

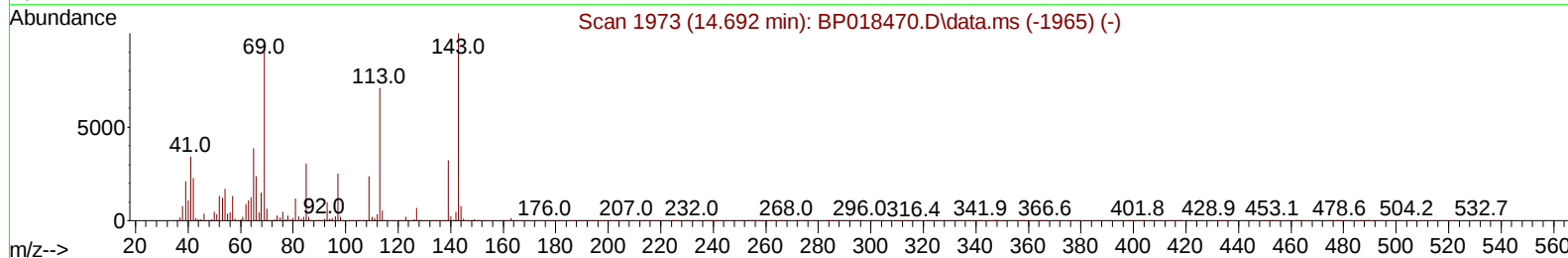
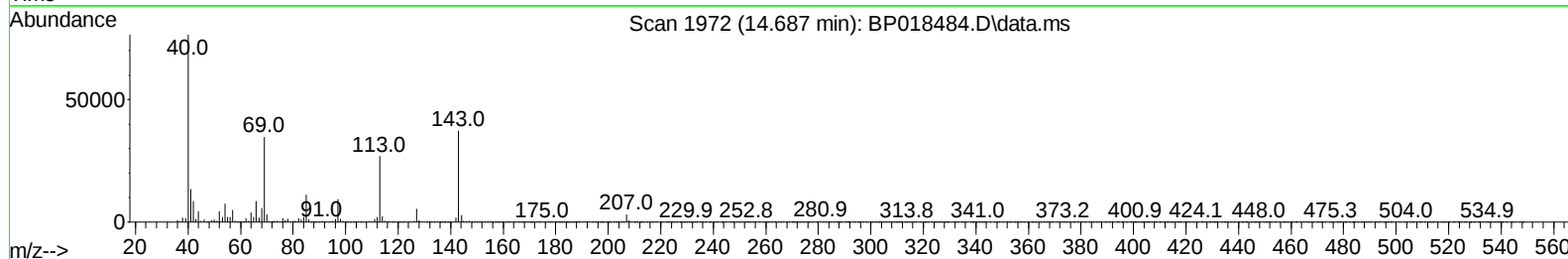
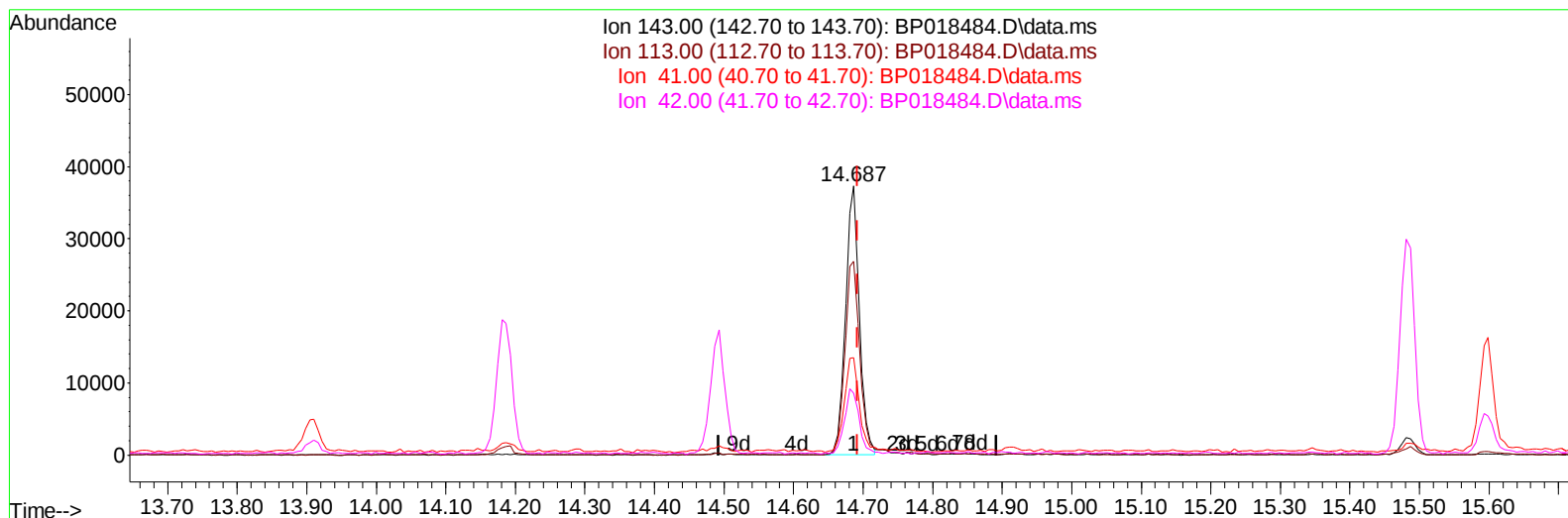
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120123\
 Data File : BP018484.D
 Acq On : 01 Dec 2023 19:37
 Operator : MA/JU
 Sample : 05332-08
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 YCGA7

Manual IntegrationsAPPROVED

Quant Time: Dec 01 22:37:16 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 01 21:56:43 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 12/04/2023
 Supervised By :mohammad ahmed 12/04/2023



TIC: BP018484.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.687min (-0.006) 3.73 ng/u1

response 51860

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	70.80	71.99
41.00	34.60	36.27
42.00	23.00	23.03

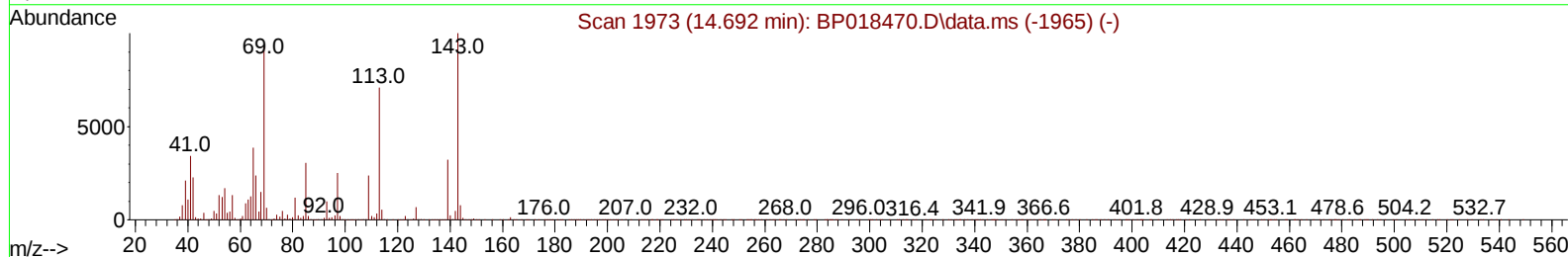
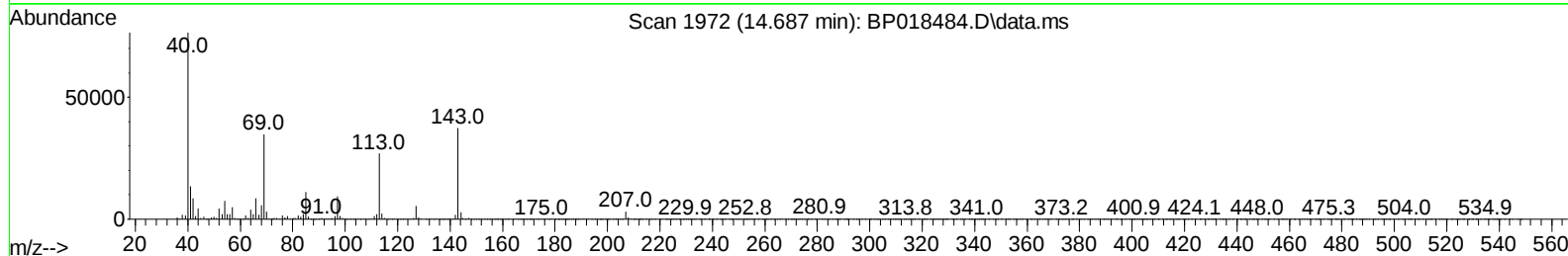
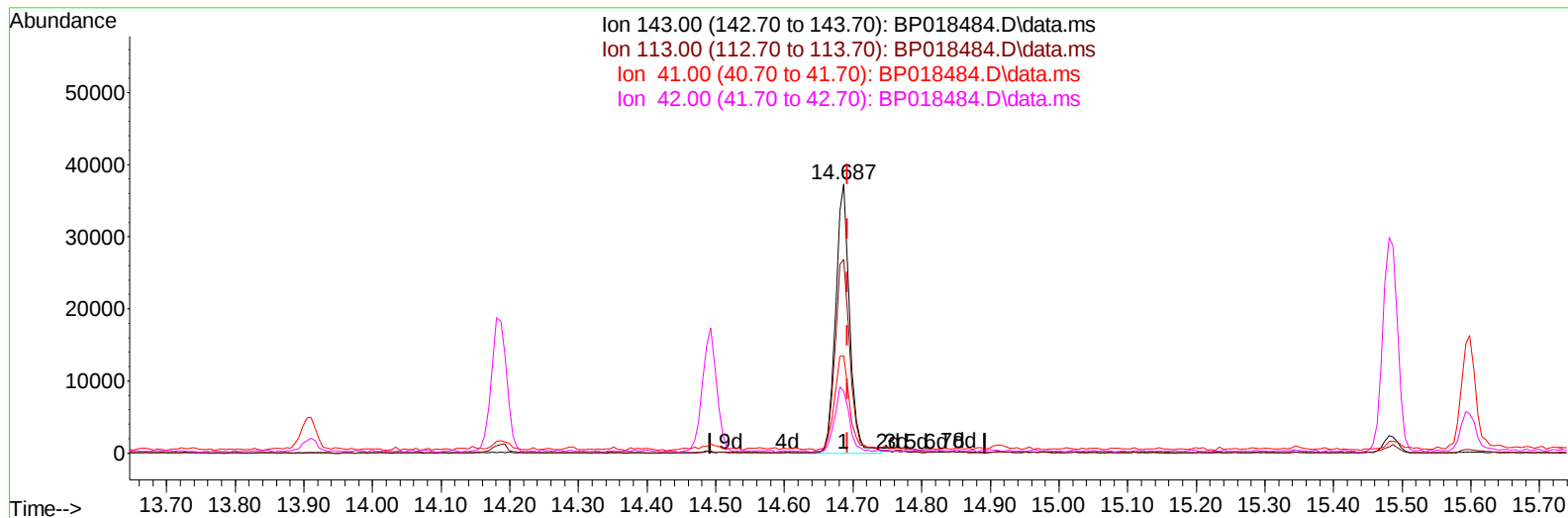
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TIC: BP018484.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.687min (-0.006) 3.80 ng/ul m

response 52897

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	70.80	71.99
41.00	34.60	36.27
42.00	23.00	23.03

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 12/04/2023
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Quant Time: Dec 01 22:55:46 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.864	152	348180	20.000	ng/ul	0.00
20) Naphthalene-d8	10.663	136	1450921	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.493	164	949830	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.240	188	1878494	20.000	ng/ul	0.00
79) Chrysene-d12	21.322	240	1280739	20.000	ng/ul	0.00
88) Perylene-d12	23.698	264	1485543	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.276	96	32971	3.227	ng/uL	0.00
4) Pyridine-d5	3.693	84	215845	7.887	ng/ul	0.00
7) Phenol-d5	7.028	99	177762	5.087	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.199	67	553838	26.744	ng/ul	0.00
11) 2-Chlorophenol-d4	7.393	132	508399	18.647	ng/ul	0.00
15) 4-Methylphenol-d8	8.569	113	345045	12.202	ng/ul	-0.01
21) Nitrobenzene-d5	9.028	128	345107	26.724	ng/ul	0.00
24) 2-Nitrophenol-d4	9.746	143	340765	25.452	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.281	165	653819	23.882	ng/ul	0.00
31) 4-Chloroaniline-d4	10.799	131	723088	20.210	ng/ul	0.00
46) Dimethylphthalate-d6	13.910	166	2597510	30.641	ng/ul	0.00
49) Acenaphthylene-d8	14.187	160	2662882	28.527	ng/ul	0.00
54) 4-Nitrophenol-d4	14.687	143	52897m	3.801	ng/ul	0.00
60) Fluorene-d10	15.487	176	2122986	29.835	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.598	200	282595	25.089	ng/ul	0.00
73) Anthracene-d10	17.340	188	3189131	32.293	ng/ul	0.00
81) Pyrene-d10	19.575	212	3174120	31.838	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.545	264	2790787	32.220	ng/ul	0.00

Target Compounds Qvalue

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

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