

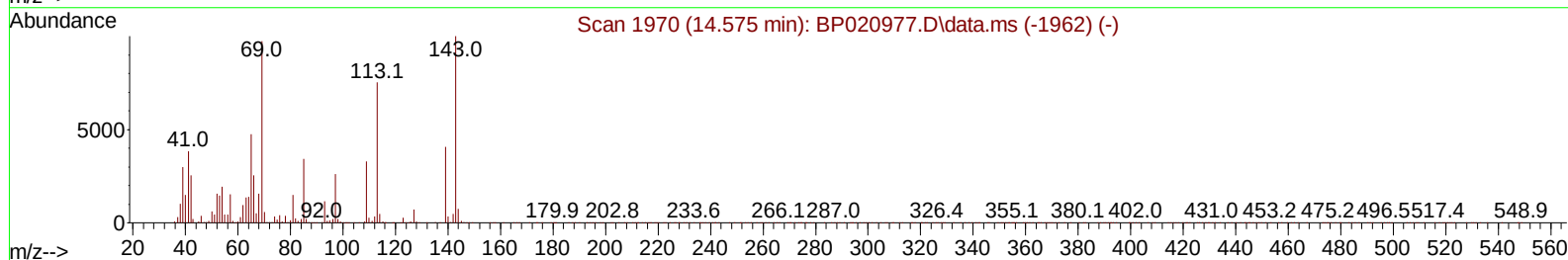
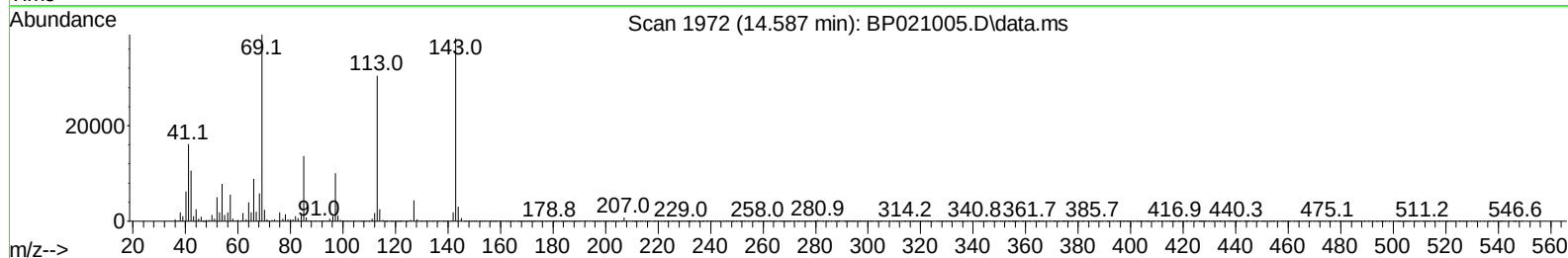
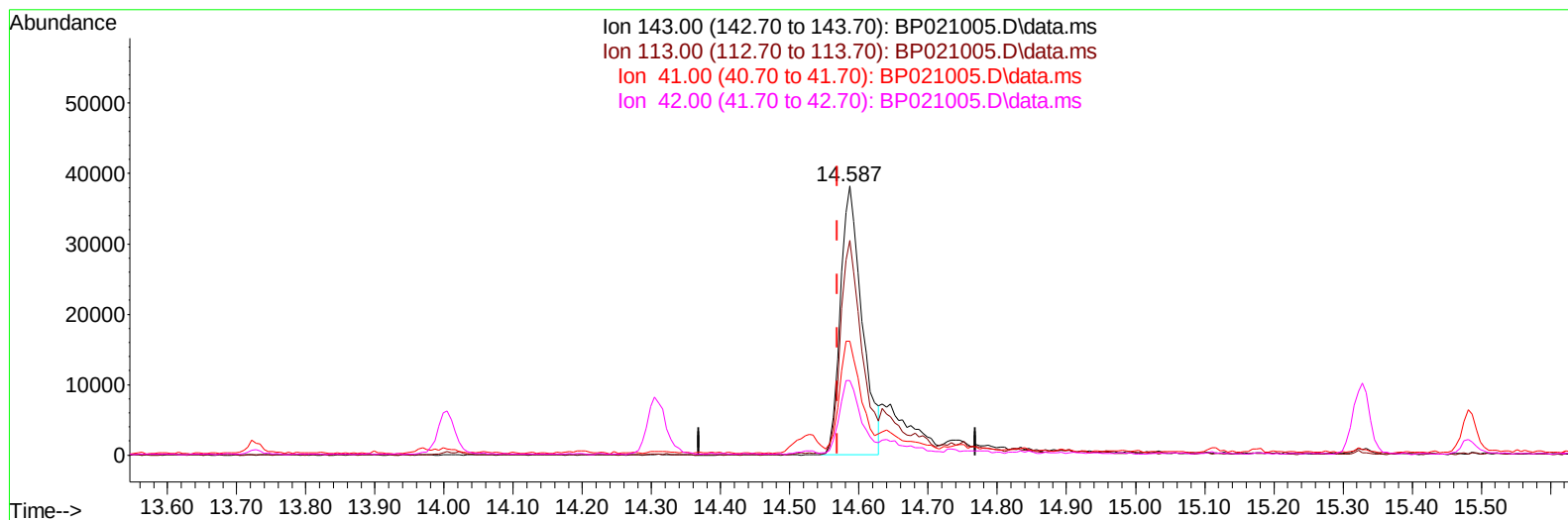
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071524\
Data File : BP021005.D
Acq On : 17 Jul 2024 01:58
Operator : MA/JU
Sample : P3178-07
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4BT2

Manual IntegrationsAPPROVED

Quant Time: Jul 17 02:59:53 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jul 15 12:22:22 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 07/18/2024
Supervised By :mohammad ahmed 07/18/2024



TIC: BP021005.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.587min (+ 0.018) 15.13 ng/ul

response 82971

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	79.82
41.00	37.90	42.22
42.00	27.10	27.68

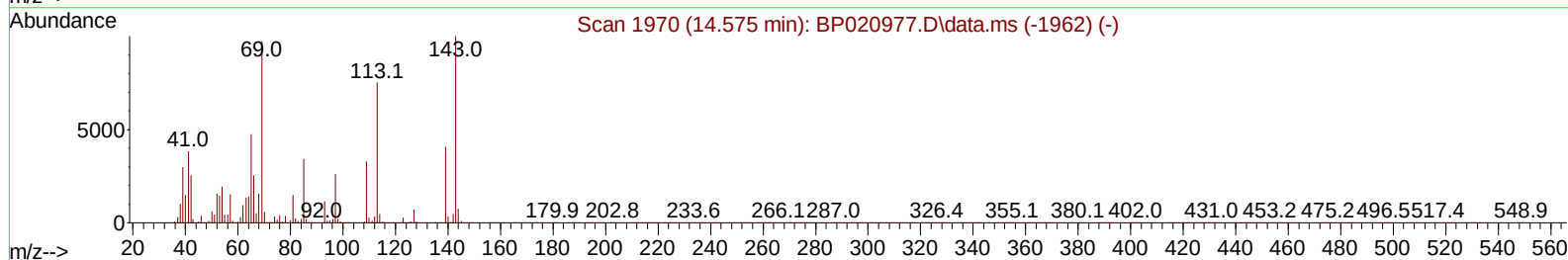
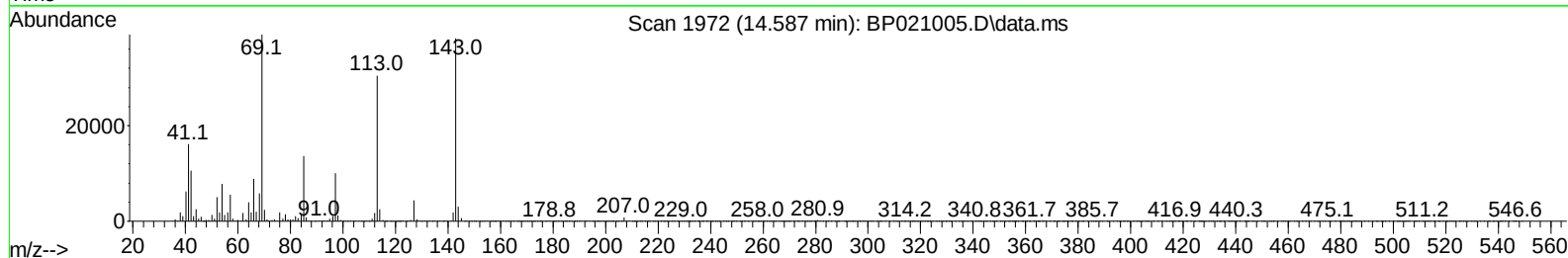
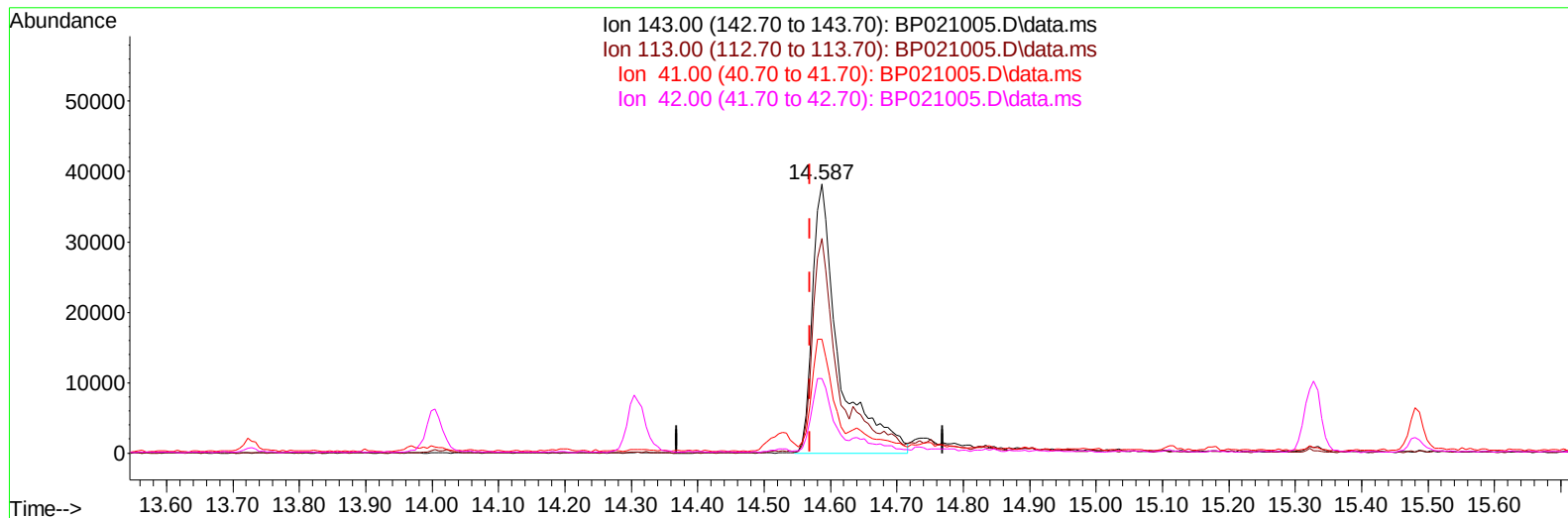
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TIC: BP021005.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.587min (+ 0.018) 19.21 ng/ul m

response 105349

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	79.82
41.00	37.90	42.22
42.00	27.10	27.68

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Manual IntegrationsAPPROVED

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Quant Time: Jul 17 03:22:17 2024
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 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jul 15 12:22:22 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.699	152	198319	20.000	ng/ul	0.01
20) Naphthalene-d8	10.452	136	792619	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.304	164	530219	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.116	188	1183625	20.000	ng/ul	0.00
79) Chrysene-d12	21.569	240	1236355	20.000	ng/ul	0.02
88) Perylene-d12	24.863	264	1428737	20.000	ng/ul	0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.240	96	13934	3.197	ng/uL	0.00
4) Pyridine-d5	3.658	84	193326	15.323	ng/ul	0.00
7) Phenol-d5	6.899	99	297972	18.429	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.046	67	183669	18.745	ng/ul	0.01
11) 2-Chlorophenol-d4	7.246	132	251610	18.888	ng/ul	0.01
15) 4-Methylphenol-d8	8.411	113	245009	18.245	ng/ul	0.02
21) Nitrobenzene-d5	8.852	128	120538	19.579	ng/ul	0.02
24) 2-Nitrophenol-d4	9.558	143	137588	19.526	ng/ul	0.02
28) 2,4-Dichlorophenol-d3	10.099	165	243499	19.111	ng/ul	0.01
31) 4-Chloroaniline-d4	10.605	131	325452	19.281	ng/ul	0.02
46) Dimethylphthalate-d6	13.728	166	813705	21.111	ng/ul	0.02
49) Acenaphthylene-d8	14.004	160	907940	20.978	ng/ul	0.00
54) 4-Nitrophenol-d4	14.587	143	105349m	19.210	ng/ul	0.02
60) Fluorene-d10	15.328	176	702217	22.207	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.481	200	102829	14.358	ng/ul	0.02
73) Anthracene-d10	17.222	188	1177391	22.396	ng/ul	0.00
81) Pyrene-d10	19.610	212	1469319	19.223	ng/ul	0.01
92) Benzo(a)pyrene-d12	24.627	264	1557748	22.107	ng/ul	0.02

Target Compounds Qvalue

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

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