

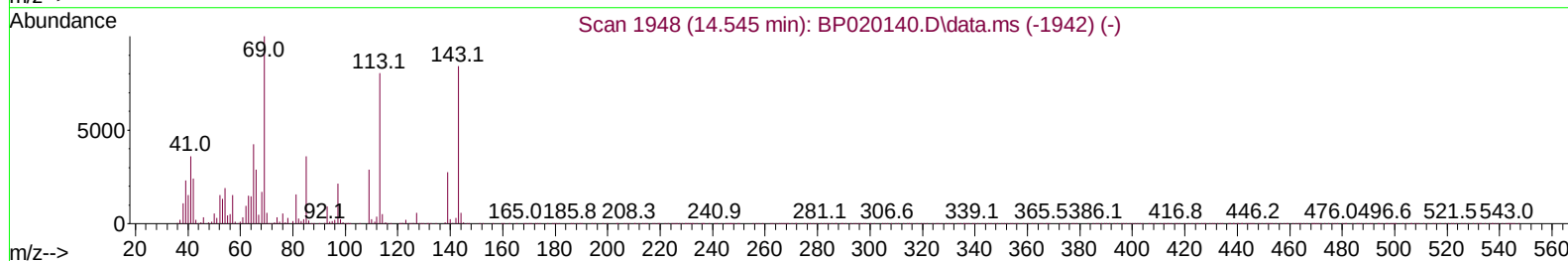
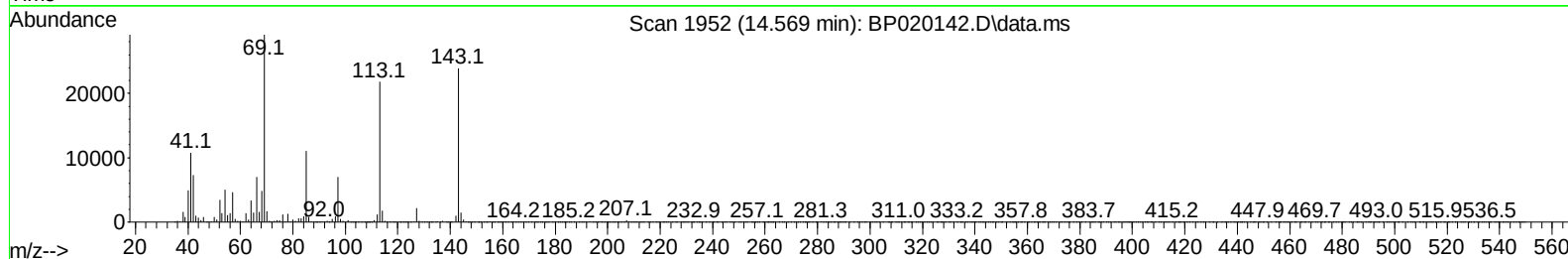
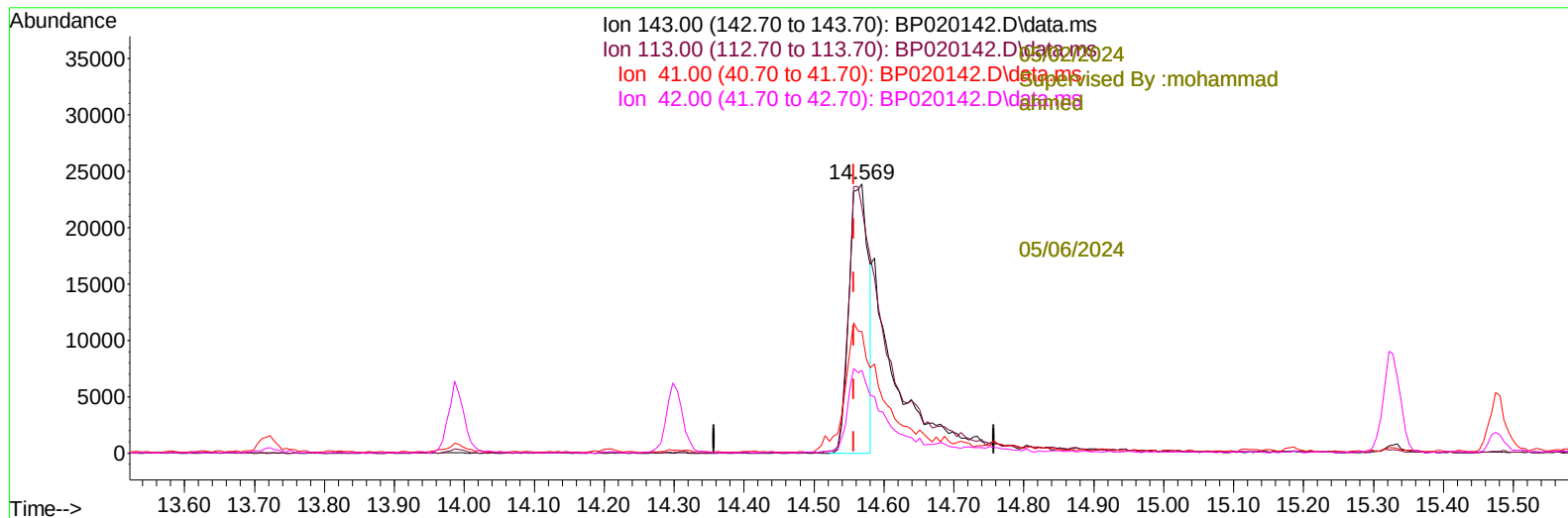
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050224\
Data File : BP020142.D
Acq On : 02 May 2024 10:45
Operator : MA/JU
Sample : P2105-08
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AD3

Manual IntegrationsAPPROVED

Quant Time: May 02 11:21:43 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed May 01 17:34:36 2024
Response via : Initial Calibration

Reviewed By :Jagrut
Upadhyay



TIC: BP020142.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.569min (+ 0.012) 15.81 ng/ul

response 46132

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	101.40	91.25
41.00	50.30	45.26
42.00	32.50	30.69

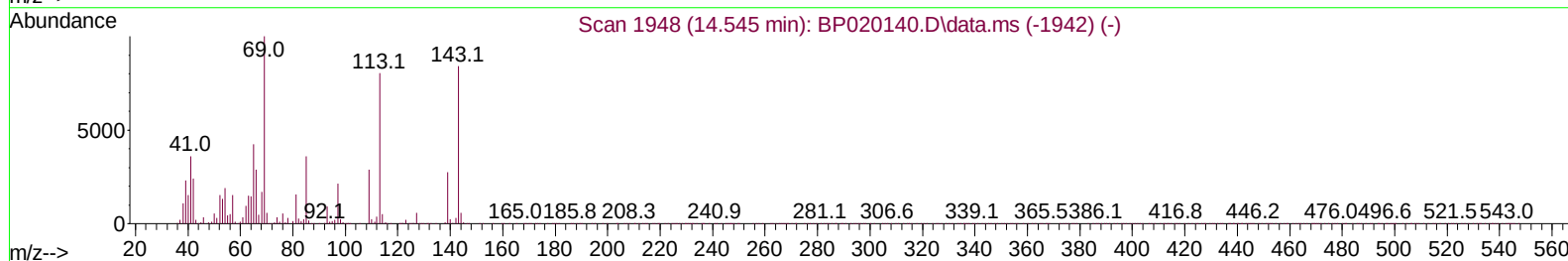
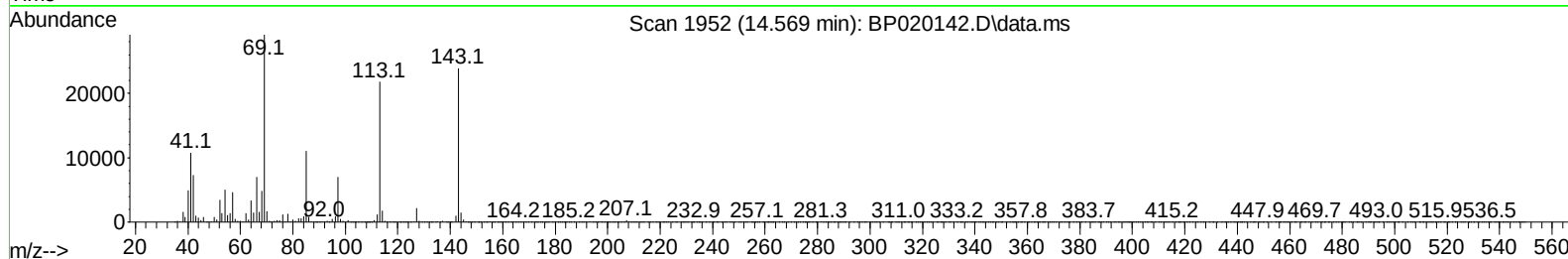
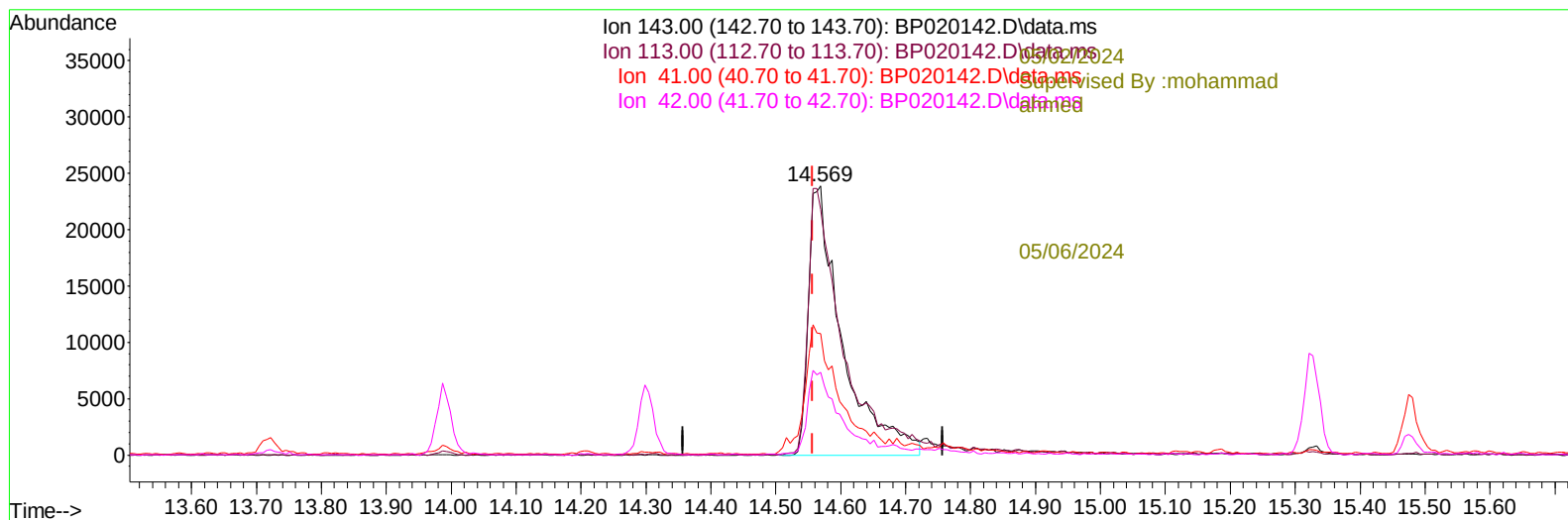
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Upadhyay



TIC: BP020142.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.569min (+ 0.012) 29.62 ng/ul m

response 86450

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	101.40	91.25
41.00	50.30	45.26
42.00	32.50	30.69

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Instrument :
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Manual IntegrationsAPPROVED

Quant Time: May 02 13:02:31 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed May 01 17:34:36 2024
 Response via : Initial Calibration

Reviewed By :Jagrut
 Upadhyay

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
-----05/02/2024						
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.628	152	82726	20.000	ng/ul	0.00
20) Naphthalene-d8	10.398	136	354482	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.298	164	260022	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.139	188	602585	20.000	ng/ul	0.00
79) Chrysene-d12	21.639	240	638711	20.000	ng/ul	0.00
88) Perylene-d12	25.004	264	686969	20.000	ng/ul	-0.01
-----06/2024						
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.222	96	9643	4.923	ng/uL	0.00
4) Pyridine-d5	3.628	84	79661	13.996	ng/ul	0.00
7) Phenol-d5	6.816	99	191565	26.964	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.987	67	123445	28.604	ng/ul	0.00
11) 2-Chlorophenol-d4	7.169	132	143452	28.245	ng/ul	0.00
15) 4-Methylphenol-d8	8.340	113	157760	27.466	ng/ul	0.00
21) Nitrobenzene-d5	8.787	128	75218	28.336	ng/ul	0.00
24) 2-Nitrophenol-d4	9.504	143	89672	29.500	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.040	165	152437	27.480	ng/ul	0.00
31) 4-Chloroaniline-d4	10.569	131	99800	12.855	ng/ul	0.01
46) Dimethylphthalate-d6	13.722	166	596043	31.390	ng/ul	0.00
49) Acenaphthylene-d8	13.986	160	626448	29.865	ng/ul	0.00
54) 4-Nitrophenol-d4	14.569	143	86450m	29.623	ng/ul	0.01
60) Fluorene-d10	15.322	176	501614	31.672	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.481	200	91528	23.936	ng/ul	0.00
73) Anthracene-d10	17.239	188	811119	31.115	ng/ul	0.00
81) Pyrene-d10	19.639	212	1043296	27.944	ng/ul	-0.01
92) Benzo(a)pyrene-d12	24.762	264	1005769	30.280	ng/ul	-0.02

Target Compounds Qvalue

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

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