

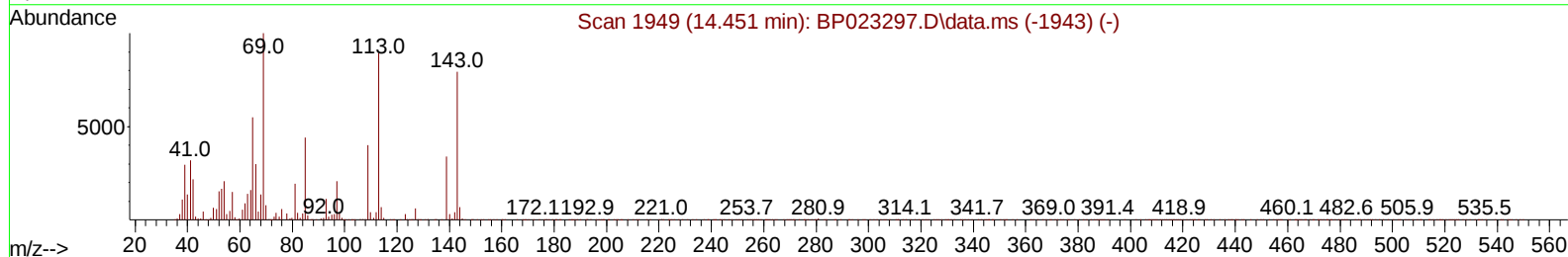
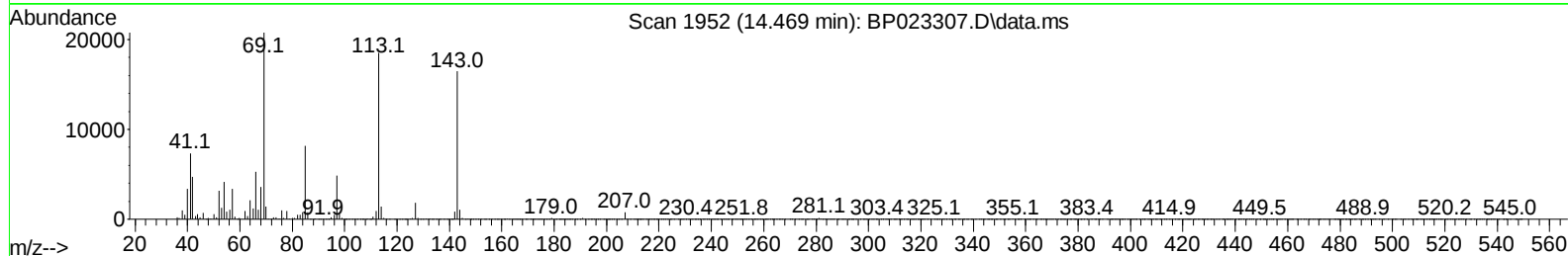
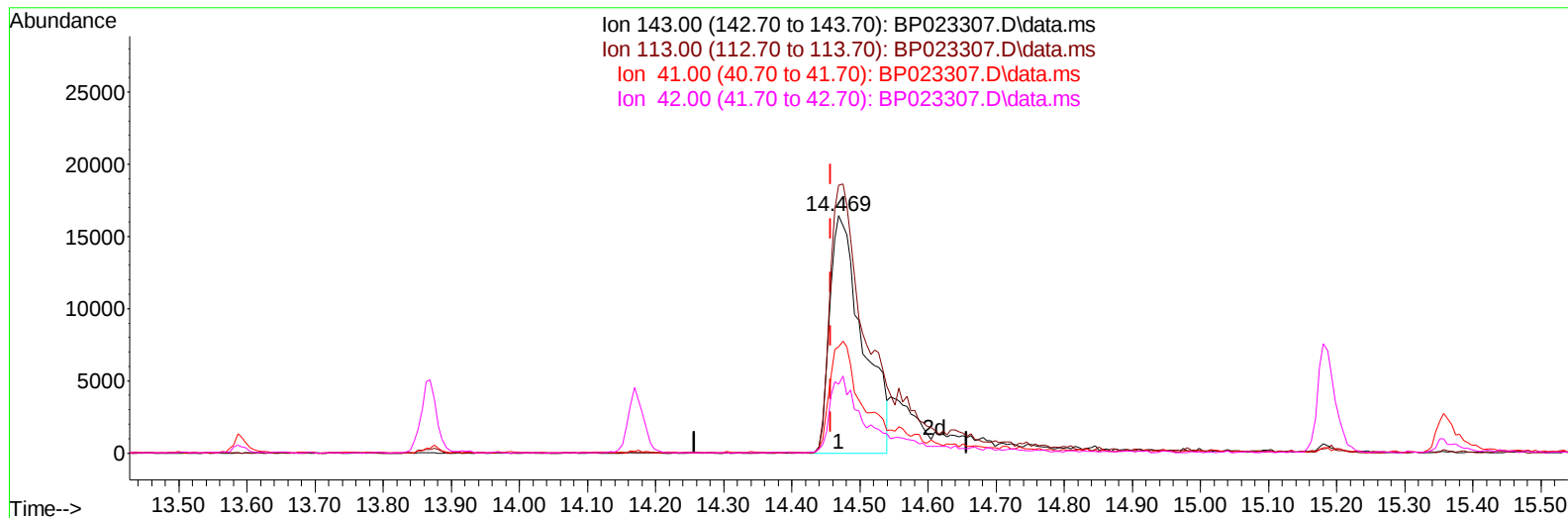
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112224\
 Data File : BP023307.D
 Acq On : 22 Nov 2024 17:24
 Operator : RC/JU
 Sample : PB165141BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK141

Manual IntegrationsAPPROVED

Quant Time: Nov 22 23:49:40 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 20 04:34:35 2024
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 11/25/2024
 Supervised By :mohammad ahmed 11/27/2024



TIC: BP023307.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.469min (+ 0.012) 24.47 ng/u1

response 54723

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	112.69
41.00	39.20	44.64
42.00	25.90	28.92

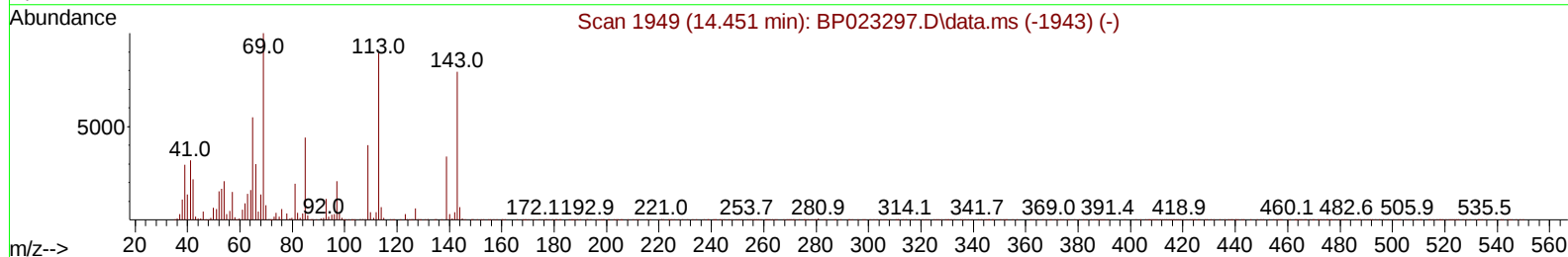
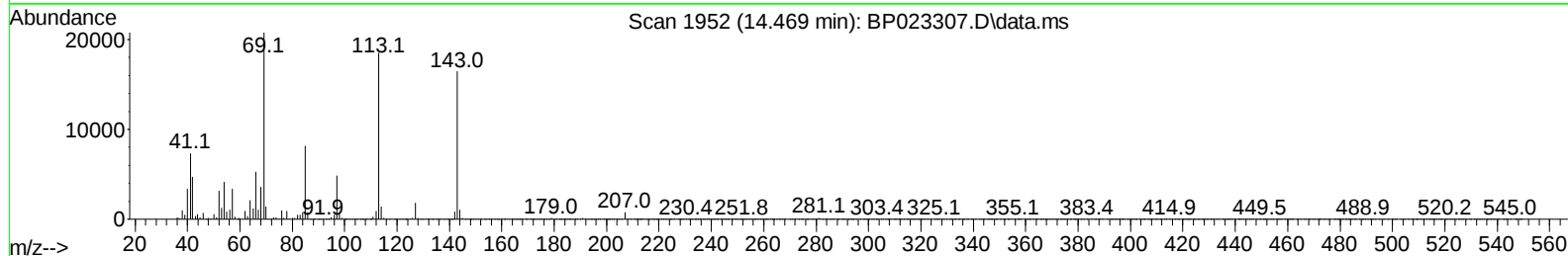
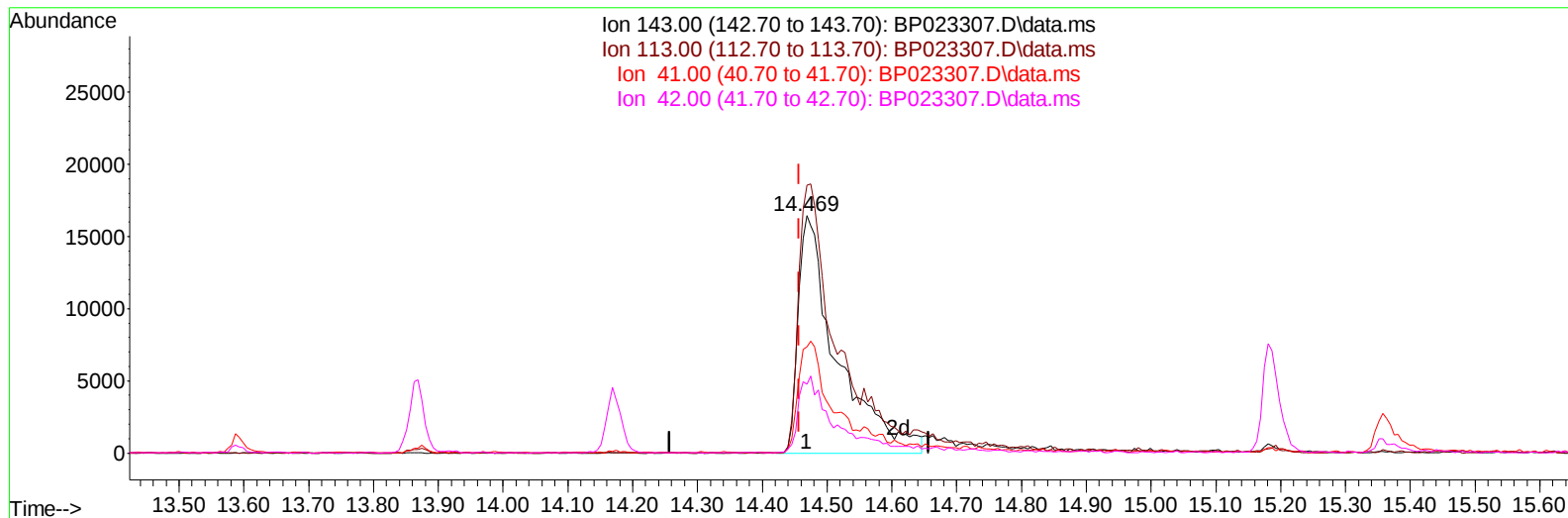
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TIC: BP023307.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.469min (+ 0.012) 30.53 ng/ul m

response 68281

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	112.69
41.00	39.20	44.64
42.00	25.90	28.92

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Manual IntegrationsAPPROVED

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Quant Time: Nov 22 23:51:30 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.563	152	73986	20.000	ng/ul	0.00
20) Naphthalene-d8	10.304	136	277339	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.169	164	206507	20.000	ng/ul	-0.02
64) Phenanthrene-d10	16.986	188	497291	20.000	ng/ul	-0.01
79) Chrysene-d12	21.445	240	534355	20.000	ng/ul	0.02
88) Perylene-d12	24.668	264	640956	20.000	ng/ul	0.04
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.134	96	12821	8.330	ng/uL	0.00
4) Pyridine-d5	3.546	84	160403	35.076	ng/ul	0.00
7) Phenol-d5	6.769	99	191349	34.982	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.910	67	126646	39.408	ng/ul	0.00
11) 2-Chlorophenol-d4	7.110	132	143538	35.830	ng/ul	0.00
15) 4-Methylphenol-d8	8.275	113	155532	34.553	ng/ul	0.00
21) Nitrobenzene-d5	8.704	128	71039	36.793	ng/ul	0.00
24) 2-Nitrophenol-d4	9.416	143	82951	36.037	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.969	165	160401	33.275	ng/ul	0.00
31) 4-Chloroaniline-d4	10.463	131	183551	31.918	ng/ul	0.00
46) Dimethylphthalate-d6	13.592	166	592482	39.808	ng/ul	0.00
49) Acenaphthylene-d8	13.869	160	589387	37.952	ng/ul	0.00
54) 4-Nitrophenol-d4	14.469	143	68281m	30.527	ng/ul	0.01
60) Fluorene-d10	15.186	176	500562	38.360	ng/ul	-0.02
65) 4,6-Dinitro-2-methylph...	15.357	200	85245	28.615	ng/ul	0.00
73) Anthracene-d10	17.098	188	833546	39.588	ng/ul	0.00
81) Pyrene-d10	19.480	212	1067206	43.360	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.433	264	1196408	39.186	ng/ul	0.01

Target Compounds Qvalue

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

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