

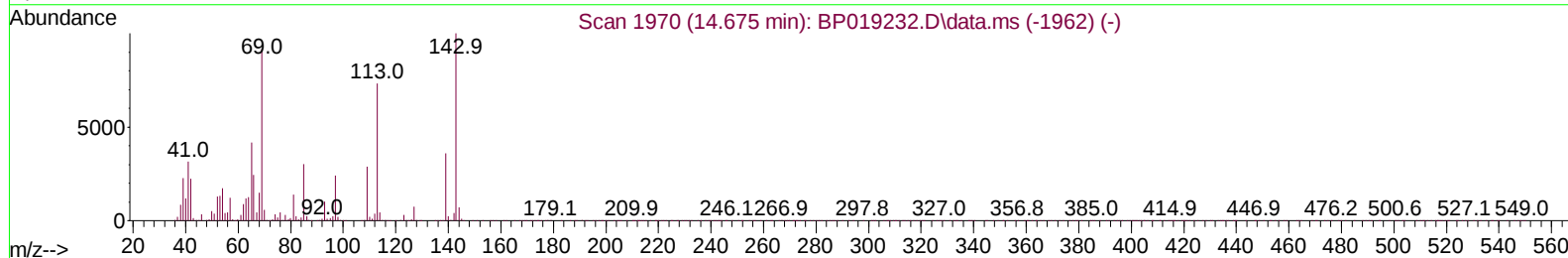
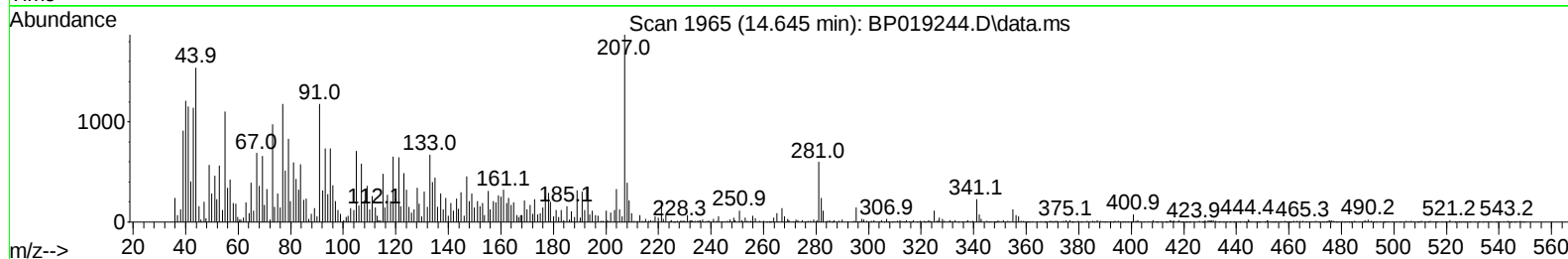
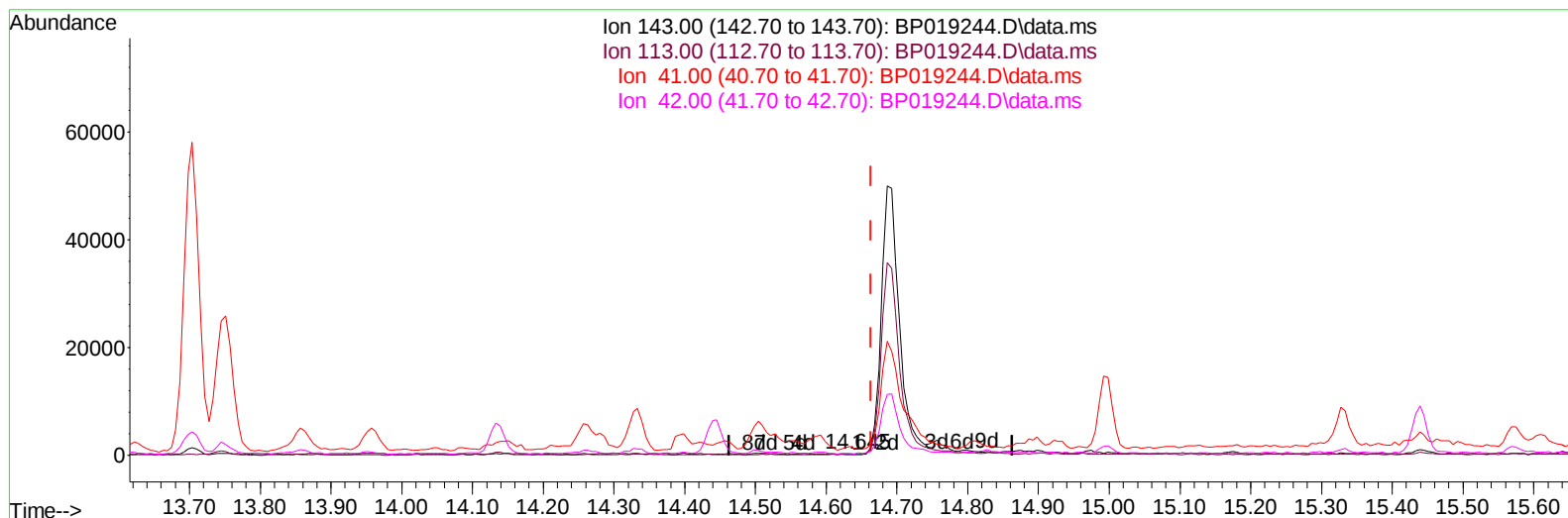
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010224\
 Data File : BP019244.D
 Acq On : 03 Jan 2024 12:10
 Operator : MA/JU
 Sample : 05952-20
 Misc :
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCFA3

Manual IntegrationsAPPROVED

Quant Time: Jan 03 13:11:03 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 29 03:48:11 2023
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 01/04/2024
 Supervised By :mohammad ahmed 01/05/2024



TIC: BP019244.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.645min (-0.018) 0.02 ng/u1

response 126

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	73.50	34.04#
41.00	33.40	491.06#
42.00	23.00	173.62#

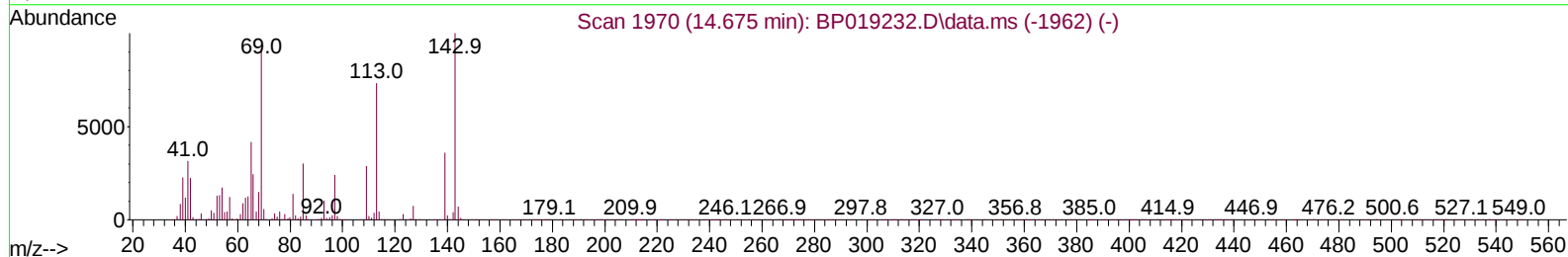
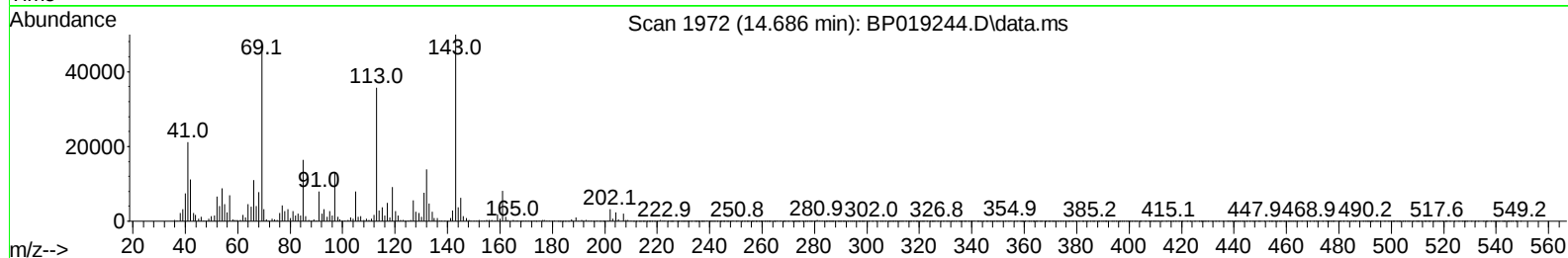
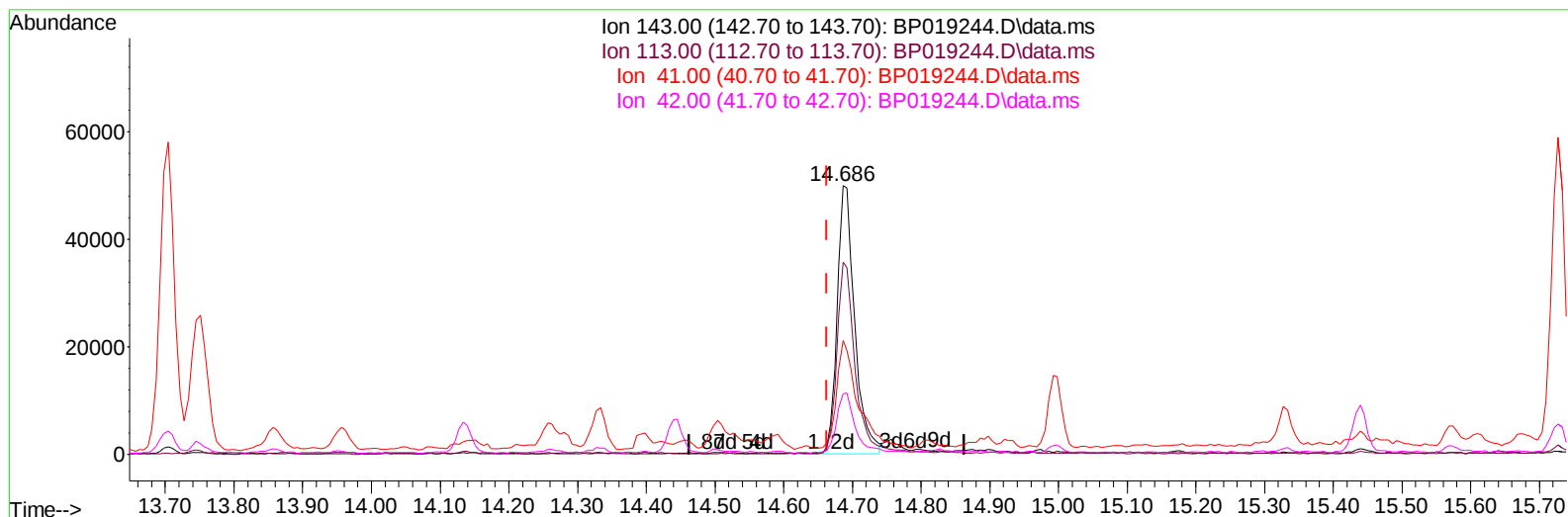
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TIC: BP019244.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.686min (+ 0.023) 16.05 ng/ul m

response 88123

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	73.50	71.56
41.00	33.40	42.34#
42.00	23.00	22.53

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Quant Time: Jan 03 13:12:20 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.798	152	157945	20.000	ng/ul	0.00
20) Naphthalene-d8	10.592	136	586285	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.445	164	368098	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.198	188	879991	20.000	ng/ul	0.00
79) Chrysene-d12	21.304	240	901730	20.000	ng/ul	0.00
88) Perylene-d12	23.674	264	1092867	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.240	96	17344	3.750	ng/uL	0.00
4) Pyridine-d5	3.669	84	85319	7.350	ng/ul	0.02
7) Phenol-d5	6.992	99	288159	20.344	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.140	67	166808	20.357	ng/ul	0.00
11) 2-Chlorophenol-d4	7.334	132	235869	20.656	ng/ul	0.00
15) 4-Methylphenol-d8	8.528	113	173765	15.758	ng/ul	0.00
21) Nitrobenzene-d5	8.963	128	112783	22.253	ng/ul	0.00
24) 2-Nitrophenol-d4	9.686	143	126831	23.156	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.233	165	228870	20.693	ng/ul	0.00
31) 4-Chloroaniline-d4	10.763	131	5130	0.377	ng/ul	0.02
46) Dimethylphthalate-d6	13.863	166	679621	20.903	ng/ul	0.00
49) Acenaphthylene-d8	14.133	160	806411	22.858	ng/ul	0.00
54) 4-Nitrophenol-d4	14.686	143	88123m	16.051	ng/ul	0.02
60) Fluorene-d10	15.439	176	636719	22.668	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.574	200	78835	13.665	ng/ul	0.00
73) Anthracene-d10	17.298	188	1002142	21.807	ng/ul	0.00
81) Pyrene-d10	19.545	212	1313910	23.909	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.521	264	1380721	21.980	ng/ul	0.00
Target Compounds						
					Qvalue	
6) Benzaldehyde	6.945	77	7443	1.132	ng/ul	90
8) Phenol	7.016	94	16306	1.166	ng/ul	97
71) Pentachlorophenol	16.857	266	208285	26.541	ng/ul	96

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

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