

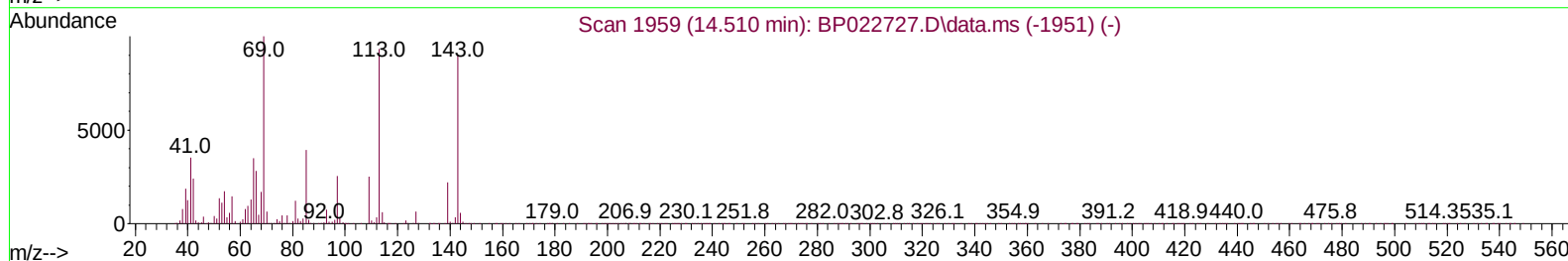
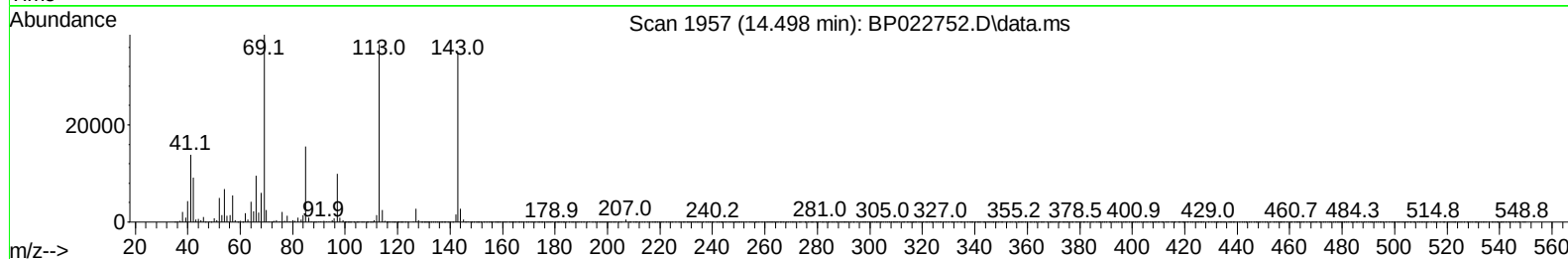
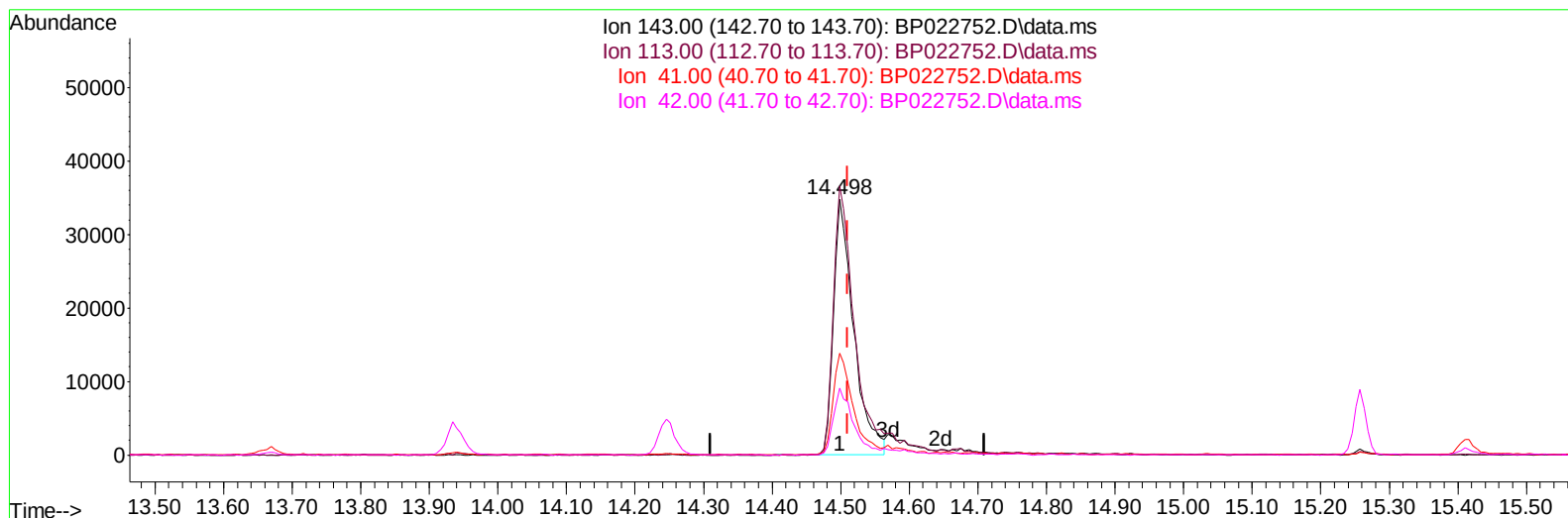
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103024\
Data File : BP022752.D
Acq On : 31 Oct 2024 02:37
Operator : RC/JU
Sample : PB164506BL
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK506

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/04/2024
Supervised By :mohammad ahmed 11/05/2024

Quant Time: Oct 31 12:38:36 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: BP022752.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (-0.012) 18.38 ng/u1

response 71038

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	105.10
41.00	35.90	39.90
42.00	23.70	26.20

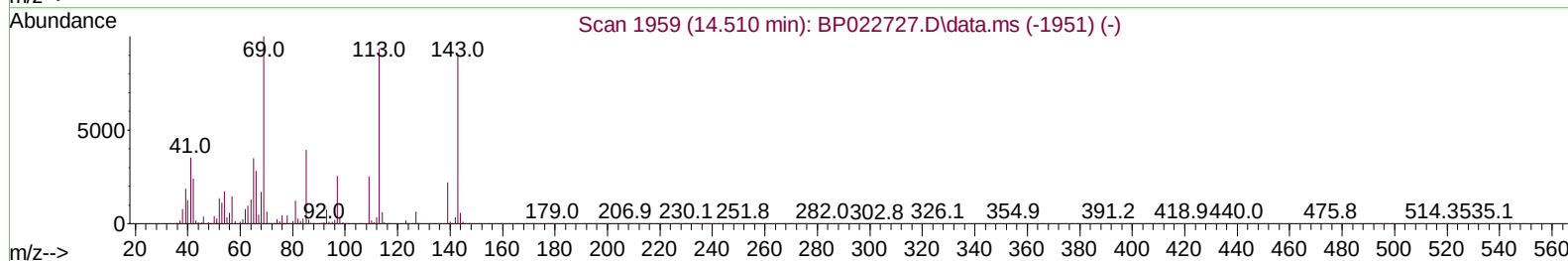
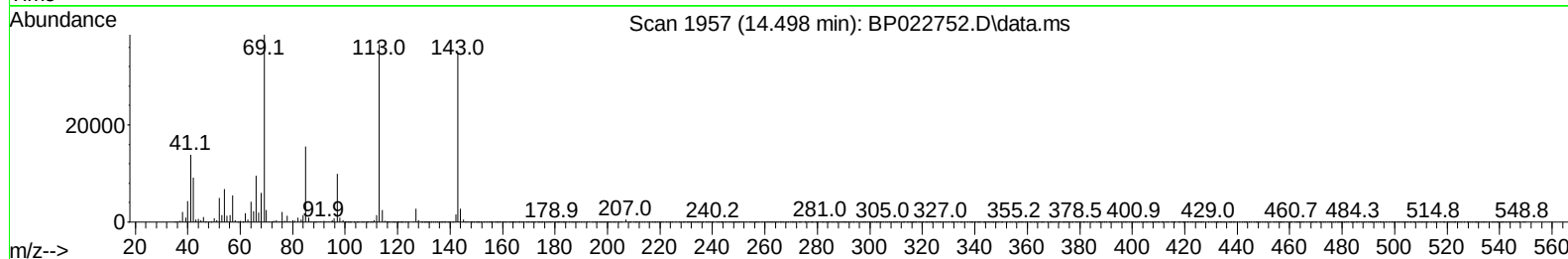
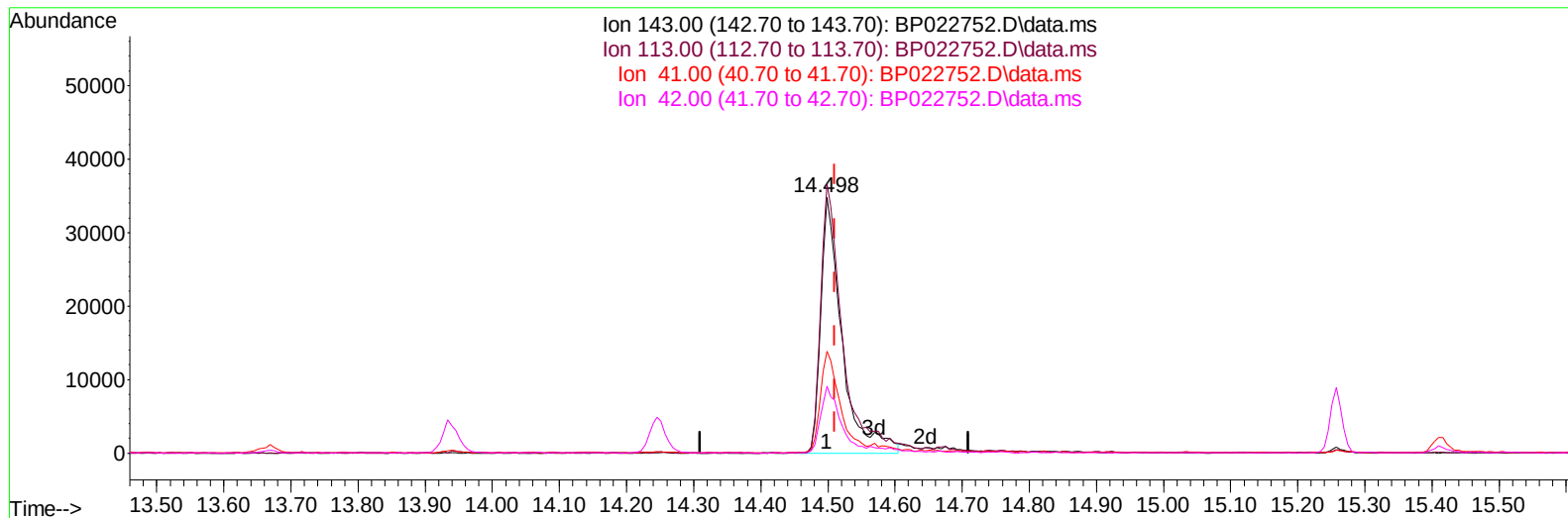
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TIC: BP022752.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (-0.012) 19.68 ng/ul m

response 76073

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	105.10
41.00	35.90	39.90
42.00	23.70	26.20

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/04/2024
 Supervised By :mohammad ahmed 11/05/2024

Quant Time: Nov 01 00:22:48 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

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.634	152	97500	20.000	ng/ul	0.00
20) Naphthalene-d8	10.387	136	369404	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.246	164	289968	20.000	ng/ul	-0.02
64) Phenanthrene-d10	17.057	188	734559	20.000	ng/ul	0.00
79) Chrysene-d12	21.498	240	782464	20.000	ng/ul	0.00
88) Perylene-d12	24.757	264	876153	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.181	96	10603	4.226	ng/uL	0.00
4) Pyridine-d5	3.593	84	153455	22.279	ng/ul	0.00
7) Phenol-d5	6.834	99	191542	23.338	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.981	67	111195	23.051	ng/ul	0.00
11) 2-Chlorophenol-d4	7.181	132	143891	24.708	ng/ul	0.00
15) 4-Methylphenol-d8	8.340	113	154050	23.520	ng/ul	0.00
21) Nitrobenzene-d5	8.781	128	71578	25.200	ng/ul	0.00
24) 2-Nitrophenol-d4	9.493	143	84557	25.714	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.028	165	170342	24.373	ng/ul	0.00
31) 4-Chloroaniline-d4	10.534	131	195231	23.640	ng/ul	0.00
46) Dimethylphthalate-d6	13.669	166	557595	24.203	ng/ul	0.00
49) Acenaphthylene-d8	13.934	160	607744	24.837	ng/ul	-0.02
54) 4-Nitrophenol-d4	14.498	143	76073m	19.682	ng/ul	-0.01
60) Fluorene-d10	15.257	176	502067	25.125	ng/ul	-0.02
65) 4,6-Dinitro-2-methylph...	15.410	200	61504	12.731	ng/ul	0.00
73) Anthracene-d10	17.157	188	854206	24.720	ng/ul	-0.01
81) Pyrene-d10	19.545	212	1083046	26.174	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.521	264	1202308	25.695	ng/ul	0.00

Target Compounds Qvalue

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

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