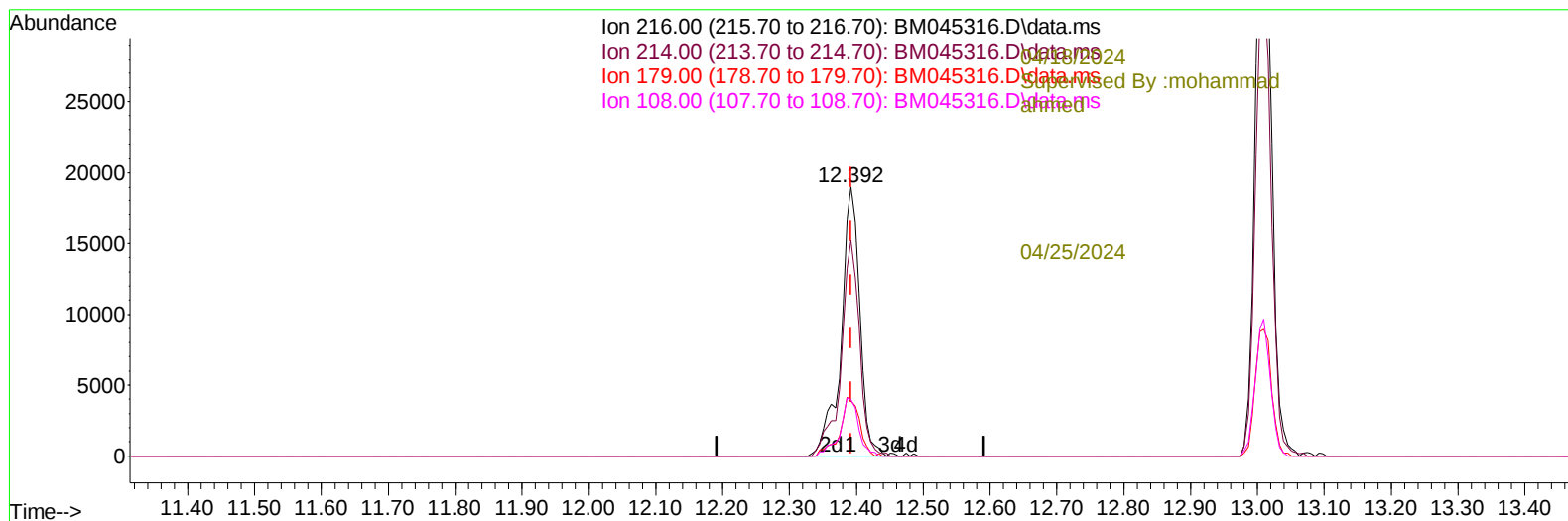
ClientSampleId :

C0RH9DL

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 Patel



TIC: BM045316.D\data.ms

(39) 1,2,4,5-Tetrachlorobenzene

12.392min (+ 0.000) 12.65 ng/ul m

response 37013

Ion	Exp%	Act%
216.00	100.00	100.00
214.00	79.40	80.03
179.00	20.80	20.42
108.00	18.70	20.78

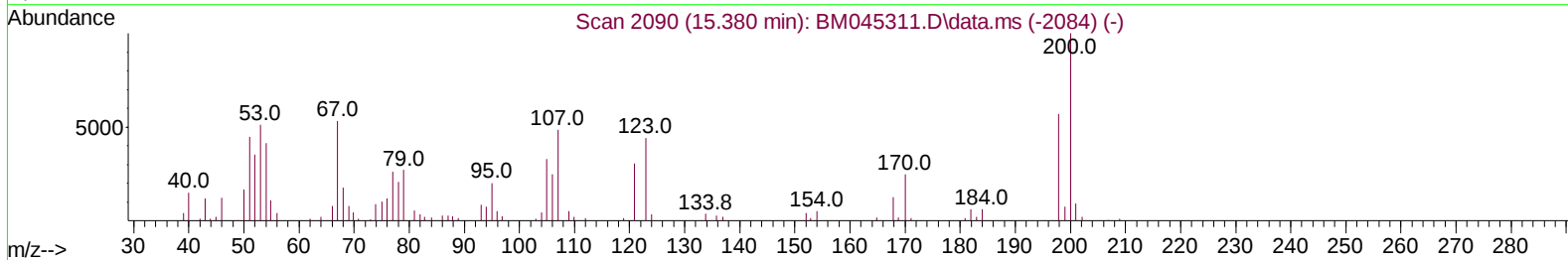
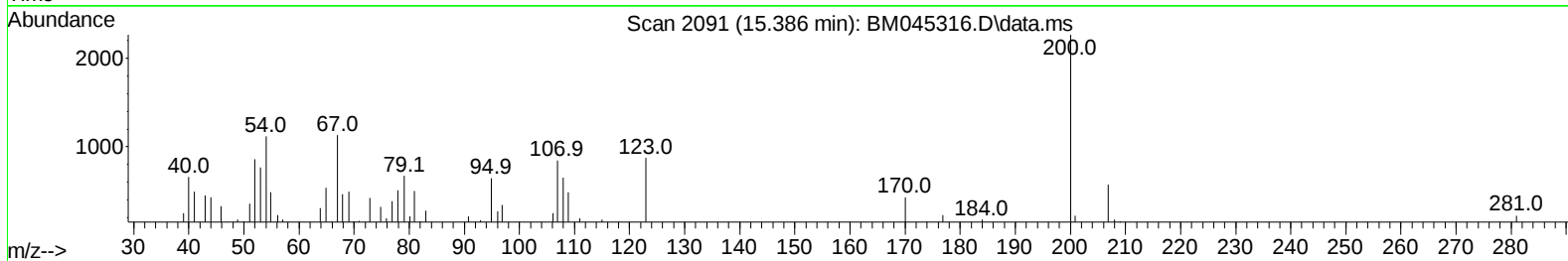
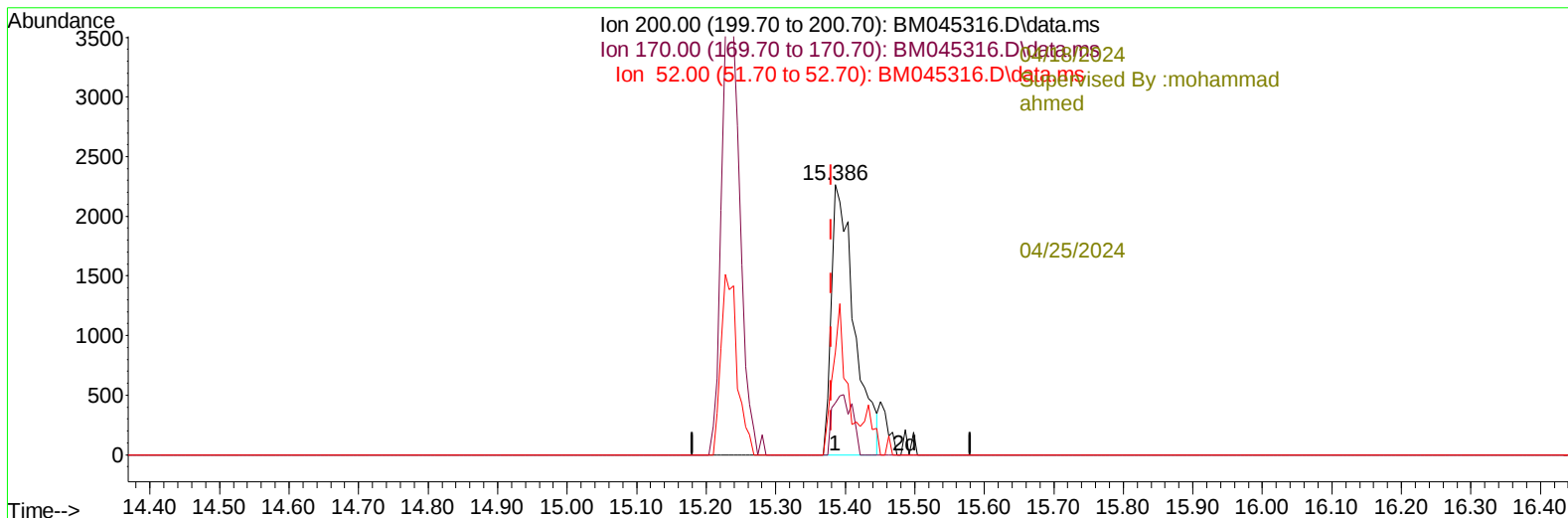
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Patel



TIC: BM045316.D\data.ms

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

15.386min (+ 0.006) 4.00 ng/u1

response 5157

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	18.30	19.18
52.00	40.20	38.00
0.00	0.00	0.00

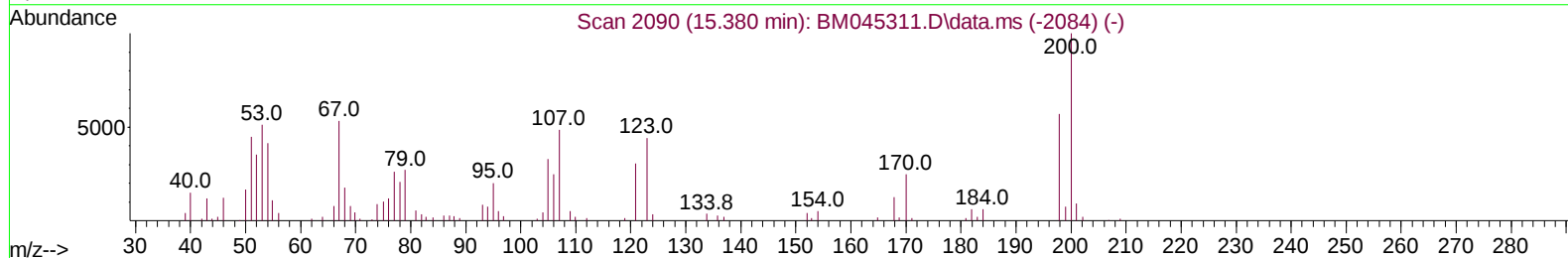
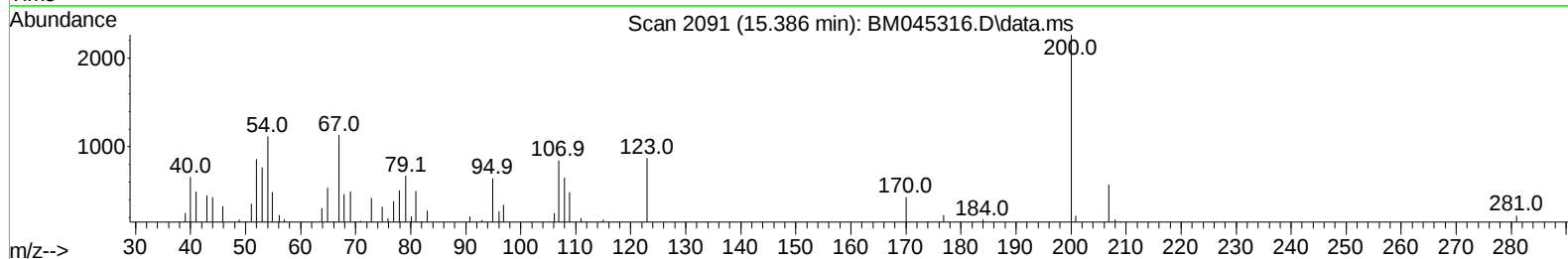
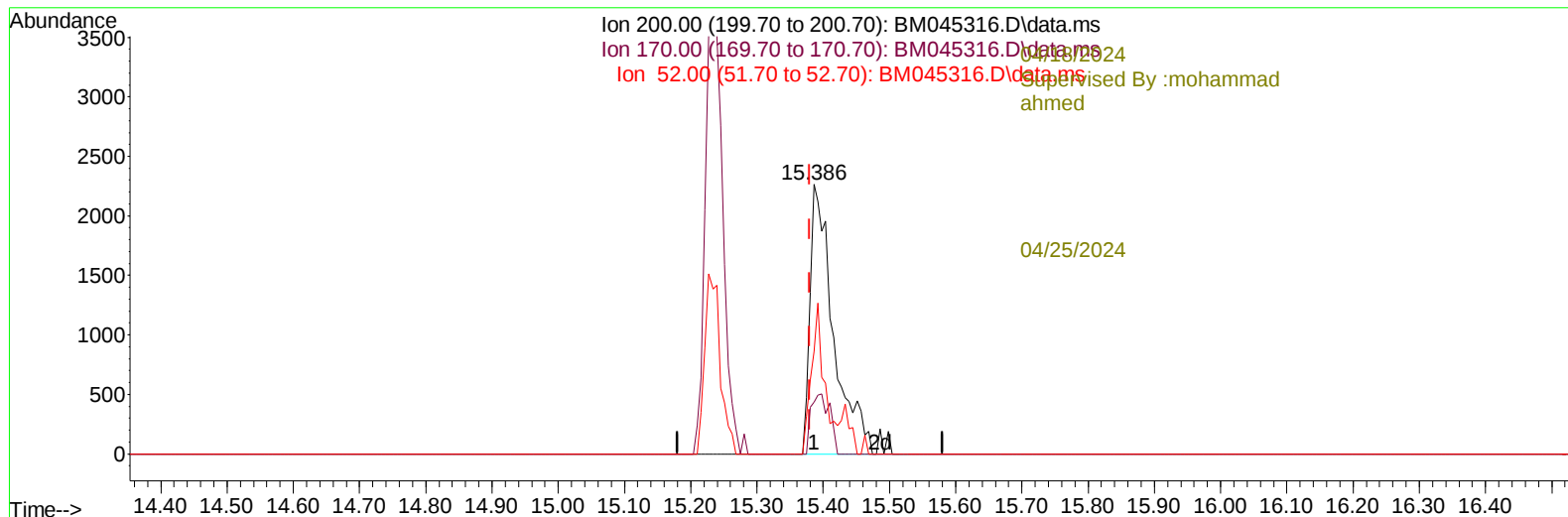
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041724\
Data File : BM045316.D
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Operator : MA/JU
Sample : P2051-12DL 5X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0RH9DL

Manual IntegrationsAPPROVED

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Reviewed By :Yogesh
Patel



TIC: BM045316.D\data.ms

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

15.386min (+ 0.006) 4.42 ng/ul m

response 5704

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	18.30	19.18
52.00	40.20	38.00
0.00	0.00	0.00

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 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
 CORH9DL

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 Patel

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
-----04/18/2024						
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.581	152	55718	20.000	ng/ul	0.00
20) Naphthalene-d8	10.351	136	176346	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.227	164	107006	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.986	188	217670	20.000	ng/ul	0.00
79) Chrysene-d12	21.203	240	228874	20.000	ng/ul	0.00
88) Perylene-d12	23.397	264	298867	20.000	ng/ul	0.00
-----04/25/2024						
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.193	96	1350	0.905	ng/uL	0.02
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	6.763	99	4421	0.936	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.940	67	28214	10.026	ng/ul	0.00
11) 2-Chlorophenol-d4	7.116	132	17067	4.612	ng/ul	0.00
15) 4-Methylphenol-d8	8.298	113	8376	2.265	ng/ul	0.01
21) Nitrobenzene-d5	8.740	128	11992	8.853	ng/ul	0.00
24) 2-Nitrophenol-d4	9.451	143	11276	7.887	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.987	165	18471	7.046	ng/ul	0.01
31) 4-Chloroaniline-d4	10.504	131	27866	6.677	ng/ul	0.00
46) Dimethylphthalate-d6	13.663	166	67062	9.001	ng/ul	0.00
49) Acenaphthylene-d8	13.916	160	75511	8.579	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
60) Fluorene-d10	15.233	176	60812	9.155	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.386	200	5704m	4.420	ng/ul	0.00
73) Anthracene-d10	17.092	188	91252	9.319	ng/ul	0.00
81) Pyrene-d10	19.398	212	111182	9.866	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.256	264	133959	9.188	ng/ul	0.00

Target Compounds						Qvalue
8) Phenol	6.793	94	7303	1.493	ng/ul	93
22) Nitrobenzene	8.793	77	1768437	532.126	ng/ul	94
32) 4-Chloroaniline	10.534	127	11842m	2.916	ng/ul	
39) 1,2,4,5-Tetrachloroben...	12.392	216	37013m	12.648	ng/ul	
75) 1,2,3,4-Tetrachloroben...	13.010	216	71991	26.241	ng/uL	98

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

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