

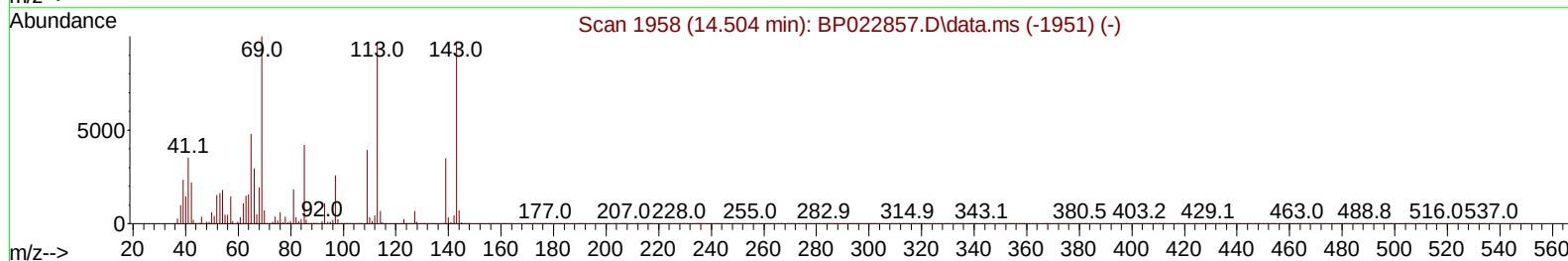
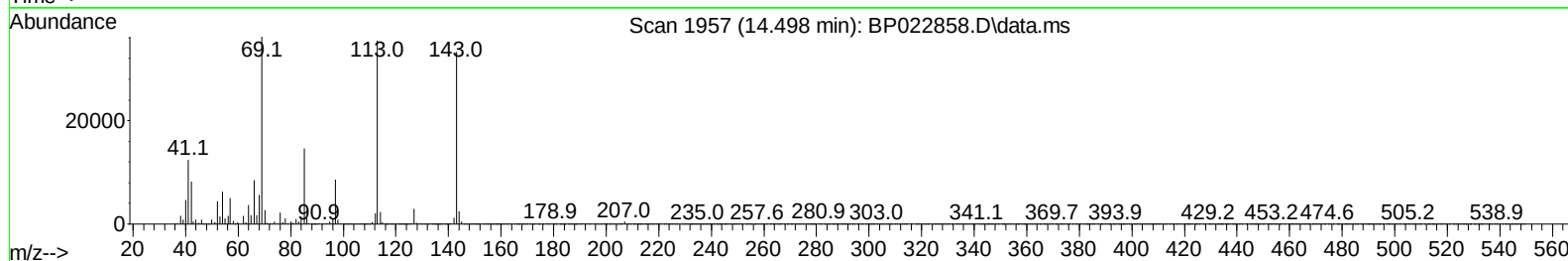
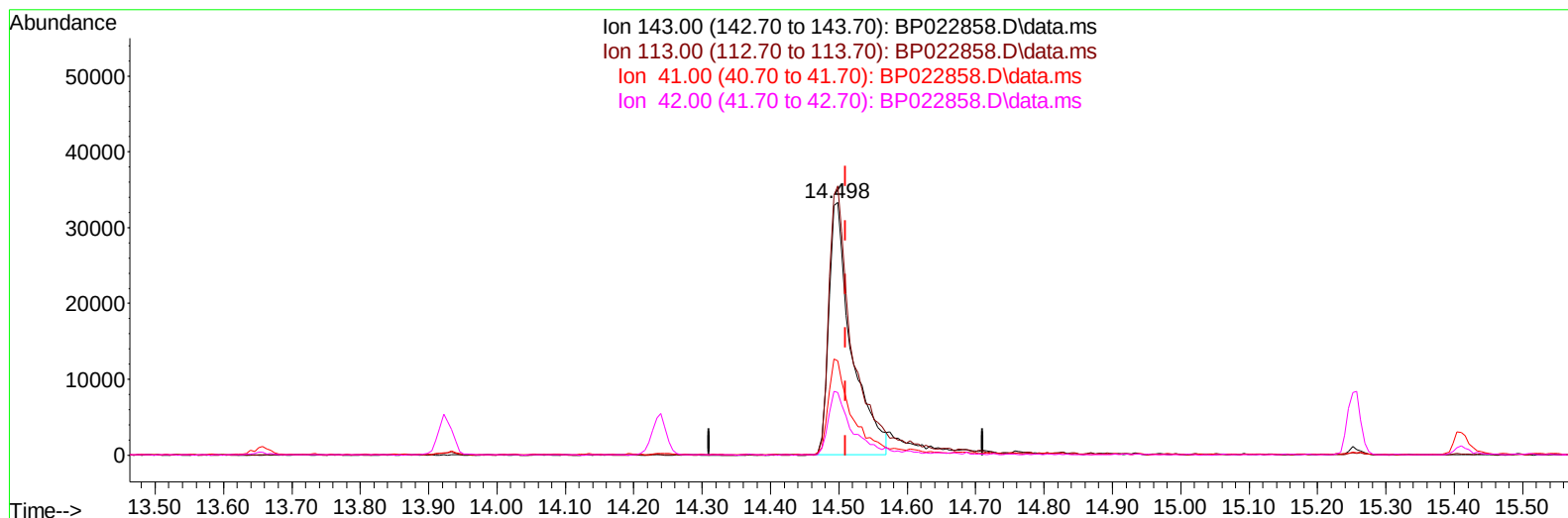
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110424\
Data File : BP022858.D
Acq On : 05 Nov 2024 12:35
Operator : RC/JU
Sample : PB164615BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK615

Manual IntegrationsAPPROVED

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

Quant Time: Nov 05 13:08:45 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: BP022858.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (-0.012) 19.50 ng/u1

response 76407

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	106.46
41.00	35.90	37.25
42.00	23.70	24.79

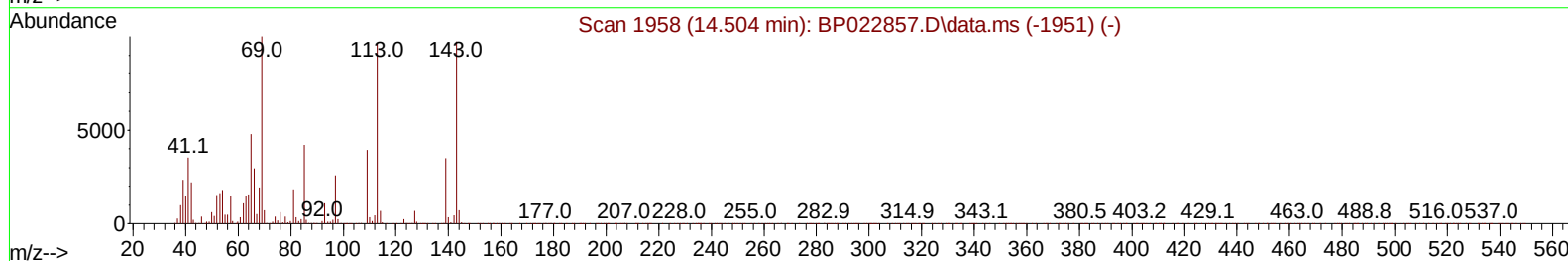
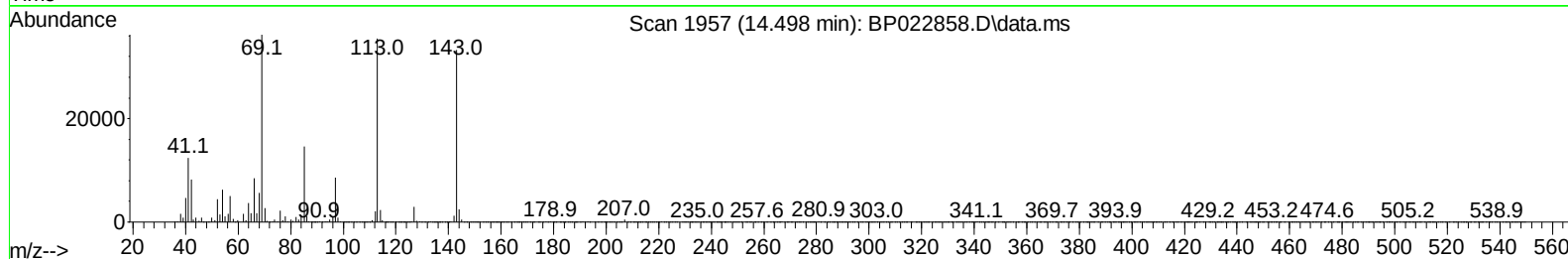
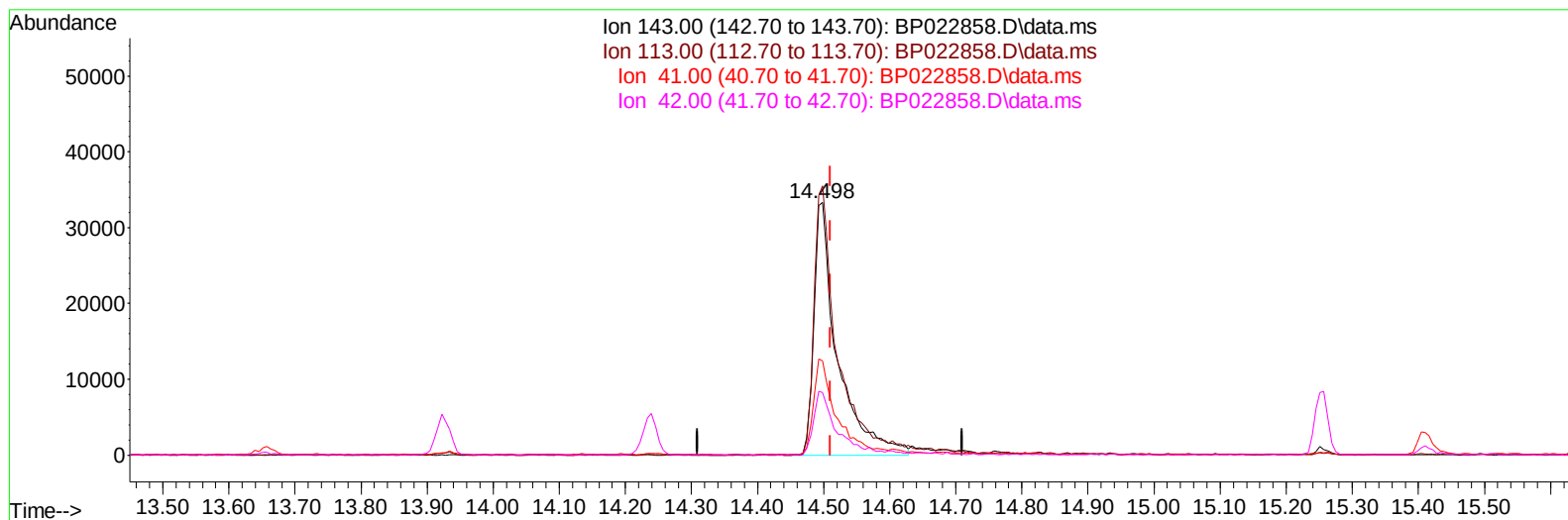
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TIC: BP022858.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (-0.012) 21.07 ng/ul m

response 82593

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	106.46
41.00	35.90	37.25
42.00	23.70	24.79

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 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	88660	20.000	ng/ul	-0.02
20) Naphthalene-d8	10.381	136	352195	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.239	164	294035	20.000	ng/ul	-0.02
64) Phenanthrene-d10	17.051	188	727569	20.000	ng/ul	-0.01
79) Chrysene-d12	21.492	240	669428	20.000	ng/ul	0.00
88) Perylene-d12	24.739	264	741418	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.175	96	8782	3.849	ng/uL	0.00
4) Pyridine-d5	3.587	84	136493	21.792	ng/ul	0.00
7) Phenol-d5	6.822	99	180668	24.208	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	6.969	67	106078	24.183	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.169	132	140845	26.596	ng/ul	-0.01
15) 4-Methylphenol-d8	8.334	113	153386	25.753	ng/ul	-0.01
21) Nitrobenzene-d5	8.769	128	71232	26.304	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.481	143	88603	28.261	ng/ul	-0.02
28) 2,4-Dichlorophenol-d3	10.016	165	171214	25.694	ng/ul	-0.02
31) 4-Chloroaniline-d4	10.522	131	204912	26.025	ng/ul	-0.02
46) Dimethylphthalate-d6	13.657	166	649638	27.808	ng/ul	0.00
49) Acenaphthylene-d8	13.928	160	667347	26.895	ng/ul	-0.02
54) 4-Nitrophenol-d4	14.498	143	82593m	21.074	ng/ul	-0.01
60) Fluorene-d10	15.257	176	551340	27.209	ng/ul	-0.02
65) 4,6-Dinitro-2-methylph...	15.410	200	82949	17.335	ng/ul	0.00
73) Anthracene-d10	17.163	188	951433	27.798	ng/ul	0.00
81) Pyrene-d10	19.533	212	1108844	31.323	ng/ul	-0.01
92) Benzo(a)pyrene-d12	24.521	264	1103463	27.868	ng/ul	0.00

Target Compounds Qvalue

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

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