

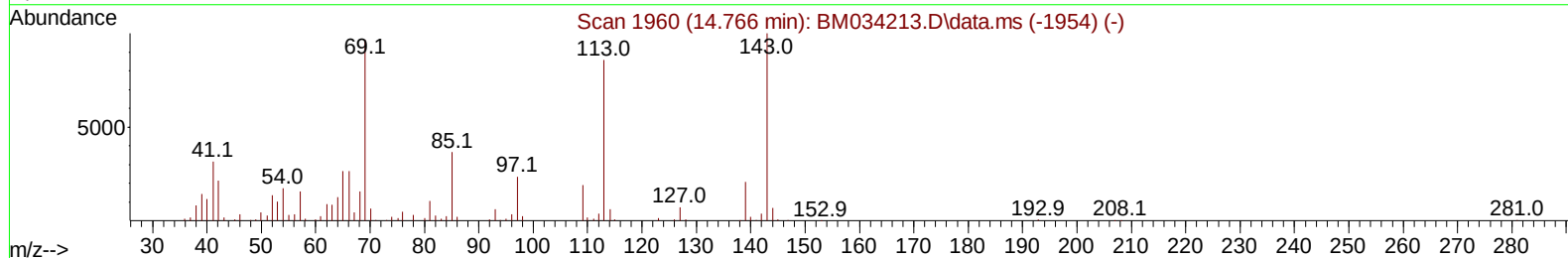
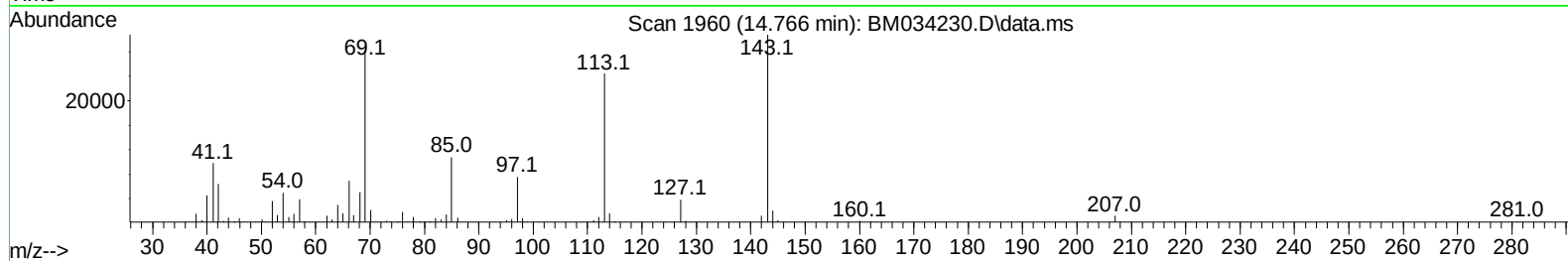
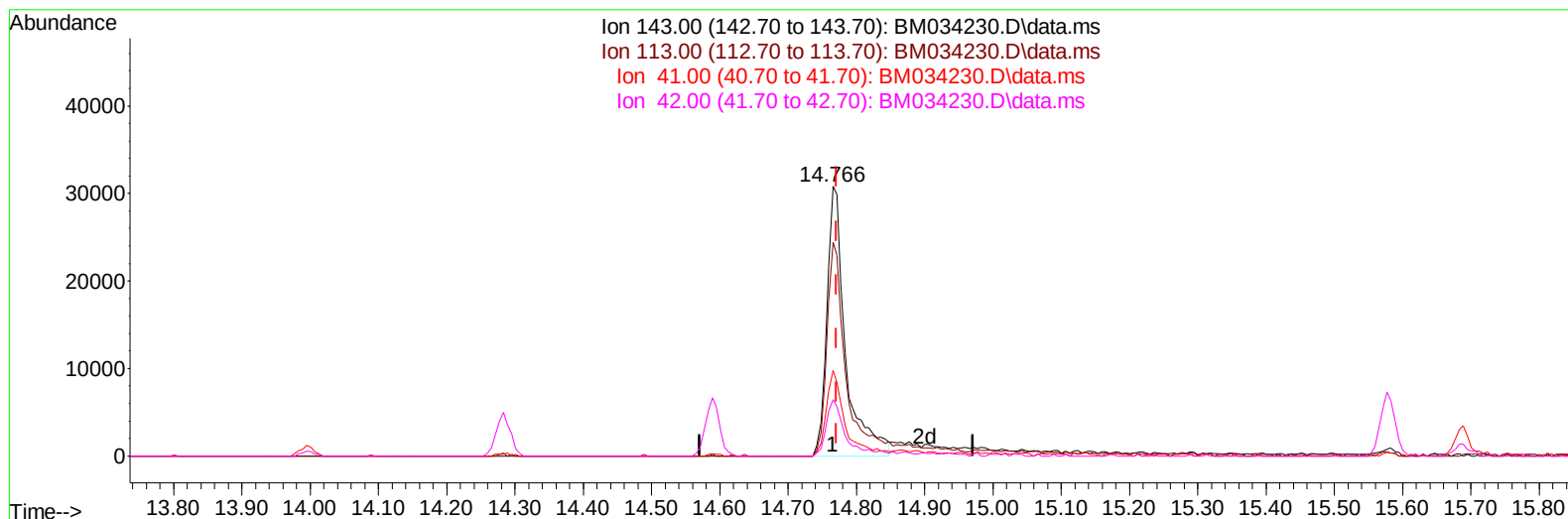
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM031122\
 Data File : BM034230.D
 Acq On : 11 Mar 2022 20:25
 Operator : CG/JU
 Sample : PB143172BL
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK172

Manual IntegrationsAPPROVED

Quant Time: Mar 11 23:42:23 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM031022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Mar 11 02:00:05 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 03/14/2022
 Supervised By :mohammad ahmed 03/14/2022



TIC: BM034230.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.766min (-0.006) 14.17 ng/ul

response 58766

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	77.80	79.42
41.00	31.90	31.78
42.00	19.10	20.77

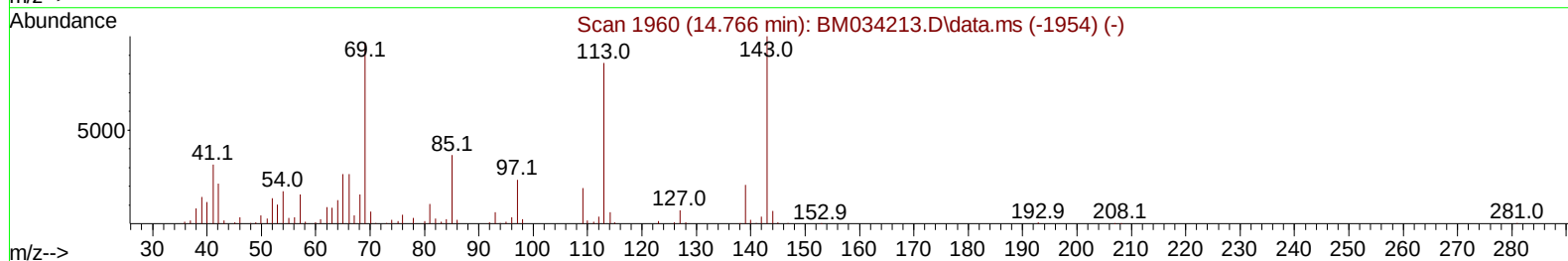
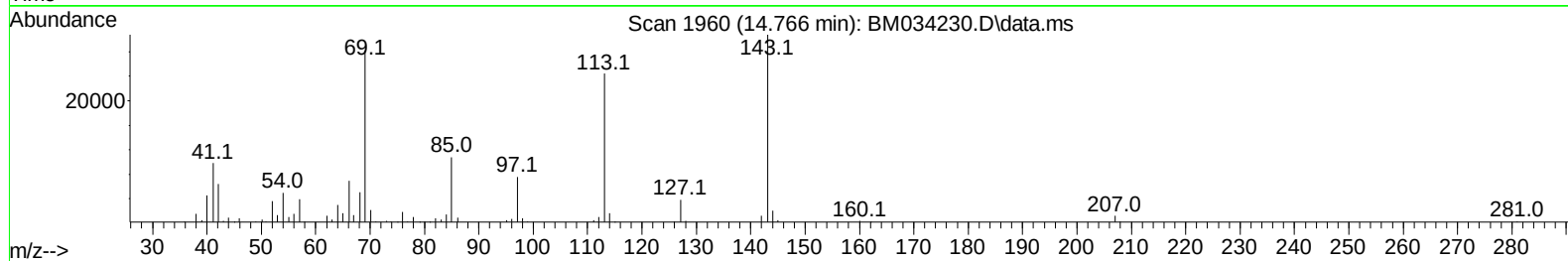
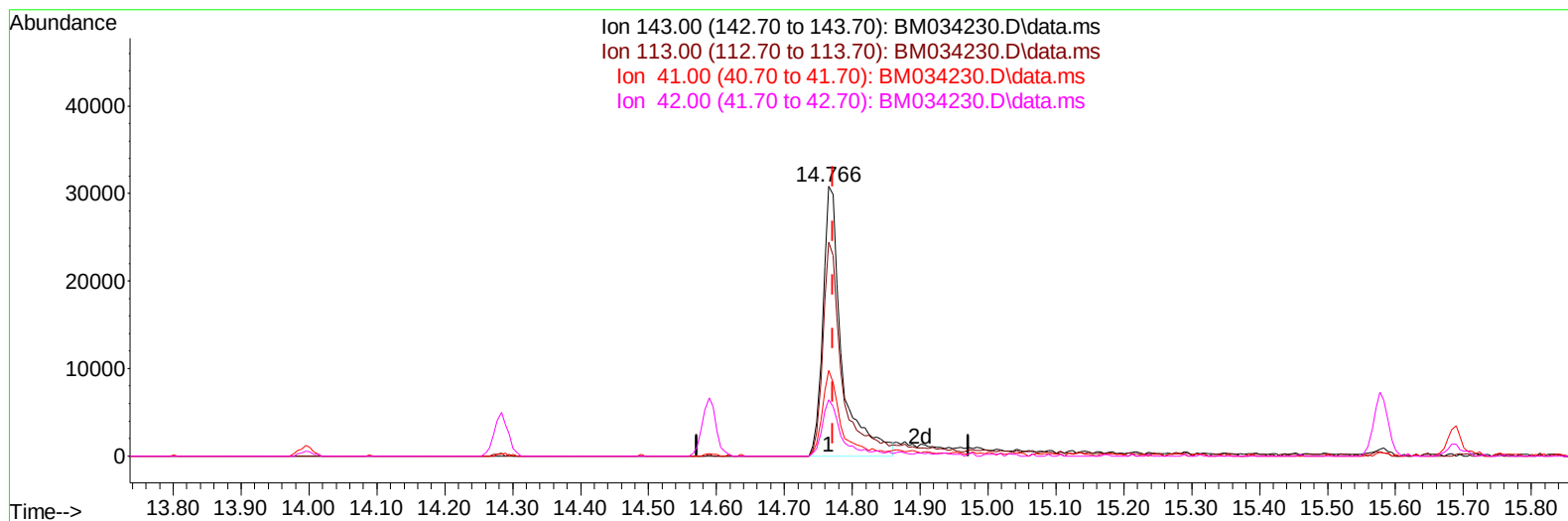
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TIC: BM034230.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.766min (-0.006) 14.43 ng/ul m

response 59855

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	77.80	79.42
41.00	31.90	31.78
42.00	19.10	20.77

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.948	152	143649	20.000	ng/ul	0.00
20) Naphthalene-d8	10.766	136	565569	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.589	164	365896	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.330	188	714743	20.000	ng/ul	0.00
79) Chrysene-d12	21.500	240	631221	20.000	ng/ul	0.00
88) Perylene-d12	23.918	264	678913	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.319	96	14607	4.439	ng/uL	0.00
4) Pyridine-d5	3.743	84	175313	22.376	ng/ul	0.00
7) Phenol-d5	7.101	99	199048	20.706	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.272	67	129229	22.915	ng/ul	0.00
11) 2-Chlorophenol-d4	7.478	132	192160	23.130	ng/ul	0.00
15) 4-Methylphenol-d8	8.654	113	171674	20.881	ng/ul	0.00
21) Nitrobenzene-d5	9.113	128	91194	22.479	ng/ul	0.00
24) 2-Nitrophenol-d4	9.836	143	98833	21.970	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.378	165	186218	20.649	ng/ul	0.00
31) 4-Chloroaniline-d4	10.895	131	225643	20.119	ng/ul	0.00
46) Dimethylphthalate-d6	13.995	166	629619	23.649	ng/ul	0.00
49) Acenaphthylene-d8	14.283	160	766743	22.675	ng/ul	0.00
54) 4-Nitrophenol-d4	14.766	143	59855m	14.434	ng/ul	0.00
60) Fluorene-d10	15.577	176	541676	23.242	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.689	200	80375	17.774	ng/ul	0.00
73) Anthracene-d10	17.430	188	791486	23.445	ng/ul	0.00
81) Pyrene-d10	19.712	212	920908	26.798	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.759	264	812673	24.263	ng/ul	0.00

Target Compounds Qvalue

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

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