

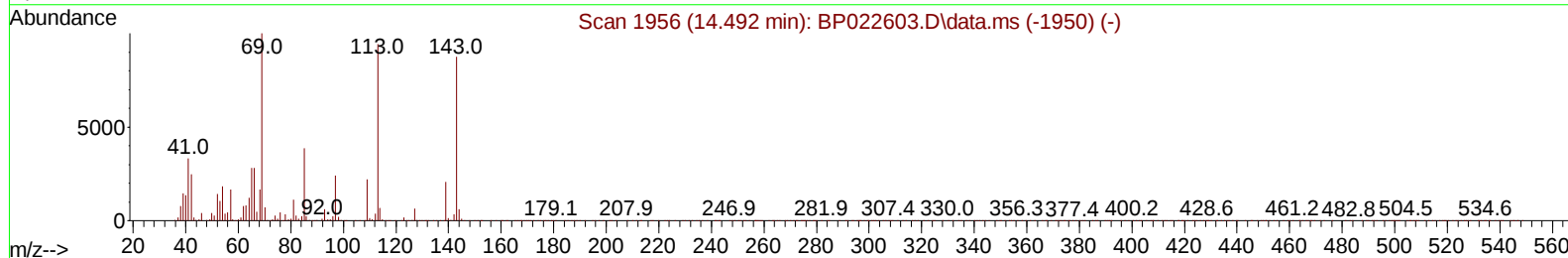
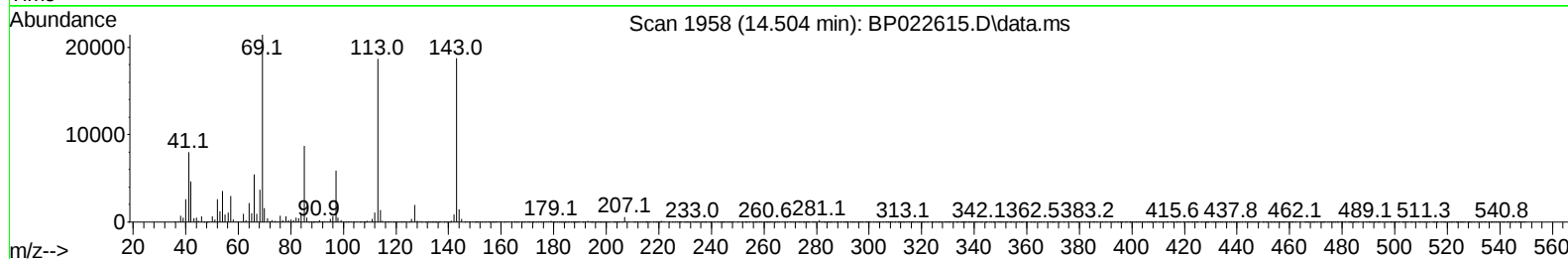
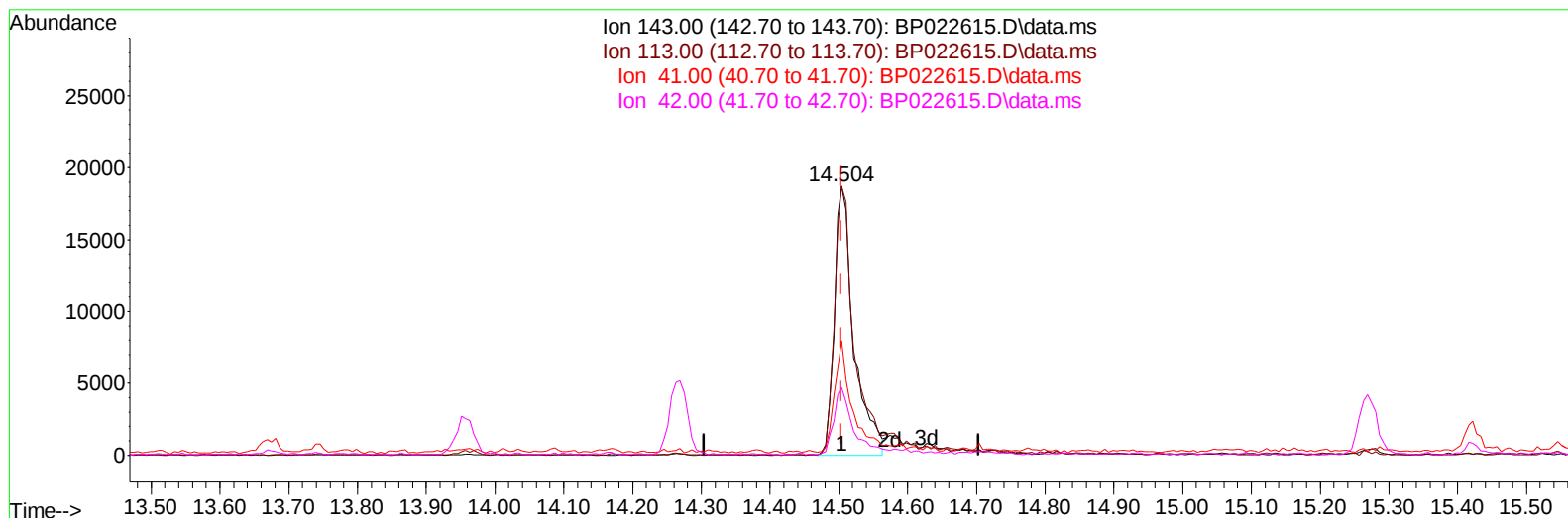
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102524\
Data File : BP022615.D
Acq On : 25 Oct 2024 17:17
Operator : RC/JU
Sample : P4436-12
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4F36

Manual IntegrationsAPPROVED

Quant Time: Oct 25 17:57:46 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Oct 22 05:59:59 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/26/2024
Supervised By :mohammad ahmed 10/29/2024



TIC: BP022615.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.504min (+ 0.000) 7.42 ng/u1

response 36643

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	99.91
41.00	35.90	42.56
42.00	23.70	25.02

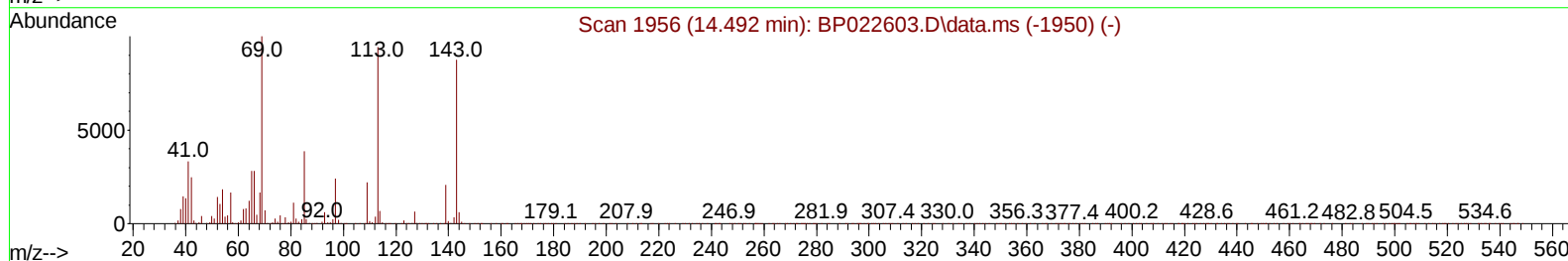
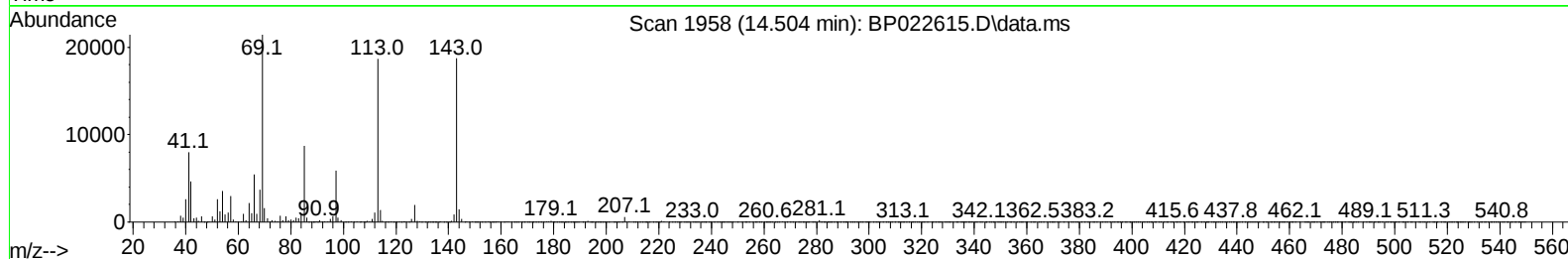
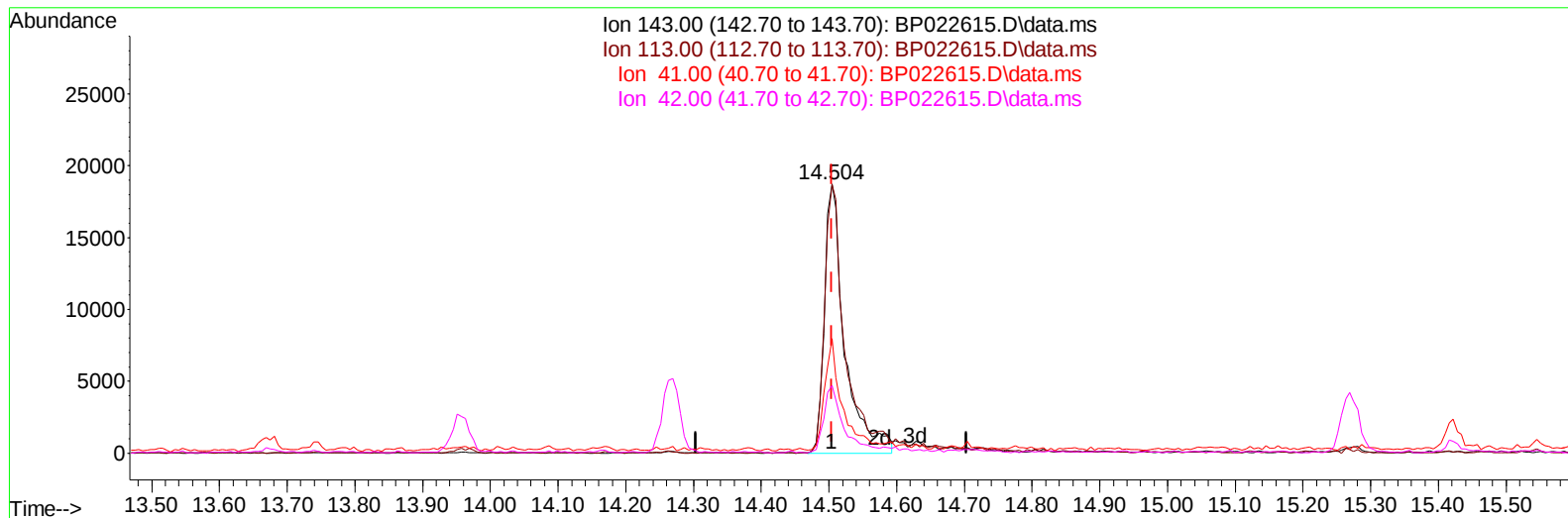
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TIC: BP022615.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.504min (+ 0.000) 7.84 ng/ul m

response 38703

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	99.91
41.00	35.90	42.56
42.00	23.70	25.02

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.652	152	127363	20.000	ng/ul	0.00
20) Naphthalene-d8	10.410	136	483795	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.269	164	370339	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.069	188	843900	20.000	ng/ul	-0.01
79) Chrysene-d12	21.516	240	856882	20.000	ng/ul	0.02
88) Perylene-d12	24.774	264	1002090	20.000	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.193	96	5593	1.706	ng/uL	0.00
4) Pyridine-d5	3.605	84	77707	8.637	ng/ul	0.00
7) Phenol-d5	6.846	99	115662	10.788	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.993	67	69930	11.098	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.199	132	91989	12.092	ng/ul	0.00
15) 4-Methylphenol-d8	8.357	113	93223	10.896	ng/ul	0.00
21) Nitrobenzene-d5	8.793	128	45288	12.175	ng/ul	0.00
24) 2-Nitrophenol-d4	9.504	143	55731	12.941	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.046	165	113028	12.348	ng/ul	0.00
31) 4-Chloroaniline-d4	10.546	131	123182	11.389	ng/ul	-0.01
46) Dimethylphthalate-d6	13.669	166	413439	14.051	ng/ul	-0.01
49) Acenaphthylene-d8	13.957	160	427435	13.677	ng/ul	0.00
54) 4-Nitrophenol-d4	14.504	143	38703m	7.841	ng/ul	0.00
60) Fluorene-d10	15.275	176	376344	14.746	ng/ul	-0.02
65) 4,6-Dinitro-2-methylph...	15.422	200	48739	8.781	ng/ul	0.00
73) Anthracene-d10	17.175	188	625310	15.751	ng/ul	0.00
81) Pyrene-d10	19.551	212	771934	17.035	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.545	264	874546	16.342	ng/ul	0.02

Target Compounds Qvalue

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

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