

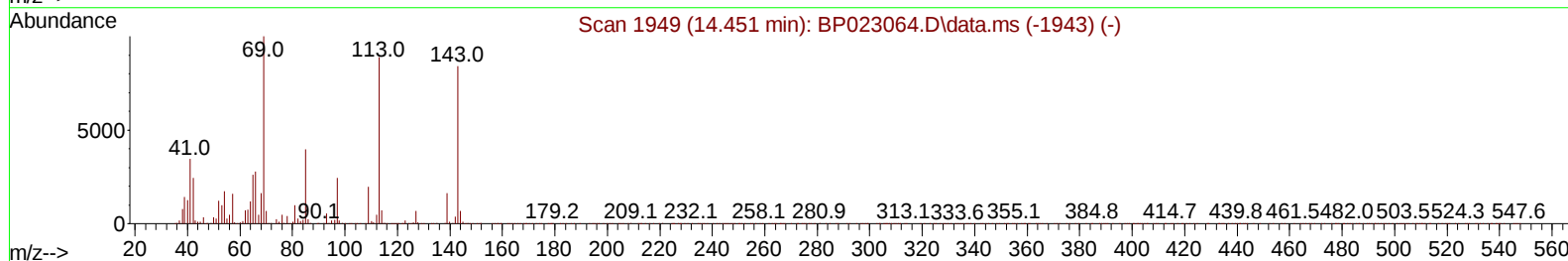
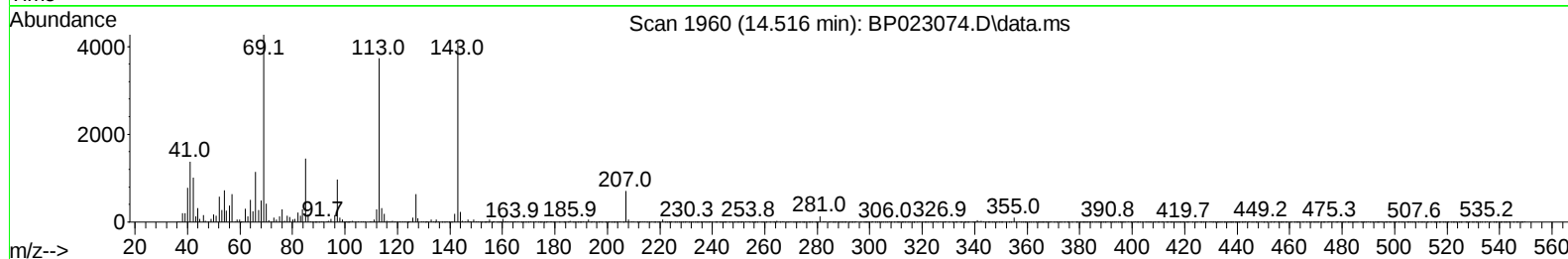
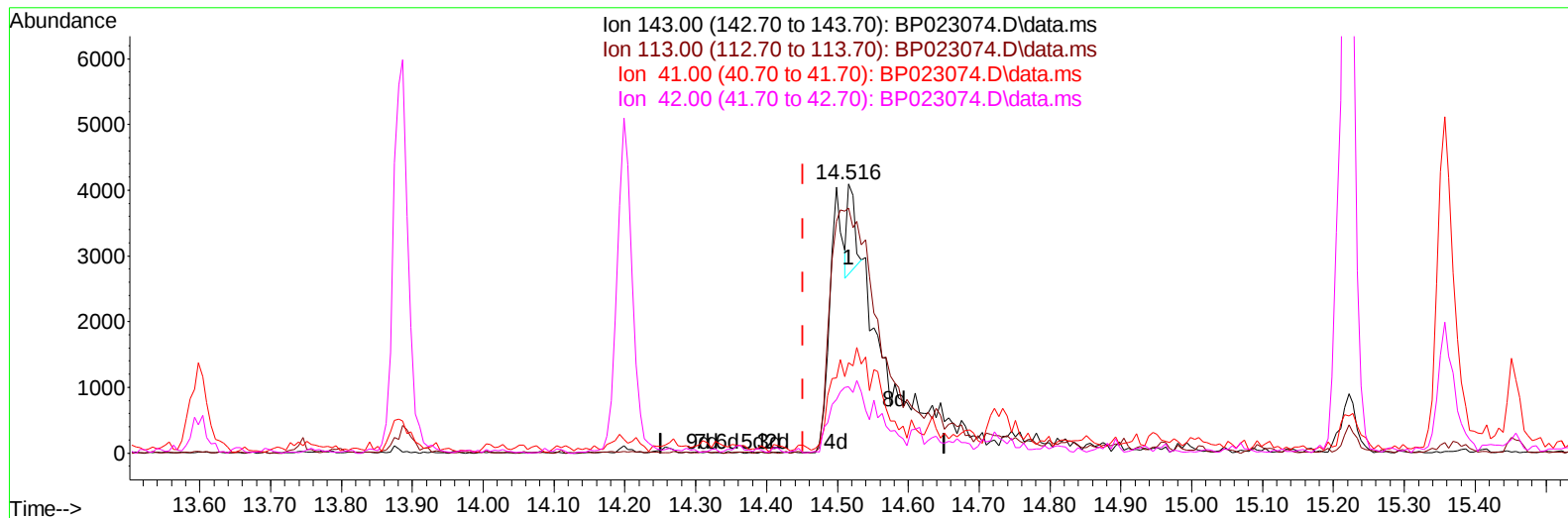
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111224\
 Data File : BP023074.D
 Acq On : 13 Nov 2024 10:07
 Operator : RC/JU
 Sample : P4687-12
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0CP3

Manual IntegrationsAPPROVED

Quant Time: Nov 14 00:41:10 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 13 04:49:01 2024
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/14/2024
 Supervised By :mohammad ahmed 11/14/2024



TIC: BP023074.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.516min (+ 0.065) 0.38 ng/ul

response 988

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	91.20
41.00	39.20	33.68
42.00	25.90	24.66

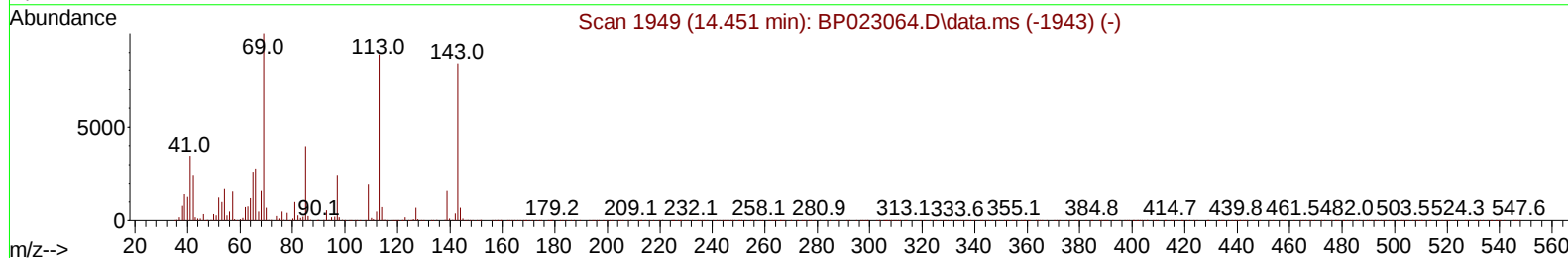
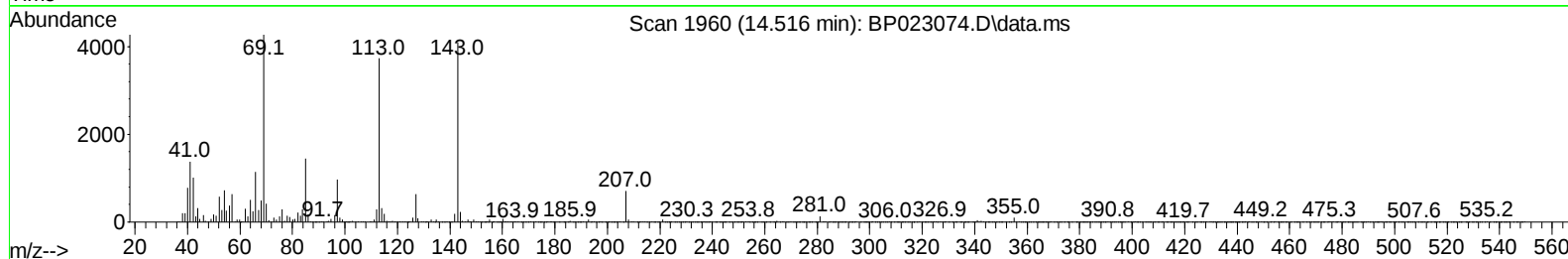
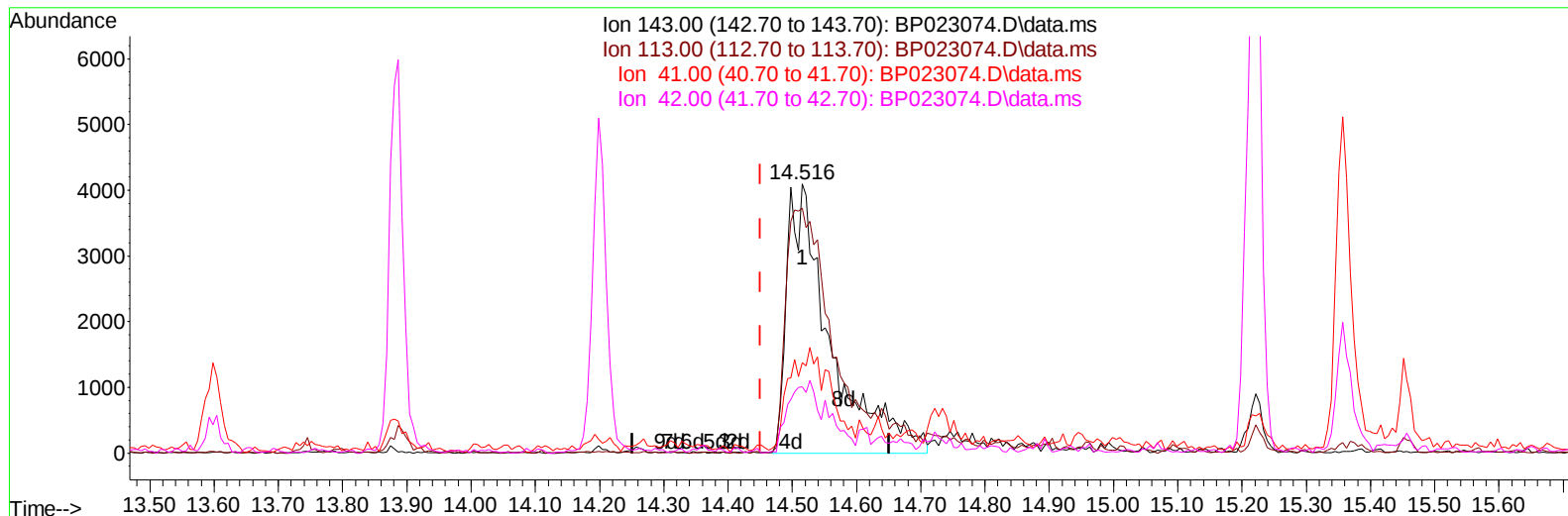
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ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Manual IntegrationsAPPROVED

Quant Time: Nov 14 00:41:10 2024
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TIC: BP023074.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.516min (+ 0.065) 7.51 ng/ul m

response 19327

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	91.20
41.00	39.20	33.68
42.00	25.90	24.66

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111224\
 Data File : BP023074.D
 Acq On : 13 Nov 2024 10:07
 Operator : RC/JU
 Sample : P4687-12
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0CP3

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/14/2024
 Supervised By :mohammad ahmed 11/14/2024

Quant Time: Nov 14 01:01:41 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 13 04:49:01 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.587	152	73665	20.000	ng/ul	0.00
20) Naphthalene-d8	10.334	136	286556	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.198	164	237552	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.010	188	606810	20.000	ng/ul	0.00
79) Chrysene-d12	21.433	240	848394	20.000	ng/ul	-0.01
88) Perylene-d12	24.651	264	1017330	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.140	96	7100	4.633	ng/uL	0.00
4) Pyridine-d5	3.564	84	41190	9.046	ng/ul	0.02
7) Phenol-d5	6.805	99	43419	7.972	ng/ul	0.02
9) Bis-(2-Chloroethyl)eth...	6.940	67	114991	35.937	ng/ul	0.01
11) 2-Chlorophenol-d4	7.140	132	123417	30.942	ng/ul	0.01
15) 4-Methylphenol-d8	8.299	113	95772	21.369	ng/ul	0.00
21) Nitrobenzene-d5	8.734	128	71823	36.003	ng/ul	0.01
24) 2-Nitrophenol-d4	9.440	143	88007	37.004	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.975	165	179510	36.042	ng/ul	0.00
31) 4-Chloroaniline-d4	10.481	131	190342	32.034	ng/ul	0.00
46) Dimethylphthalate-d6	13.598	166	685108	40.015	ng/ul	-0.01
49) Acenaphthylene-d8	13.887	160	645814	36.150	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.516	143	19327m	7.512	ng/ul	0.06
60) Fluorene-d10	15.222	176	591503	39.405	ng/ul	0.01
65) 4,6-Dinitro-2-methylph...	15.357	200	123275	33.912	ng/ul	0.00
73) Anthracene-d10	17.104	188	1092218	42.510	ng/ul	-0.01
81) Pyrene-d10	19.480	212	1420244	36.344	ng/ul	-0.01
92) Benzo(a)pyrene-d12	24.421	264	2146248	44.290	ng/ul	-0.02

Target Compounds Qvalue

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

