

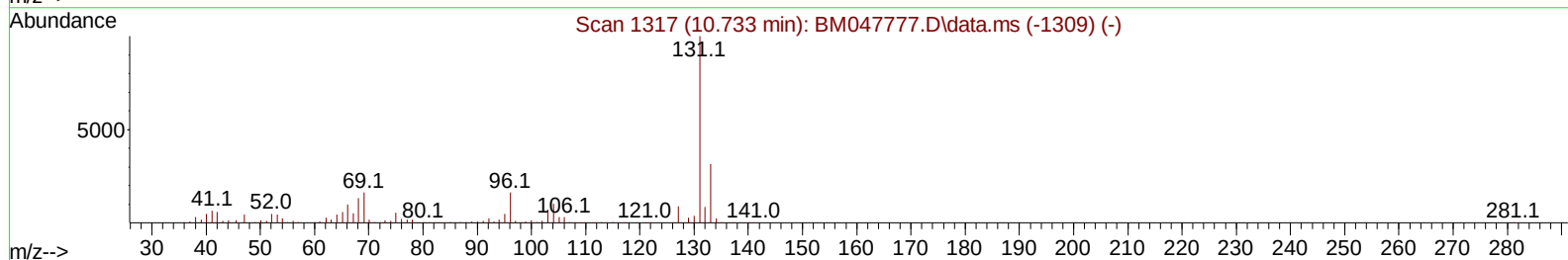
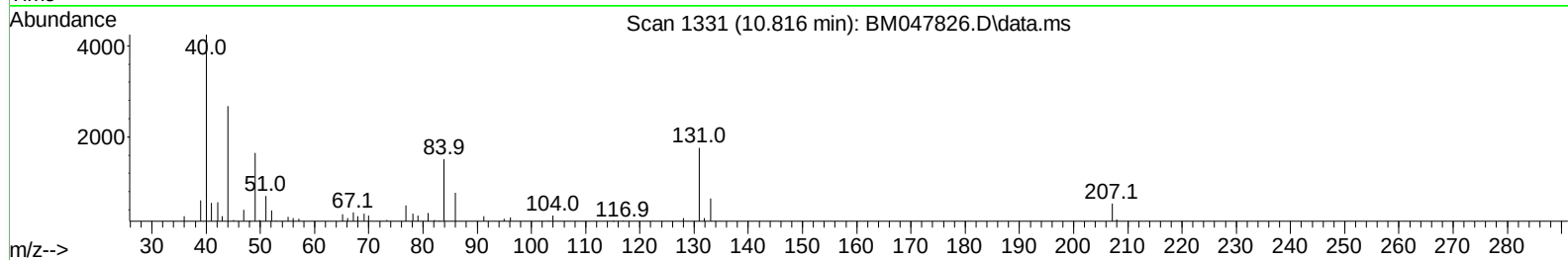
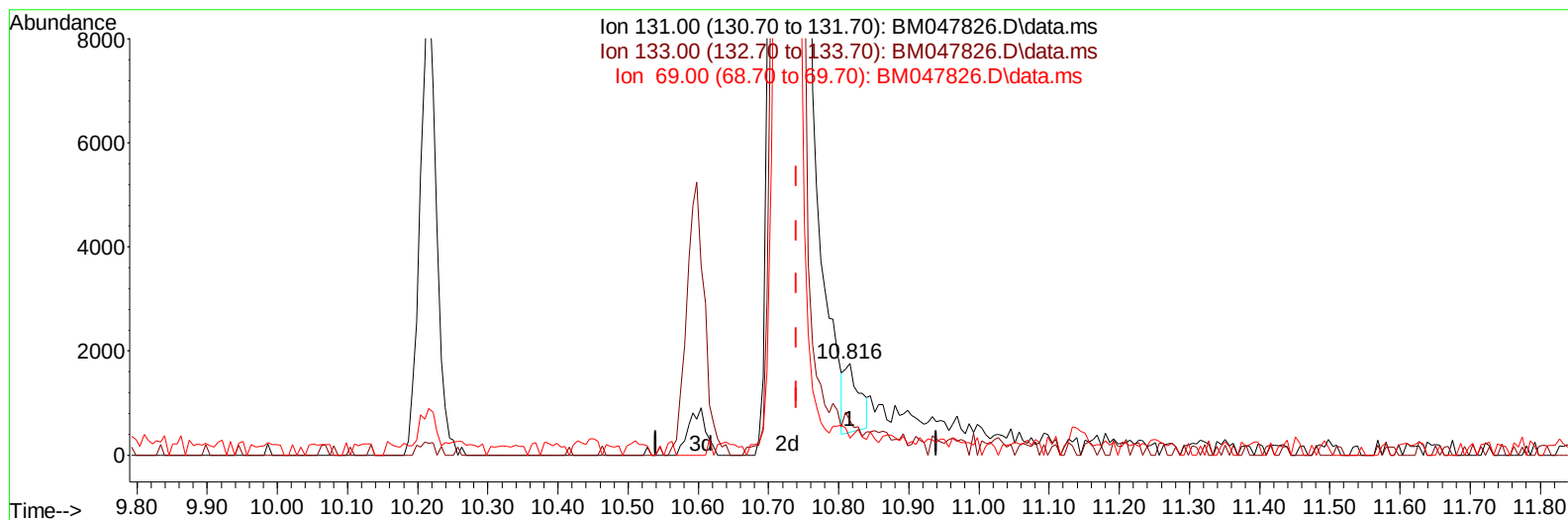
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100324\
Data File : BM047826.D
Acq On : 04 Oct 2024 03:11
Operator : RC/JU
Sample : P4084-17
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCYM1

Manual IntegrationsAPPROVED

Quant Time: Oct 04 03:59:50 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Sep 28 02:30:36 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 10/04/2024
Supervised By :mohammad ahmed 10/06/2024



TIC: BM047826.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.816min (+ 0.076) 0.07 ng/ul

response 1933

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	34.00	37.27
69.00	20.90	18.04
0.00	0.00	0.00

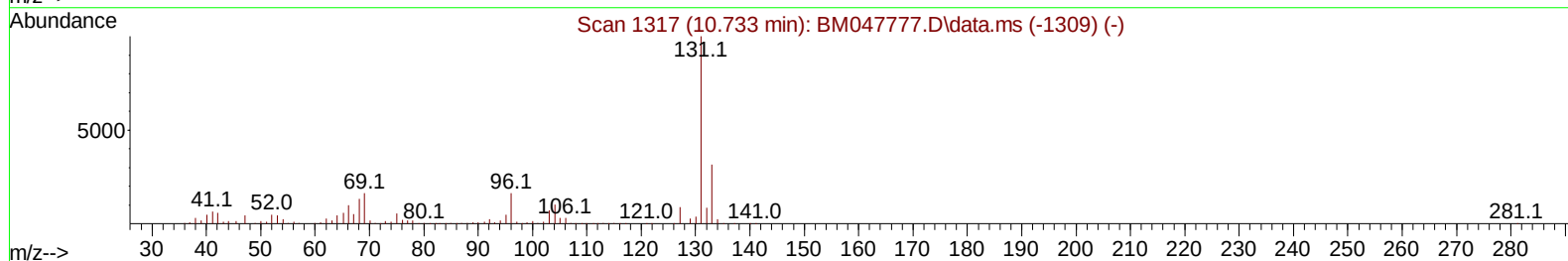
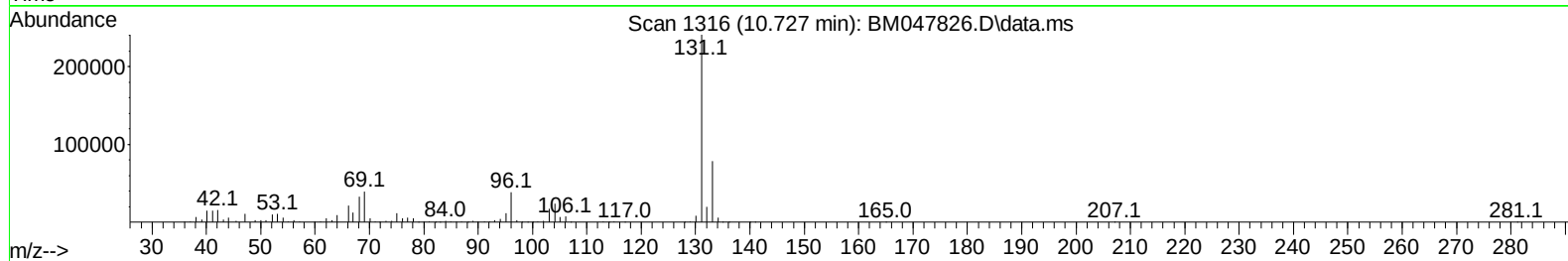
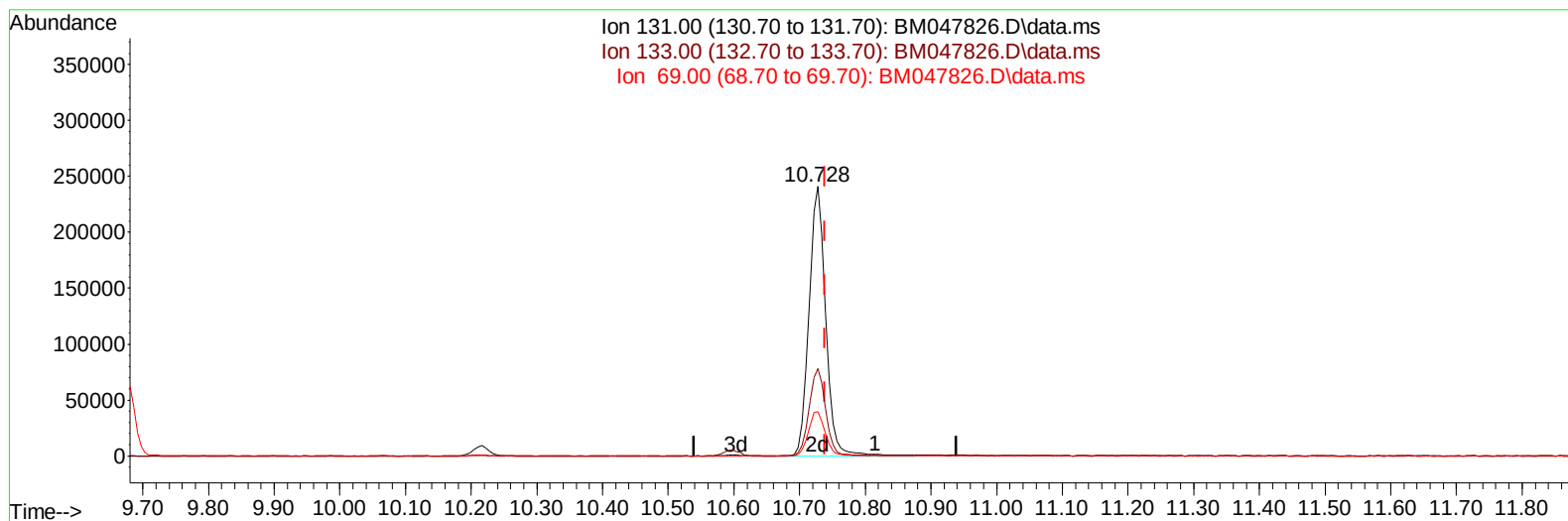
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(31) 4-Chloroaniline-d4 (S)

10.727min (-0.012) 16.30 ng/ul m

response 424424

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	34.00	32.51
69.00	20.90	16.39#
0.00	0.00	0.00

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 Sample : P4084-17
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 ALS Vial : 25 Sample Multiplier: 1

Instrument :
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Manual IntegrationsAPPROVED

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Quant Time: Oct 04 04:00:14 2024
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 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Sep 28 02:30:36 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.804	152	280264	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.592	136	1104900	20.000	ng/ul	-0.02
38) Acenaphthene-d10	14.439	164	676162	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.174	188	1351199	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.397	240	1295566	20.000	ng/ul	-0.01
88) Perylene-d12	24.397	264	1506053	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.252	96	26838	3.095	ng/uL	0.00
4) Pyridine-d5	3.681	84	27160	1.070	ng/ul	0.00
7) Phenol-d5	6.969	99	561242	21.178	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.134	67	364654	19.391	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.334	132	463377	24.032	ng/ul	-0.01
15) 4-Methylphenol-d8	8.504	113	459394	23.315	ng/ul	-0.01
21) Nitrobenzene-d5	8.957	128	238288	27.087	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.675	143	227664	27.005	ng/ul	-0.02
28) 2,4-Dichlorophenol-d3	10.216	165	445087	26.429	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.727	131	424424m	16.301	ng/ul	-0.01
46) Dimethylphthalate-d6	13.845	166	1219727	24.881	ng/ul	-0.01
49) Acenaphthylene-d8	14.127	160	1663969	28.628	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.627	143	200461	20.538	ng/ul	0.00
60) Fluorene-d10	15.427	176	1259386	29.659	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.539	200	144479	18.272	ng/ul	-0.01
73) Anthracene-d10	17.274	188	1961877	29.529	ng/ul	-0.01
81) Pyrene-d10	19.556	212	2287946	31.116	ng/ul	-0.01
92) Benzo(a)pyrene-d12	24.191	264	2377454	30.425	ng/ul	-0.02

Target Compounds Qvalue

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

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