

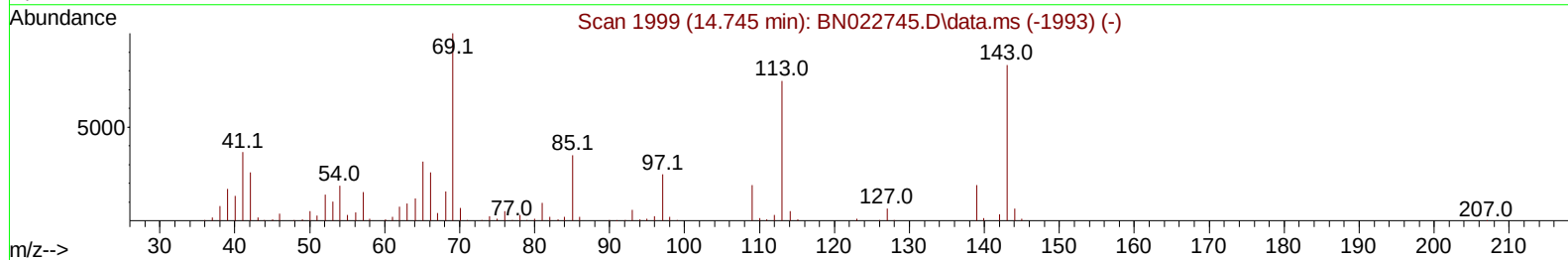
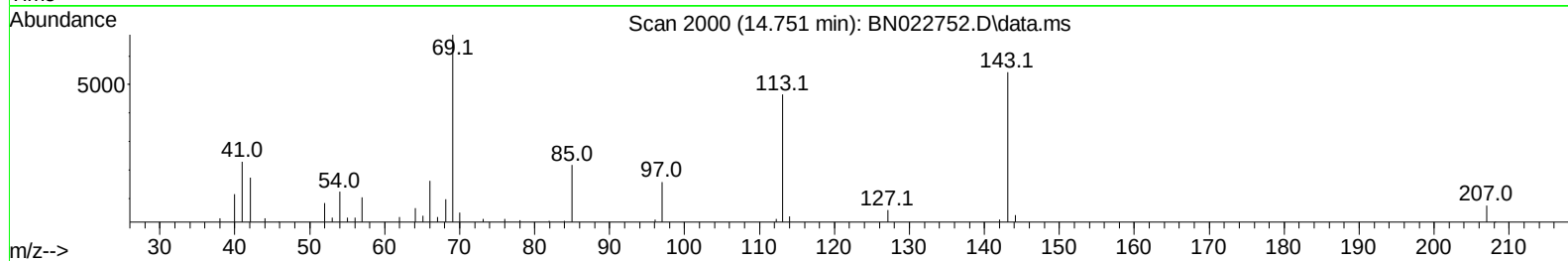
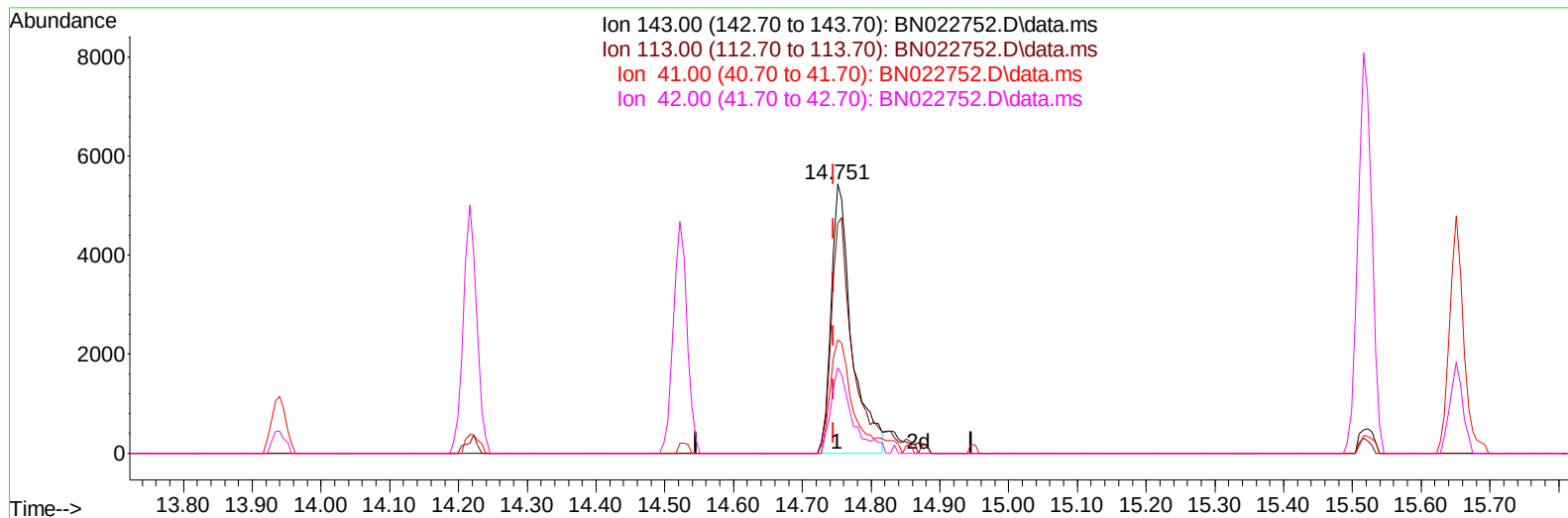
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111722\
Data File : BN022752.D
Acq On : 18 Nov 2022 12:10
Operator : CG/JU
Sample : N5689-03
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BGCC6

Manual IntegrationsAPPROVED

Quant Time: Nov 18 21:47:57 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN111522.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Nov 18 05:17:17 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 11/21/2022
Supervised By :mohammad ahmed 11/21/2022



TIC: BN022752.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.751min (+ 0.006) 4.46 ng/ul

response 11228

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	95.00	85.55
41.00	46.30	42.16
42.00	31.30	31.73

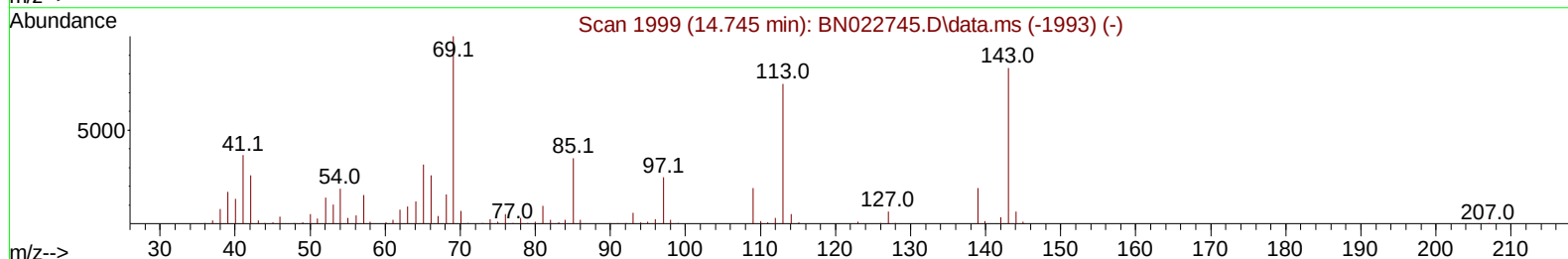
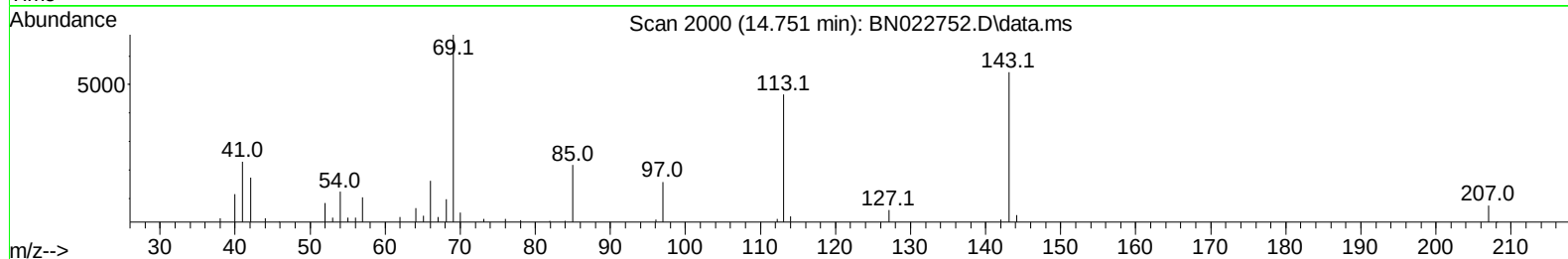
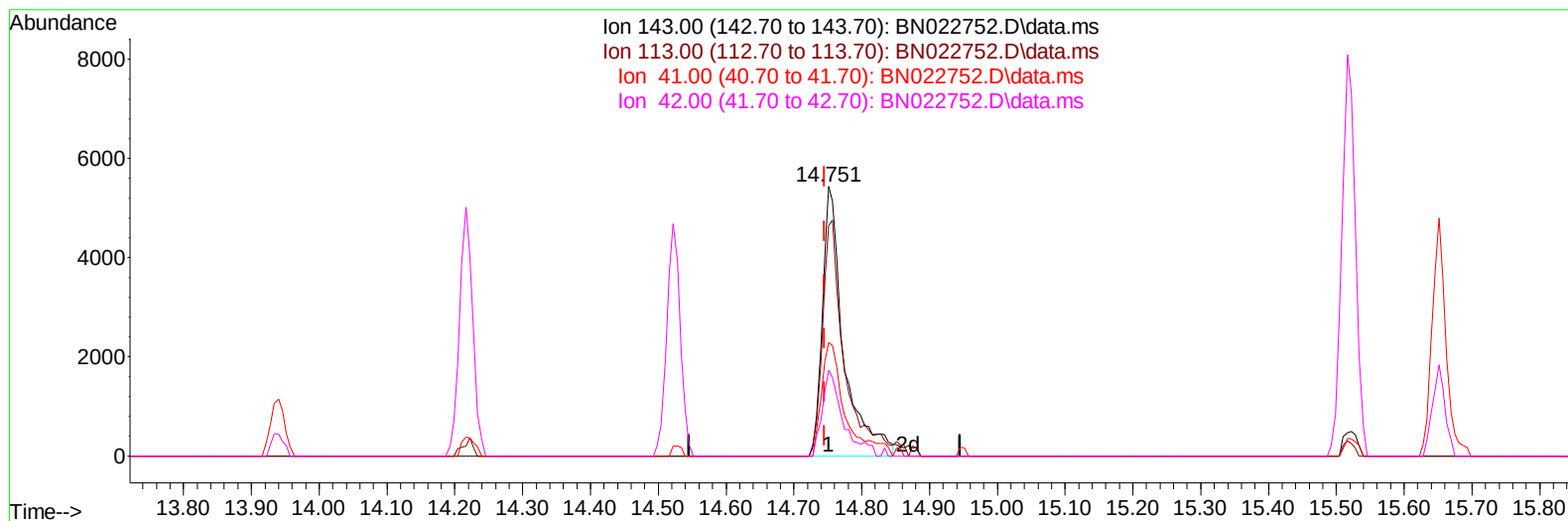
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TIC: BN022752.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.751min (+ 0.006) 4.71 ng/ul m

response 11873

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	95.00	85.55
41.00	46.30	42.16
42.00	31.30	31.73

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.852	152	51646	20.000	ng/ul	0.00
20) Naphthalene-d8	10.669	136	249776	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.522	164	173719	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.281	188	375797	20.000	ng/ul	0.00
79) Chrysene-d12	21.486	240	406216	20.000	ng/ul	0.00
88) Perylene-d12	23.886	264	422476	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.181	96	5310	3.510	ng/uL	0.00
4) Pyridine-d5	3.628	84	22055	5.155	ng/ul	0.01
7) Phenol-d5	7.022	99	27981	5.736	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.187	67	100782	34.460	ng/ul	0.00
11) 2-Chlorophenol-d4	7.381	132	85504	23.936	ng/ul	0.00
15) 4-Methylphenol-d8	8.581	113	58692	15.241	ng/ul	0.00
21) Nitrobenzene-d5	9.034	128	61227	31.669	ng/ul	0.00
24) 2-Nitrophenol-d4	9.758	143	64738	30.009	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.299	165	100654	27.369	ng/ul	0.00
31) 4-Chloroaniline-d4	10.822	131	132131	23.587	ng/ul	0.00
46) Dimethylphthalate-d6	13.940	166	435140	34.438	ng/ul	0.00
49) Acenaphthylene-d8	14.216	160	459398	32.997	ng/ul	0.00
54) 4-Nitrophenol-d4	14.751	143	11873m	4.714	ng/ul	0.00
60) Fluorene-d10	15.522	176	367681	35.479	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.651	200	64608	27.048	ng/ul	0.00
73) Anthracene-d10	17.381	188	609503	37.040	ng/ul	0.00
81) Pyrene-d10	19.681	212	697947	33.154	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.733	264	787600	37.086	ng/ul	0.00

Target Compounds Qvalue

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

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