

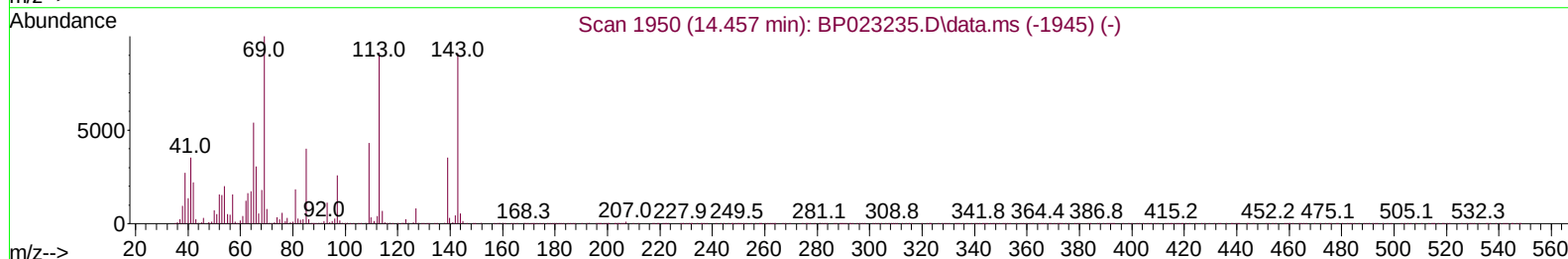
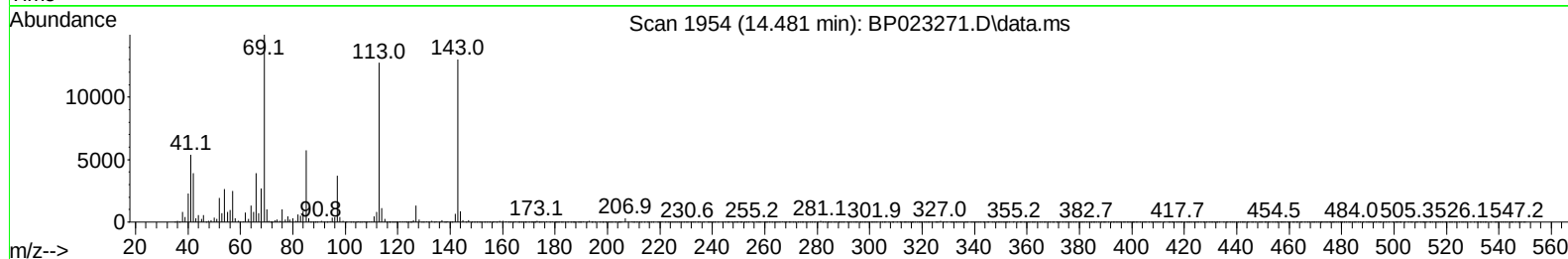
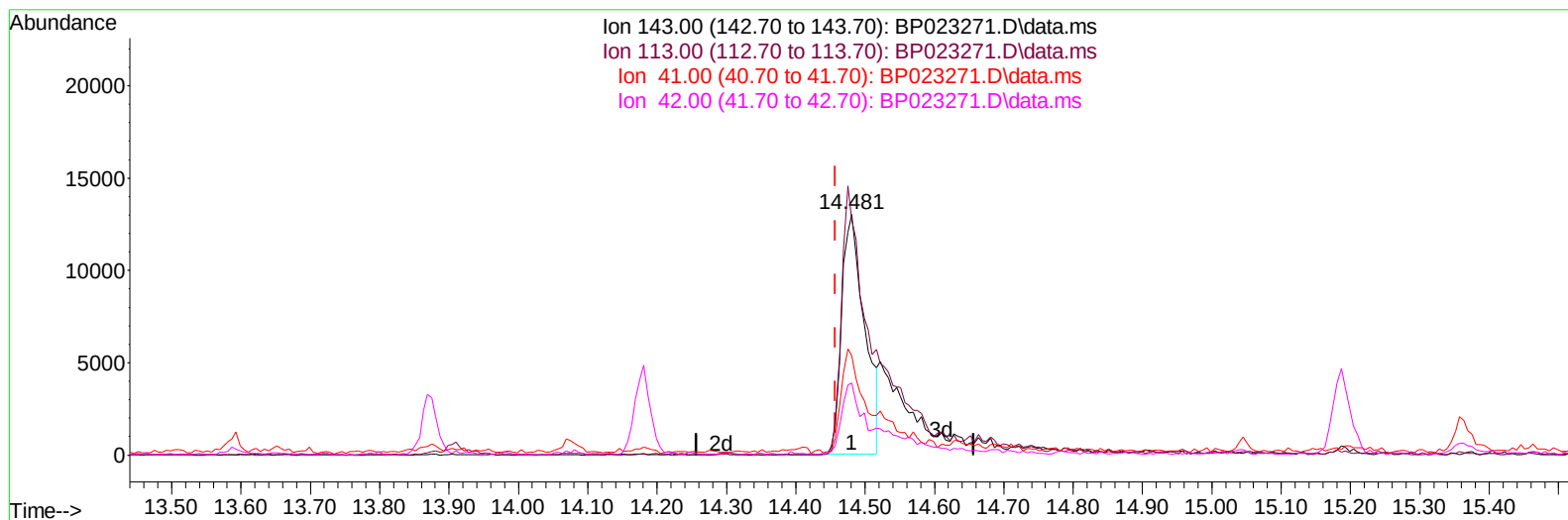
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111924\
Data File : BP023271.D
Acq On : 21 Nov 2024 05:13
Operator : RC/JU
Sample : P4872-10
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29X8

Manual IntegrationsAPPROVED

Quant Time: Nov 21 05:39:23 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 :Yogesh Patel 11/22/2024
Supervised By :mohammad ahmed 11/22/2024



TIC: BP023271.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.481min (+ 0.024) 10.52 ng/ul

response 29653

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	97.72
41.00	39.20	41.30
42.00	25.90	29.90

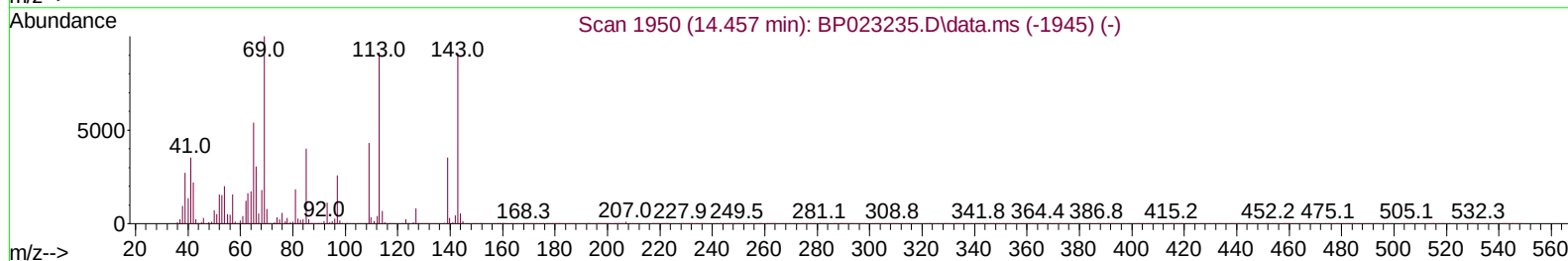
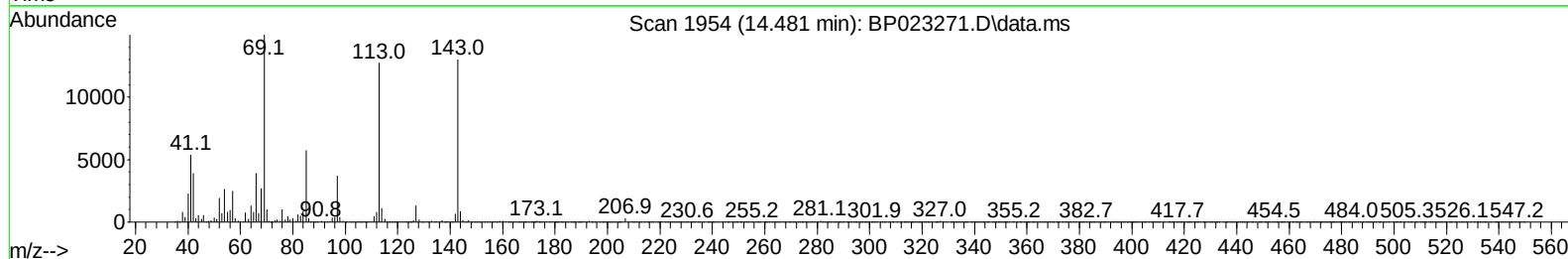
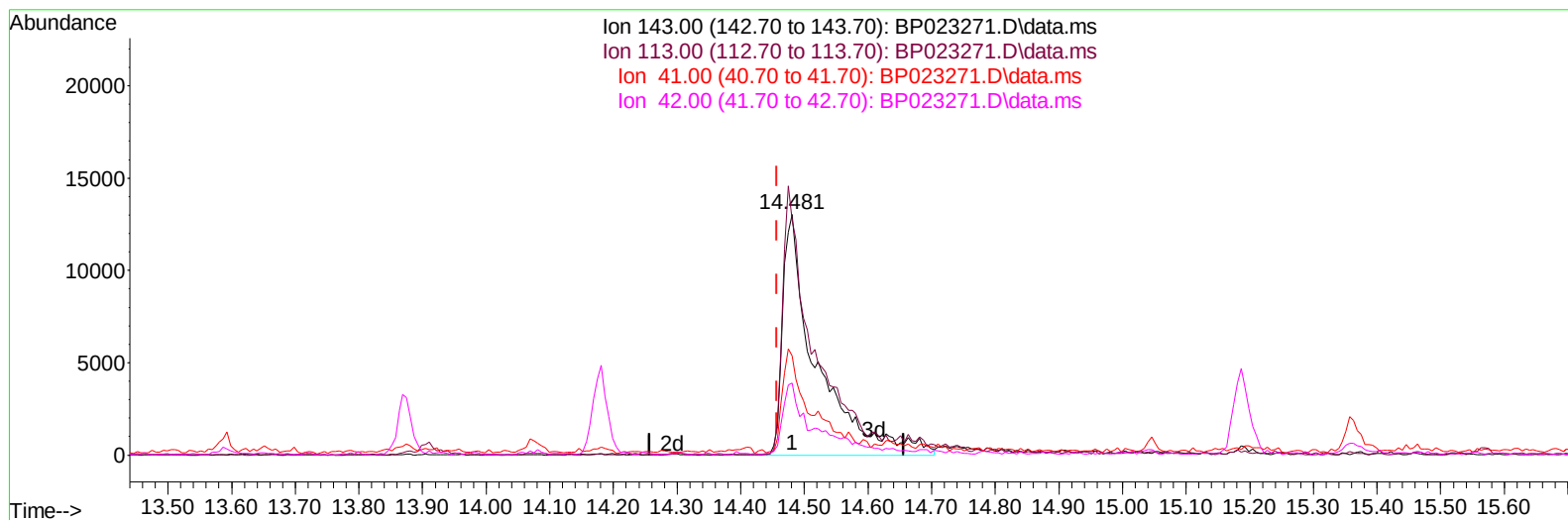
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TIC: BP023271.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.481min (+ 0.024) 17.11 ng/ul m

response 48238

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	97.72
41.00	39.20	41.30
42.00	25.90	29.90

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Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/22/2024
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Quant Time: Nov 21 05:45:23 2024
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 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 20 04:34:35 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.558	152	94717	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.310	136	346189	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.181	164	260243	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.987	188	629670	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.433	240	749311	20.000	ng/ul	0.01
88) Perylene-d12	24.645	264	912883	20.000	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.134	96	6984	3.544	ng/uL	0.00
4) Pyridine-d5	3.546	84	81322	13.891	ng/ul	0.00
7) Phenol-d5	6.769	99	130352	18.615	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.911	67	76490	18.592	ng/ul	0.00
11) 2-Chlorophenol-d4	7.111	132	99360	19.374	ng/ul	0.00
15) 4-Methylphenol-d8	8.269	113	105216	18.258	ng/ul	-0.01
21) Nitrobenzene-d5	8.705	128	49191	20.411	ng/ul	0.00
24) 2-Nitrophenol-d4	9.416	143	56634	19.711	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.969	165	114290	18.994	ng/ul	0.00
31) 4-Chloroaniline-d4	10.463	131	124641	17.363	ng/ul	0.00
46) Dimethylphthalate-d6	13.587	166	394555	21.036	ng/ul	0.00
49) Acenaphthylene-d8	13.875	160	413808	21.144	ng/ul	0.00
54) 4-Nitrophenol-d4	14.481	143	48238m	17.113	ng/ul	0.02
60) Fluorene-d10	15.187	176	354261	21.543	ng/ul	-0.02
65) 4,6-Dinitro-2-methylph...	15.363	200	48176	12.772	ng/ul	0.01
73) Anthracene-d10	17.098	188	593410	22.258	ng/ul	0.00
81) Pyrene-d10	19.475	212	771261	22.346	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.439	264	994979	22.881	ng/ul	0.02

Target Compounds Qvalue

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

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