

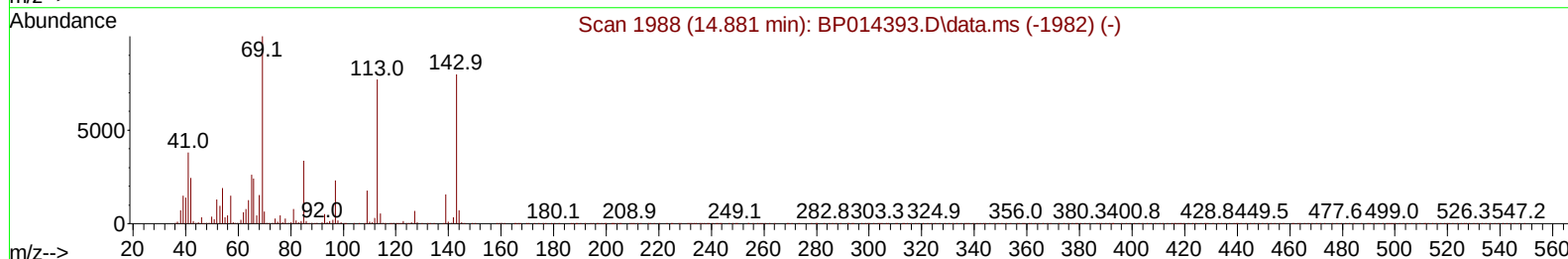
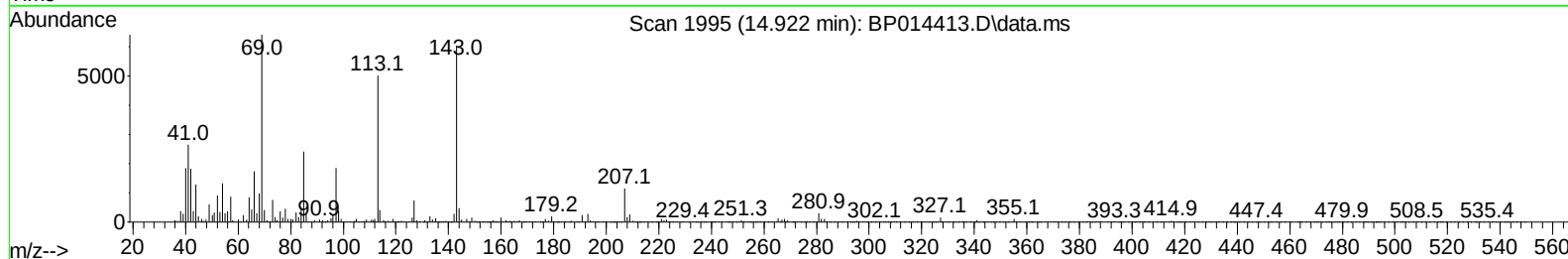
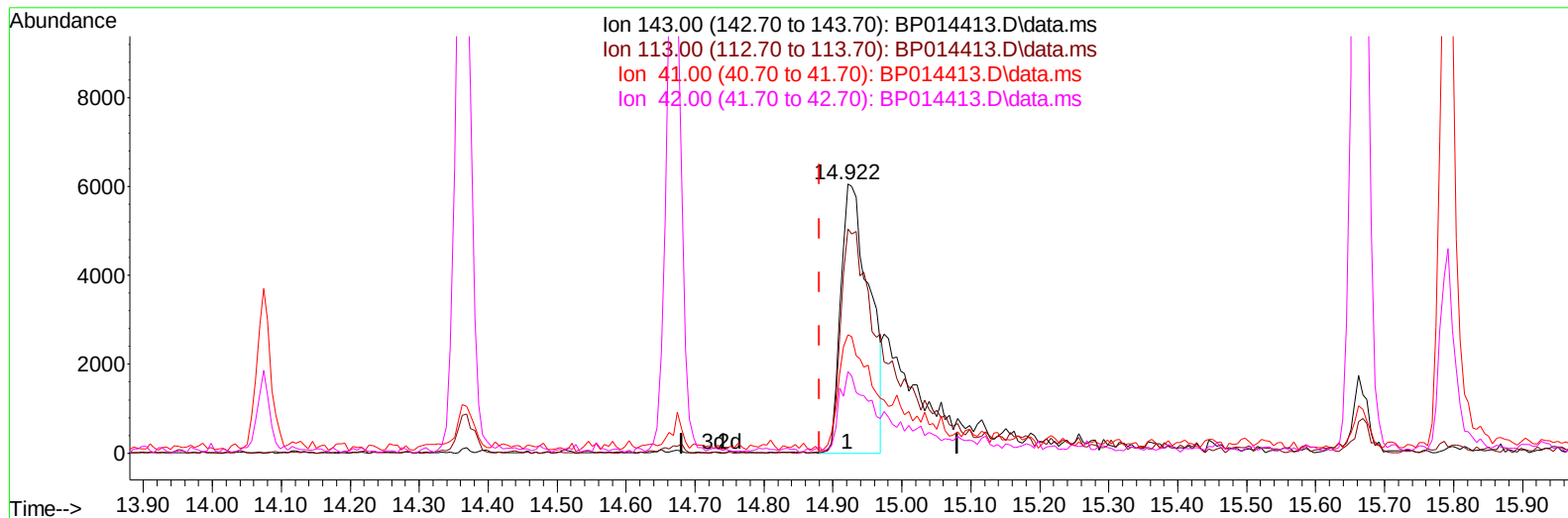
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
Data File : BP014413.D
Acq On : 30 Mar 2023 21:29
Operator : CG/JU
Sample : 02038-08
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EYF73

Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 03/31/2023
Supervised By :Jagrut Upadhyay 03/31/2023

Quant Time: Mar 30 22:44:31 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Mar 30 12:32:14 2023
Response via : Initial Calibration



TIC: BP014413.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.922min (+ 0.041) 2.74 ng/u1

response 17269

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	83.28
41.00	44.20	44.07
42.00	31.50	30.25

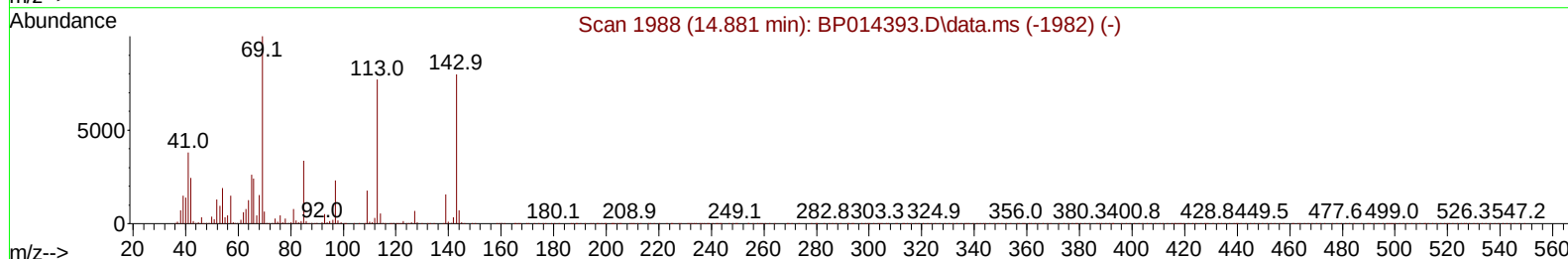
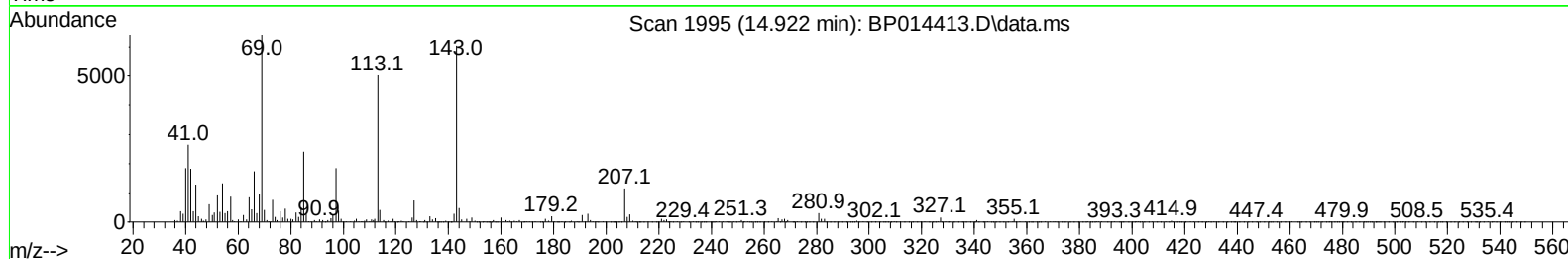
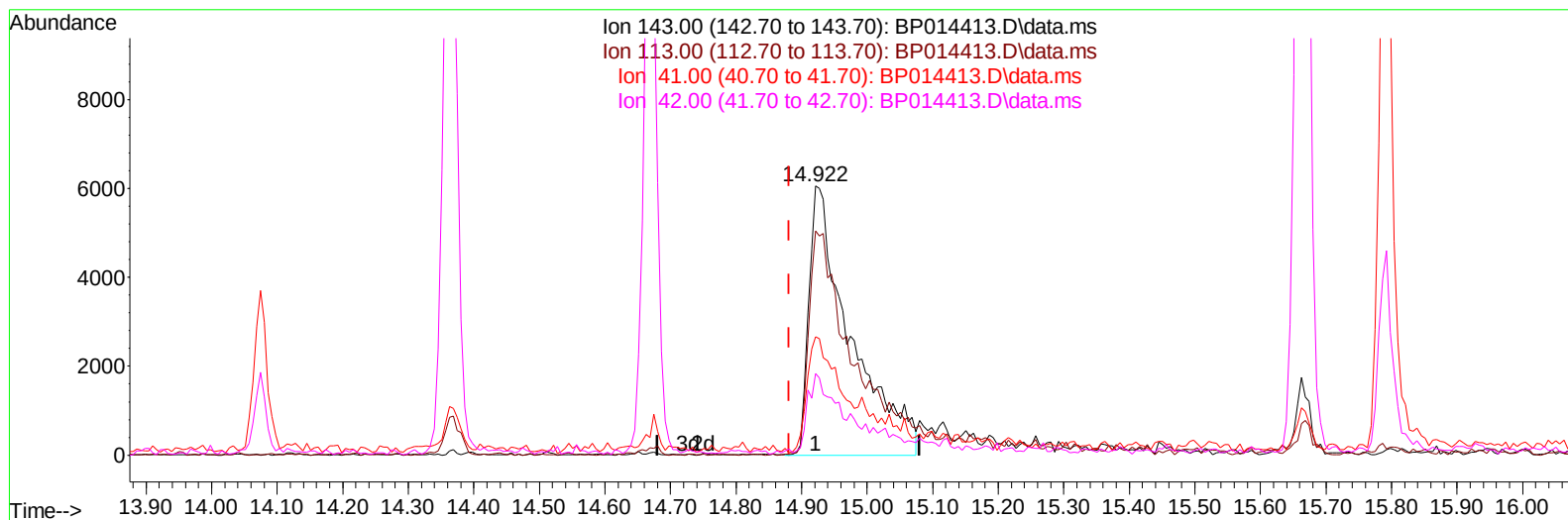
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TIC: BP014413.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.922min (+ 0.041) 4.21 ng/ul m

response 26459

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	83.28
41.00	44.20	44.07
42.00	31.50	30.25

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.016	152	173464	20.000	ng/ul	0.00
20) Naphthalene-d8	10.839	136	773833	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.669	164	513653	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.433	188	1067452	20.000	ng/ul	0.00
79) Chrysene-d12	21.521	240	807362	20.000	ng/ul	0.00
88) Perylene-d12	24.039	264	848290	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	21211	4.742	ng/uL	0.00
4) Pyridine-d5	3.828	84	96147	8.388	ng/ul	0.00
7) Phenol-d5	7.181	99	101422	6.340	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.340	67	343984	33.513	ng/ul	0.00
11) 2-Chlorophenol-d4	7.545	132	285771	24.715	ng/ul	0.00
15) 4-Methylphenol-d8	8.734	113	199979	15.497	ng/ul	0.00
21) Nitrobenzene-d5	9.192	128	199755	31.976	ng/ul	0.00
24) 2-Nitrophenol-d4	9.916	143	225958	31.588	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.463	165	348184	26.727	ng/ul	0.00
31) 4-Chloroaniline-d4	10.981	131	474115	27.629	ng/ul	0.00
46) Dimethylphthalate-d6	14.074	166	1438496	34.539	ng/ul	0.00
49) Acenaphthylene-d8	14.363	160	1541633	33.360	ng/ul	0.00
54) 4-Nitrophenol-d4	14.922	143	26459m	4.205	ng/ul	0.04
60) Fluorene-d10	15.663	176	1192564	34.337	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.792	200	227392	28.951	ng/ul	0.00
73) Anthracene-d10	17.533	188	1850581	35.822	ng/ul	0.00
81) Pyrene-d10	19.762	212	1942313	35.816	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.874	264	1675109	37.167	ng/ul	0.00

Target Compounds Qvalue

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

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