

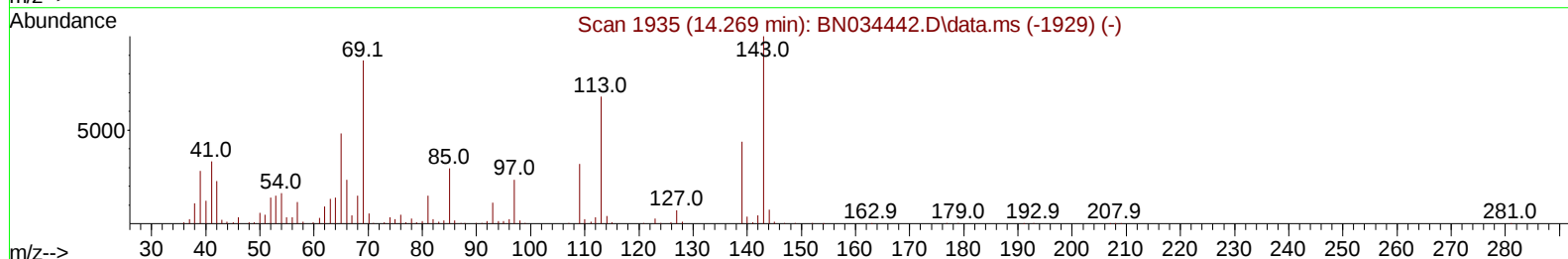
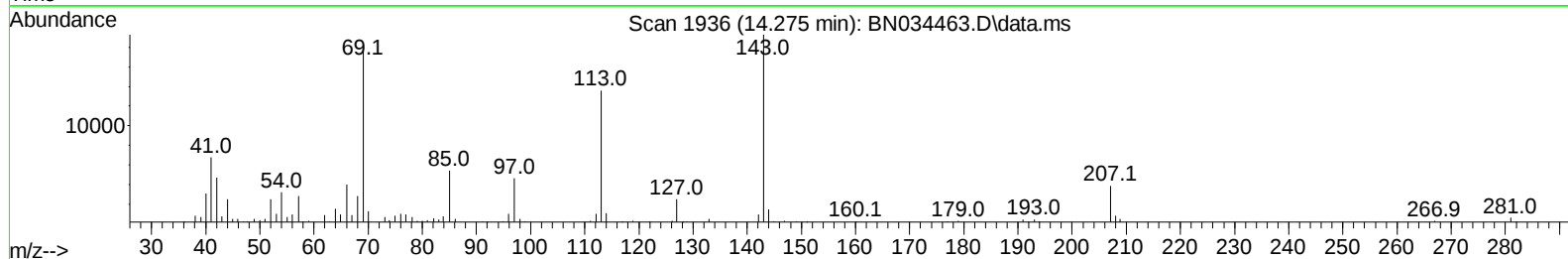
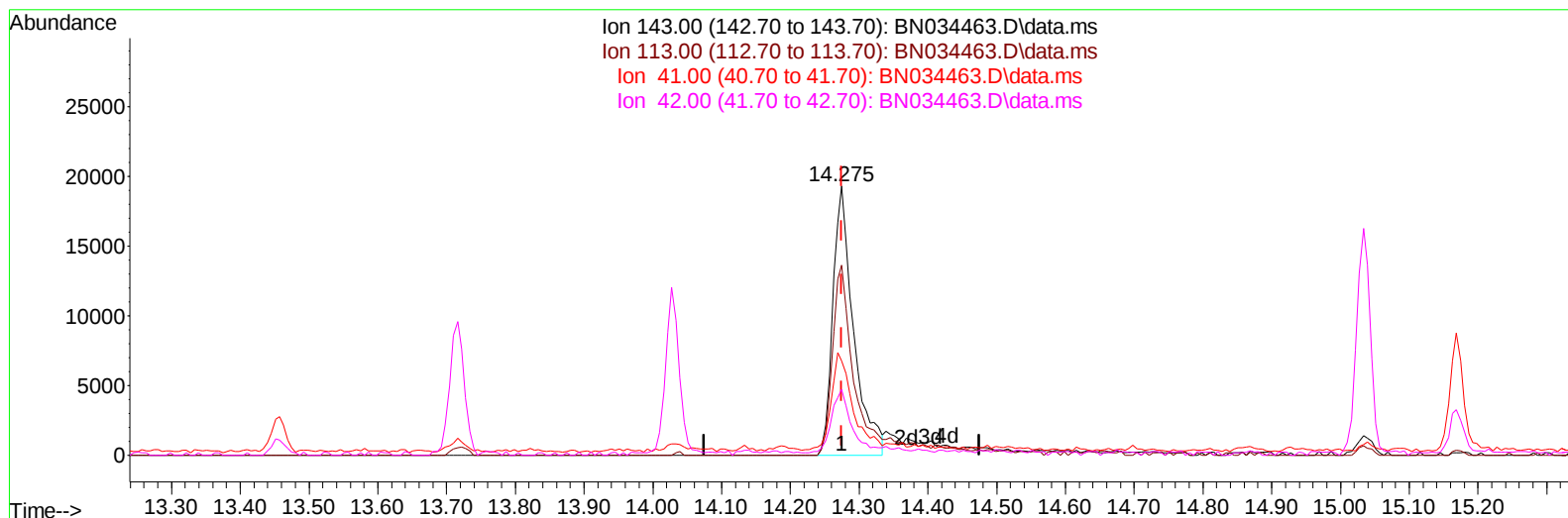
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100724\
Data File : BN034463.D
Acq On : 08 Oct 2024 08:01
Operator : RC/JU
Sample : P4187-02
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
EY046

Manual IntegrationsAPPROVED

Quant Time: Oct 08 08:30:08 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 :Yogesh Patel 10/09/2024
Supervised By :mohammad ahmed 10/09/2024



TIC: BN034463.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.275min (-0.000) 4.26 ng/u1

response 40271

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	76.90	70.54
41.00	41.70	35.02
42.00	28.50	24.31

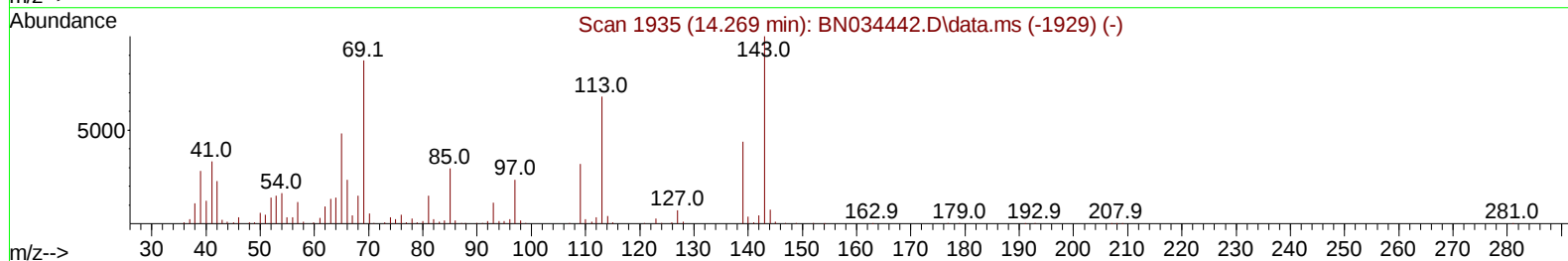
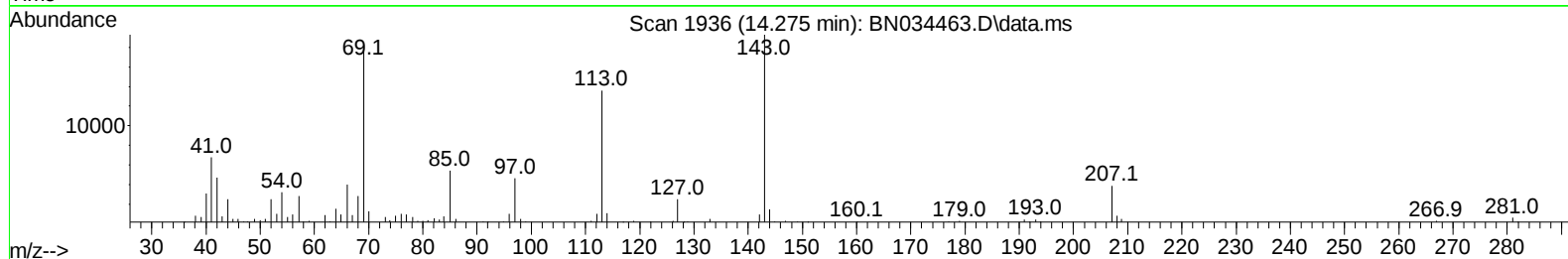
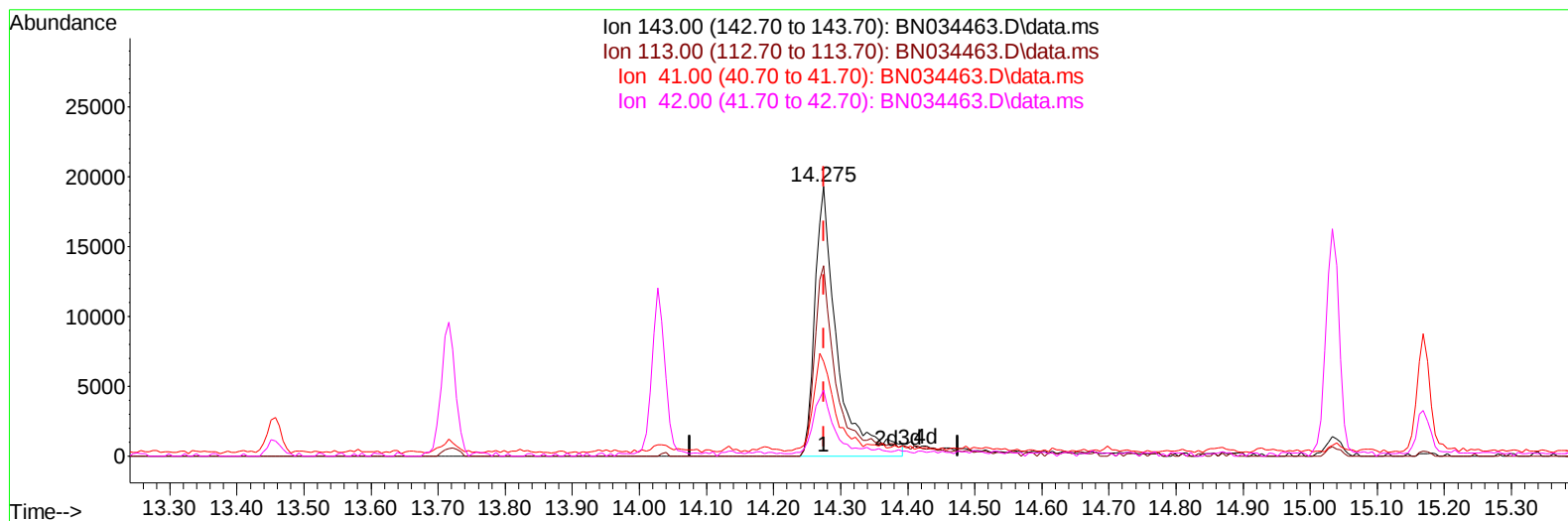
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TIC: BN034463.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.275min (-0.000) 4.69 ng/ul m

response 44347

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	76.90	70.54
41.00	41.70	35.02
42.00	28.50	24.31

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Manual IntegrationsAPPROVED

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 Supervised By :mohammad ahmed 10/09/2024

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 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	7.375	152	242523	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.134	136	989679	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.028	164	615868	20.000	ng/ul	-0.01
64) Phenanthrene-d10	16.786	188	1165418	20.000	ng/ul	-0.01
79) Chrysene-d12	21.021	240	825870	20.000	ng/ul	-0.01
88) Perylene-d12	23.733	264	992240	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.993	96	20375	3.428	ng/uL	0.00
4) Pyridine-d5	3.387	84	81635	4.621	ng/ul	0.00
7) Phenol-d5	6.581	99	137451	6.386	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.734	67	281220	21.058	ng/ul	0.00
11) 2-Chlorophenol-d4	6.916	132	381091	22.493	ng/ul	-0.01
15) 4-Methylphenol-d8	8.093	113	253513	14.314	ng/ul	-0.01
21) Nitrobenzene-d5	8.522	128	199667	25.312	ng/ul	0.00
24) 2-Nitrophenol-d4	9.234	143	222810	27.468	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.763	165	390720	25.577	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.281	131	408898	17.680	ng/ul	-0.01
46) Dimethylphthalate-d6	13.457	166	1310785	28.479	ng/ul	-0.01
49) Acenaphthylene-d8	13.716	160	1315016	25.397	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.275	143	44347m	4.695	ng/ul	0.00
60) Fluorene-d10	15.033	176	1051152	26.948	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.169	200	150392	21.643	ng/ul	0.00
73) Anthracene-d10	16.886	188	1661153	30.318	ng/ul	-0.01
81) Pyrene-d10	19.198	212	1678230	36.471	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.545	264	1606982	31.529	ng/ul	-0.01

Target Compounds Qvalue

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

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