

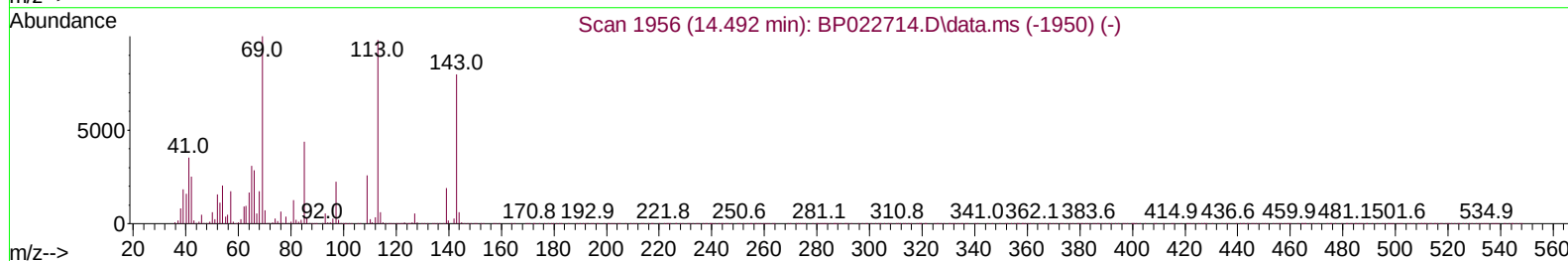
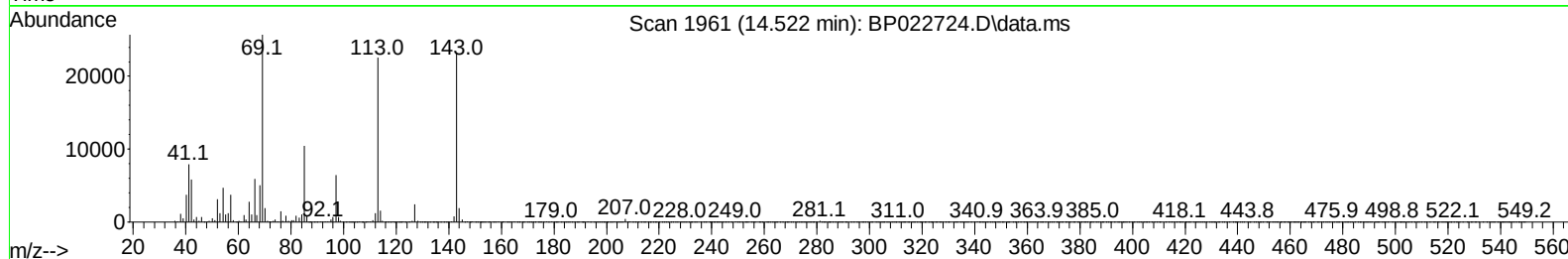
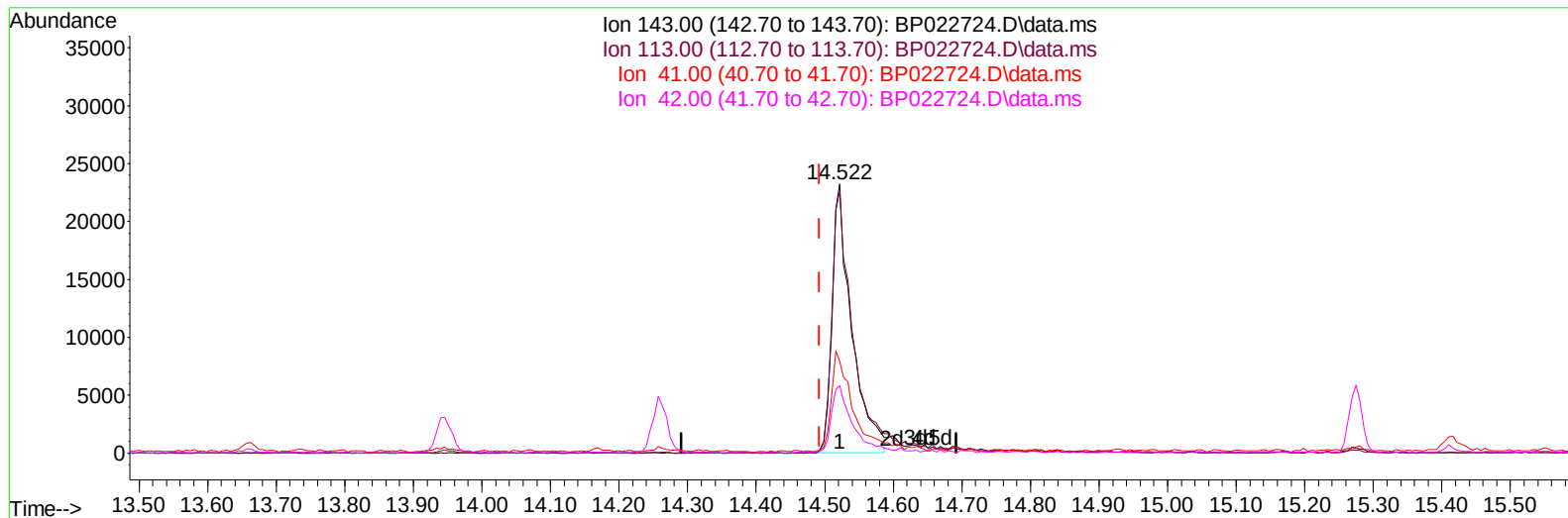
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102924\
 Data File : BP022724.D
 Acq On : 30 Oct 2024 00:11
 Operator : RC/JU
 Sample : P4492-12
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A4ER0

Manual IntegrationsAPPROVED

Quant Time: Oct 30 01:16:32 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Oct 29 18:08:42 2024
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 10/30/2024
 Supervised By :mohammad ahmed 11/04/2024



TIC: BP022724.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.522min (+ 0.030) 13.48 ng/ul

response 45966

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	97.28
41.00	35.90	34.08
42.00	23.70	25.15

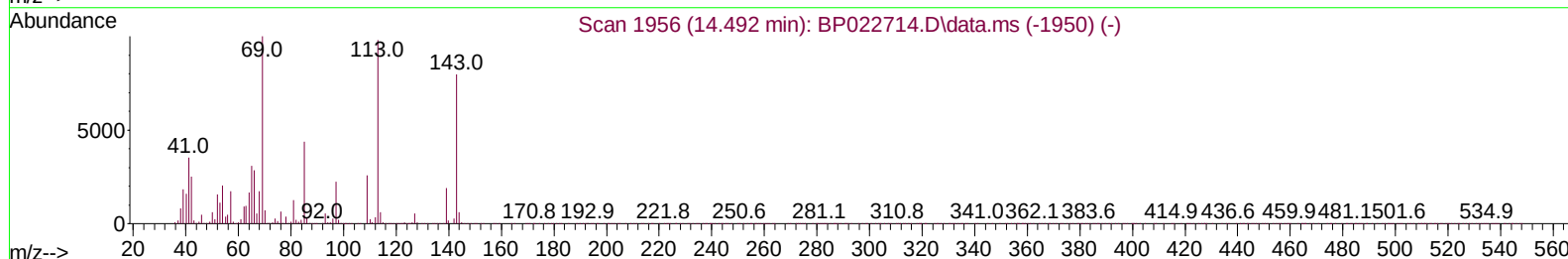
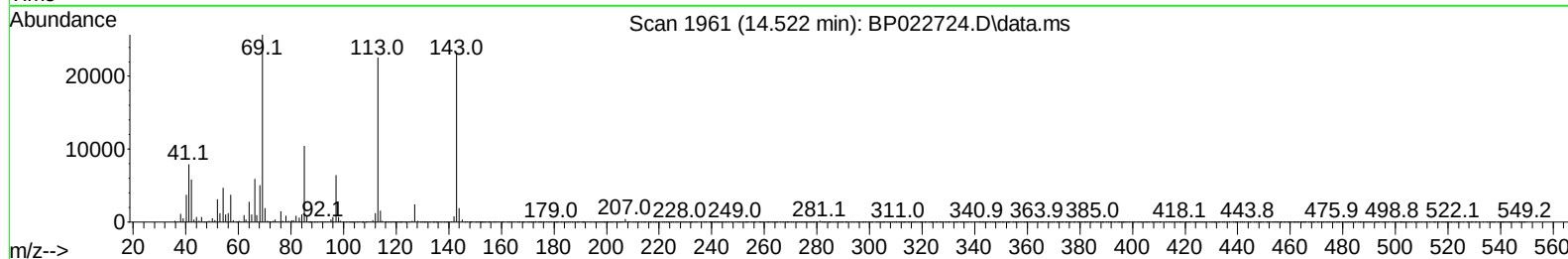
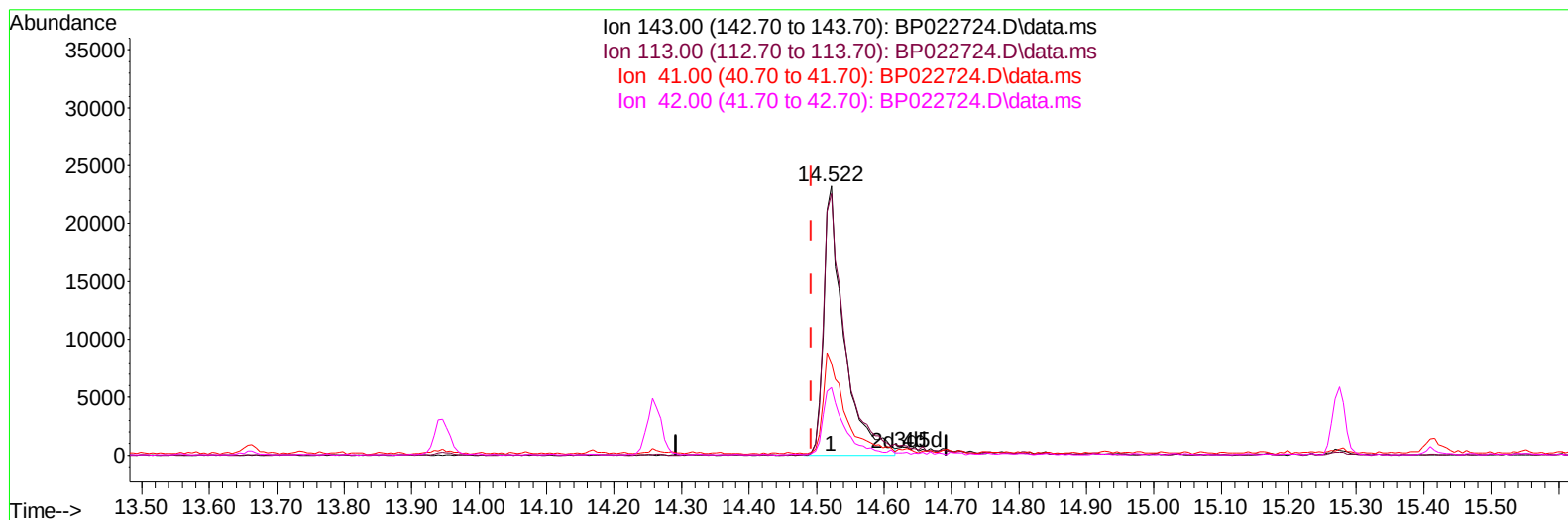
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TIC: BP022724.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.522min (+ 0.030) 14.06 ng/ul m

response 47966

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	97.28
41.00	35.90	34.08
42.00	23.70	25.15

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.640	152	84148	20.000	ng/ul	0.00
20) Naphthalene-d8	10.393	136	324641	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.263	164	255895	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.063	188	599325	20.000	ng/ul	0.00
79) Chrysene-d12	21.498	240	677178	20.000	ng/ul	0.00
88) Perylene-d12	24.745	264	833382	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.181	96	5896	2.723	ng/uL	0.00
4) Pyridine-d5	3.599	84	80087	13.472	ng/ul	0.00
7) Phenol-d5	6.834	99	117029	16.522	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.987	67	65500	15.733	ng/ul	0.00
11) 2-Chlorophenol-d4	7.187	132	86229	17.156	ng/ul	0.00
15) 4-Methylphenol-d8	8.346	113	95554	16.904	ng/ul	0.00
21) Nitrobenzene-d5	8.781	128	42607	17.069	ng/ul	0.00
24) 2-Nitrophenol-d4	9.493	143	51696	17.888	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.034	165	106440	17.329	ng/ul	0.00
31) 4-Chloroaniline-d4	10.540	131	118223	16.289	ng/ul	0.00
46) Dimethylphthalate-d6	13.663	166	372318	18.312	ng/ul	0.00
49) Acenaphthylene-d8	13.945	160	399758	18.512	ng/ul	0.00
54) 4-Nitrophenol-d4	14.522	143	47966m	14.063	ng/ul	0.03
60) Fluorene-d10	15.275	176	342063	19.397	ng/ul	0.01
65) 4,6-Dinitro-2-methylph...	15.416	200	37802	9.590	ng/ul	0.00
73) Anthracene-d10	17.163	188	556058	19.723	ng/ul	0.00
81) Pyrene-d10	19.545	212	738362	20.619	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.527	264	932862	20.960	ng/ul	0.01

Target Compounds Qvalue

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

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