

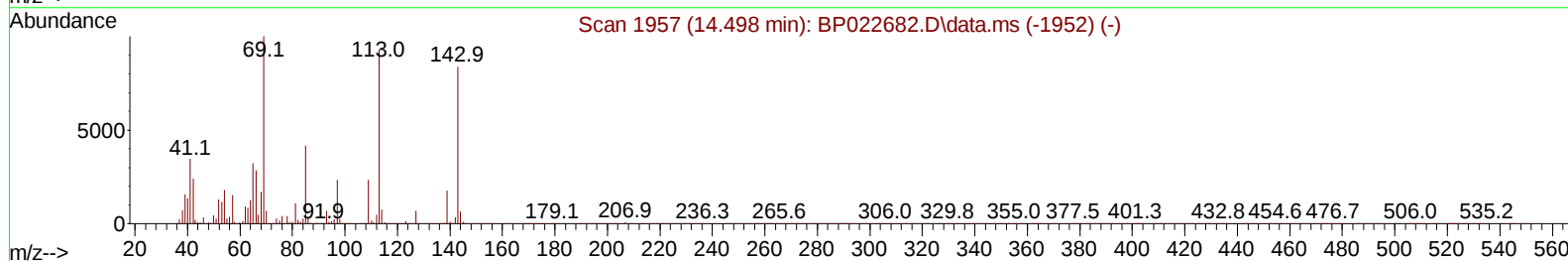
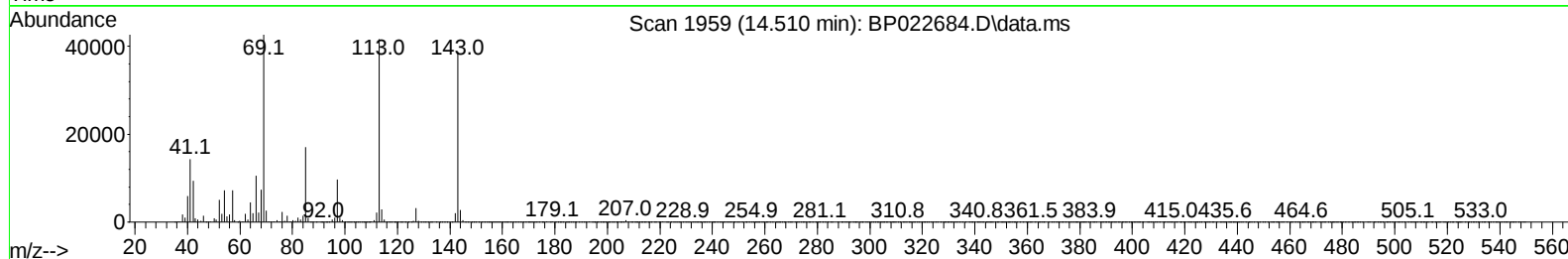
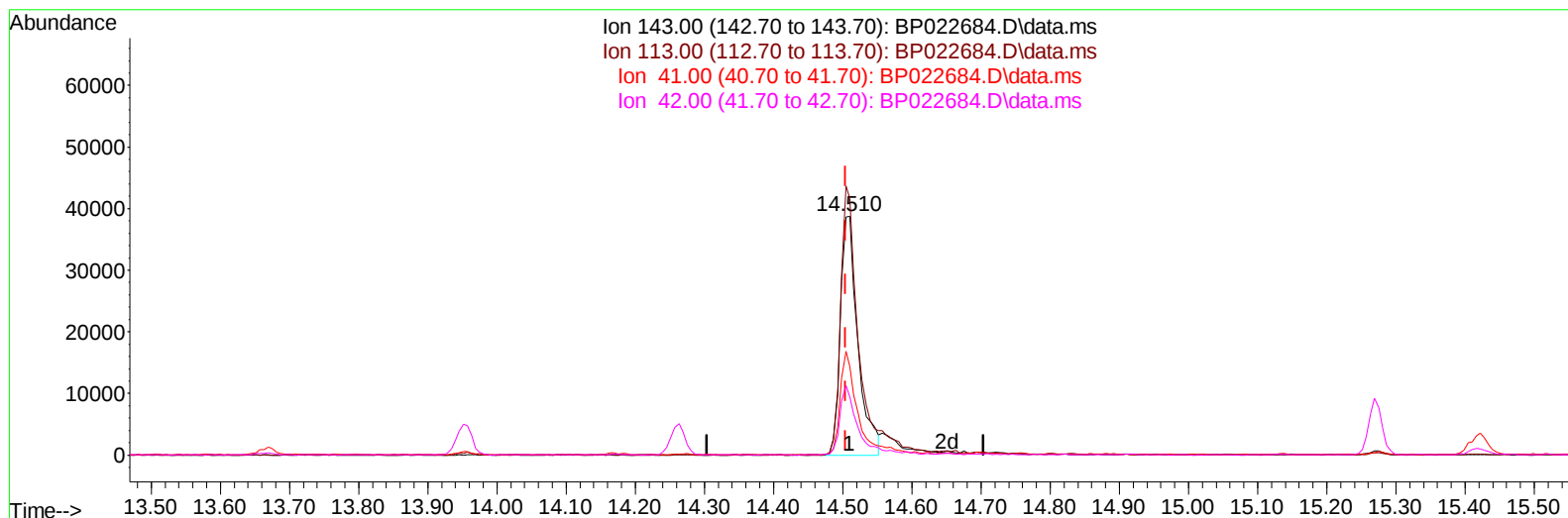
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102824\
 Data File : BP022684.D
 Acq On : 28 Oct 2024 17:39
 Operator : RC/JU
 Sample : P4530-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 F7H26

Manual IntegrationsAPPROVED

Quant Time: Oct 28 18:09:20 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 10/29/2024
 Supervised By :mohammad ahmed 10/30/2024



TIC: BP022684.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.510min (+ 0.006) 22.36 ng/u1

response 68111

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	107.54
41.00	35.90	37.09
42.00	23.70	24.38

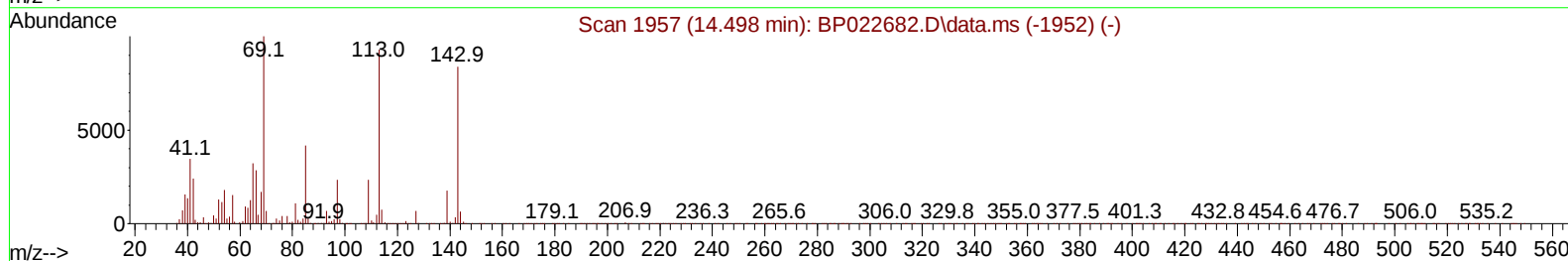
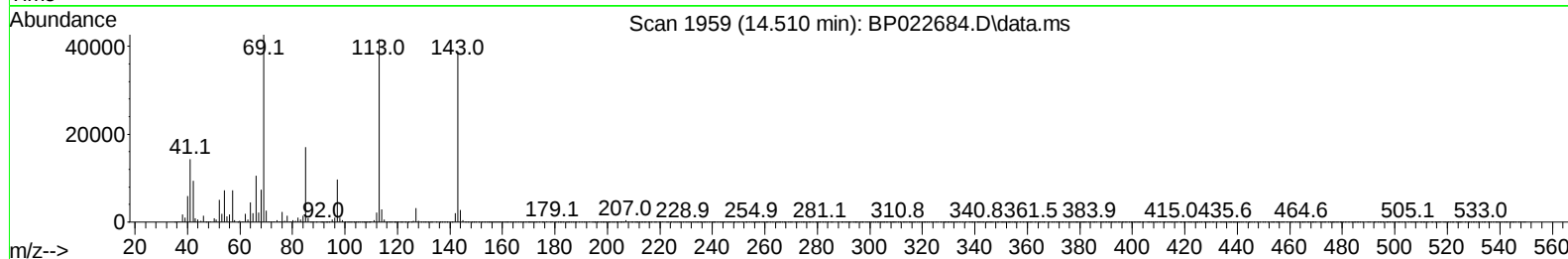
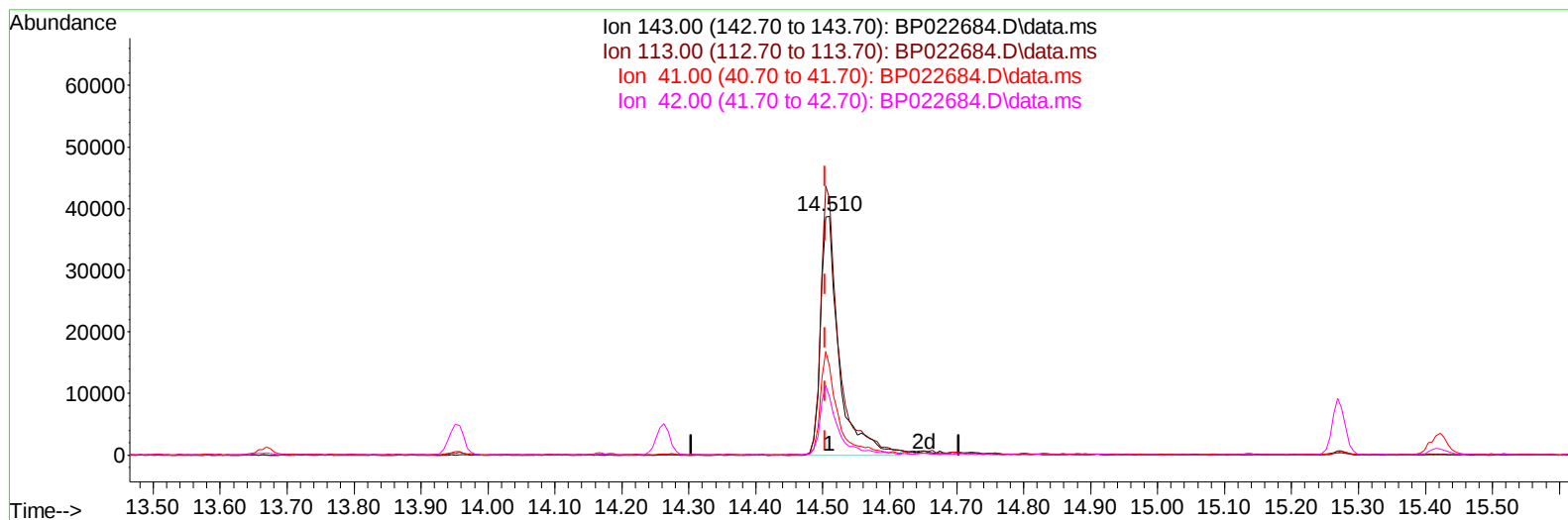
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TIC: BP022684.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.510min (+ 0.006) 24.61 ng/ul m

response 74968

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	107.54
41.00	35.90	37.09
42.00	23.70	24.38

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Instrument :
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Reviewed By :Yogesh Patel 10/29/2024
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Quant Time: Oct 28 18:10:03 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.646	152	75424	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.405	136	288821	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.263	164	228503	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.057	188	572569	20.000	ng/ul	-0.02
79) Chrysene-d12	21.492	240	645765	20.000	ng/ul	# 0.00
88) Perylene-d12	24.745	264	775973	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.181	96	8874	4.572	ng/uL	-0.02
4) Pyridine-d5	3.593	84	114423	21.475	ng/ul	-0.02
7) Phenol-d5	6.840	99	165311	26.037	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	6.987	67	99258	26.600	ng/ul	-0.02
11) 2-Chlorophenol-d4	7.187	132	121285	26.922	ng/ul	-0.02
15) 4-Methylphenol-d8	8.352	113	135262	26.696	ng/ul	-0.01
21) Nitrobenzene-d5	8.787	128	60625	27.299	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.499	143	75683	29.437	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.046	165	151884	27.795	ng/ul	0.00
31) 4-Chloroaniline-d4	10.540	131	172060	26.647	ng/ul	-0.02
46) Dimethylphthalate-d6	13.669	166	582982	32.111	ng/ul	-0.01
49) Acenaphthylene-d8	13.951	160	580834	30.122	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.510	143	74968m	24.614	ng/ul	0.00
60) Fluorene-d10	15.275	176	500805	31.803	ng/ul	-0.02
65) 4,6-Dinitro-2-methylph...	15.422	200	92469	24.556	ng/ul	0.00
73) Anthracene-d10	17.169	188	888157	32.974	ng/ul	0.00
81) Pyrene-d10	19.545	212	1095921	32.092	ng/ul	-0.01
92) Benzo(a)pyrene-d12	24.516	264	1362548	32.879	ng/ul	-0.01

Target Compounds Qvalue

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

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