

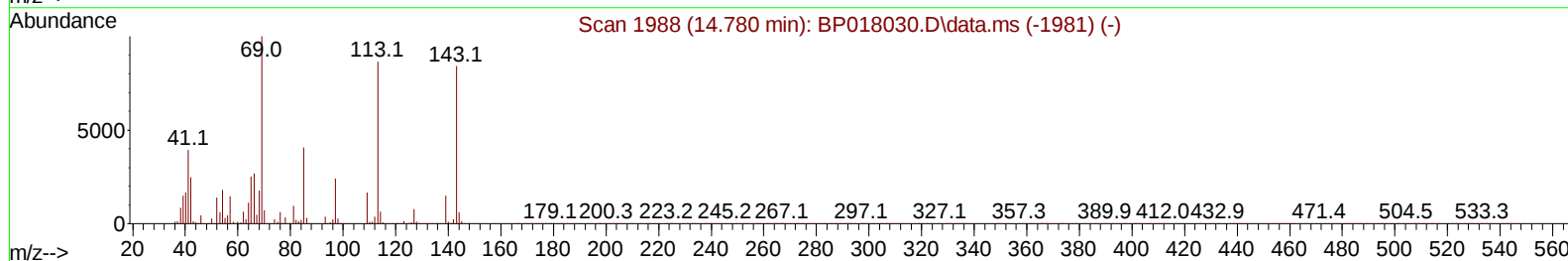
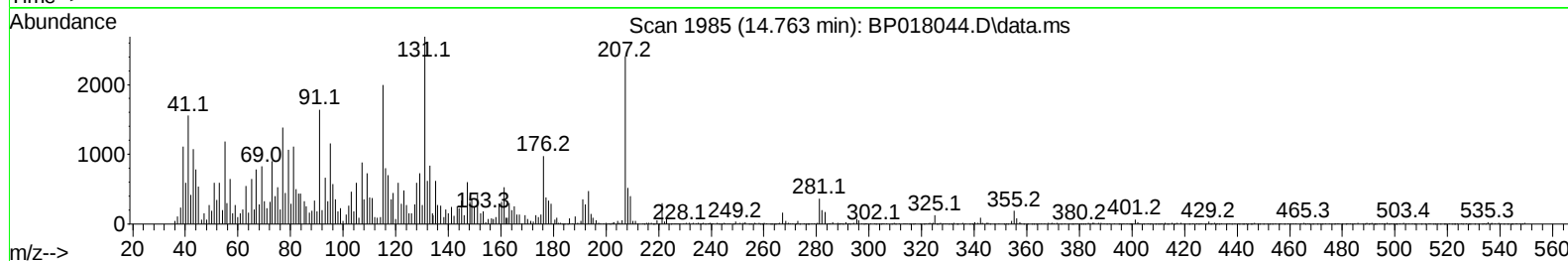
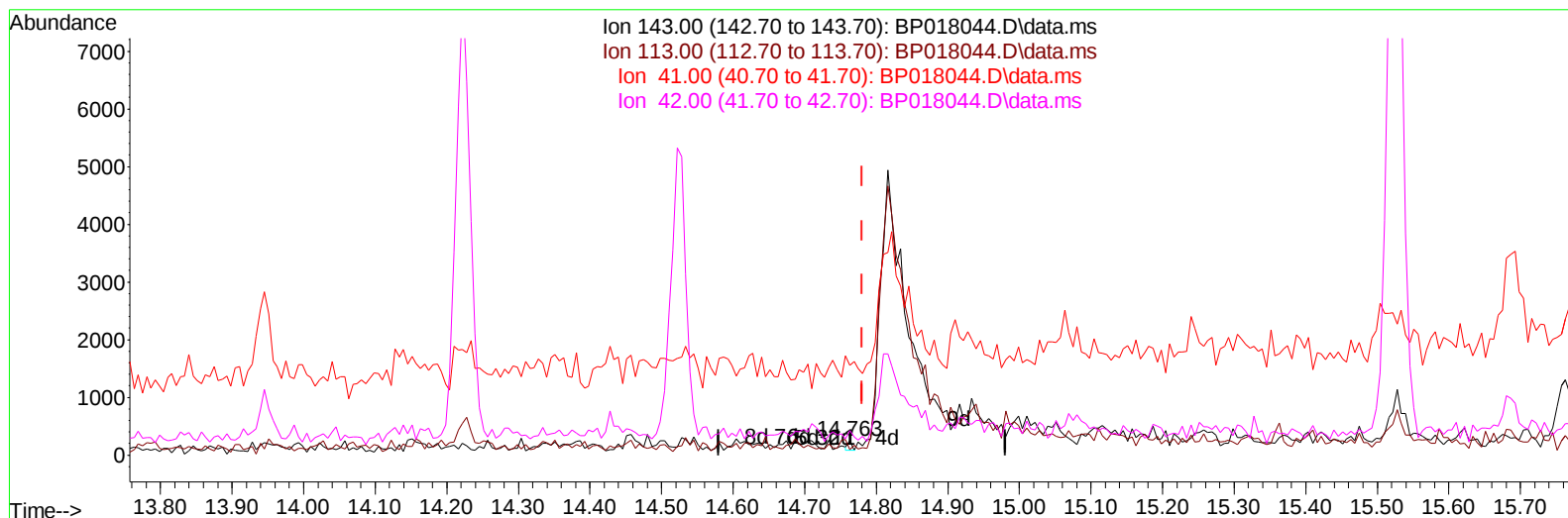
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
Data File : BP018044.D
Acq On : 11 Nov 2023 02:14
Operator : MA/JU
Sample : 05044-08
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHEE8

Manual IntegrationsAPPROVED

Quant Time: Nov 11 02:47:40 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Nov 10 21:55:18 2023
Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/11/2023
Supervised By :mohammad ahmed 11/15/2023



TIC: BP018044.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (-0.017) 0.02 ng/u1

response 58

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	37.30#
41.00	47.70	617.86#
42.00	30.10	166.27#

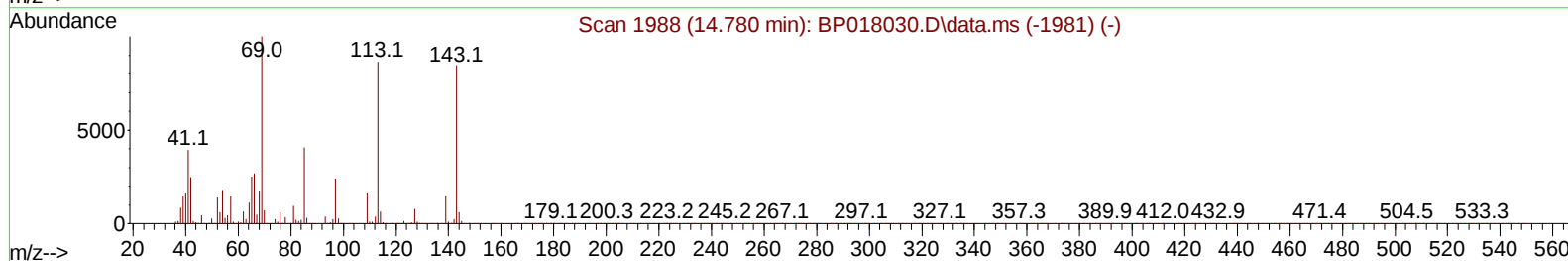
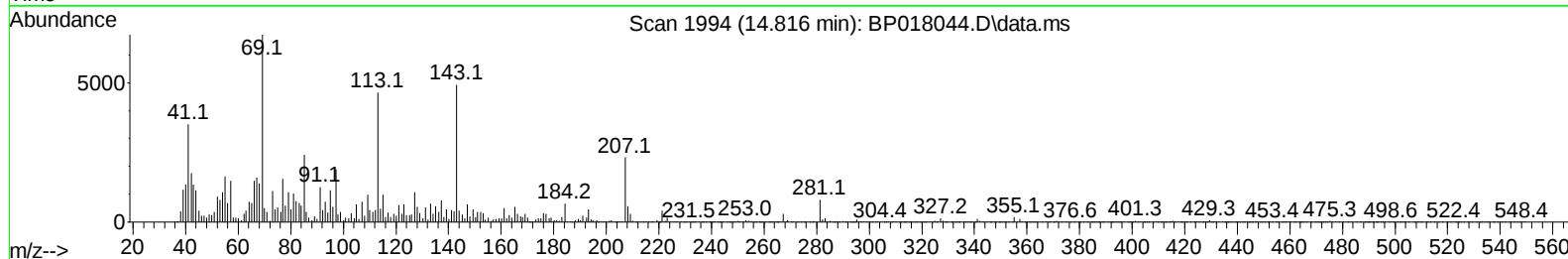
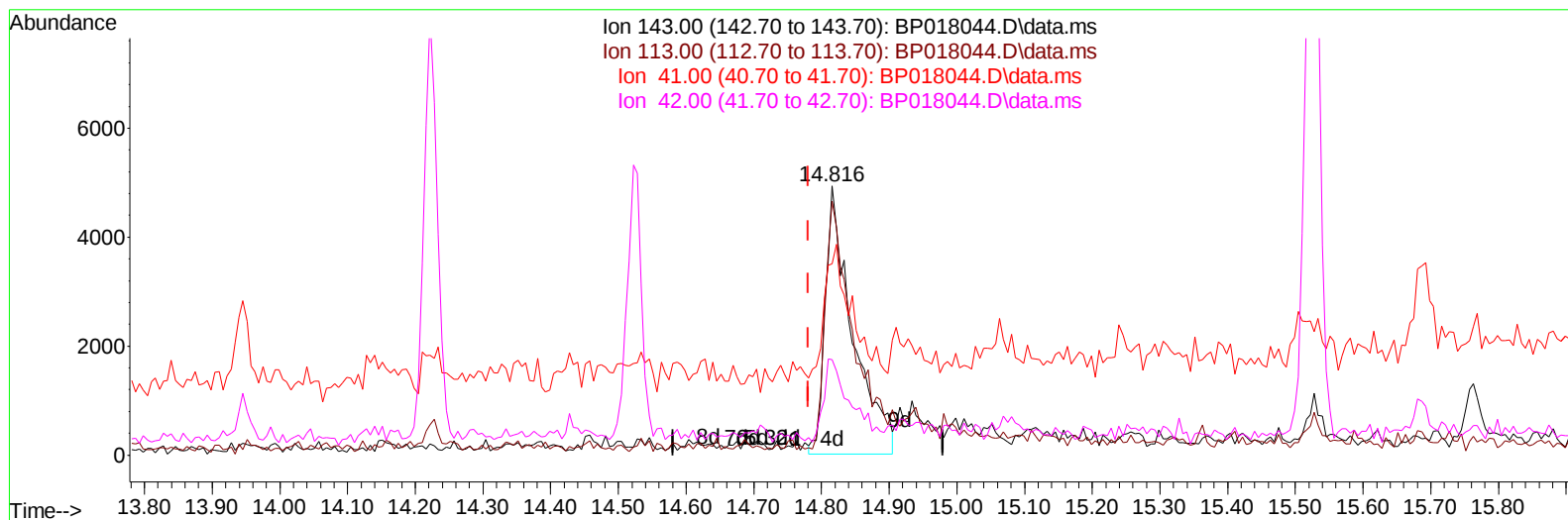
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TIC: BP018044.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.816min (+ 0.036) 5.72 ng/ul m

response 13795

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	94.45
41.00	47.70	71.19#
42.00	30.10	35.47

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 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHEE8

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/11/2023
 Supervised By :mohammad ahmed 11/15/2023

Quant Time: Nov 11 02:48:16 2023
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 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 10 21:55:18 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.852	152	62027	20.000	ng/ul	0.00
20) Naphthalene-d8	10.669	136	265811	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.522	164	175031	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.304	188	398082	20.000	ng/ul	0.00
79) Chrysene-d12	21.427	240	376725	20.000	ng/ul	0.00
88) Perylene-d12	23.933	264	415388	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	7628	4.338	ng/uL	0.00
4) Pyridine-d5	3.670	84	63841	13.925	ng/ul	0.00
7) Phenol-d5	7.022	99	50668	8.461	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.199	67	129740	36.867	ng/ul	0.00
11) 2-Chlorophenol-d4	7.375	132	139839	29.489	ng/ul	0.00
15) 4-Methylphenol-d8	8.575	113	101039	20.591	ng/ul	0.00
21) Nitrobenzene-d5	9.057	128	86095	35.951	ng/ul	0.00
24) 2-Nitrophenol-d4	9.769	143	97131	35.801	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.299	165	173563	35.893	ng/ul	0.00
31) 4-Chloroaniline-d4	10.852	131	215900	32.558	ng/ul	0.00
46) Dimethylphthalate-d6	13.945	166	623429	38.603	ng/ul	0.00
49) Acenaphthylene-d8	14.222	160	669293	38.050	ng/ul	0.00
54) 4-Nitrophenol-d4	14.816	143	13795m	5.716	ng/ul	0.04
60) Fluorene-d10	15.528	176	526312	39.473	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.687	200	22797	7.922	ng/ul	0.00
73) Anthracene-d10	17.404	188	864753	40.165	ng/ul	0.00
81) Pyrene-d10	19.657	212	1045383	41.543	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.768	264	999509	41.381	ng/ul	0.00

Target Compounds Qvalue

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

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