

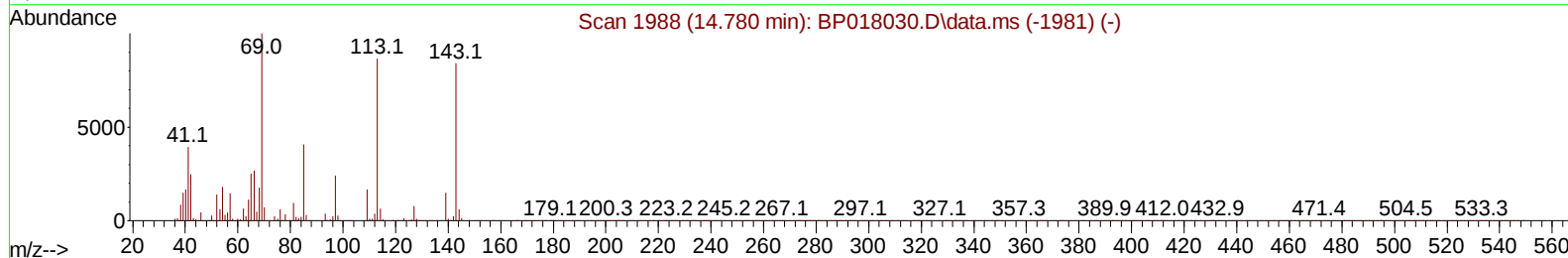
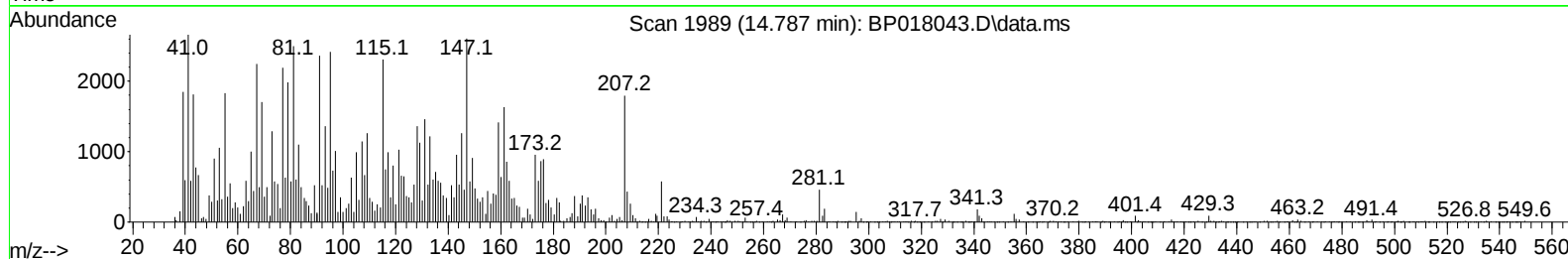
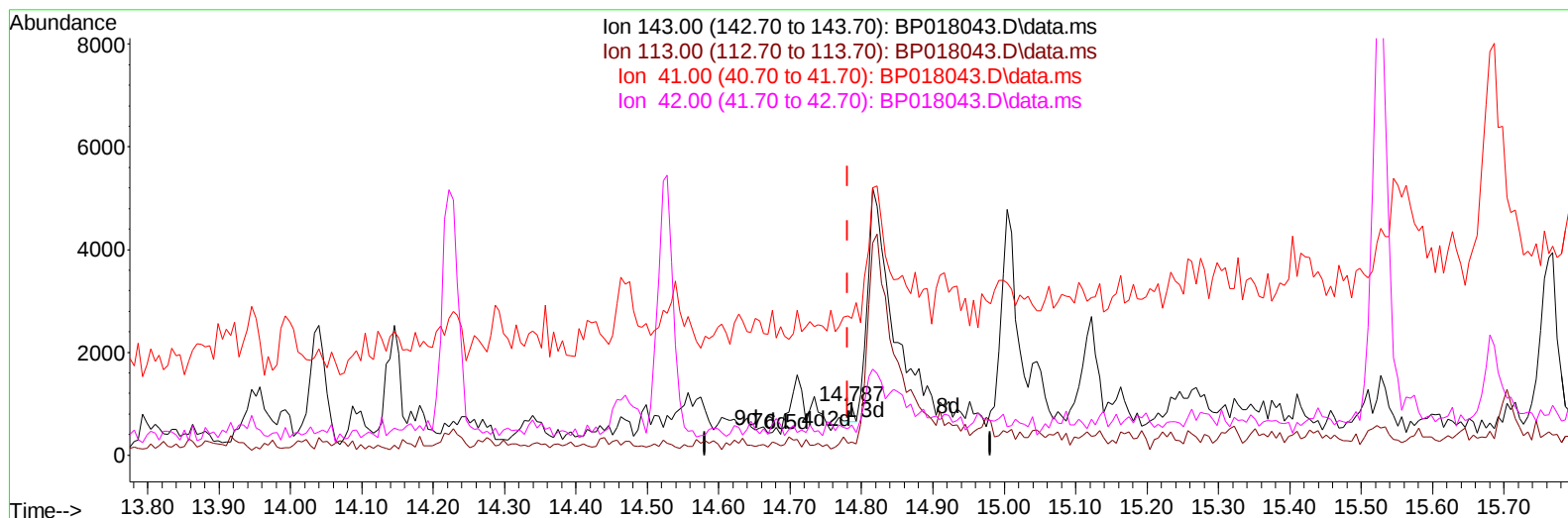
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
Data File : BP018043.D
Acq On : 11 Nov 2023 01:40
Operator : MA/JU
Sample : 05044-07
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHEE5

Manual IntegrationsAPPROVED

Quant Time: Nov 11 02:27:24 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: BP018043.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.787min (+ 0.006) 0.05 ng/u1

response 120

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	26.75#
41.00	47.70	278.06#
42.00	30.10	61.55#

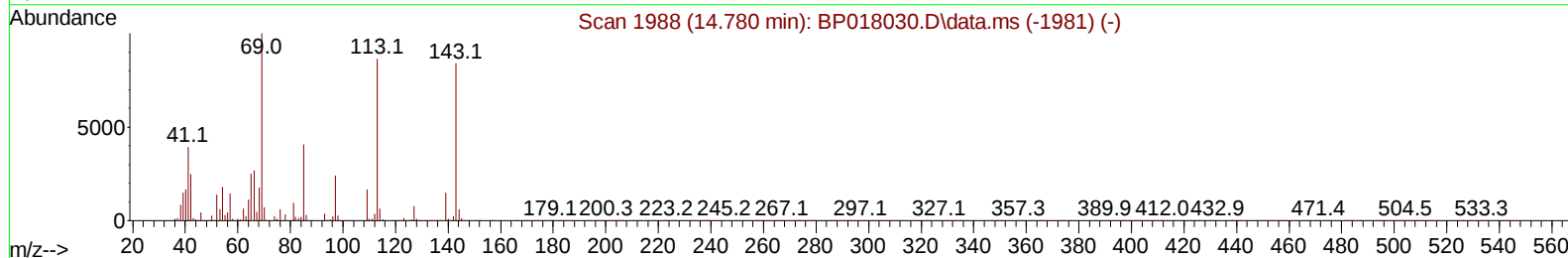
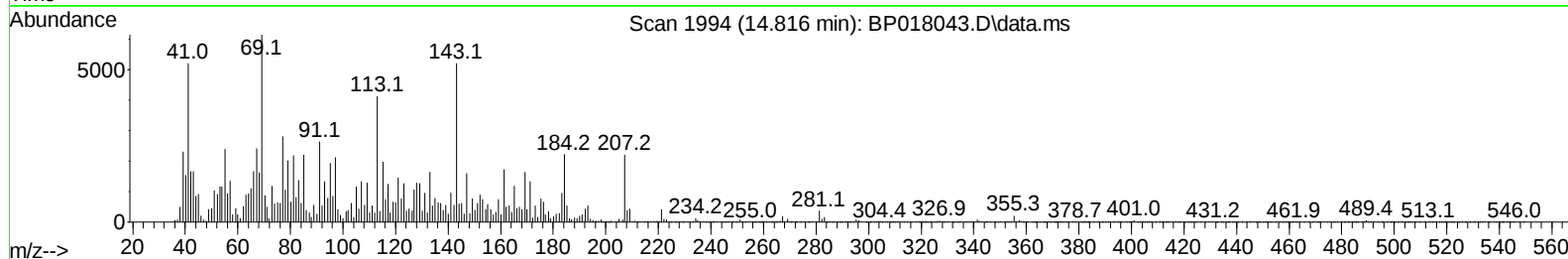
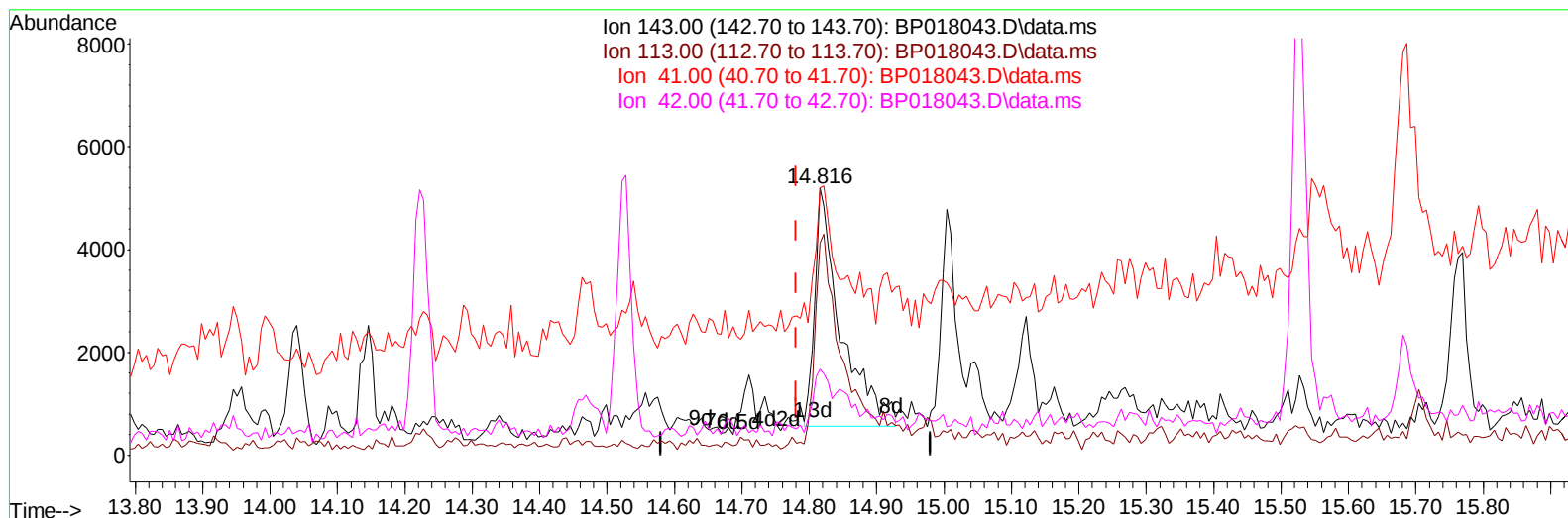
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TIC: BP018043.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.816min (+ 0.036) 5.01 ng/ul m

response 12475

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	79.21#
41.00	47.70	99.83#
42.00	30.10	32.02

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 Operator : MA/JU
 Sample : 05044-07
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHEE5

Manual IntegrationsAPPROVED

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

Quant Time: Nov 11 02:28:22 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

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.852	152	65648	20.000	ng/ul	0.00
20) Naphthalene-d8	10.669	136	272258	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.528	164	180703	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.304	188	412638	20.000	ng/ul	0.00
79) Chrysene-d12	21.433	240	410014	20.000	ng/ul	0.00
88) Perylene-d12	23.939	264	477678	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.228	96	6790	3.649	ng/uL	0.00
4) Pyridine-d5	3.670	84	54593	11.251	ng/ul	0.00
7) Phenol-d5	7.022	99	35311	5.571	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.199	67	96980	26.038	ng/ul	0.00
11) 2-Chlorophenol-d4	7.375	132	103302	20.582	ng/ul	0.00
15) 4-Methylphenol-d8	8.575	113	70245	13.526	ng/ul	0.00
21) Nitrobenzene-d5	9.057	128	63602	25.929	ng/ul	0.00
24) 2-Nitrophenol-d4	9.769	143	69988	25.185	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.299	165	119945	24.217	ng/ul	0.00
31) 4-Chloroaniline-d4	10.851	131	139339	20.515	ng/ul	0.00
46) Dimethylphthalate-d6	13.945	166	434137	26.038	ng/ul	0.00
49) Acenaphthylene-d8	14.228	160	476439	26.236	ng/ul	0.00
54) 4-Nitrophenol-d4	14.816	143	12475m	5.007	ng/ul	0.04
60) Fluorene-d10	15.528	176	381681	27.727	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.681	200	58947	19.761	ng/ul	0.00
73) Anthracene-d10	17.404	188	640916	28.719	ng/ul	0.00
81) Pyrene-d10	19.657	212	796082	29.068	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.768	264	801205	28.846	ng/ul	0.00

Target Compounds Qvalue

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

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