

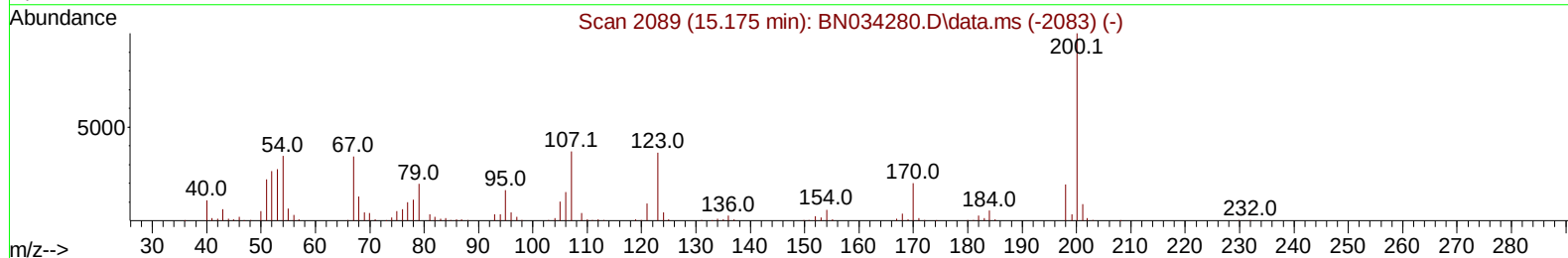
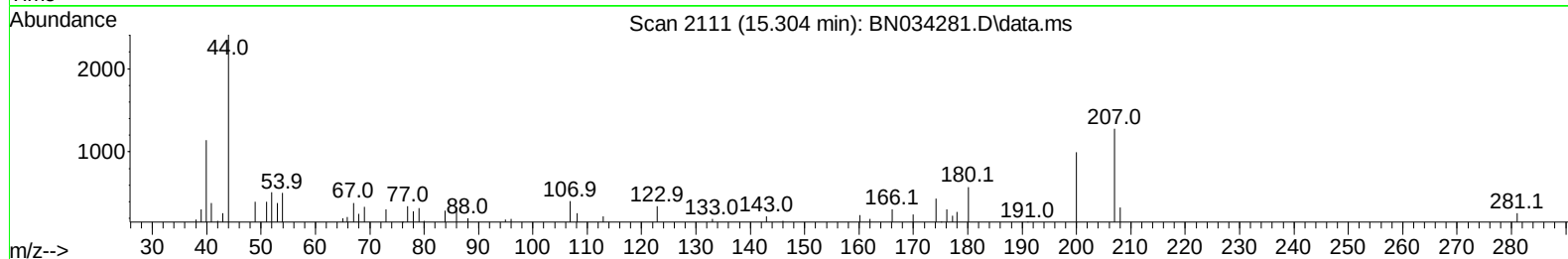
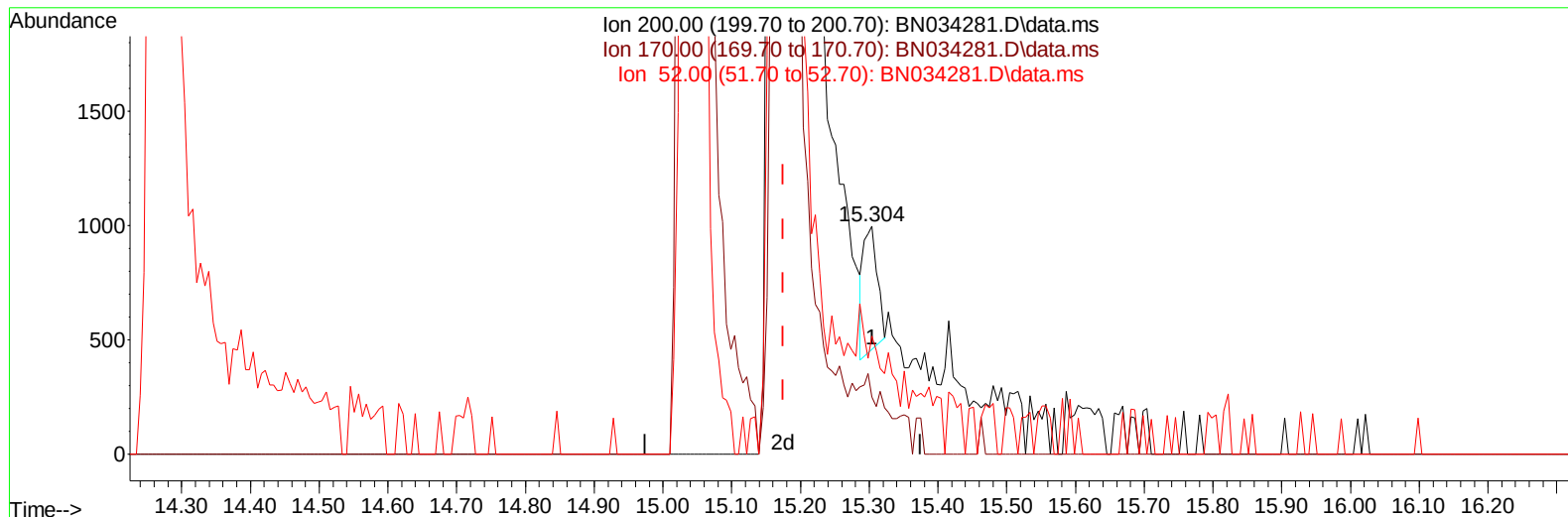
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN093024\
Data File : BN034281.D
Acq On : 30 Sep 2024 17:29
Operator : RC/JU
Sample : PB163394BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
SBLK394

Manual IntegrationsAPPROVED

Quant Time: Sep 30 18:11:05 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN091124.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Sep 30 14:14:38 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 10/01/2024
Supervised By :mohammad ahmed 10/02/2024



TIC: BN034281.D\data.ms

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

15.304min (+ 0.129) 0.13 ng/u1

response 764

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	21.50	24.92
52.00	46.90	51.76
0.00	0.00	0.00

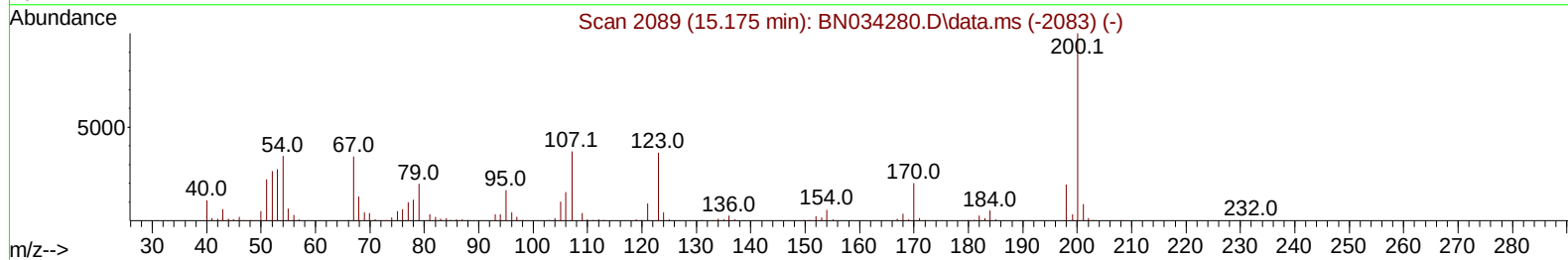
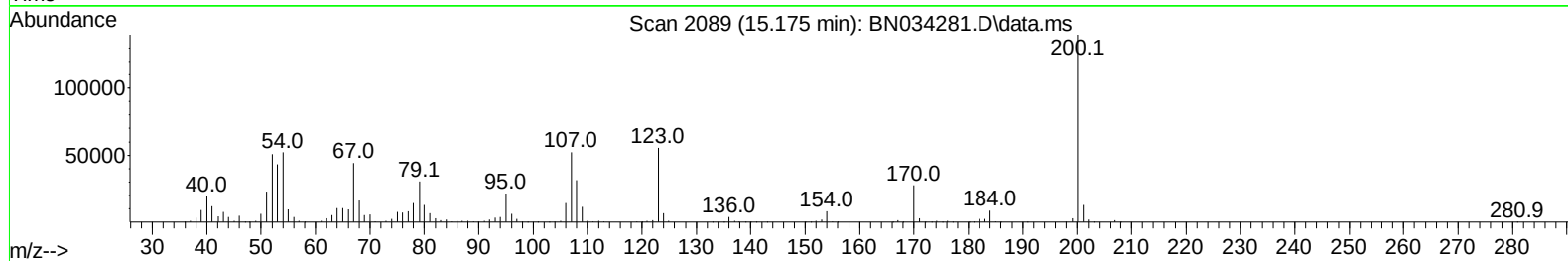
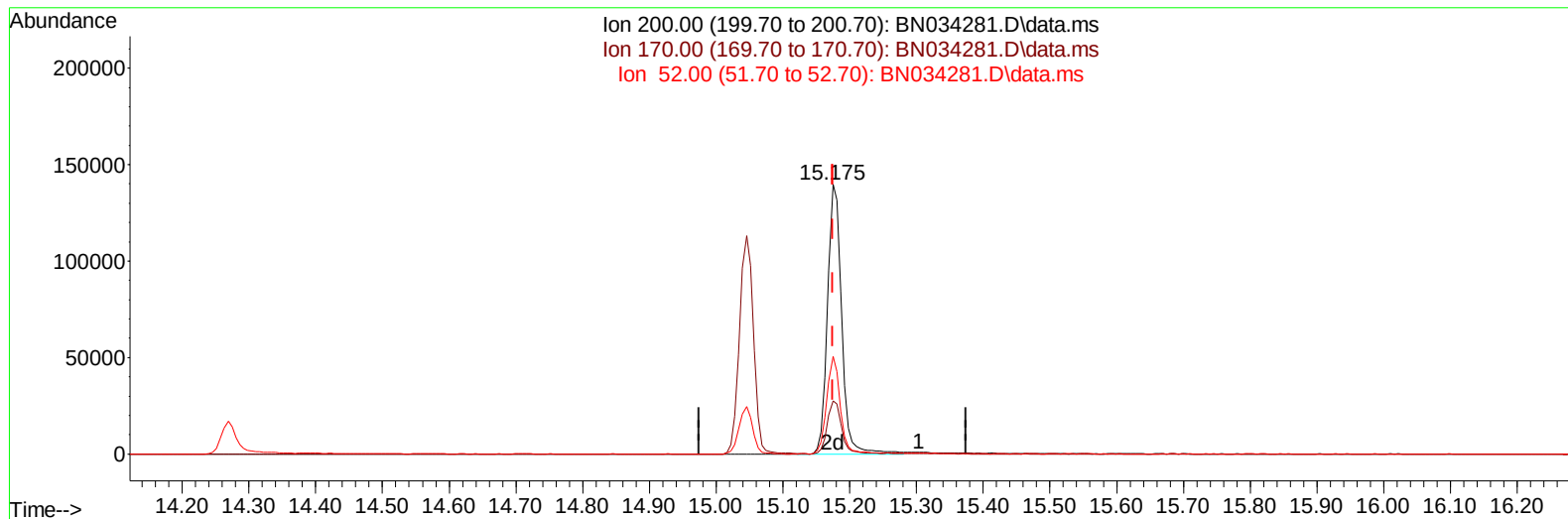
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TIC: BN034281.D\data.ms

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

15.175min (-0.000) 36.06 ng/ul m

response 205302

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	21.50	19.74
52.00	46.90	36.23#
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 10/01/2024
 Supervised By :mohammad ahmed 10/02/2024

Quant Time: Sep 30 18:16:14 2024
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 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Sep 30 14:14:38 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.387	152	208921	20.000	ng/ul	0.00
20) Naphthalene-d8	10.146	136	825309	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.040	164	493845	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.798	188	954784	20.000	ng/ul	0.00
79) Chrysene-d12	21.027	240	862609	20.000	ng/ul	0.00
88) Perylene-d12	23.733	264	1004473	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.999	96	44670	8.725	ng/uL	0.00
4) Pyridine-d5	3.387	84	488828	32.122	ng/ul	0.00
7) Phenol-d5	6.587	99	690363	37.233	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.746	67	457220	39.744	ng/ul	0.00
11) 2-Chlorophenol-d4	6.928	132	622017	42.618	ng/ul	0.00
15) 4-Methylphenol-d8	8.105	113	557311	36.529	ng/ul	0.00
21) Nitrobenzene-d5	8.534	128	301091	45.771	ng/ul	0.00
24) 2-Nitrophenol-d4	9.246	143	310242	45.864	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.775	165	541328	42.493	ng/ul	0.00
31) 4-Chloroaniline-d4	10.293	131	727981	37.746	ng/ul	0.00
46) Dimethylphthalate-d6	13.469	166	1664852	45.110	ng/ul	0.00
49) Acenaphthylene-d8	13.728	160	1922711	46.309	ng/ul	0.00
54) 4-Nitrophenol-d4	14.269	143	234692	30.986	ng/ul	0.00
60) Fluorene-d10	15.045	176	1422960	45.493	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.175	200	205302m	36.064	ng/ul	0.00
73) Anthracene-d10	16.898	188	2182666	48.624	ng/ul	0.00
81) Pyrene-d10	19.204	212	2457422	51.130	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.557	264	2480737	48.079	ng/ul	0.00

Target Compounds Qvalue

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

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