

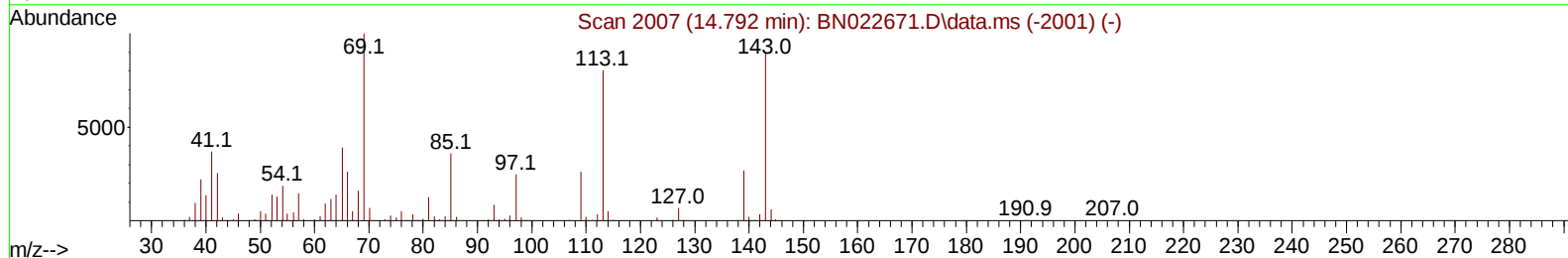
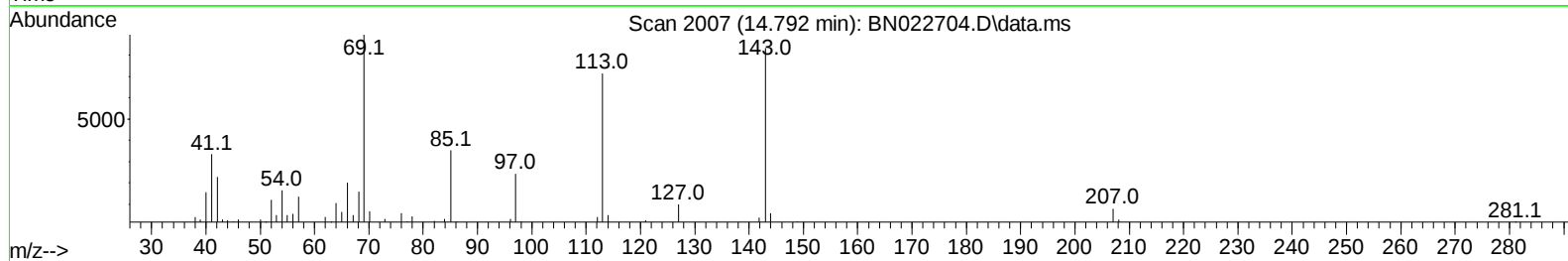
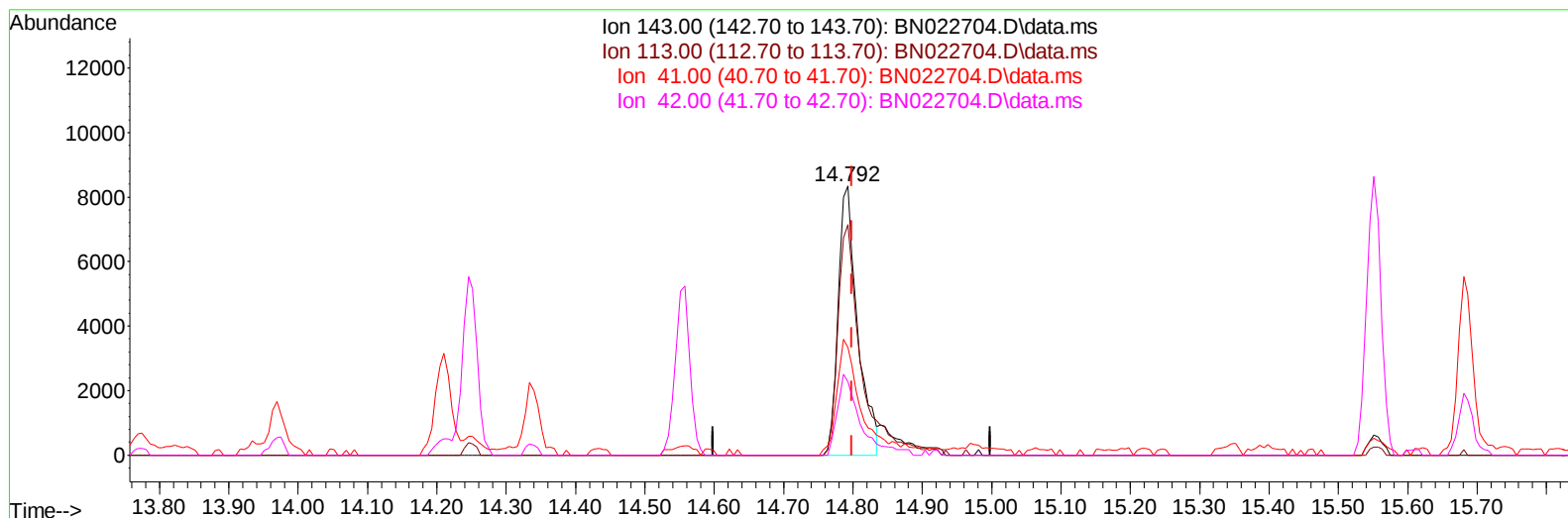
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111522\
Data File : BN022704.D
Acq On : 17 Nov 2022 02:10
Operator : CG/JU
Sample : N5577-05
Misc :
ALS Vial : 64 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
C9818

Manual IntegrationsAPPROVED

Quant Time: Nov 17 03:27:06 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN111522.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Nov 15 22:53:43 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 11/17/2022
Supervised By :mohammad ahmed 11/18/2022



TIC: BN022704.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.792min (-0.006) 5.07 ng/u1

response 16089

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	95.00	85.61
41.00	46.30	40.01
42.00	31.30	27.30

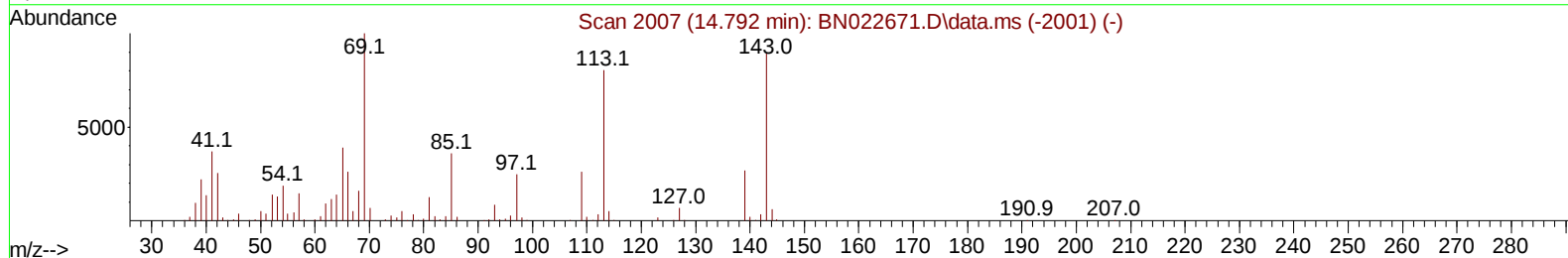
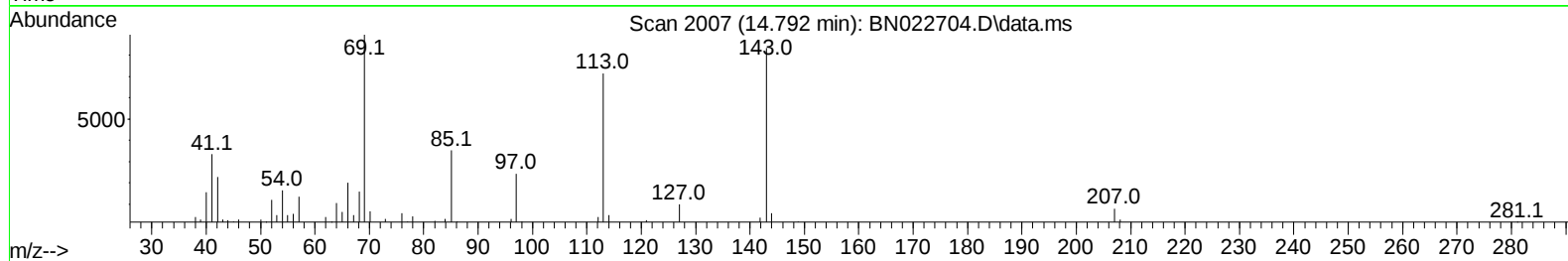
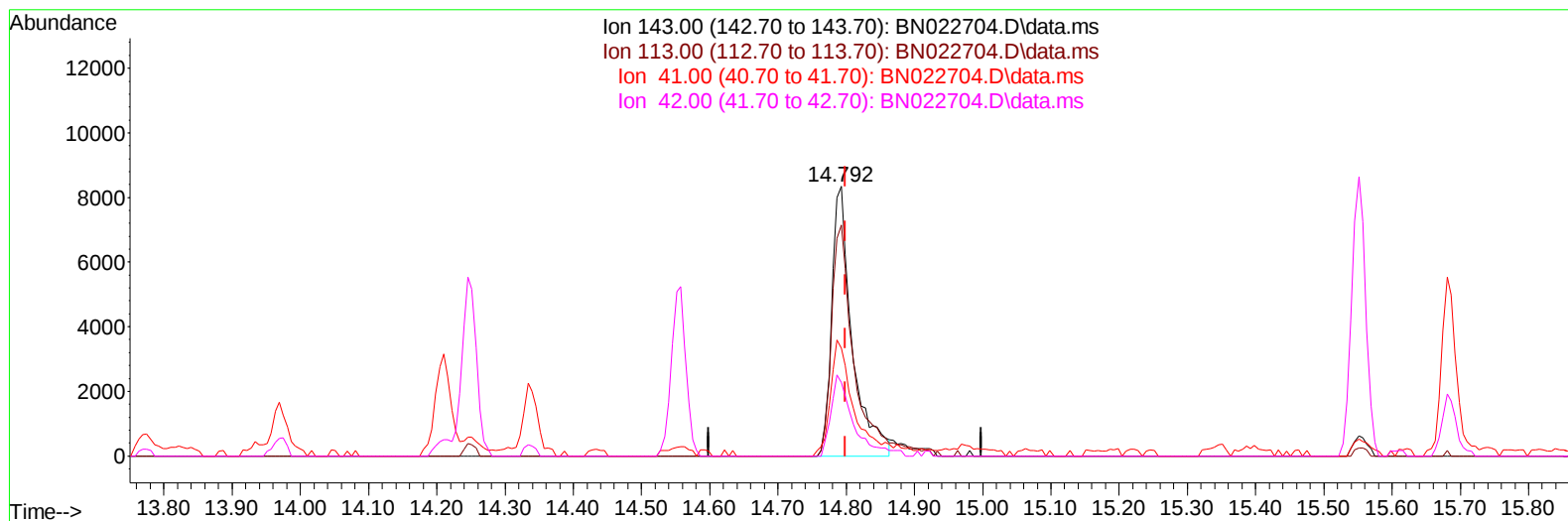
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TIC: BN022704.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.792min (-0.006) 5.47 ng/ul m

response 17345

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	95.00	85.61
41.00	46.30	40.01
42.00	31.30	27.30

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.869	152	78557	20.000	ng/ul	0.00
20) Naphthalene-d8	10.693	136	342924	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.557	164	218715	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.316	188	451994	20.000	ng/ul	0.00
79) Chrysene-d12	21.522	240	415268	20.000	ng/ul	-0.01
88) Perylene-d12	23.951	264	375423	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.187	96	9853	4.282	ng/uL	0.00
4) Pyridine-d5	3.634	84	53125	8.164	ng/ul	0.00
7) Phenol-d5	7.040	99	47794	6.442	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.199	67	131435	29.546	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.399	132	123786	22.781	ng/ul	0.00
15) 4-Methylphenol-d8	8.599	113	82997	14.169	ng/ul	-0.01
21) Nitrobenzene-d5	9.052	128	82844	31.210	ng/ul	0.00
24) 2-Nitrophenol-d4	9.781	143	83640	28.240	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.322	165	135508	26.838	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.846	131	117121	15.228	ng/ul	0.00
46) Dimethylphthalate-d6	13.969	166	518097	32.567	ng/ul	0.00
49) Acenaphthylene-d8	14.245	160	542742	30.963	ng/ul	0.00
54) 4-Nitrophenol-d4	14.792	143	17345m	5.469	ng/ul	0.00
60) Fluorene-d10	15.551	176	423539	32.461	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.687	200	75678	26.341	ng/ul	0.00
73) Anthracene-d10	17.416	188	642565	32.466	ng/ul	0.00
81) Pyrene-d10	19.722	212	754363	35.053	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.798	264	657154	34.822	ng/ul	0.00

Target Compounds Qvalue

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

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