

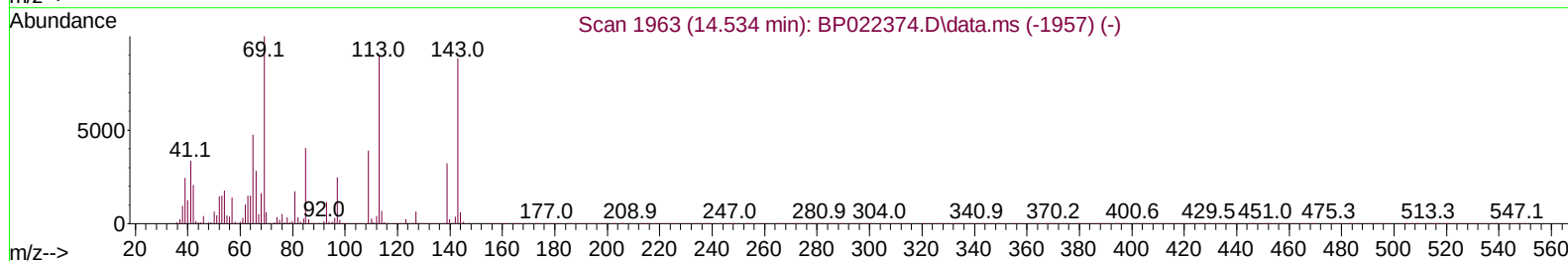
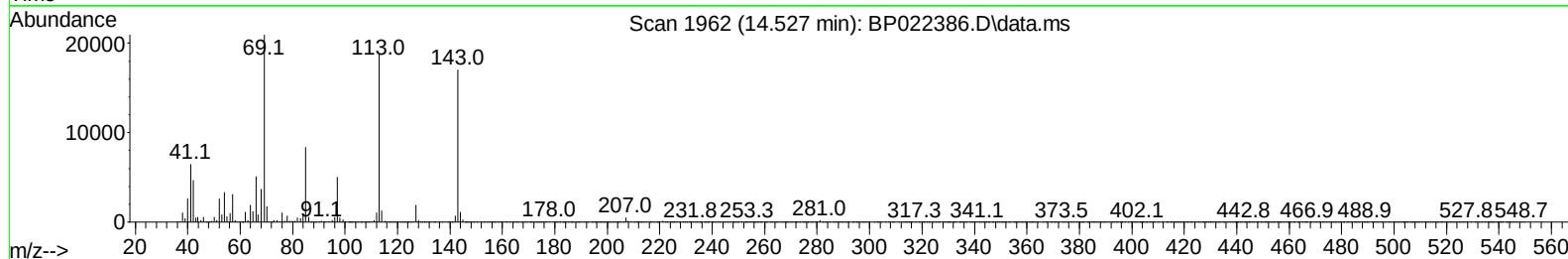
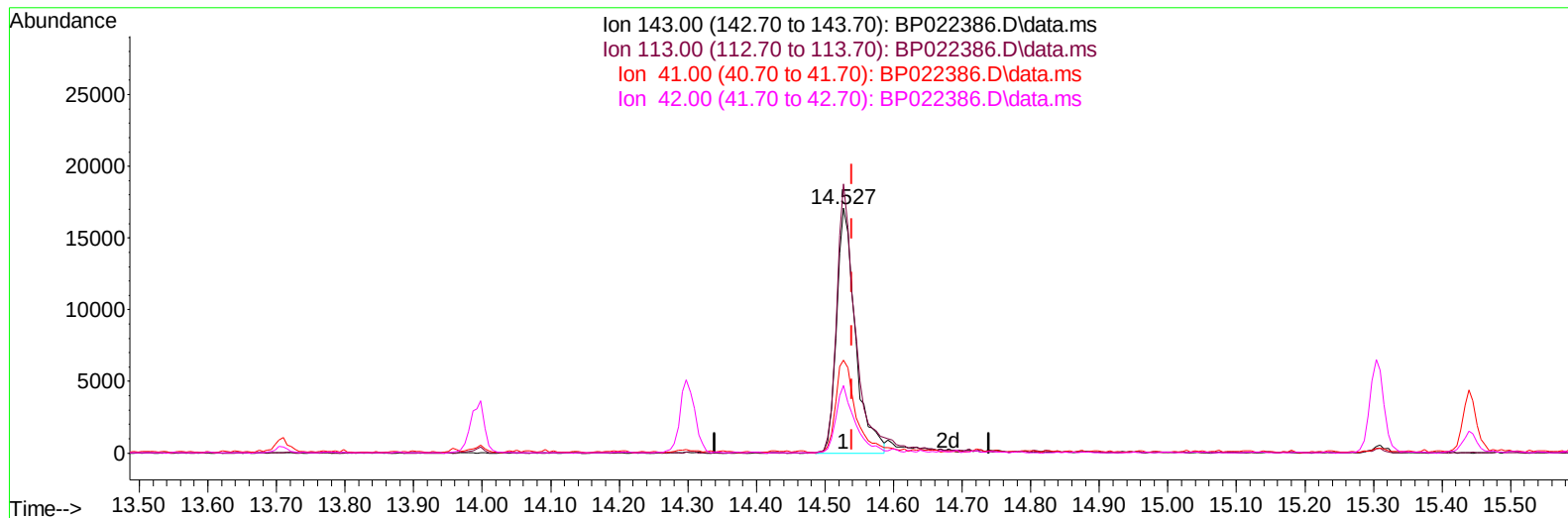
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101424\
Data File : BP022386.D
Acq On : 15 Oct 2024 12:24
Operator : RC/JU
Sample : P4349-08
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EOAG6

Manual IntegrationsAPPROVED

Quant Time: Oct 16 01:25:09 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

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



TIC: BP022386.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.527min (-0.012) 9.45 ng/u1

response 32302

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	109.79
41.00	35.90	37.90
42.00	23.70	27.58

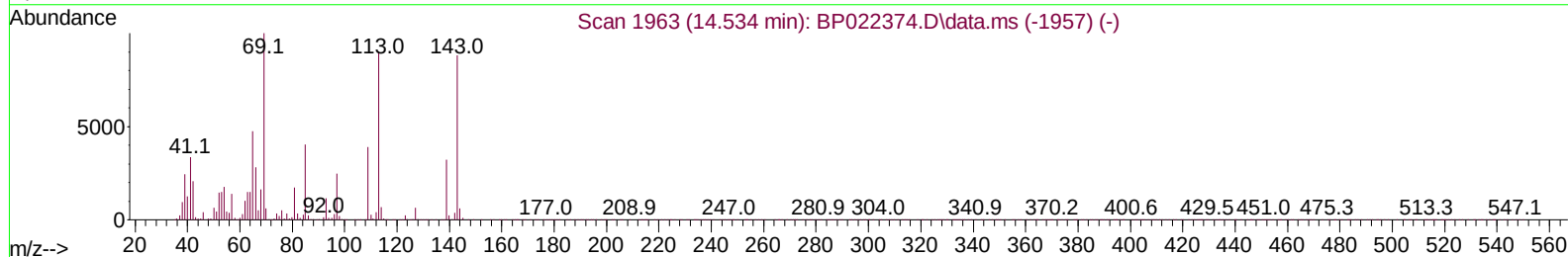
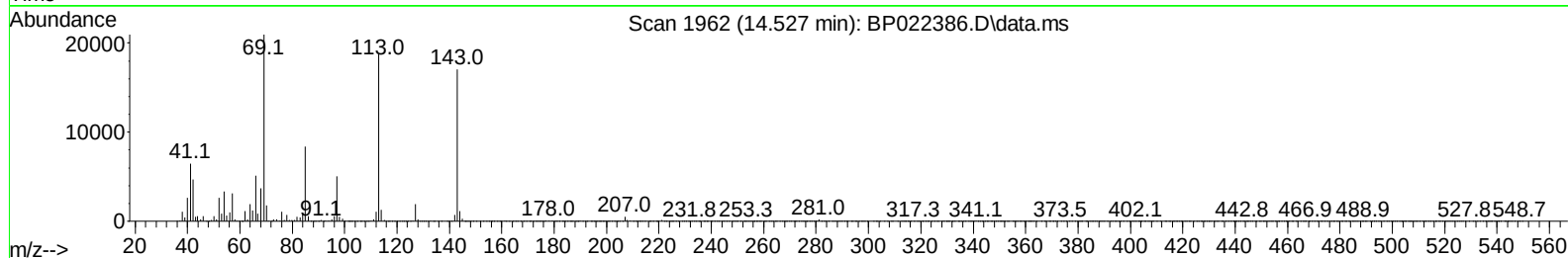
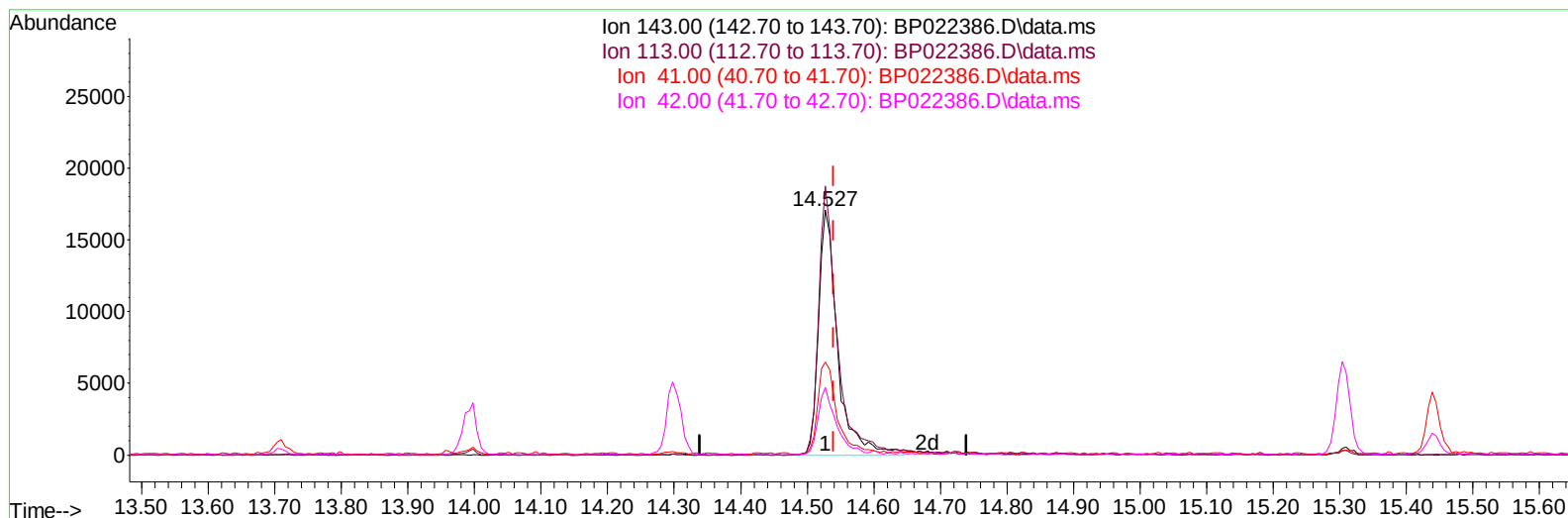
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TIC: BP022386.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.527min (-0.012) 9.93 ng/ul m

response 33945

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	109.79
41.00	35.90	37.90
42.00	23.70	27.58

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 10/16/2024
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Quant Time: Oct 16 01:26:58 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

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.698	152	83660	20.000	ng/ul	0.00
20) Naphthalene-d8	10.451	136	314983	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.298	164	256362	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.104	188	660665	20.000	ng/ul	0.00
79) Chrysene-d12	21.545	240	763336	20.000	ng/ul	# 0.00
88) Perylene-d12	24.821	264	909076	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.222	96	8296	3.853	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	6.887	99	63818	9.062	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.040	67	88024	21.267	ng/ul	0.00
11) 2-Chlorophenol-d4	7.245	132	115077	23.029	ng/ul	0.00
15) 4-Methylphenol-d8	8.398	113	92738	16.501	ng/ul	0.00
21) Nitrobenzene-d5	8.839	128	57161	23.602	ng/ul	0.00
24) 2-Nitrophenol-d4	9.551	143	69196	24.678	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.081	165	143718	24.116	ng/ul	0.00
31) 4-Chloroaniline-d4	10.598	131	13566	1.926	ng/ul	0.01
46) Dimethylphthalate-d6	13.710	166	472429	23.194	ng/ul	0.00
49) Acenaphthylene-d8	13.992	160	441831	20.423	ng/ul	0.00
54) 4-Nitrophenol-d4	14.527	143	33945m	9.934	ng/ul	-0.01
60) Fluorene-d10	15.304	176	433997	24.565	ng/ul	-0.02
65) 4,6-Dinitro-2-methylph...	15.445	200	91070	20.959	ng/ul	0.00
73) Anthracene-d10	17.204	188	773780	24.897	ng/ul	0.00
81) Pyrene-d10	19.580	212	1056409	26.170	ng/ul	-0.02
92) Benzo(a)pyrene-d12	24.603	264	1148425	23.655	ng/ul	-0.03

Target Compounds Qvalue

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

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