

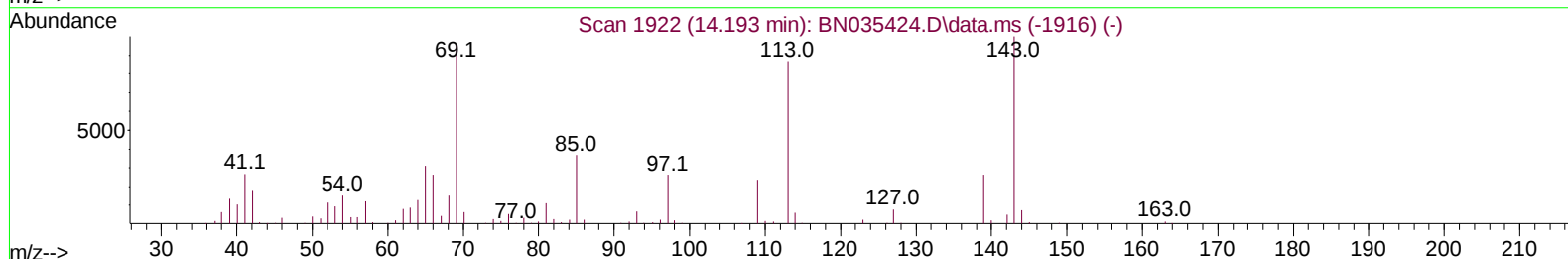
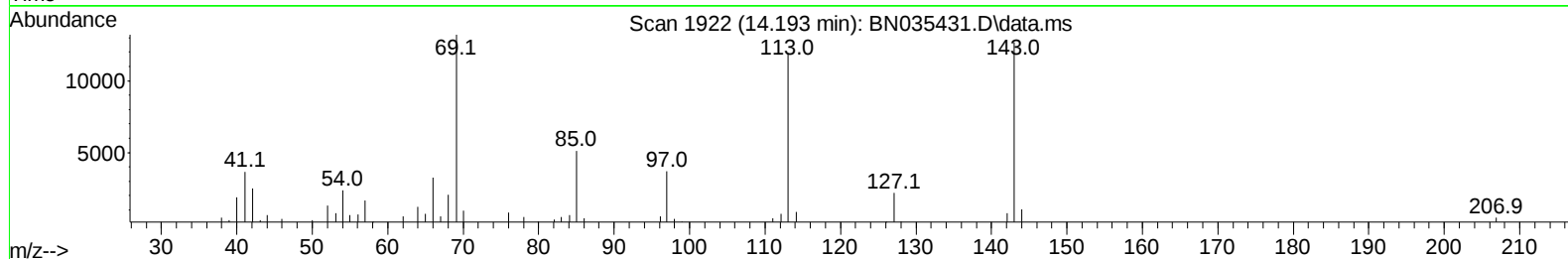
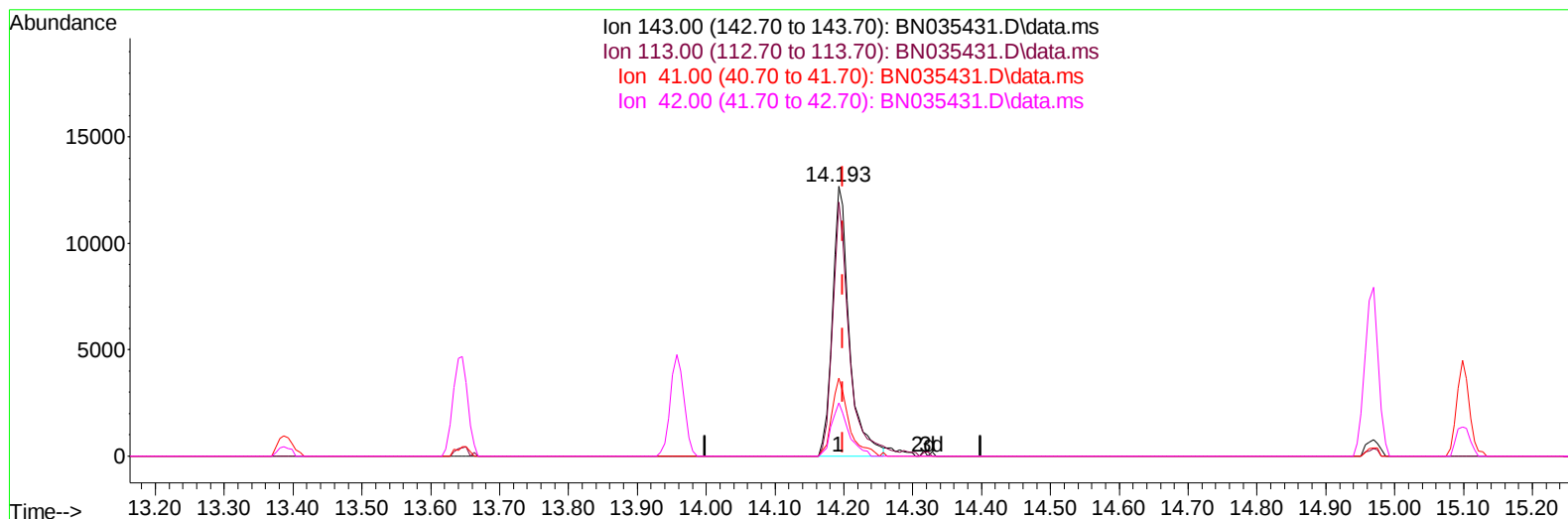
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN120624\
Data File : BN035431.D
Acq On : 05 Dec 2024 17:45
Operator : RC/JU
Sample : P5104-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BHJ63

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 12/07/2024
Supervised By :mohammad ahmed 12/13/2024

Quant Time: Dec 06 00:00:21 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN120224.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Dec 02 18:25:21 2024
Response via : Initial Calibration



TIC: BN035431.D\data.ms

(54) 4-Nitrophenol-d4 (S)**14.193min (-0.006) 6.32 ng/u1****response 21608**

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	84.60	94.30
41.00	26.90	28.91
42.00	17.10	19.73

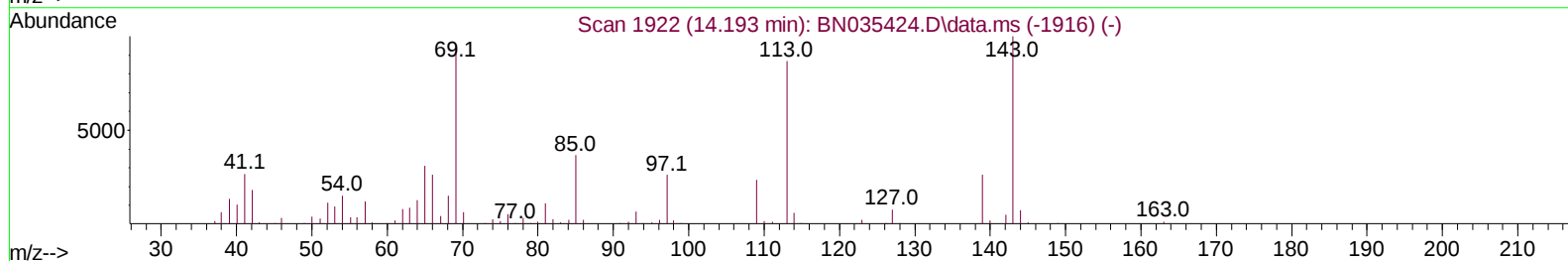
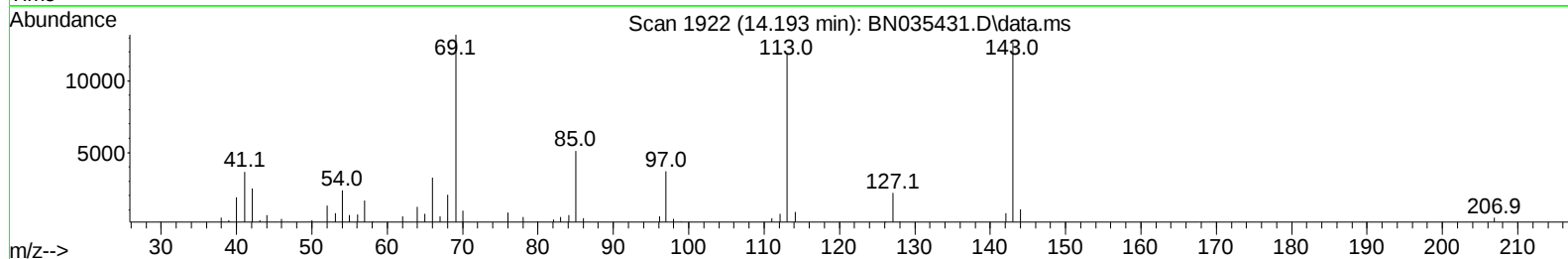
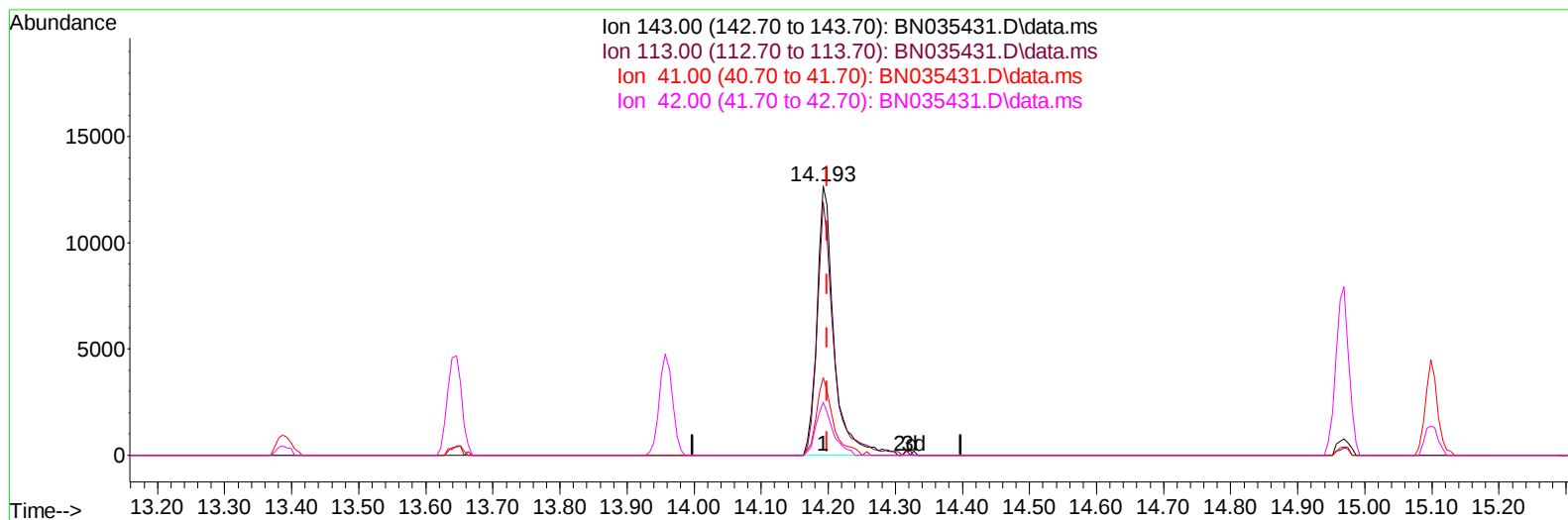
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TIC: BN035431.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.193min (-0.006) 6.50 ng/u1 m

response 22234

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	84.60	94.30
41.00	26.90	28.91
42.00	17.10	19.73

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Reviewed By :Yogesh Patel 12/07/2024
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 QLast Update : Mon Dec 02 18:25:21 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.299	152	90540	20.000	ng/ul	0.00
20) Naphthalene-d8	10.046	136	378673	20.000	ng/ul	0.00
38) Acenaphthene-d10	13.957	164	289270	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.722	188	667534	20.000	ng/ul	0.00
79) Chrysene-d12	20.980	240	713510	20.000	ng/ul	0.00
88) Perylene-d12	23.651	264	814717	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.964	96	8252	4.307	ng/uL	0.00
4) Pyridine-d5	3.358	84	7089	1.404	ng/ul	0.02
7) Phenol-d5	6.499	99	50690	7.859	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.658	67	115020	31.976	ng/ul	0.00
11) 2-Chlorophenol-d4	6.840	132	147497	26.986	ng/ul	0.00
15) 4-Methylphenol-d8	8.011	113	108875	19.369	ng/ul	0.00
21) Nitrobenzene-d5	8.440	128	82661	33.599	ng/ul	0.00
24) 2-Nitrophenol-d4	9.146	143	97542	34.254	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.681	165	190480	30.605	ng/ul	0.00
31) 4-Chloroaniline-d4	10.193	131	175157	22.792	ng/ul	0.00
46) Dimethylphthalate-d6	13.387	166	749822	36.100	ng/ul	0.00
49) Acenaphthylene-d8	13.646	160	752028	32.877	ng/ul	0.00
54) 4-Nitrophenol-d4	14.193	143	22234m	6.501	ng/ul	0.00
60) Fluorene-d10	14.969	176	621581	34.653	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.098	200	113443	32.224	ng/ul	0.00
73) Anthracene-d10	16.828	188	1064350	36.150	ng/ul	0.00
81) Pyrene-d10	19.145	212	1321167	35.058	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.474	264	1484389	37.626	ng/ul	0.00

Target Compounds Qvalue

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

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