

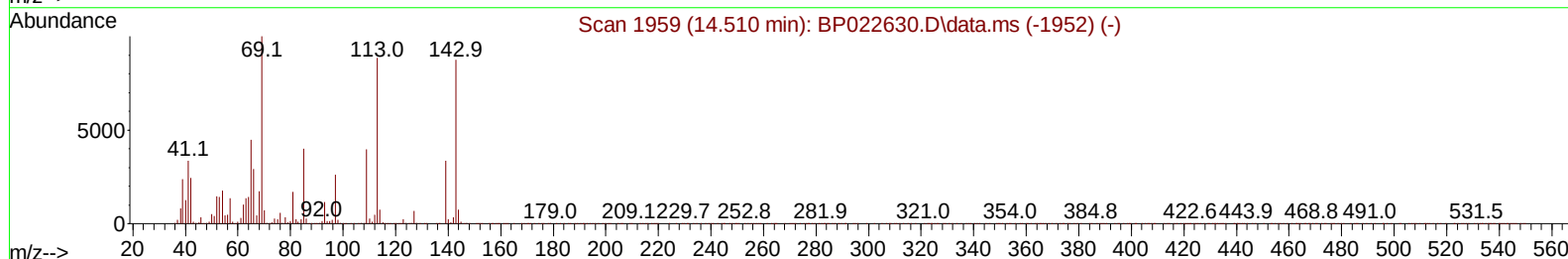
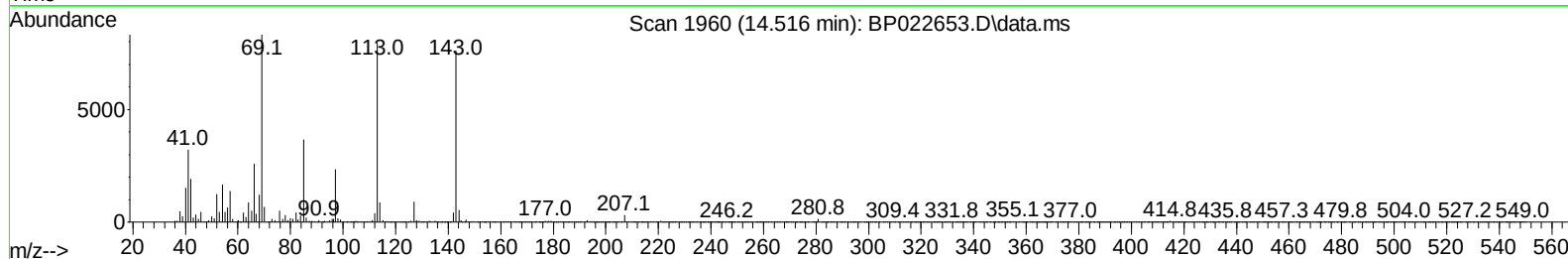
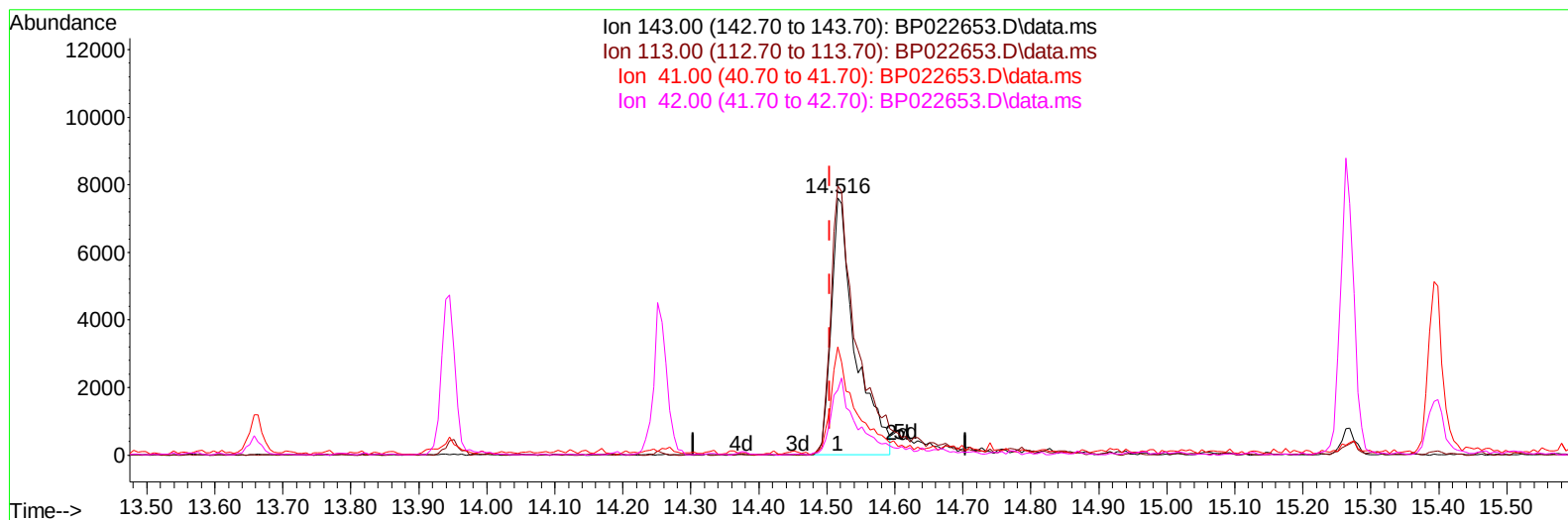
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102524\
 Data File : BP022653.D
 Acq On : 26 Oct 2024 22:05
 Operator : RC/JU
 Sample : P4559-02
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCNX9

Manual IntegrationsAPPROVED

Quant Time: Oct 28 00:32:33 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Oct 26 05:55:04 2024
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/28/2024
 Supervised By :mohammad ahmed 10/29/2024



TIC: BP022653.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.516min (+ 0.012) 6.78 ng/ul

response 18728

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	104.53
41.00	35.90	42.09
42.00	23.70	25.28

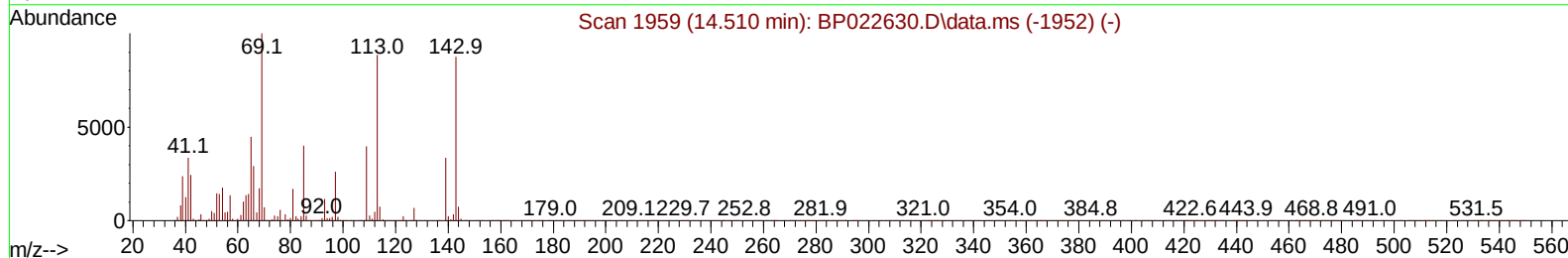
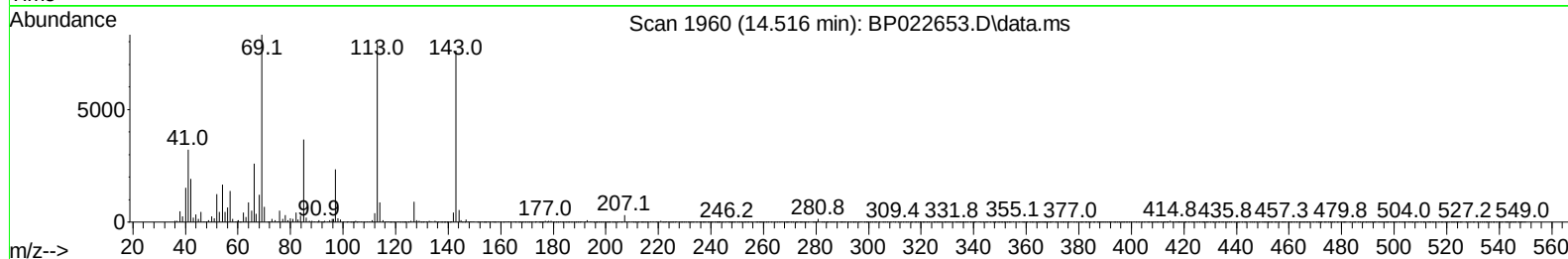
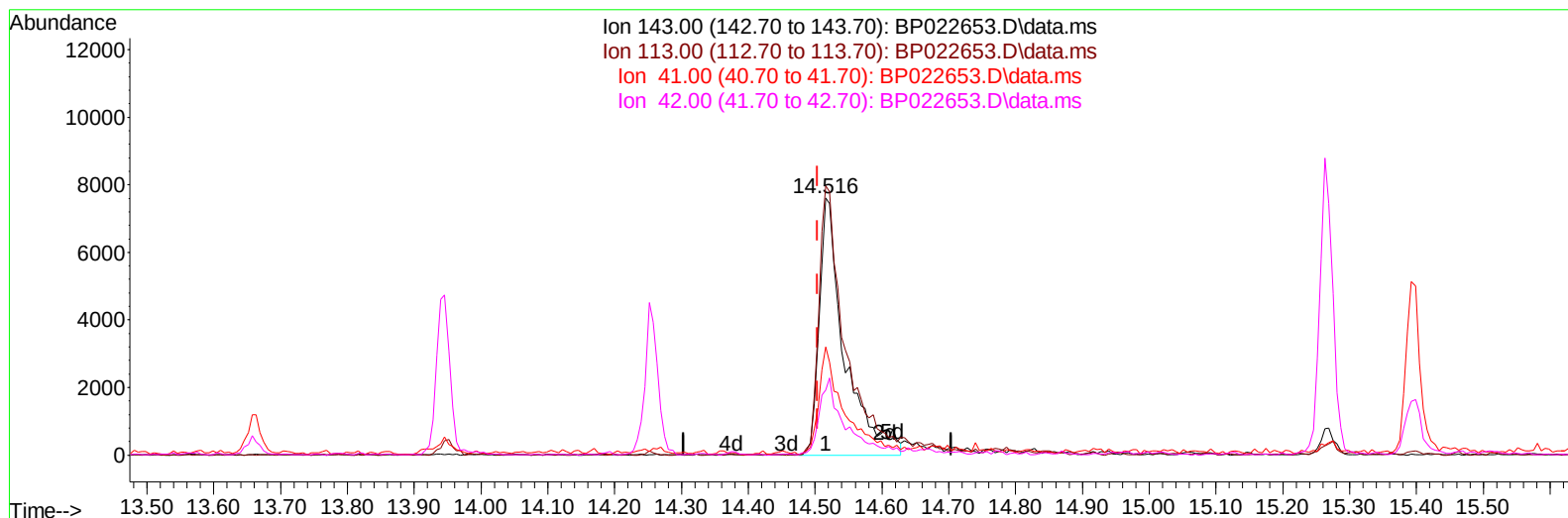
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TIC: BP022653.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.516min (+ 0.012) 7.22 ng/ul m

response 19950

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	104.53
41.00	35.90	42.09
42.00	23.70	25.28

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.646	152	67648	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.399	136	259752	20.000	ng/ul	-0.02
38) Acenaphthene-d10	14.257	164	207383	20.000	ng/ul	-0.02
64) Phenanthrene-d10	17.051	188	530164	20.000	ng/ul	-0.03
79) Chrysene-d12	21.492	240	611840	20.000	ng/ul	0.00
88) Perylene-d12	24.751	264	700375	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.187	96	8526	4.898	ng/uL	-0.01
4) Pyridine-d5	3.617	84	10096	2.113	ng/ul	0.00
7) Phenol-d5	6.846	99	44927	7.890	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.987	67	108045	32.283	ng/ul	-0.02
11) 2-Chlorophenol-d4	7.193	132	122585	30.338	ng/ul	-0.01
15) 4-Methylphenol-d8	8.352	113	92305	20.312	ng/ul	-0.01
21) Nitrobenzene-d5	8.787	128	69961	35.029	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.499	143	84943	36.736	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.040	165	161389	32.839	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.557	131	6419	1.105	ng/ul	0.00
46) Dimethylphthalate-d6	13.663	166	585389	35.528	ng/ul	-0.02
49) Acenaphthylene-d8	13.945	160	579885	33.135	ng/ul	-0.02
54) 4-Nitrophenol-d4	14.516	143	19950m	7.217	ng/ul	0.01
60) Fluorene-d10	15.263	176	522934	36.590	ng/ul	-0.03
65) 4,6-Dinitro-2-methylph...	15.398	200	112801	32.351	ng/ul	-0.02
73) Anthracene-d10	17.157	188	997506	39.996	ng/ul	-0.01
81) Pyrene-d10	19.528	212	1275999	39.437	ng/ul	-0.03
92) Benzo(a)pyrene-d12	24.521	264	1529792	40.900	ng/ul	0.00
Target Compounds						
2) 1,4-Dioxane	3.223	88	2565	1.370	ng/uL	Qvalue 93

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

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