

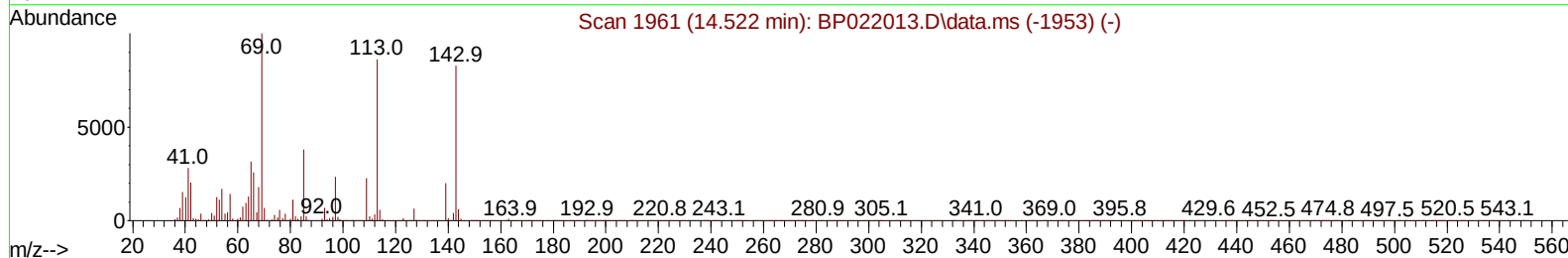
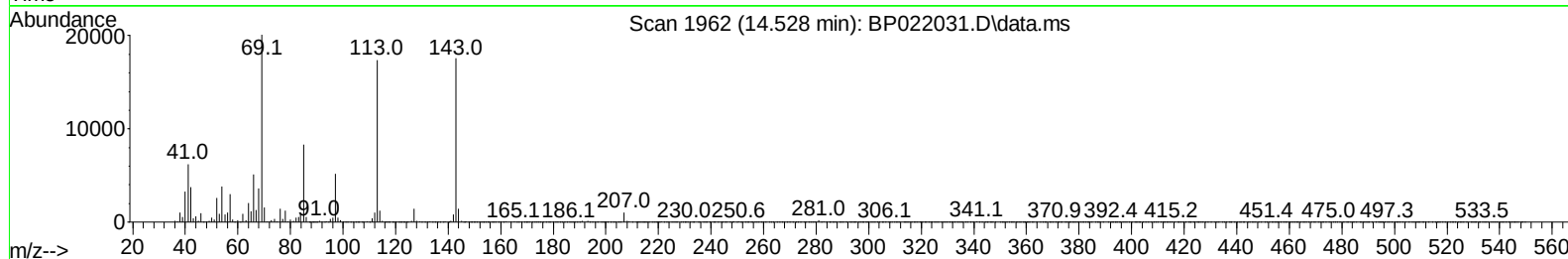
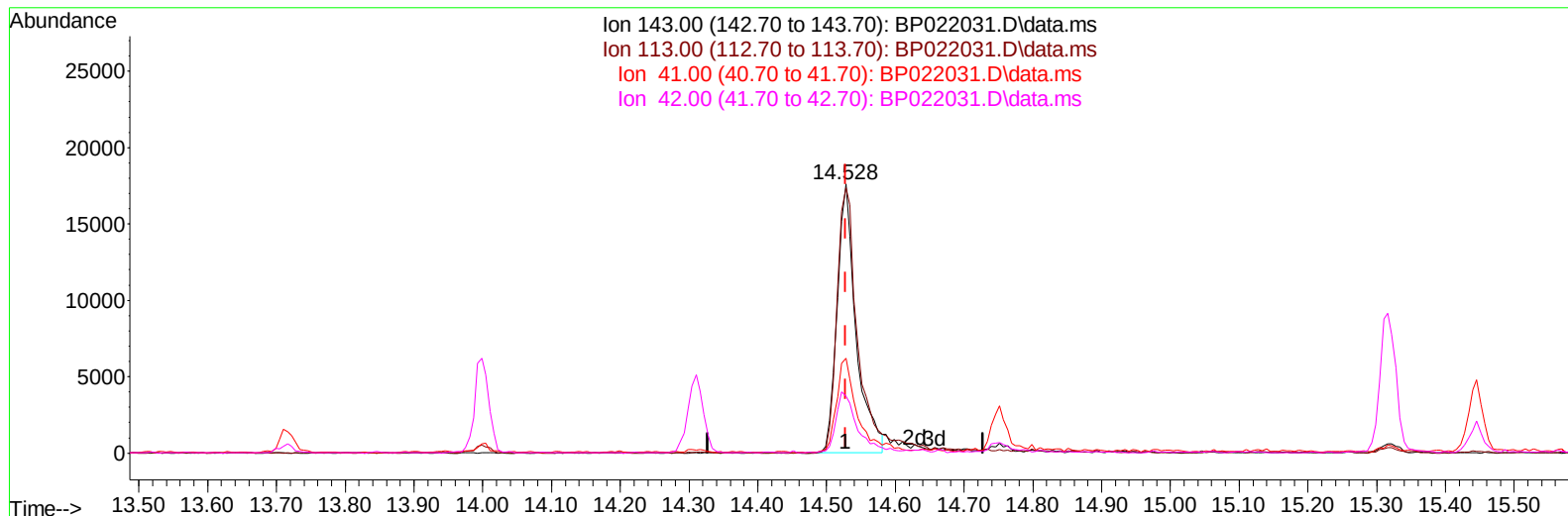
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092424\
Data File : BP022031.D
Acq On : 25 Sep 2024 16:45
Operator : RC/JU
Sample : P4107-06
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0B43

Manual IntegrationsAPPROVED

Quant Time: Sep 25 18:17:48 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Sep 24 15:23:13 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 09/26/2024
Supervised By :mohammad ahmed 09/28/2024



TIC: BP022031.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.528min (+ 0.000) 9.38 ng/ul

response 33605

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	106.70	98.77
41.00	35.80	35.23
42.00	26.20	21.16

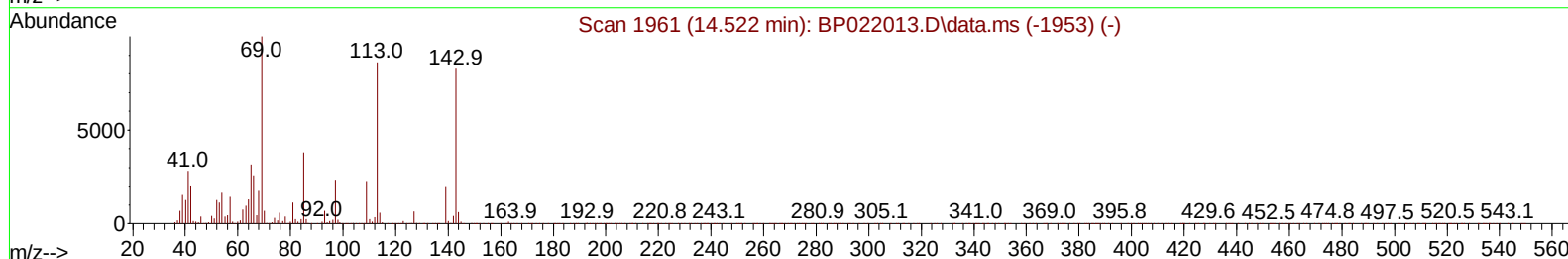
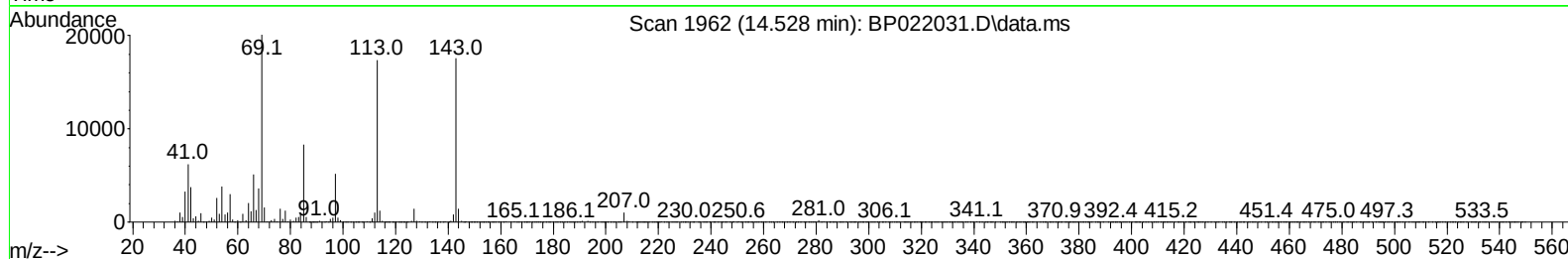
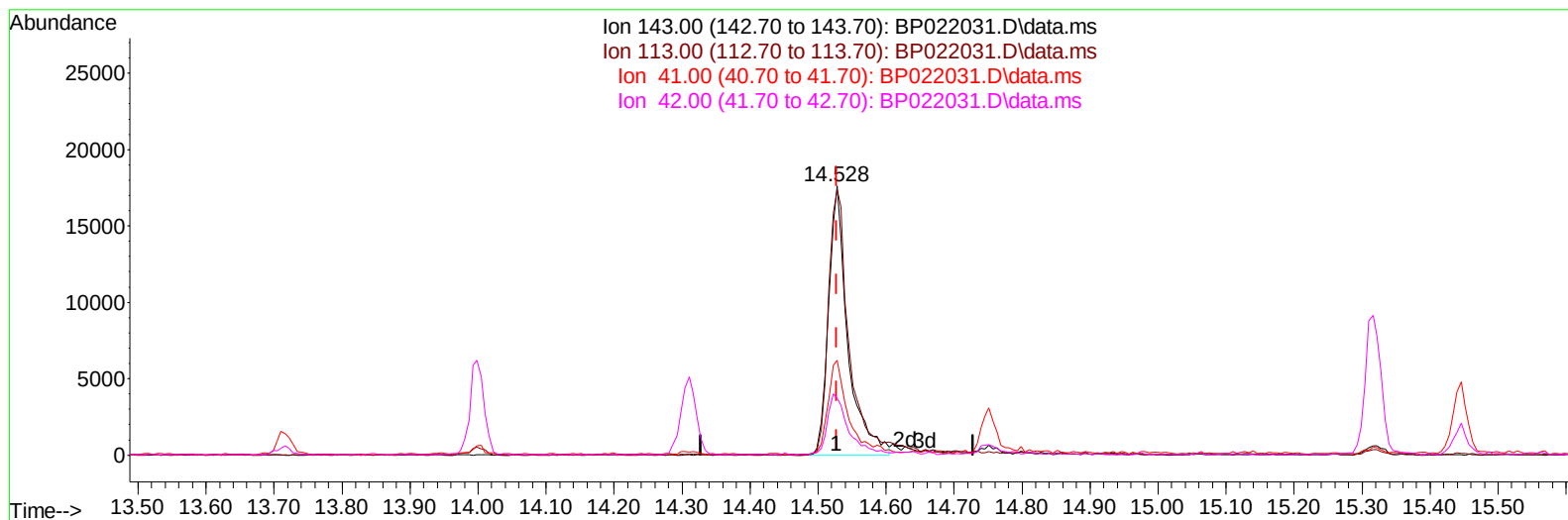
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TIC: BP022031.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.528min (+ 0.000) 9.73 ng/ul m

response 34852

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	106.70	98.77
41.00	35.80	35.23
42.00	26.20	21.16

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 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 09/26/2024
 Supervised By :mohammad ahmed 09/28/2024

Quant Time: Sep 26 02:45:41 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 24 15:23:13 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.705	152	96455	20.000	ng/ul	0.00
20) Naphthalene-d8	10.463	136	352635	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.310	164	270977	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.116	188	643668	20.000	ng/ul	0.00
79) Chrysene-d12	21.533	240	748335	20.000	ng/ul	0.00
88) Perylene-d12	24.816	264	914289	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.223	96	12876	5.229	ng/uL	0.00
4) Pyridine-d5	3.640	84	47511	6.750	ng/ul	0.00
7) Phenol-d5	6.893	99	96397	12.217	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.046	67	153746	31.645	ng/ul	0.00
11) 2-Chlorophenol-d4	7.246	132	179593	31.460	ng/ul	0.00
15) 4-Methylphenol-d8	8.405	113	154954	24.784	ng/ul	0.00
21) Nitrobenzene-d5	8.840	128	90347	34.231	ng/ul	0.00
24) 2-Nitrophenol-d4	9.558	143	105949	35.404	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.093	165	216240	33.273	ng/ul	0.00
31) 4-Chloroaniline-d4	10.599	131	178736	23.050	ng/ul	0.00
46) Dimethylphthalate-d6	13.716	166	745477	35.589	ng/ul	0.00
49) Acenaphthylene-d8	13.998	160	726829	32.773	ng/ul	0.00
54) 4-Nitrophenol-d4	14.528	143	34852m	9.727	ng/ul	0.00
60) Fluorene-d10	15.316	176	624586	33.709	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.445	200	118214	29.582	ng/ul	0.00
73) Anthracene-d10	17.222	188	1098216	36.548	ng/ul	0.00
81) Pyrene-d10	19.592	212	1447754	37.342	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.598	264	1750146	36.394	ng/ul	-0.01

Target Compounds Qvalue

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

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