

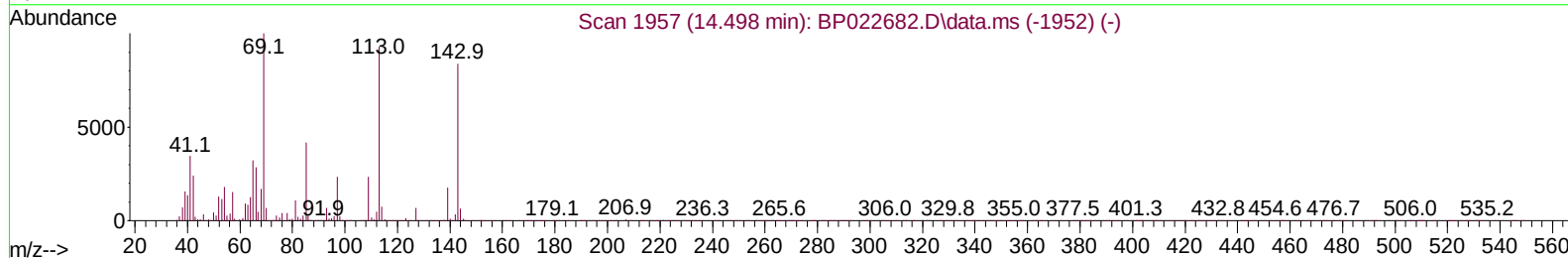
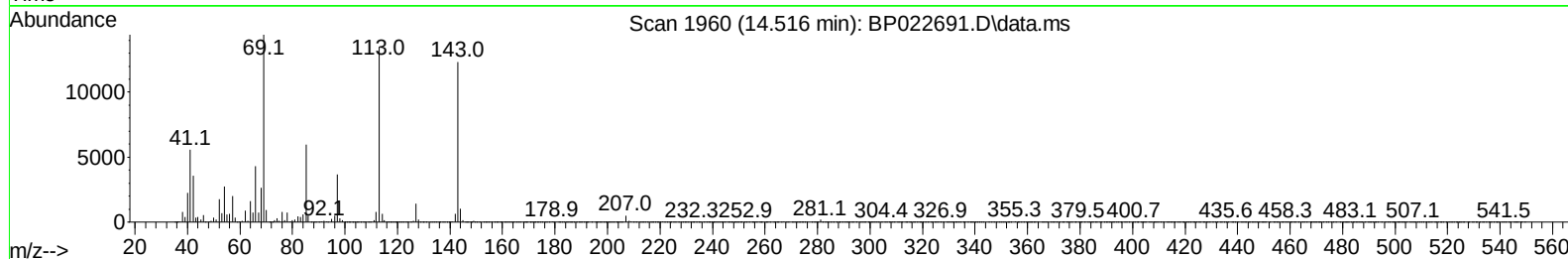
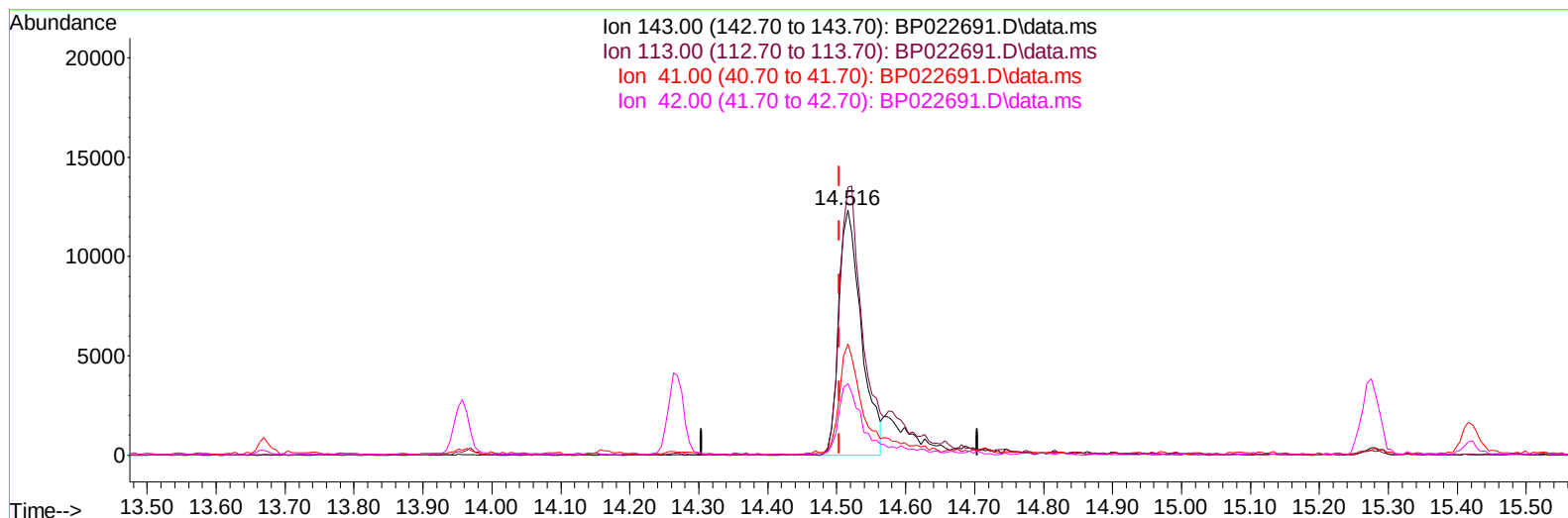
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102824\
Data File : BP022691.D
Acq On : 28 Oct 2024 22:21
Operator : RC/JU
Sample : P4530-09
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
F7H33

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 10/29/2024
Supervised By :mohammad ahmed 10/30/2024

Quant Time: Oct 29 01:23:12 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



TIC: BP022691.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.516min (+ 0.012) 10.64 ng/ul

response 28164

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	109.33
41.00	35.90	45.33#
42.00	23.70	29.02#

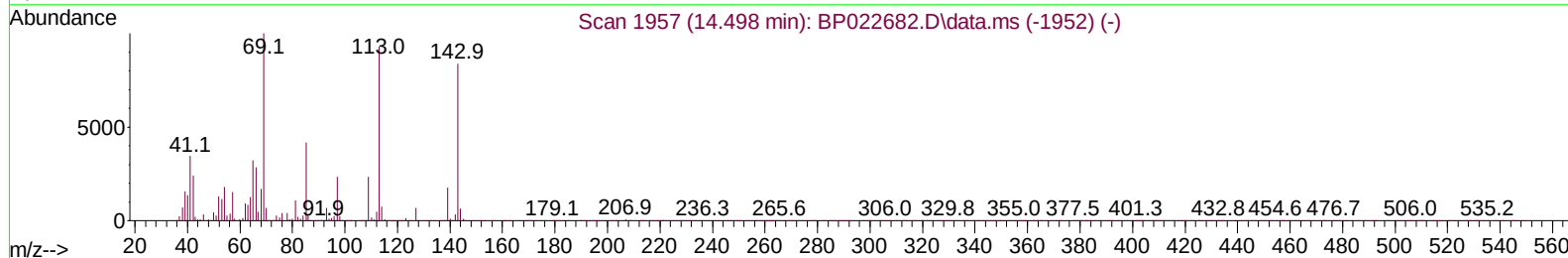
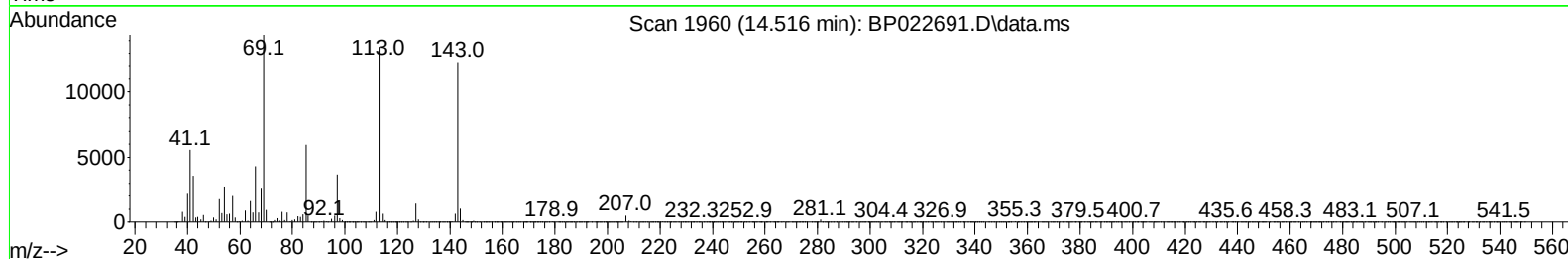
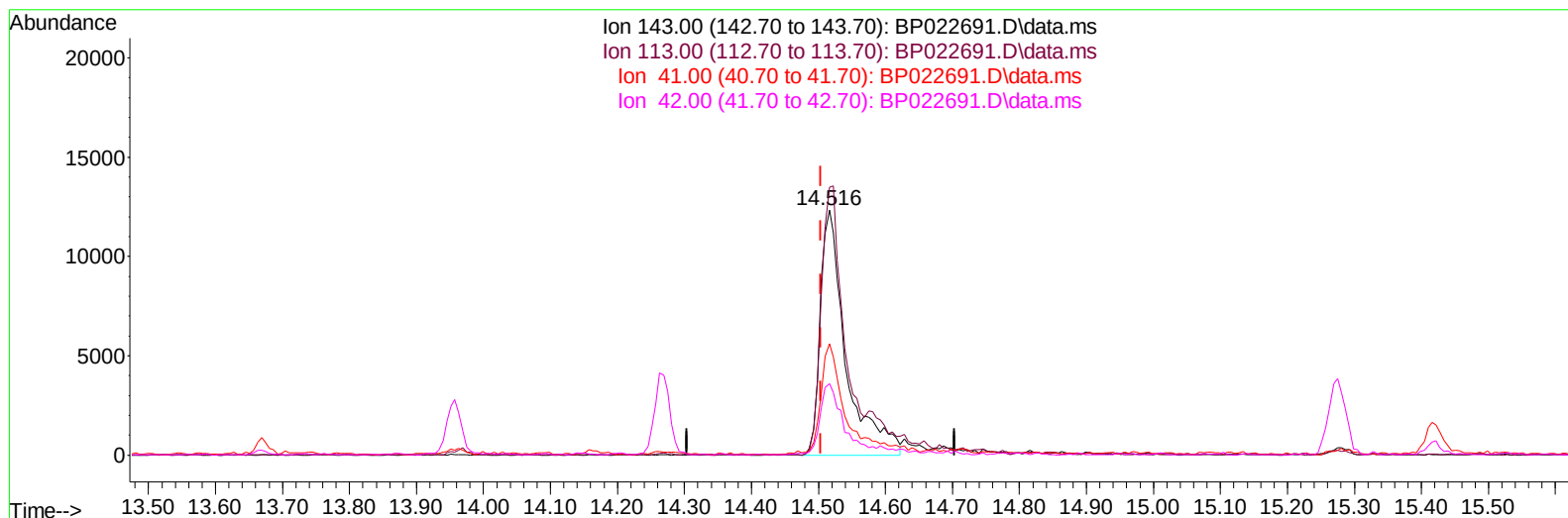
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TIC: BP022691.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.516min (+ 0.012) 12.39 ng/ul m

response 32792

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	109.33
41.00	35.90	45.33#
42.00	23.70	29.02#

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 10/29/2024
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Quant Time: Oct 29 01:23:58 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

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.652	152	66817	20.000	ng/ul	0.00
20) Naphthalene-d8	10.404	136	246291	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.269	164	198640	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.075	188	499462	20.000	ng/ul	0.00
79) Chrysene-d12	21.498	240	579251	20.000	ng/ul	0.00
88) Perylene-d12	24.745	264	733358	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.187	96	5256	3.057	ng/uL	-0.01
4) Pyridine-d5	3.599	84	69697	14.766	ng/ul	-0.01
7) Phenol-d5	6.840	99	92585	16.461	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	6.993	67	54460	16.474	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.193	132	65772	16.480	ng/ul	-0.01
15) 4-Methylphenol-d8	8.357	113	75429	16.805	ng/ul	0.00
21) Nitrobenzene-d5	8.793	128	32312	17.063	ng/ul	0.00
24) 2-Nitrophenol-d4	9.505	143	39529	18.030	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.046	165	79814	17.128	ng/ul	0.00
31) 4-Chloroaniline-d4	10.546	131	90174	16.377	ng/ul	-0.01
46) Dimethylphthalate-d6	13.669	166	307918	19.510	ng/ul	-0.01
49) Acenaphthylene-d8	13.957	160	302278	18.033	ng/ul	0.00
54) 4-Nitrophenol-d4	14.516	143	32792m	12.385	ng/ul	0.01
60) Fluorene-d10	15.275	176	269130	19.660	ng/ul	-0.02
65) 4,6-Dinitro-2-methylph...	15.422	200	42844	13.043	ng/ul	0.00
73) Anthracene-d10	17.169	188	471041	20.048	ng/ul	0.00
81) Pyrene-d10	19.557	212	604329	19.729	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.527	264	777734	19.858	ng/ul	0.00

Target Compounds Qvalue

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

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