

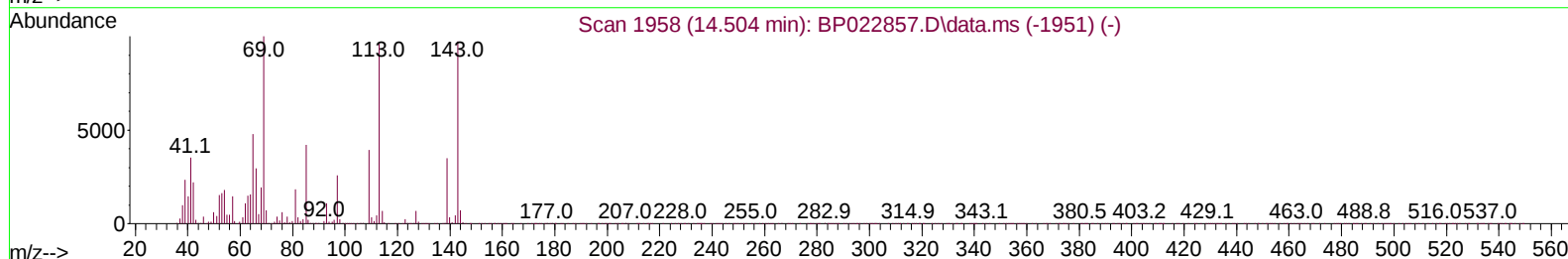
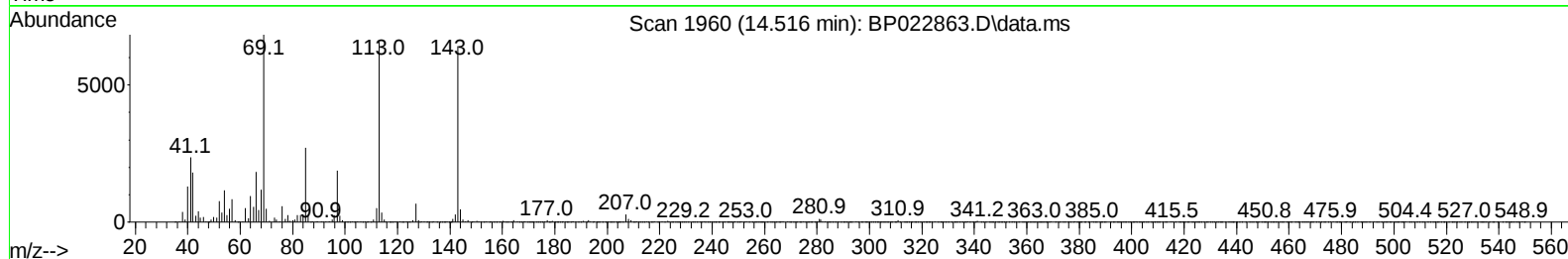
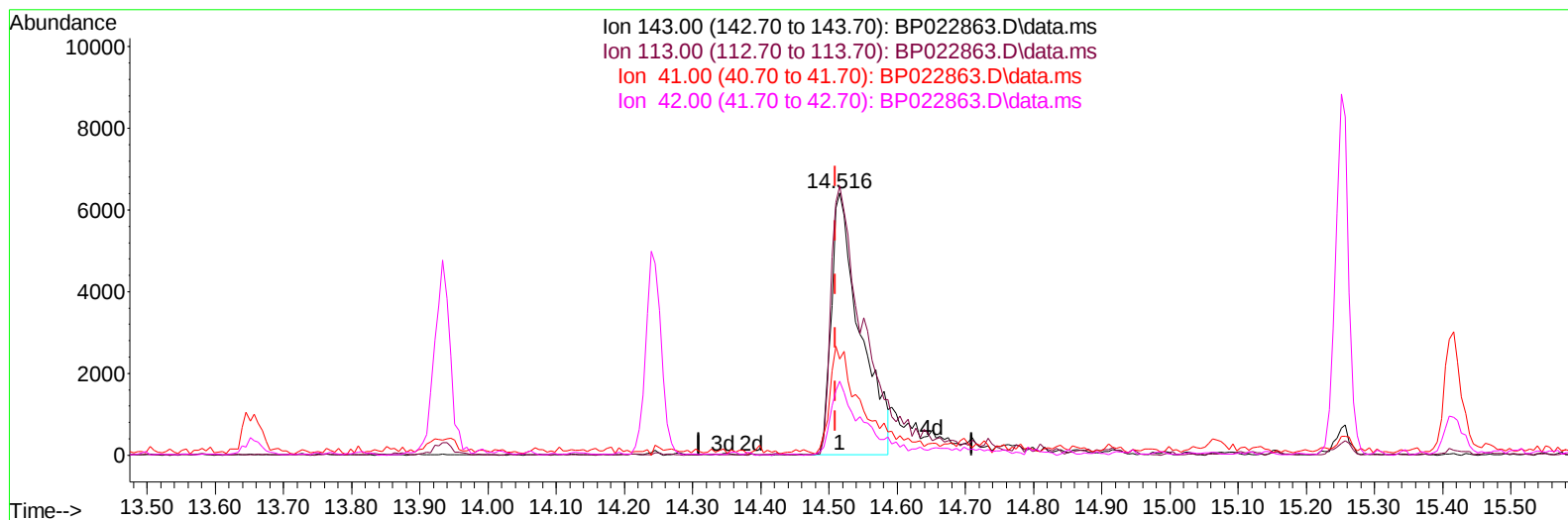
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110424\
Data File : BP022863.D
Acq On : 05 Nov 2024 16:50
Operator : RC/JU
Sample : P4671-02
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AB1

Manual IntegrationsAPPROVED

Quant Time: Nov 05 23:50:48 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

Reviewed By :Yogesh Patel 11/06/2024
Supervised By :mohammad ahmed 11/08/2024



TIC: BP022863.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.516min (+ 0.006) 5.31 ng/u1

response 18714

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	102.67
41.00	35.90	36.81
42.00	23.70	28.28

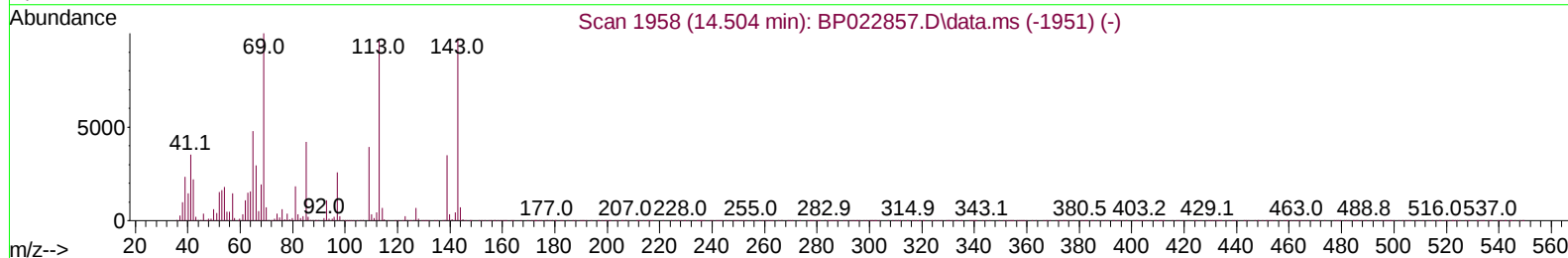
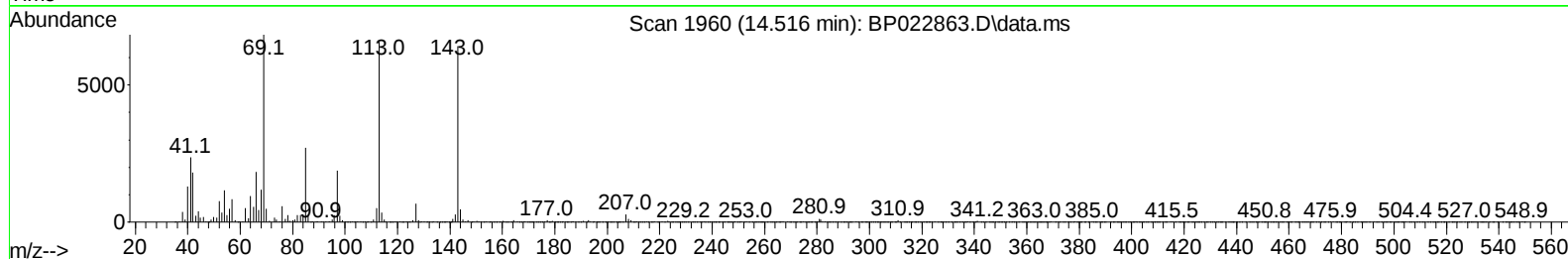
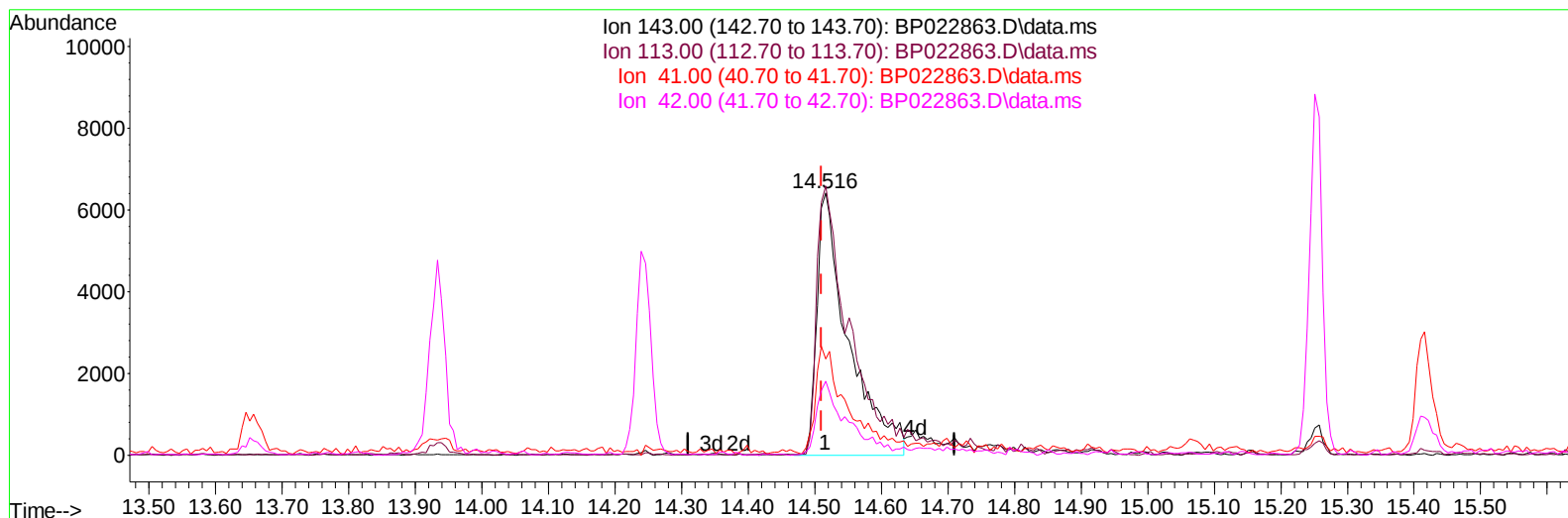
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TIC: BP022863.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.516min (+ 0.006) 5.96 ng/ul m

response 20998

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	102.67
41.00	35.90	36.81
42.00	23.70	28.28

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 ALS Vial : 8 Sample Multiplier: 1

Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/06/2024
 Supervised By :mohammad ahmed 11/08/2024

Quant Time: Nov 05 23:51:54 2024
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 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 30 10:56:39 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.622	152	84949	20.000	ng/ul	-0.02
20) Naphthalene-d8	10.381	136	332045	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.245	164	264490	20.000	ng/ul	-0.02
64) Phenanthrene-d10	17.051	188	636396	20.000	ng/ul	-0.01
79) Chrysene-d12	21.498	240	677602	20.000	ng/ul	0.00
88) Perylene-d12	24.751	264	785797	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.175	96	6917	3.164	ng/uL	0.00
4) Pyridine-d5	3.593	84	48697	8.115	ng/ul	0.00
7) Phenol-d5	6.828	99	53385	7.466	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.969	67	96937	23.065	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.169	132	119024	23.458	ng/ul	-0.01
15) 4-Methylphenol-d8	8.334	113	97016	17.000	ng/ul	-0.01
21) Nitrobenzene-d5	8.769	128	68477	26.821	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.481	143	83667	28.306	ng/ul	-0.02
28) 2,4-Dichlorophenol-d3	10.022	165	166008	26.425	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.522	131	172438	23.229	ng/ul	-0.02
46) Dimethylphthalate-d6	13.657	166	586464	27.908	ng/ul	0.00
49) Acenaphthylene-d8	13.934	160	591837	26.516	ng/ul	-0.02
54) 4-Nitrophenol-d4	14.516	143	20998m	5.956	ng/ul	0.00
60) Fluorene-d10	15.251	176	523352	28.713	ng/ul	-0.02
65) 4,6-Dinitro-2-methylph...	15.416	200	79722	19.047	ng/ul	0.01
73) Anthracene-d10	17.157	188	964764	32.226	ng/ul	-0.01
81) Pyrene-d10	19.539	212	1203422	33.584	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.527	264	1340960	31.954	ng/ul	0.00

Target Compounds Qvalue

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

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