

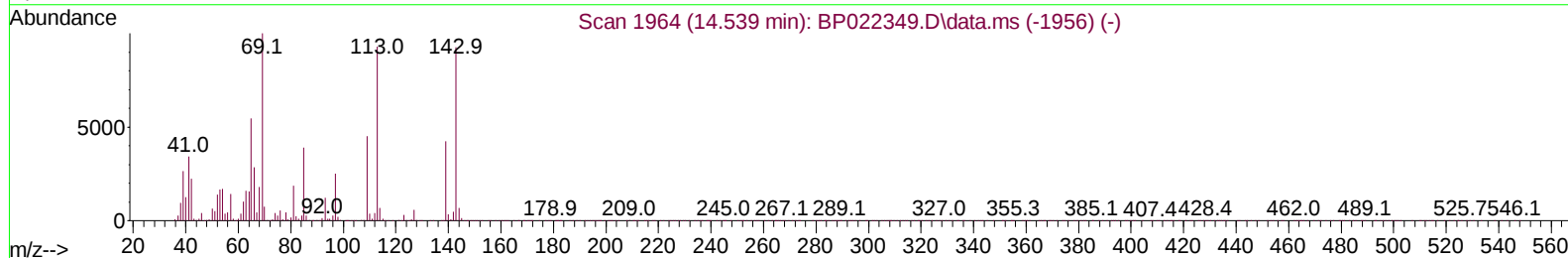
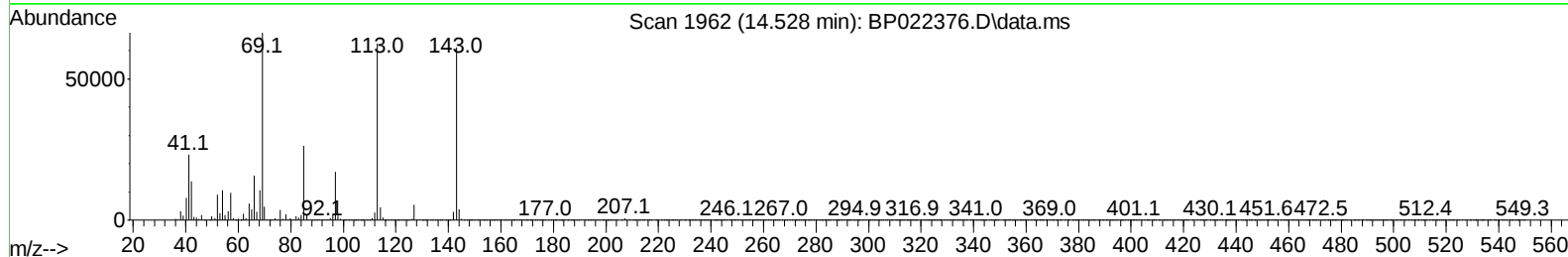
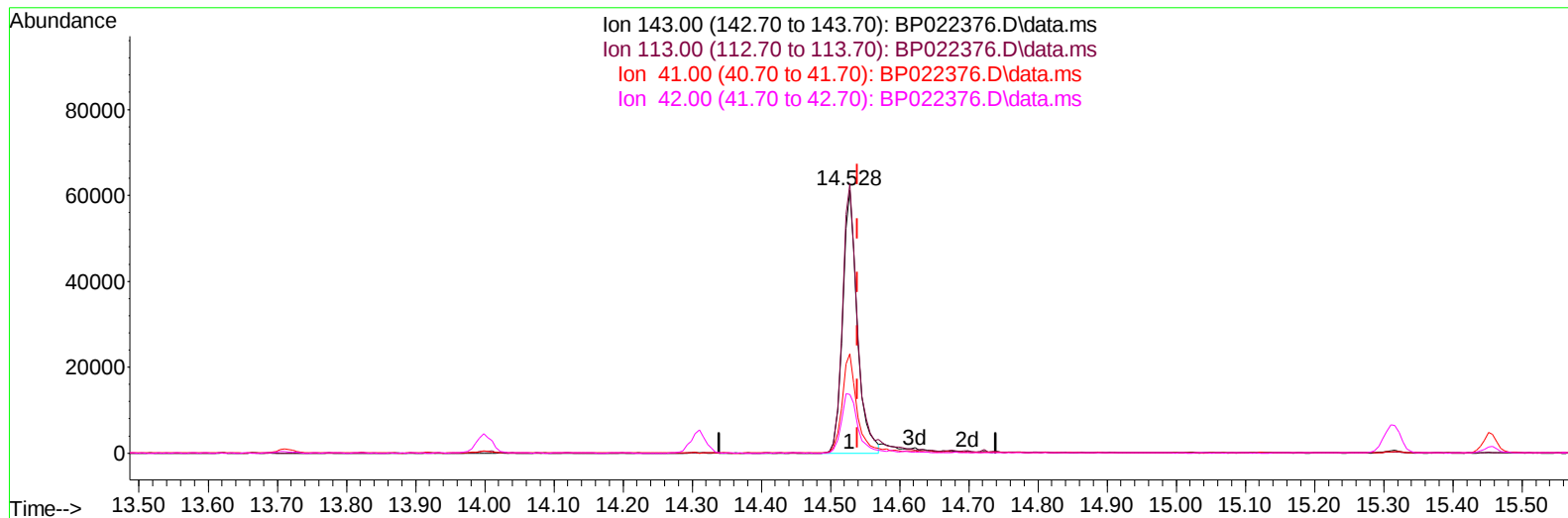
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101424\
Data File : BP022376.D
Acq On : 15 Oct 2024 05:37
Operator : RC/JU
Sample : PB164091BL
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK091

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 10/16/2024
Supervised By :mohammad ahmed 10/22/2024

Quant Time: Oct 15 06:08:54 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Oct 14 11:01:55 2024
Response via : Initial Calibration



TIC: BP022376.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.528min (-0.012) 23.10 ng/u1

response 89702

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	102.27
41.00	35.90	37.66
42.00	23.70	22.37

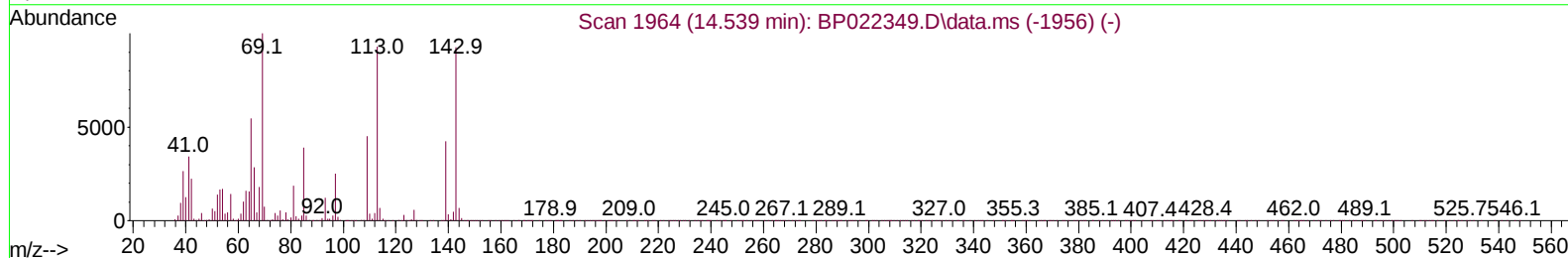
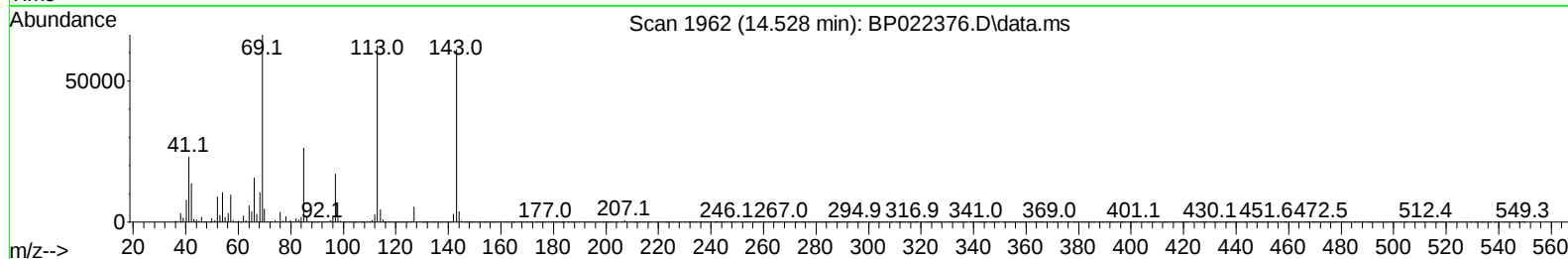
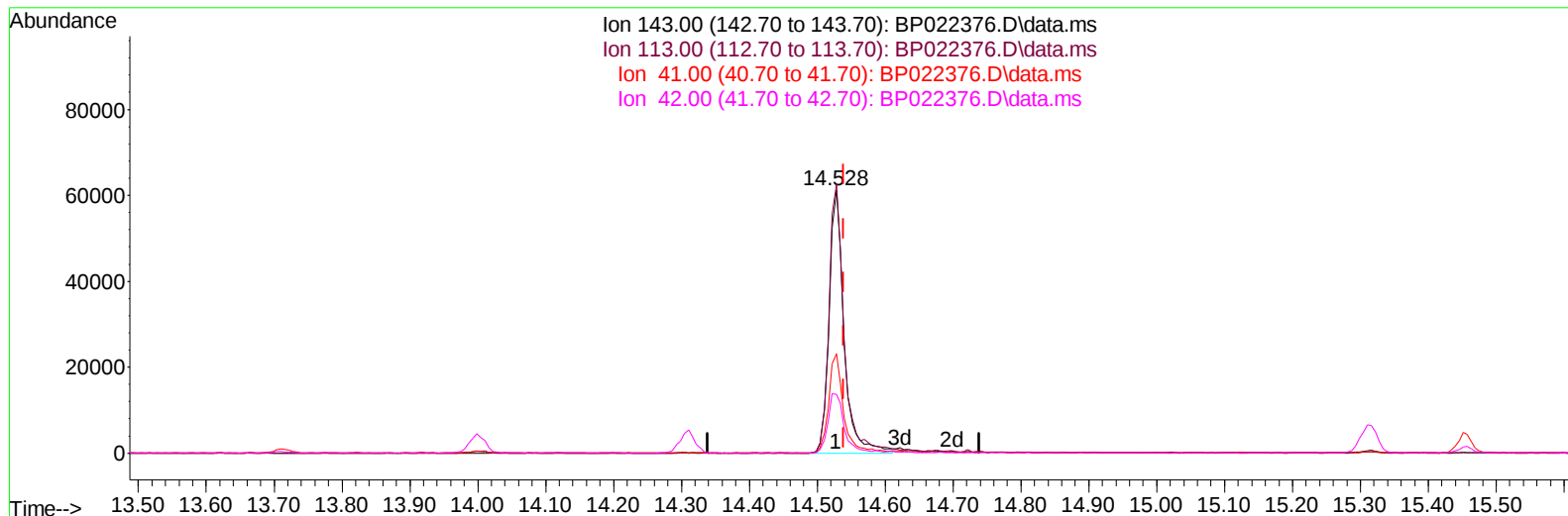
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TIC: BP022376.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.528min (-0.012) 24.02 ng/ul m

response 93304

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	102.27
41.00	35.90	37.66
42.00	23.70	22.37

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Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 10/16/2024
 Supervised By :mohammad ahmed 10/22/2024

Quant Time: Oct 15 06:10:57 2024
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 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Oct 14 11:01:55 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.699	152	96262	20.000	ng/ul	0.00
20) Naphthalene-d8	10.458	136	370031	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.310	164	291387	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.110	188	748332	20.000	ng/ul	0.00
79) Chrysene-d12	21.551	240	799306	20.000	ng/ul	0.00
88) Perylene-d12	24.863	264	915668	20.000	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.223	96	10853	4.381	ng/uL	0.00
4) Pyridine-d5	3.634	84	164243	24.152	ng/ul	0.00
7) Phenol-d5	6.887	99	199932	24.674	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.040	67	117148	24.598	ng/ul	0.00
11) 2-Chlorophenol-d4	7.240	132	153804	26.750	ng/ul	0.00
15) 4-Methylphenol-d8	8.399	113	162586	25.142	ng/ul	0.00
21) Nitrobenzene-d5	8.840	128	74012	26.013	ng/ul	0.00
24) 2-Nitrophenol-d4	9.552	143	89482	27.165	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.093	165	181225	25.886	ng/ul	0.01
31) 4-Chloroaniline-d4	10.587	131	209073	25.273	ng/ul	0.00
46) Dimethylphthalate-d6	13.716	166	581644	25.124	ng/ul	0.00
49) Acenaphthylene-d8	13.999	160	638833	25.980	ng/ul	0.00
54) 4-Nitrophenol-d4	14.528	143	93304m	24.023	ng/ul	-0.01
60) Fluorene-d10	15.310	176	530612	26.424	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.457	200	101466	20.616	ng/ul	0.01
73) Anthracene-d10	17.216	188	940762	26.724	ng/ul	0.01
81) Pyrene-d10	19.598	212	1207803	28.574	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.621	264	1358348	27.777	ng/ul	-0.01

Target Compounds Qvalue

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

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