

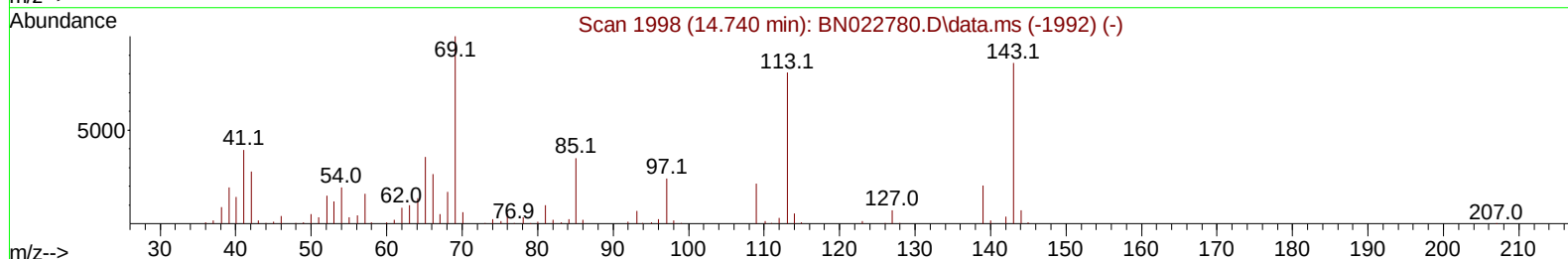
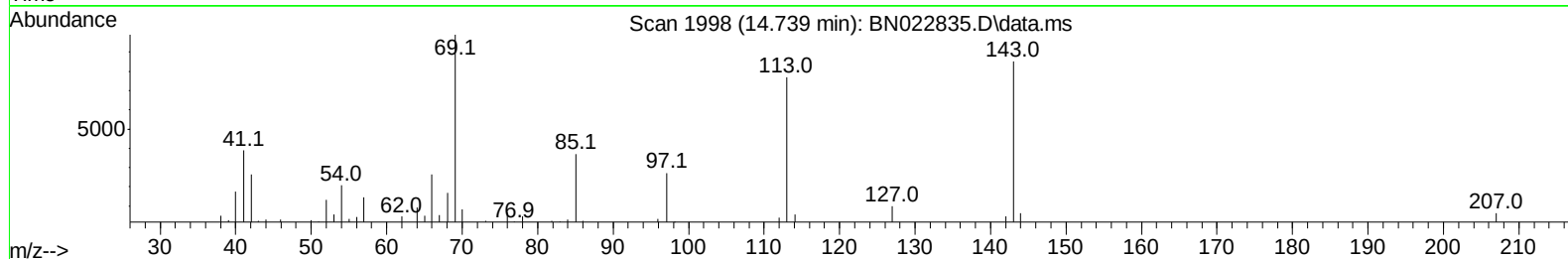
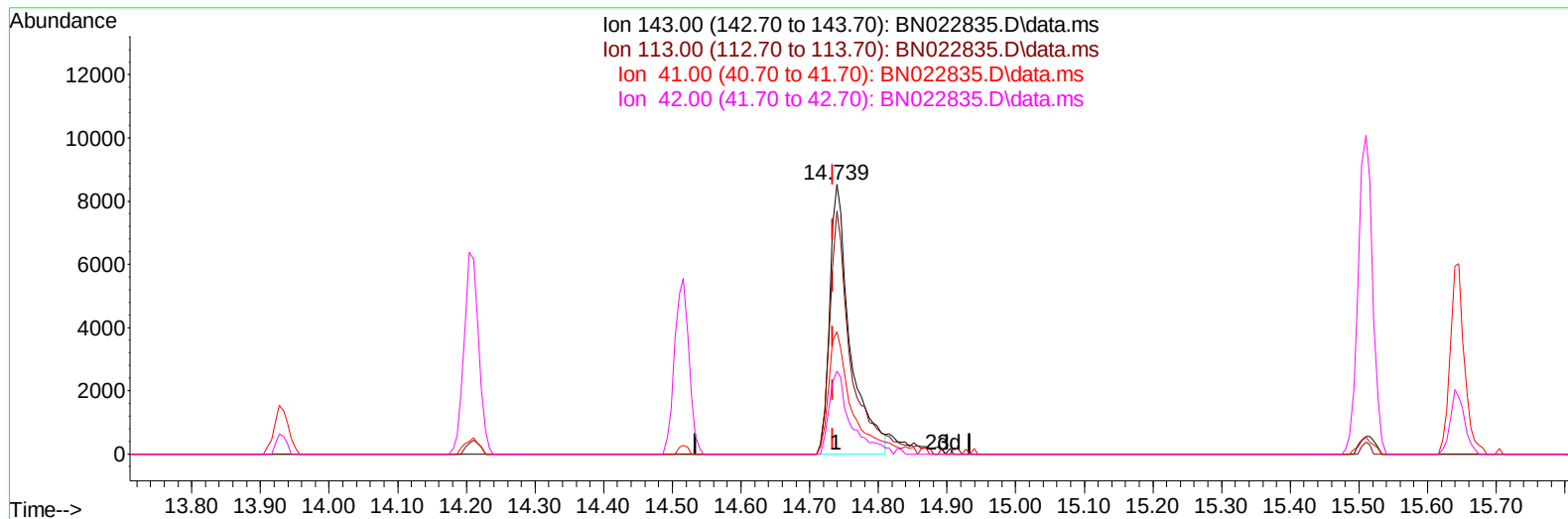
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112122\
Data File : BN022835.D
Acq On : 22 Nov 2022 21:35
Operator : CG/JU
Sample : N5592-21
Misc :
ALS Vial : 57 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
H0AE1

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/23/2022
Supervised By :mohammad ahmed 11/25/2022

Quant Time: Nov 22 22:22:39 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN11522.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Nov 22 05:15:51 2022
Response via : Initial Calibration



TIC: BN022835.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.739min (+ 0.006) 5.77 ng/ul

response 17643

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	95.00	90.29
41.00	46.30	45.48
42.00	31.30	30.87

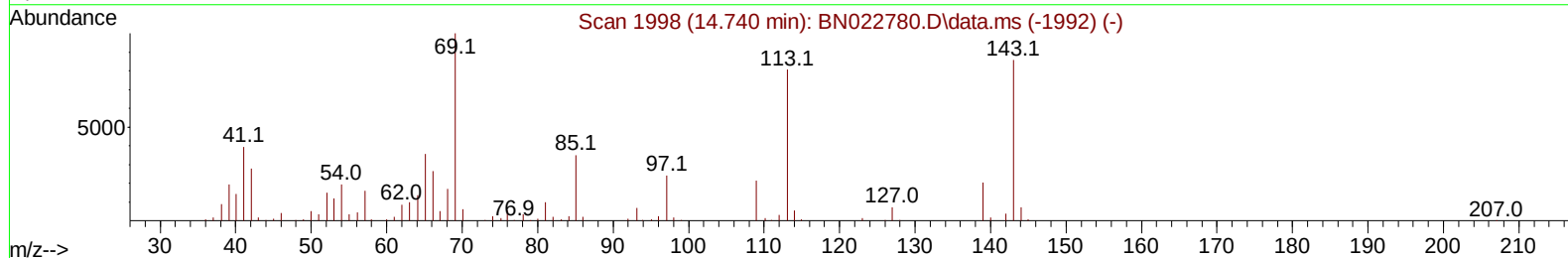
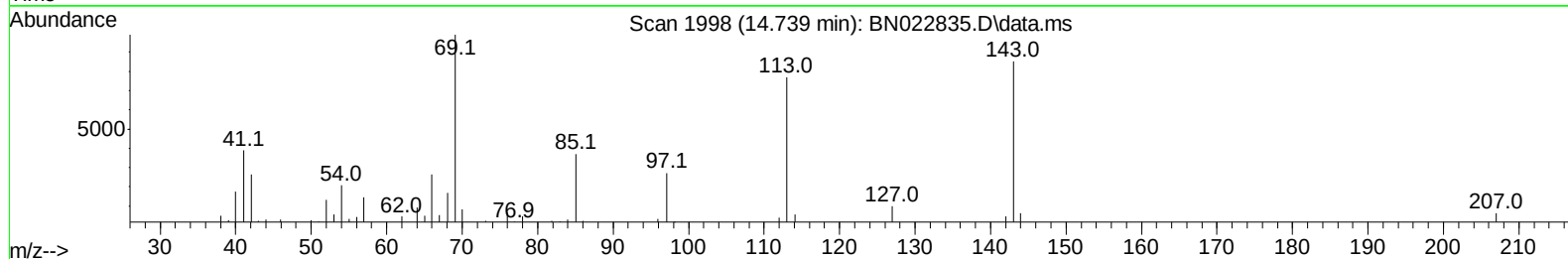
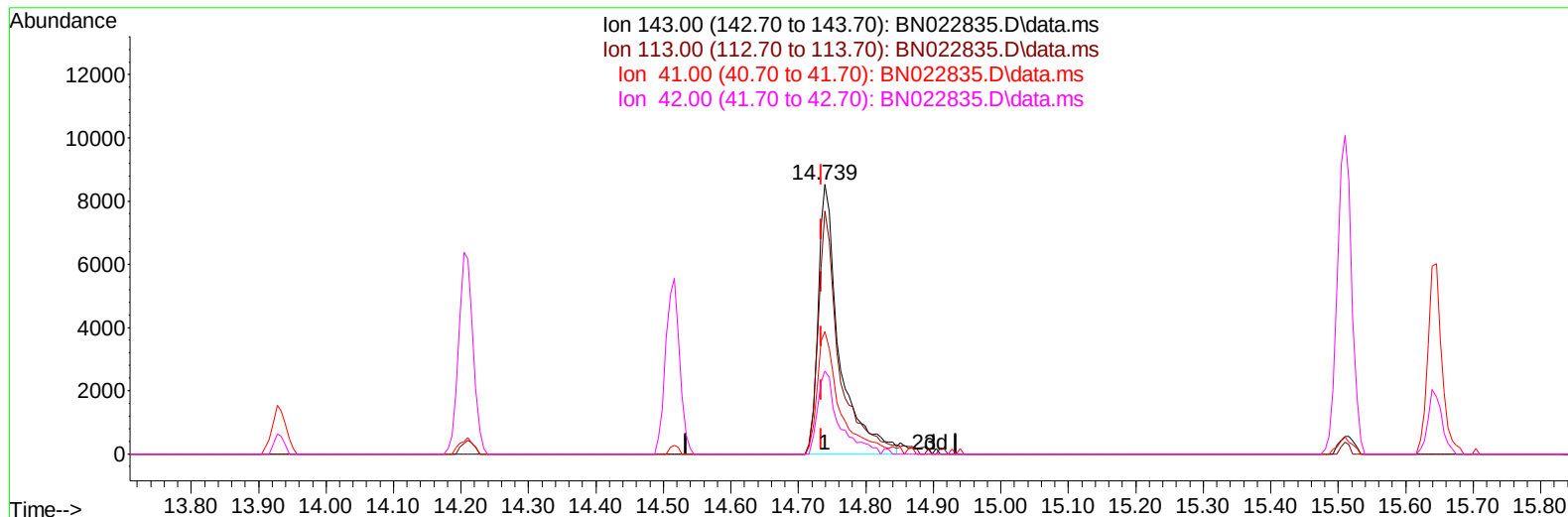
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TIC: BN022835.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.739min (+ 0.006) 6.06 ng/ul m

response 18545

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	95.00	90.29
41.00	46.30	45.48
42.00	31.30	30.87

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 Sample : N5592-21
 Misc :
 ALS Vial : 57 Sample Multiplier: 1

Instrument :
 BNA_N
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Quant Time: Nov 22 22:22:39 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN111522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 22 05:15:51 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.846	152	66373	20.000	ng/ul	0.00
20) Naphthalene-d8	10.663	136	316807	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.516	164	210880	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.269	188	425936	20.000	ng/ul	0.00
79) Chrysene-d12	21.469	240	406687	20.000	ng/ul	0.00
88) Perylene-d12	23.863	264	400966	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.175	96	9230	4.748	ng/uL	0.00
4) Pyridine-d5	3.617	84	57003	10.368	ng/ul	0.00
7) Phenol-d5	7.016	99	51088	8.150	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.181	67	143616	38.210	ng/ul	0.00
11) 2-Chlorophenol-d4	7.369	132	126346	27.521	ng/ul	0.00
15) 4-Methylphenol-d8	8.569	113	94856	19.167	ng/ul	0.00
21) Nitrobenzene-d5	9.022	128	88689	36.167	ng/ul	0.00
24) 2-Nitrophenol-d4	9.746	143	89647	32.763	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.287	165	143077	30.673	ng/ul	0.00
31) 4-Chloroaniline-d4	10.810	131	211912	29.825	ng/ul	0.00
46) Dimethylphthalate-d6	13.934	166	559059	36.448	ng/ul	0.00
49) Acenaphthylene-d8	14.210	160	612369	36.234	ng/ul	0.00
54) 4-Nitrophenol-d4	14.739	143	18545m	6.065	ng/ul	0.00
60) Fluorene-d10	15.510	176	464538	36.927	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	79334	29.303	ng/ul	0.00
73) Anthracene-d10	17.369	188	721730	38.697	ng/ul	0.00
81) Pyrene-d10	19.674	212	802190	38.062	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.710	264	772891	38.345	ng/ul	0.00

Target Compounds Qvalue

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

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