

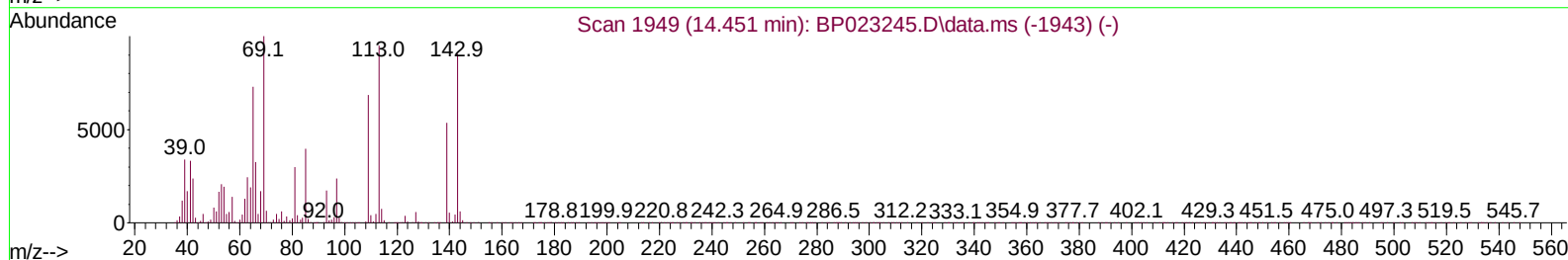
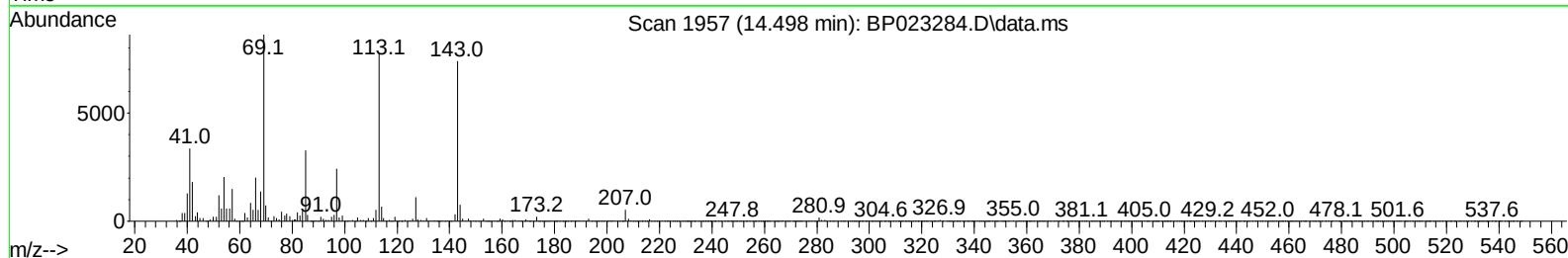
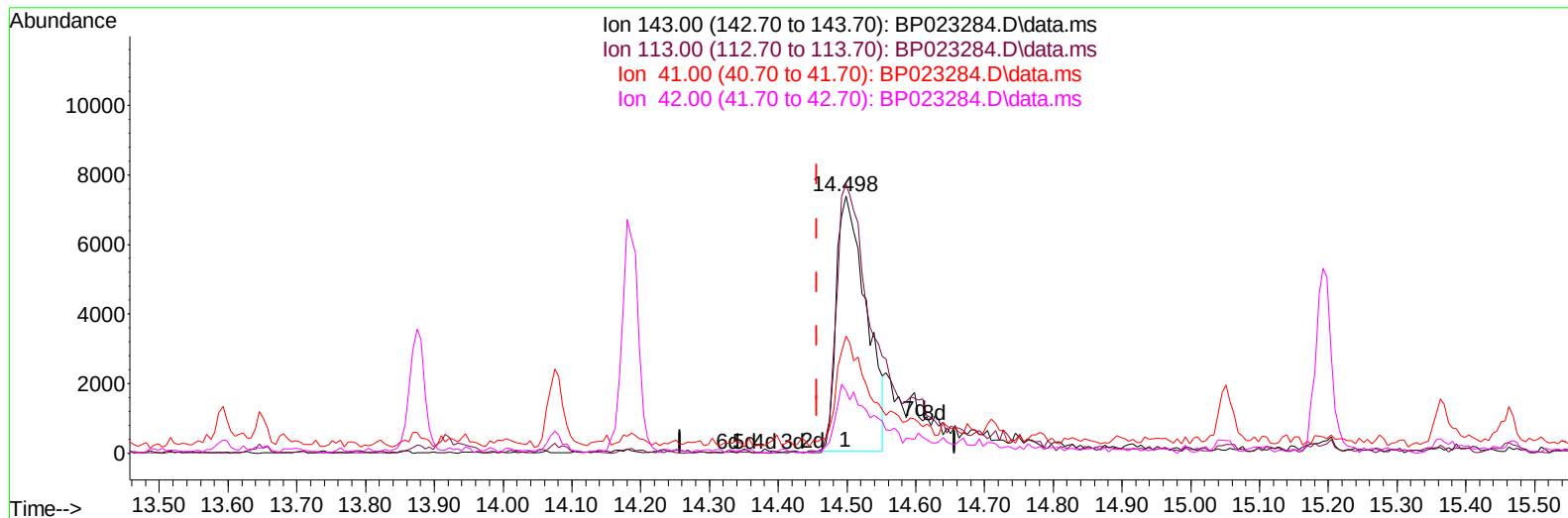
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111924\
Data File : BP023284.D
Acq On : 21 Nov 2024 14:42
Operator : RC/JU
Sample : P4872-22
Misc :
ALS Vial : 43 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29Y9

Manual IntegrationsAPPROVED

Quant Time: Nov 22 00:33:51 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 20 04:34:35 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/22/2024
Supervised By :mohammad ahmed 11/22/2024



TIC: BP023284.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (+ 0.041) 5.94 ng/ul

response 22099

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	104.50
41.00	39.20	45.45
42.00	25.90	24.35

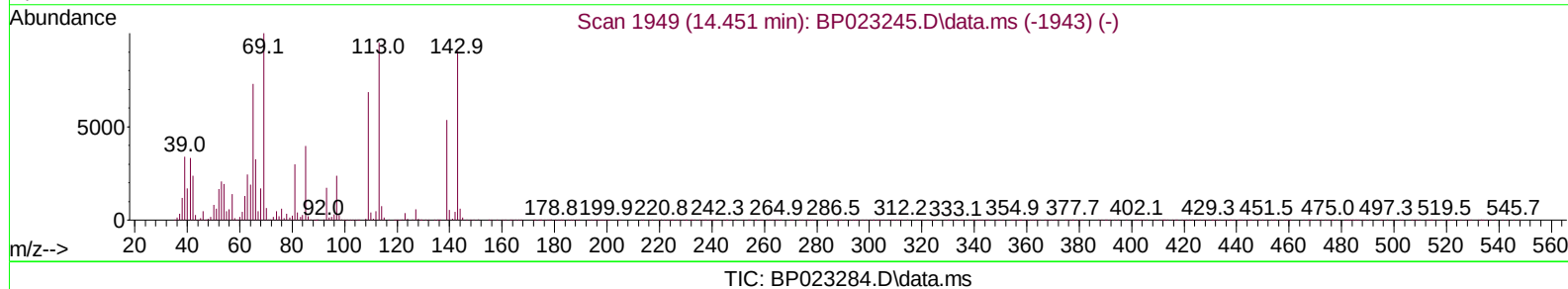
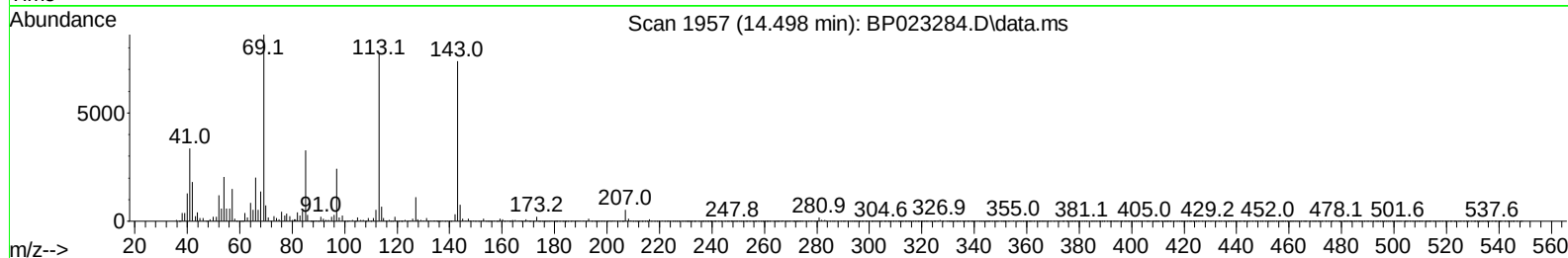
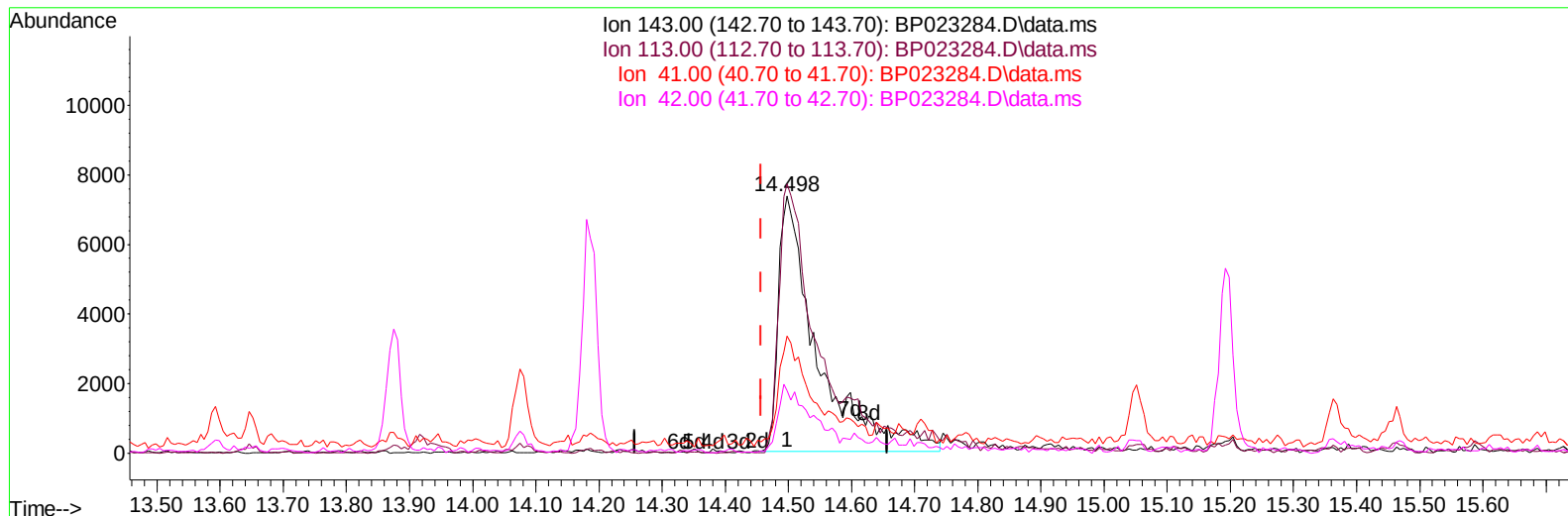
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(54) 4-Nitrophenol-d4 (S)

14.498min (+ 0.041) 8.53 ng/ul m

response 31712

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	104.50
41.00	39.20	45.45
42.00	25.90	24.35

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.563	152	124072	20.000	ng/ul	0.00
20) Naphthalene-d8	10.310	136	456527	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.187	164	343308	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.998	188	824187	20.000	ng/ul	0.00
79) Chrysene-d12	21.445	240	955648	20.000	ng/ul	0.02
88) Perylene-d12	24.674	264	1130156	20.000	ng/ul	0.04
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.134	96	6218	2.409	ng/uL	0.00
4) Pyridine-d5	3.552	84	79814	10.408	ng/ul	0.00
7) Phenol-d5	6.769	99	118249	12.891	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.916	67	72039	13.367	ng/ul	0.00
11) 2-Chlorophenol-d4	7.110	132	90010	13.398	ng/ul	0.00
15) 4-Methylphenol-d8	8.281	113	103589	13.723	ng/ul	0.00
21) Nitrobenzene-d5	8.710	128	44350	13.954	ng/ul	0.00
24) 2-Nitrophenol-d4	9.422	143	54461	14.373	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.963	165	108793	13.711	ng/ul	0.00
31) 4-Chloroaniline-d4	10.475	131	127579	13.477	ng/ul	0.00
46) Dimethylphthalate-d6	13.592	166	418790	16.925	ng/ul	0.00
49) Acenaphthylene-d8	13.875	160	416602	16.136	ng/ul	0.00
54) 4-Nitrophenol-d4	14.498	143	31712m	8.528	ng/ul	0.04
60) Fluorene-d10	15.192	176	359811	16.586	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.363	200	32439	6.570	ng/ul	0.01
73) Anthracene-d10	17.104	188	609078	17.454	ng/ul	0.00
81) Pyrene-d10	19.486	212	805852	18.307	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.462	264	968232	17.986	ng/ul	0.04

Target Compounds Qvalue

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

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