

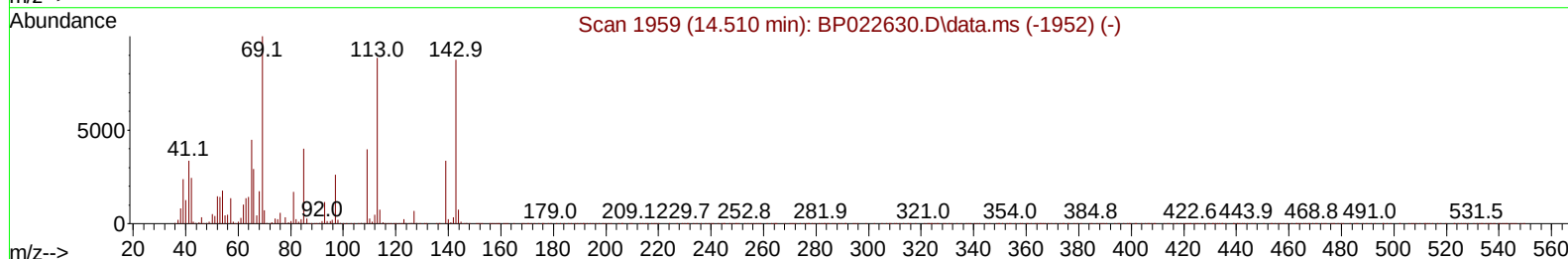
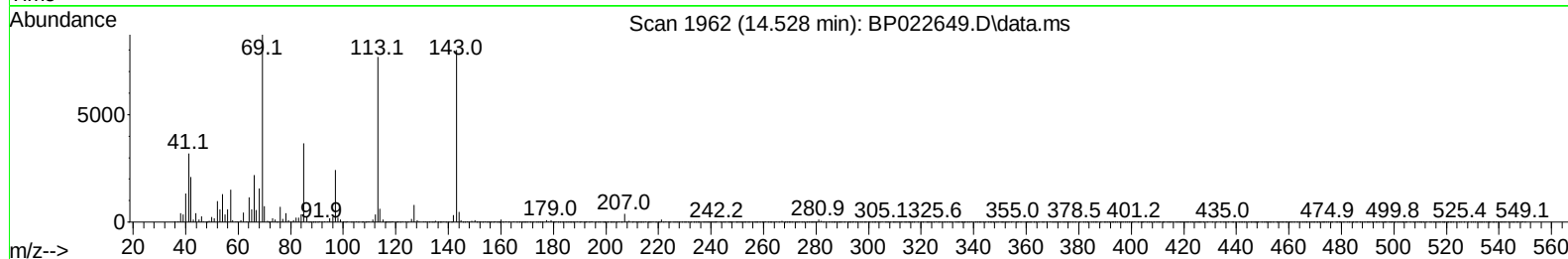
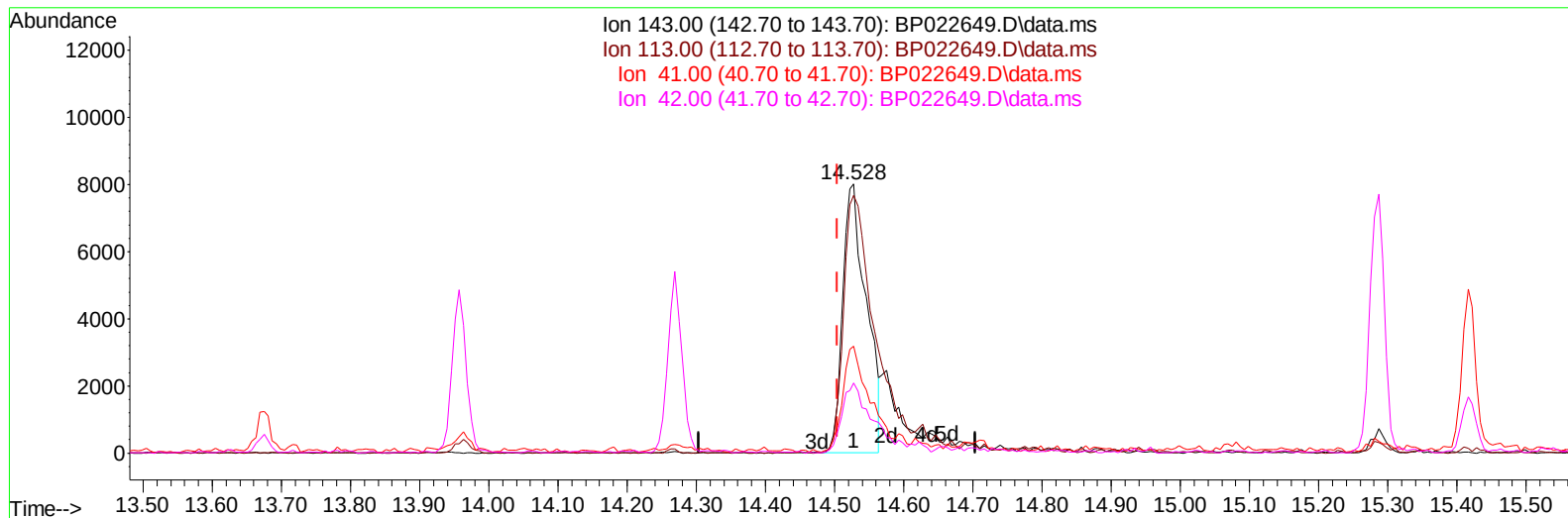
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102524\
Data File : BP022649.D
Acq On : 26 Oct 2024 19:23
Operator : RC/JU
Sample : P4532-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCNX0

Manual IntegrationsAPPROVED

Quant Time: Oct 28 00:25:28 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: BP022649.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.528min (+ 0.024) 5.56 ng/u1

response 18941

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	95.93
41.00	35.90	39.74
42.00	23.70	26.05

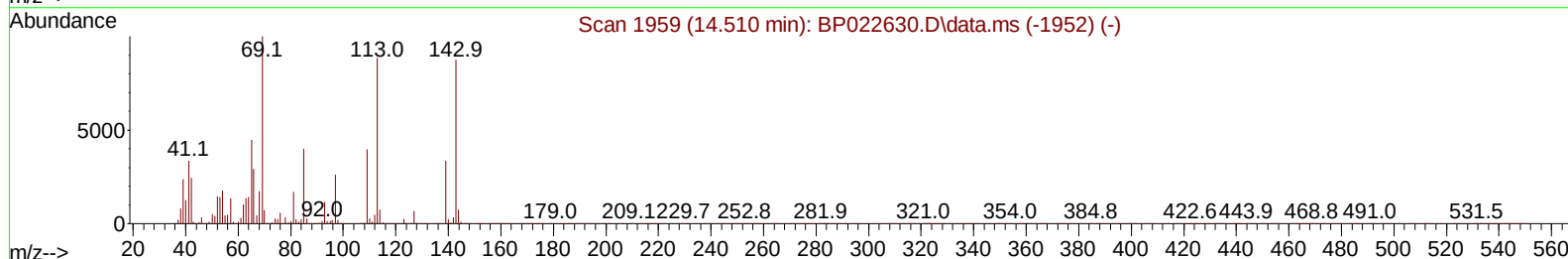
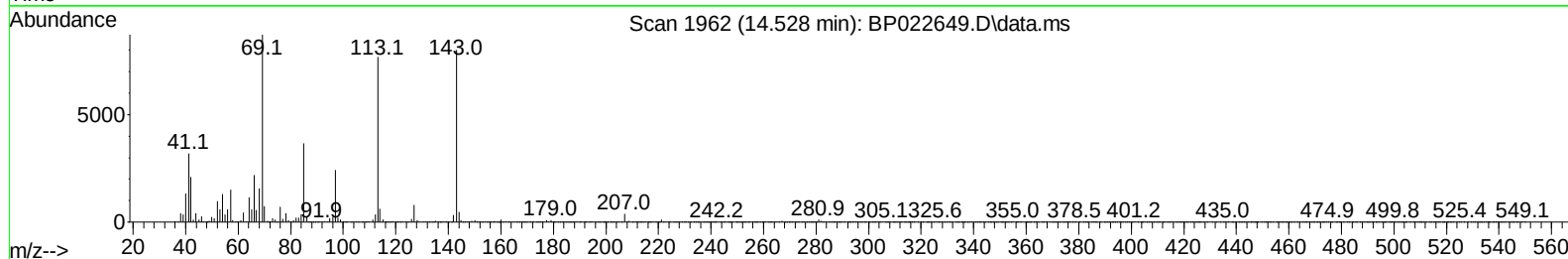
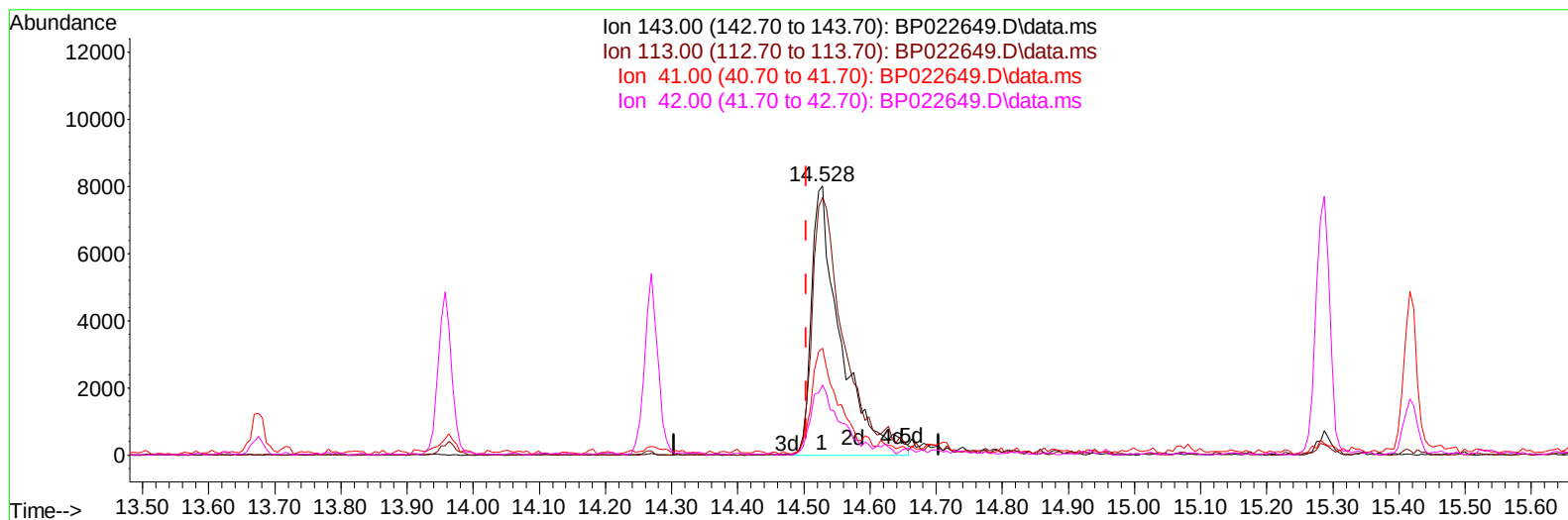
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TIC: BP022649.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.528min (+ 0.024) 7.19 ng/ul m

response 24501

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	95.93
41.00	35.90	39.74
42.00	23.70	26.05

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 Sample : P4532-09
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Instrument :
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 GCNX0

Manual IntegrationsAPPROVED

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

Quant Time: Oct 28 00:26:10 2024
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 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	87044	20.000	ng/ul	0.00
20) Naphthalene-d8	10.405	136	328066	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.269	164	255681	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.075	188	613709	20.000	ng/ul	0.00
79) Chrysene-d12	21.498	240	708434	20.000	ng/ul	0.00
88) Perylene-d12	24.757	264	828765	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.187	96	8135	3.632	ng/uL	-0.01
4) Pyridine-d5	3.611	84	38104	6.197	ng/ul	0.00
7) Phenol-d5	6.846	99	50771	6.929	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.993	67	103231	23.971	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.193	132	121829	23.433	ng/ul	-0.01
15) 4-Methylphenol-d8	8.358	113	95937	16.407	ng/ul	0.00
21) Nitrobenzene-d5	8.793	128	65958	26.148	ng/ul	0.00
24) 2-Nitrophenol-d4	9.505	143	78741	26.962	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.040	165	159679	25.726	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.546	131	157213	21.435	ng/ul	-0.01
46) Dimethylphthalate-d6	13.675	166	576396	28.374	ng/ul	0.00
49) Acenaphthylene-d8	13.957	160	581409	26.947	ng/ul	0.00
54) 4-Nitrophenol-d4	14.528	143	24501m	7.189	ng/ul	0.02
60) Fluorene-d10	15.287	176	514967	29.226	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.416	200	105370	26.106	ng/ul	0.00
73) Anthracene-d10	17.163	188	954222	33.052	ng/ul	0.00
81) Pyrene-d10	19.551	212	1245280	33.240	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.533	264	1491678	33.702	ng/ul	0.00

Target Compounds Qvalue

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

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