

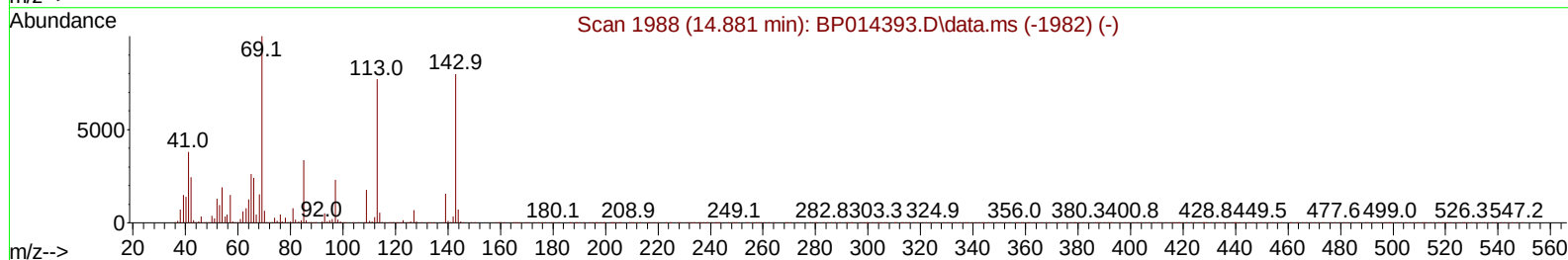
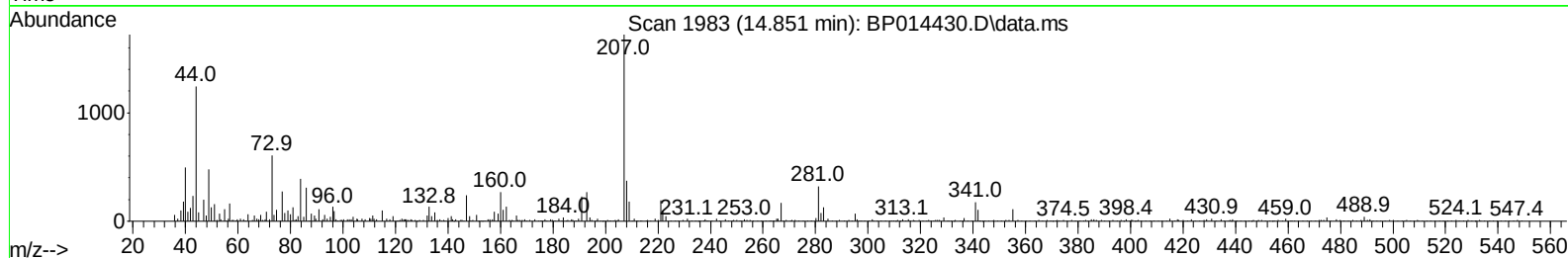
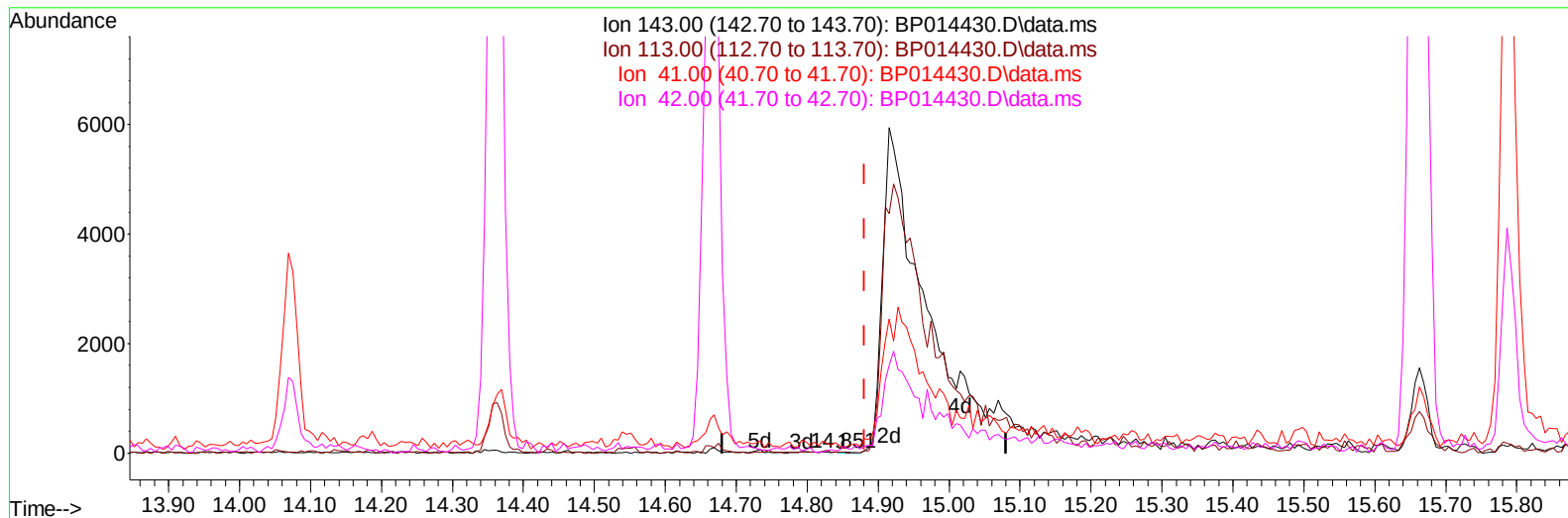
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
Data File : BP014430.D
Acq On : 31 Mar 2023 07:32
Operator : CG/JU
Sample : 02041-05
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB959

Manual IntegrationsAPPROVED

Quant Time: Mar 31 07:56:14 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Mar 31 04:41:29 2023
Response via : Initial Calibration

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



TIC: BP014430.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.851min (-0.030) 0.00 ng/u1

response 18

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	83.33
41.00	44.20	488.89#
42.00	31.50	672.22#

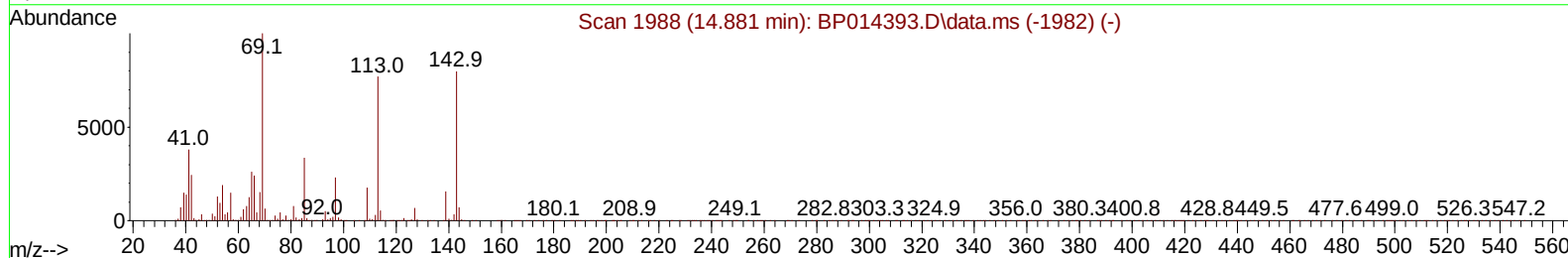
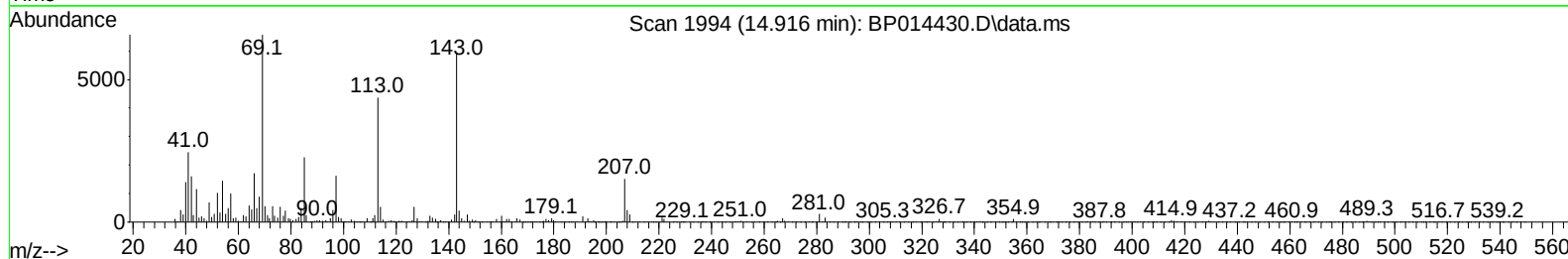
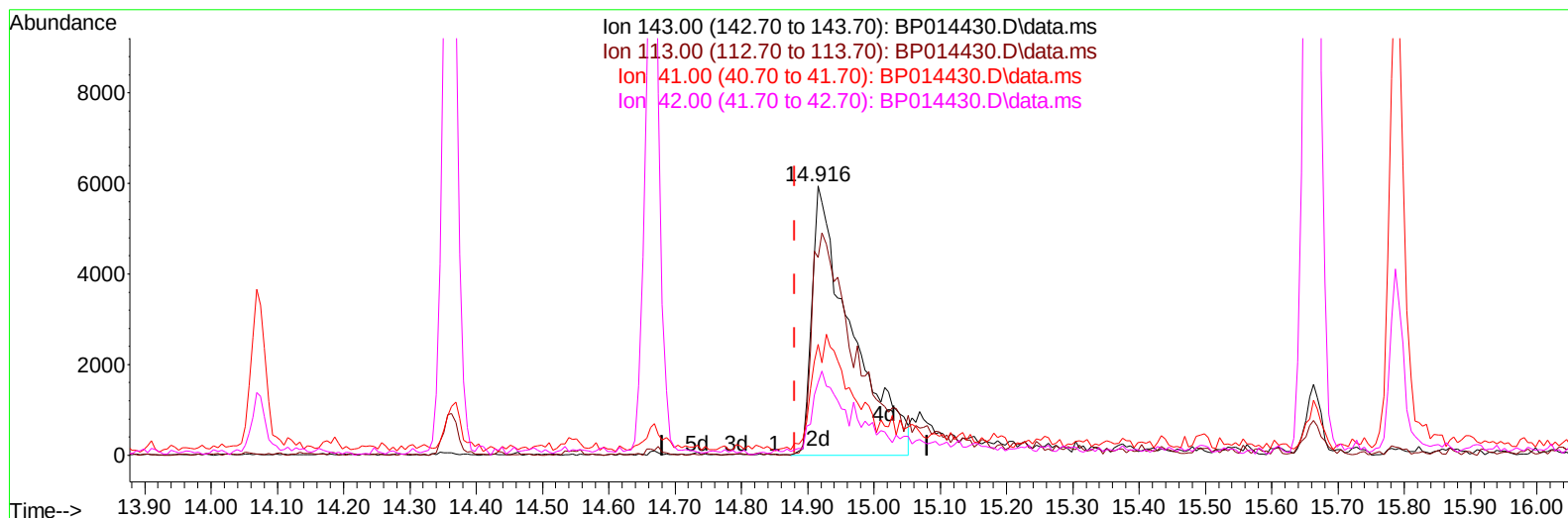
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TIC: BP014430.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.916min (+ 0.035) 4.00 ng/ul m

response 24298

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	73.56#
41.00	44.20	41.25
42.00	31.50	27.01

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.010	152	171315	20.000	ng/ul	0.00
20) Naphthalene-d8	10.834	136	760946	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.669	164	496523	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.428	188	1013459	20.000	ng/ul	0.00
79) Chrysene-d12	21.516	240	759704	20.000	ng/ul	0.00
88) Perylene-d12	24.033	264	792561	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	21381	4.840	ng/uL	0.00
4) Pyridine-d5	3.822	84	115838	10.232	ng/ul	0.00
7) Phenol-d5	7.175	99	100346	6.351	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.334	67	349438	34.471	ng/ul	0.00
11) 2-Chlorophenol-d4	7.540	132	285685	25.018	ng/ul	0.00
15) 4-Methylphenol-d8	8.728	113	202148	15.862	ng/ul	0.00
21) Nitrobenzene-d5	9.193	128	202574	32.976	ng/ul	0.00
24) 2-Nitrophenol-d4	9.916	143	225005	31.987	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.457	165	353982	27.632	ng/ul	0.00
31) 4-Chloroaniline-d4	10.981	131	456032	27.026	ng/ul	0.00
46) Dimethylphthalate-d6	14.075	166	1468793	36.484	ng/ul	0.00
49) Acenaphthylene-d8	14.363	160	1520857	34.045	ng/ul	0.00
54) 4-Nitrophenol-d4	14.916	143	24298m	3.995	ng/ul	0.04
60) Fluorene-d10	15.663	176	1213847	36.155	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.786	200	185407	24.863	ng/ul	0.00
73) Anthracene-d10	17.528	188	1824539	37.200	ng/ul	0.00
81) Pyrene-d10	19.757	212	1942067	38.058	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.874	264	1607341	38.172	ng/ul	0.00

Target Compounds Qvalue

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

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