

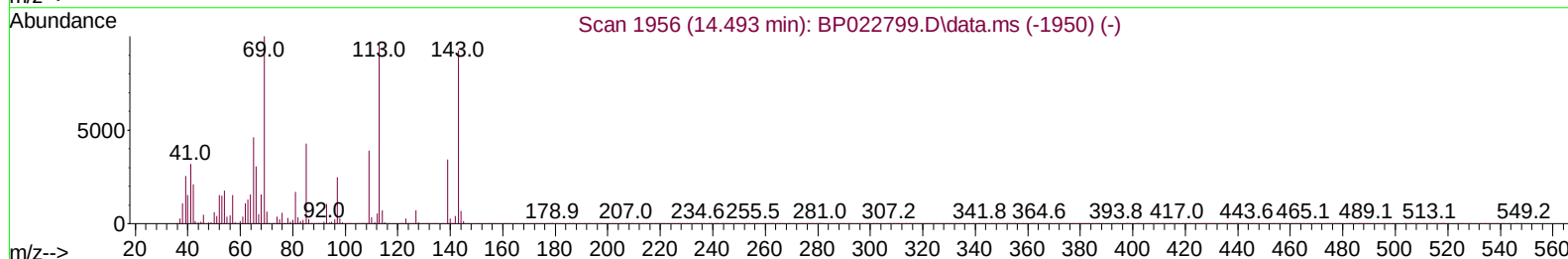
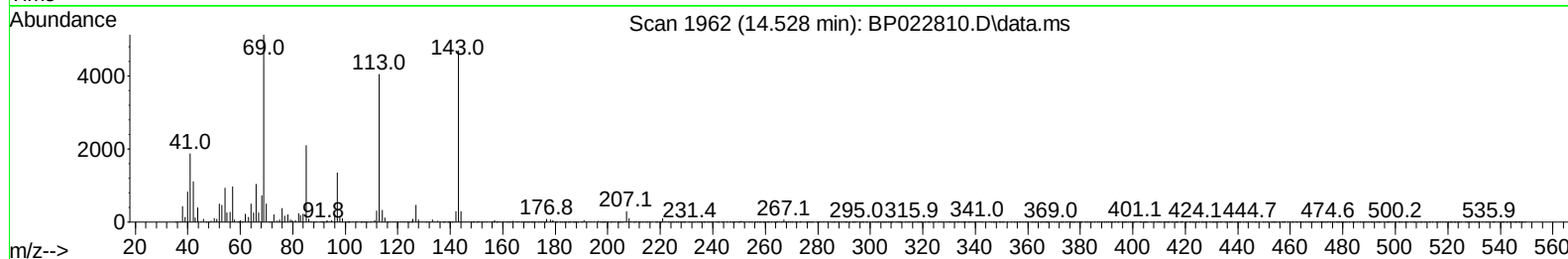
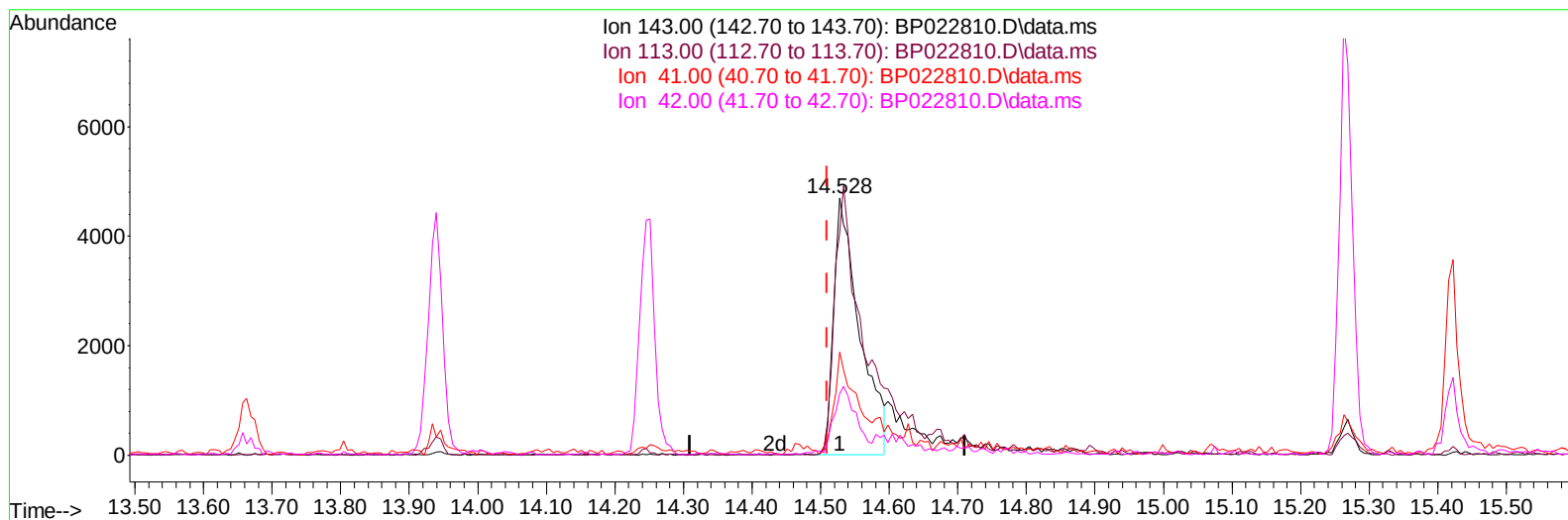
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103024\
Data File : BP022810.D
Acq On : 01 Nov 2024 18:50
Operator : RC/JU
Sample : P4627-01
Misc :
ALS Vial : 46 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0CN6

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/04/2024
Supervised By :mohammad ahmed 11/05/2024

Quant Time: Nov 04 04:36:37 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Oct 30 10:56:39 2024
Response via : Initial Calibration



TIC: BP022810.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.528min (+ 0.018) 4.84 ng/u1

response 12298

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	86.44
41.00	35.90	40.03
42.00	23.70	23.89

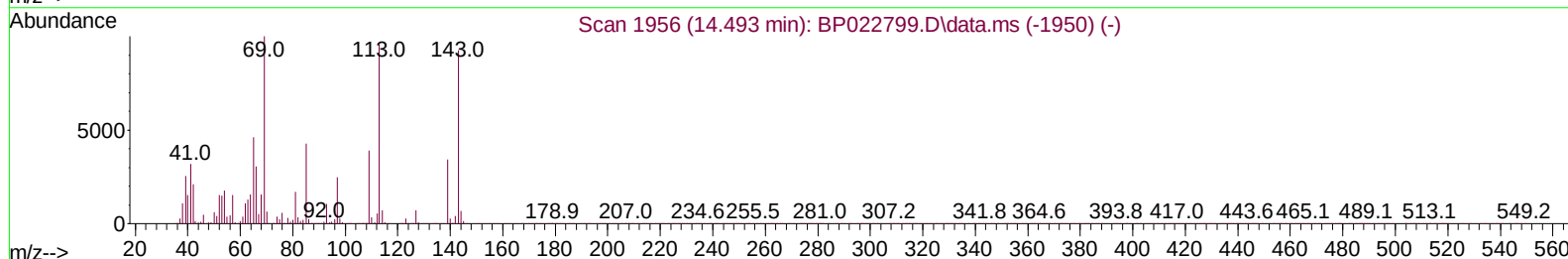
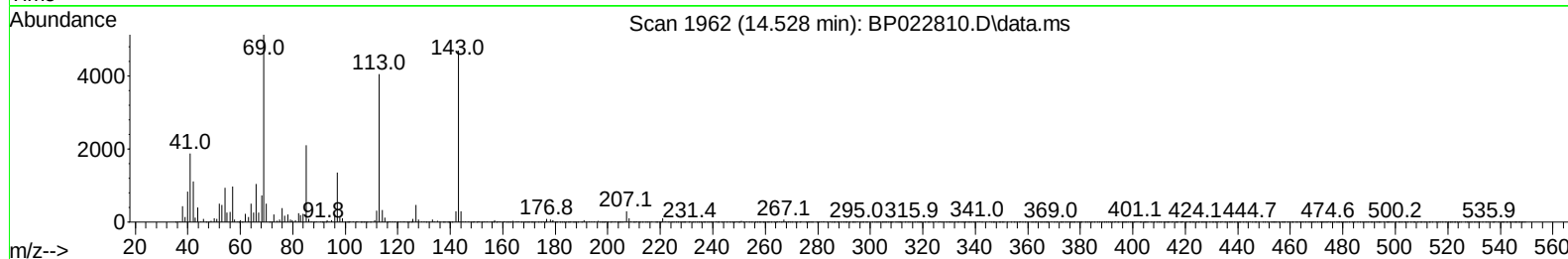
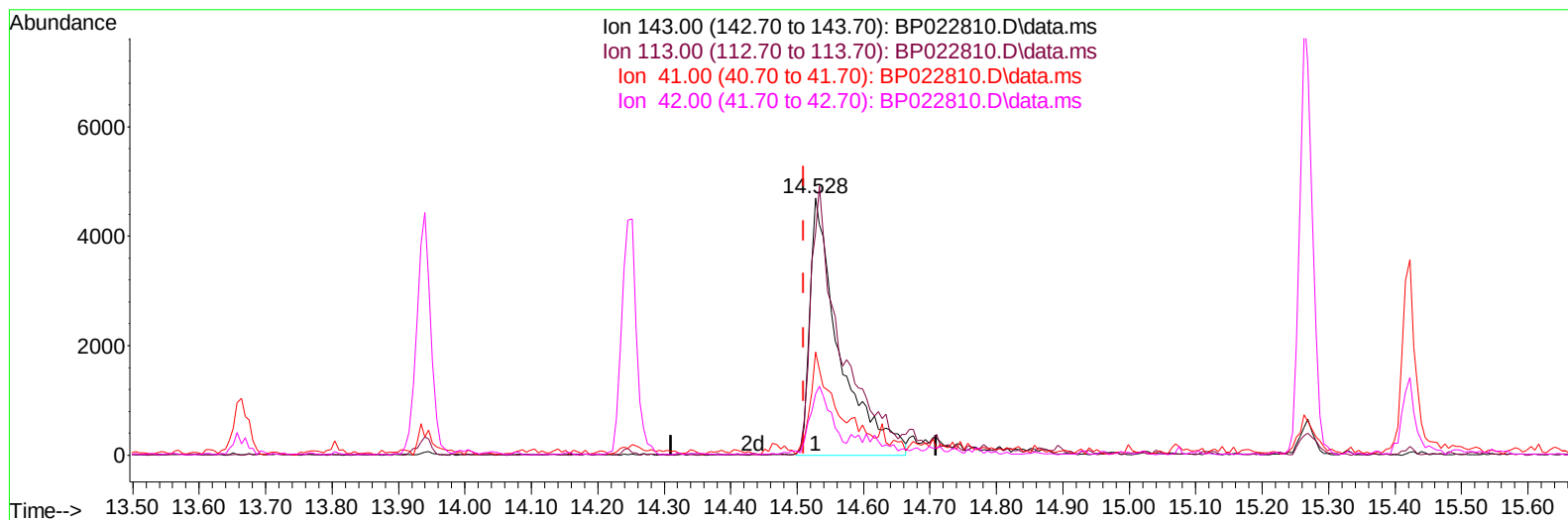
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TIC: BP022810.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.528min (+ 0.018) 5.73 ng/ul m

response 14552

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	86.44
41.00	35.90	40.03
42.00	23.70	23.89

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.628	152	67506	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.381	136	254049	20.000	ng/ul	#-0.01
38) Acenaphthene-d10	14.245	164	190616	20.000	ng/ul	-0.02
64) Phenanthrene-d10	17.063	188	461515	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.504	240	528164	20.000	ng/ul	0.00
88) Perylene-d12	24.768	264	640105	20.000	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.175	96	7716	4.442	ng/uL	0.00
4) Pyridine-d5	3.593	84	46241	9.696	ng/ul	0.00
7) Phenol-d5	6.828	99	44985	7.916	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.969	67	93777	28.078	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.169	132	108286	26.856	ng/ul	-0.01
15) 4-Methylphenol-d8	8.334	113	83762	18.471	ng/ul	-0.01
21) Nitrobenzene-d5	8.769	128	60021	30.727	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.487	143	72768	32.177	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.022	165	144301	30.022	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.528	131	146622	25.816	ng/ul	-0.01
46) Dimethylphthalate-d6	13.663	166	478106	31.569	ng/ul	0.00
49) Acenaphthylene-d8	13.940	160	500038	31.086	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.528	143	14552m	5.727	ng/ul	0.02
60) Fluorene-d10	15.269	176	424805	32.339	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.422	200	72794	23.982	ng/ul	0.02
73) Anthracene-d10	17.169	188	750779	34.581	ng/ul	0.00
81) Pyrene-d10	19.551	212	971028	34.766	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.539	264	1182723	34.598	ng/ul	0.02

Target Compounds Qvalue

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

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