

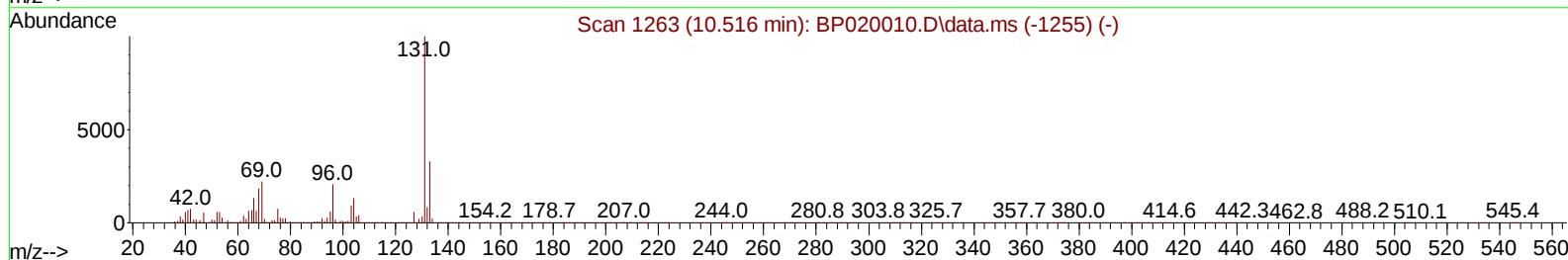
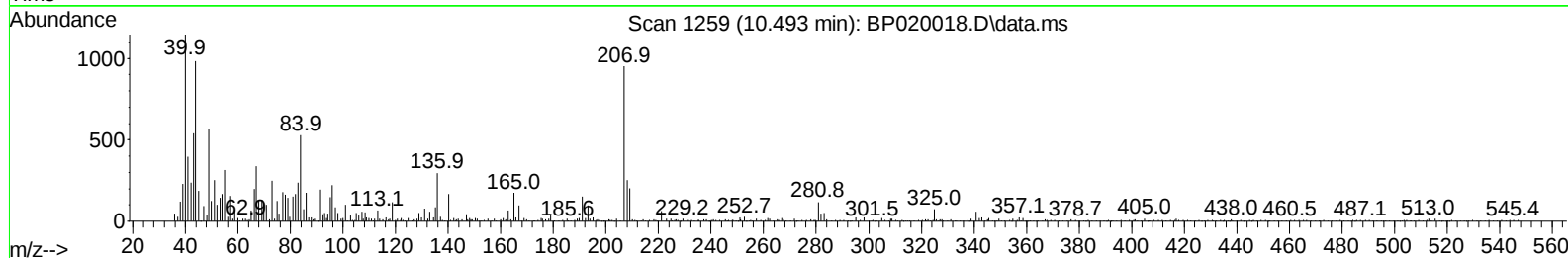
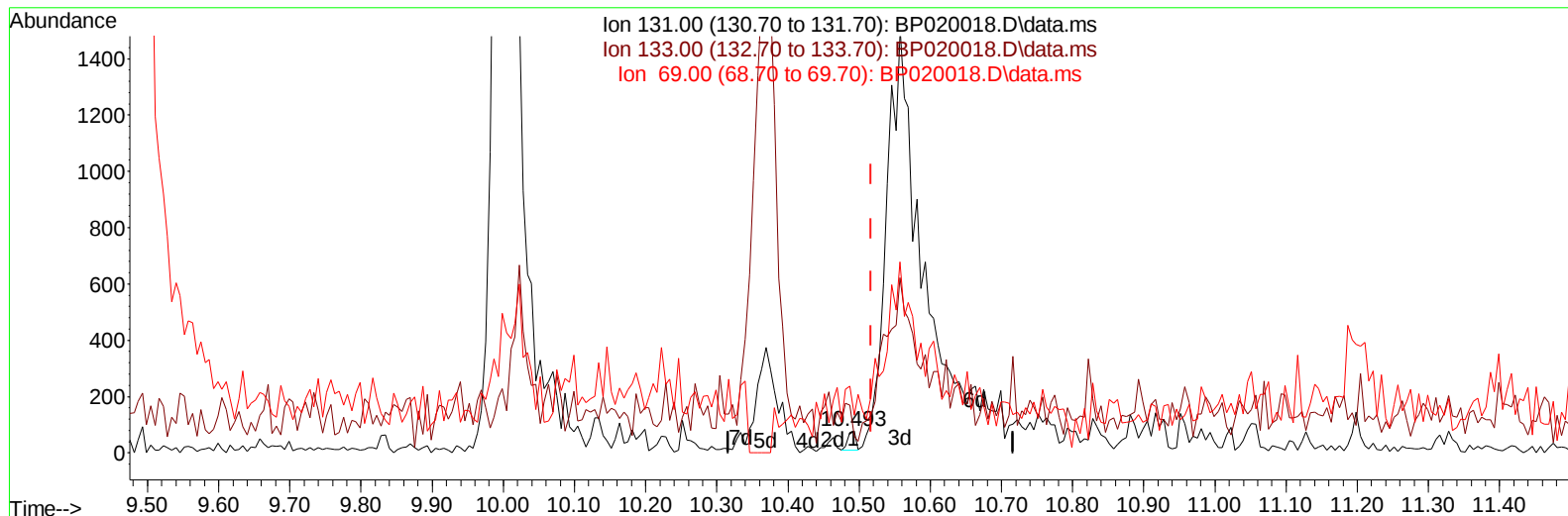
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032824\
Data File : BP020018.D
Acq On : 28 Mar 2024 20:13
Operator : MA/JU
Sample : P1856-01 10X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E0LZ0

Manual IntegrationsAPPROVED

Quant Time: Mar 28 23:10:05 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032224.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Mar 28 14:08:48 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 03/29/2024
Supervised By :Jagrut Upadhyay 03/29/2024



TIC: BP020018.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.493min (-0.024) 0.01 ng/u1

response 53

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.80	76.92#
69.00	22.40	171.79#
0.00	0.00	0.00

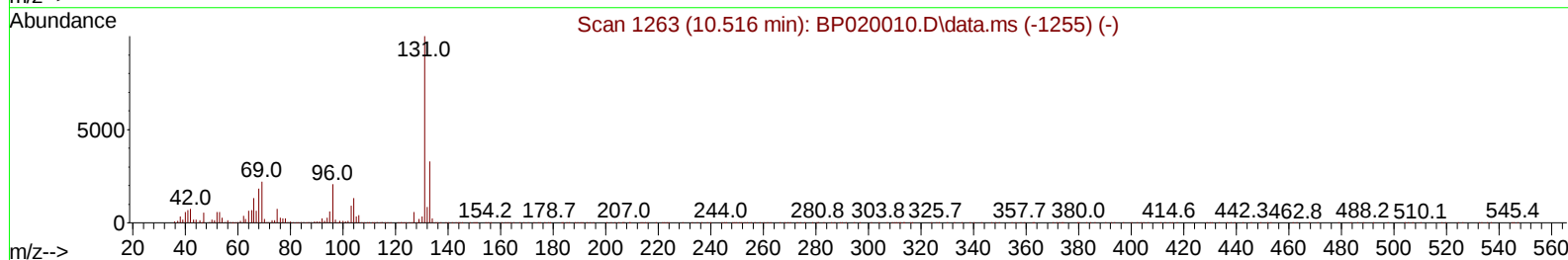
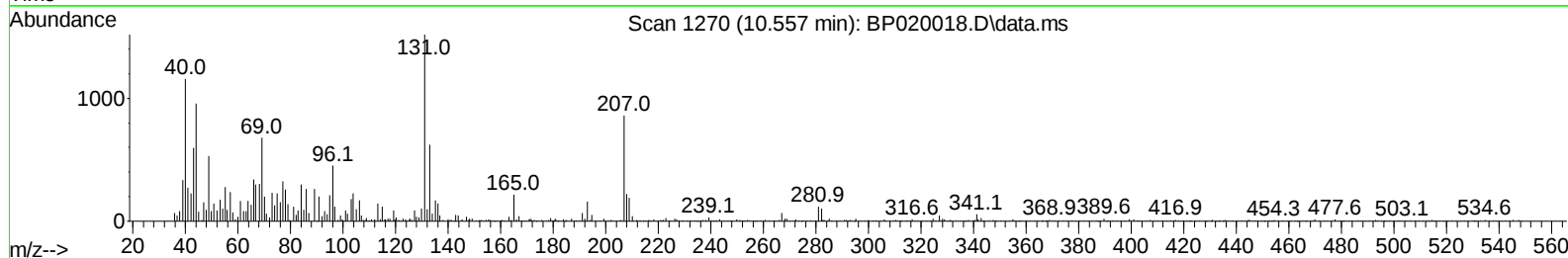
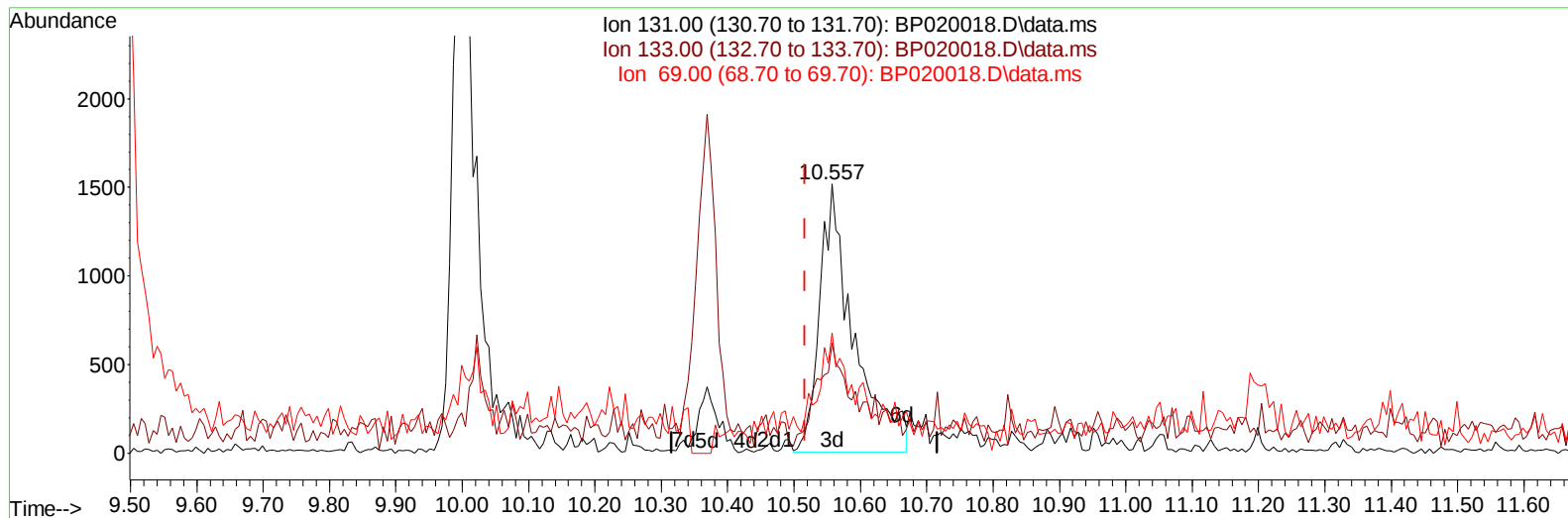
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TIC: BP020018.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.557min (+ 0.041) 0.58 ng/ul m

response 5387

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.80	40.97#
69.00	22.40	44.73#
0.00	0.00	0.00

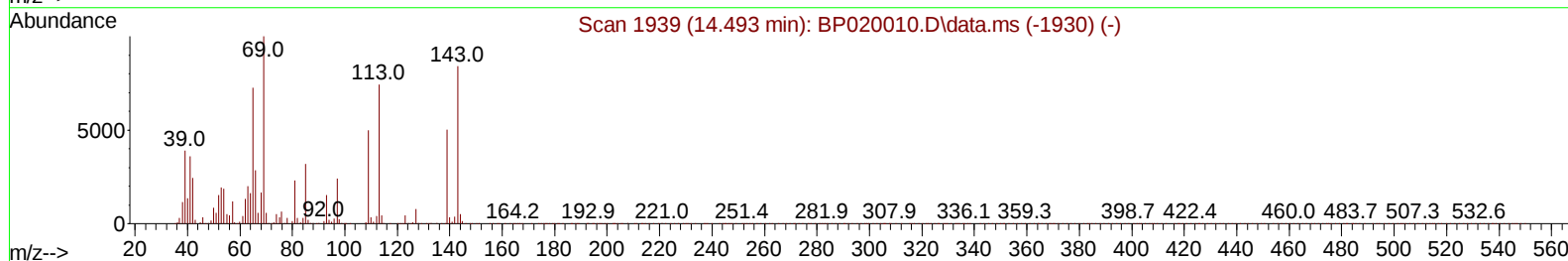
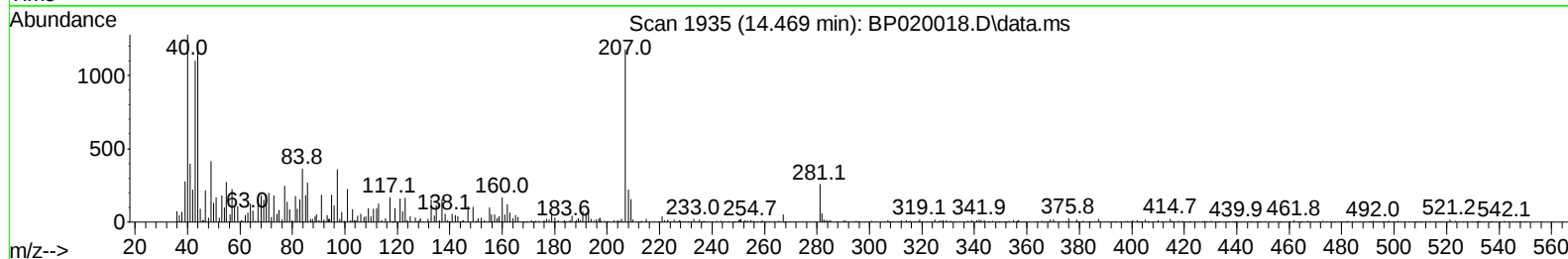
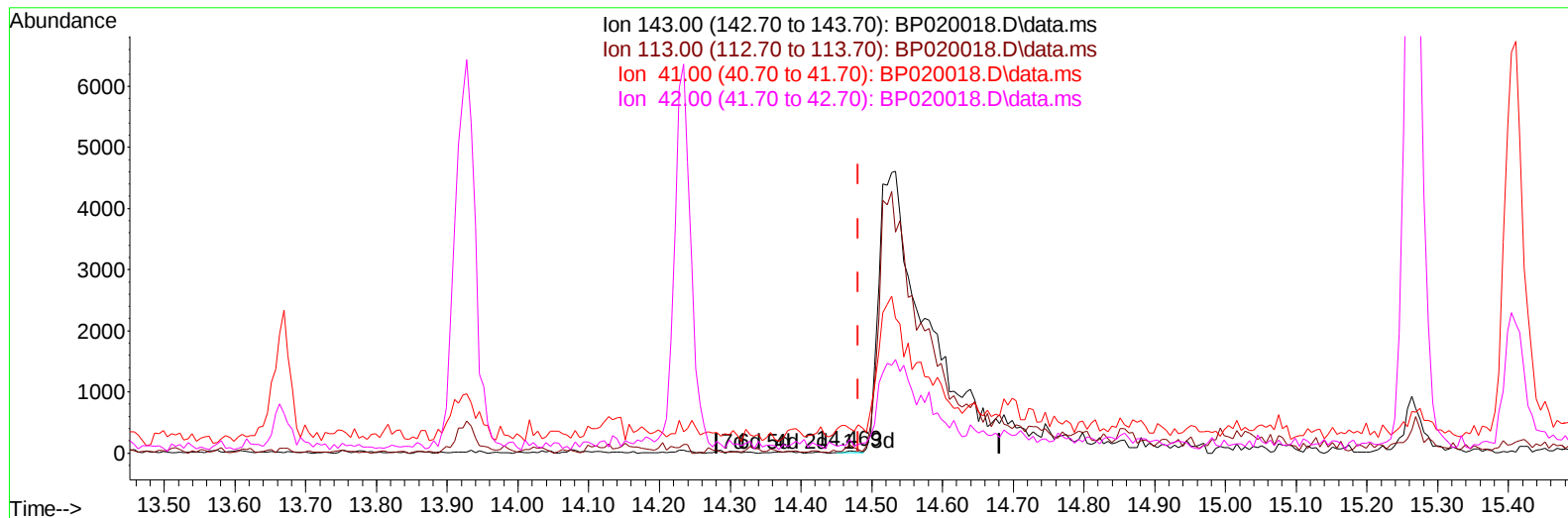
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TIC: BP020018.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.469min (-0.012) 0.01 ng/u1

response 37

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.50	334.21#
41.00	40.90	1050.00#
42.00	28.50	576.32#

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 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
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Manual IntegrationsAPPROVED

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 Supervised By :Jagrut Upadhyay 03/29/2024

Quant Time: Mar 28 23:20:19 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.605	152	110659	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.363	136	453091	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.234	164	305327	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.075	188	650352	20.000	ng/ul	0.00
79) Chrysene-d12	21.533	240	586441	20.000	ng/ul	-0.01
88) Perylene-d12	24.816	264	635290	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.240	96	7839	3.104	ng/uL	0.00
4) Pyridine-d5	3.664	84	17995	2.360	ng/ul	0.02
7) Phenol-d5	6.799	99	67824	7.267	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.969	67	234582	38.580	ng/ul	0.00
11) 2-Chlorophenol-d4	7.152	132	182700	27.259	ng/ul	0.00
15) 4-Methylphenol-d8	8.305	113	127266	17.625	ng/ul	0.00
21) Nitrobenzene-d5	8.758	128	128707	37.061	ng/ul	0.00
24) 2-Nitrophenol-d4	9.463	143	135361	34.836	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.999	165	218000	30.720	ng/ul	0.00
31) 4-Chloroaniline-d4	10.557	131	5387m	0.582	ng/ul	0.04
46) Dimethylphthalate-d6	13.669	166	888451	41.502	ng/ul	0.02
49) Acenaphthylene-d8	13.928	160	942457	37.688	ng/ul	0.00
54) 4-Nitrophenol-d4	14.534	143	22628m	6.724	ng/ul	0.05
60) Fluorene-d10	15.263	176	746890	40.957	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.404	200	136255	35.001	ng/ul	-0.02
73) Anthracene-d10	17.181	188	1203119	41.792	ng/ul	0.01
81) Pyrene-d10	19.569	212	1490453	43.318	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.586	264	1387356	45.456	ng/ul	-0.02

Target Compounds Qvalue

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

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