

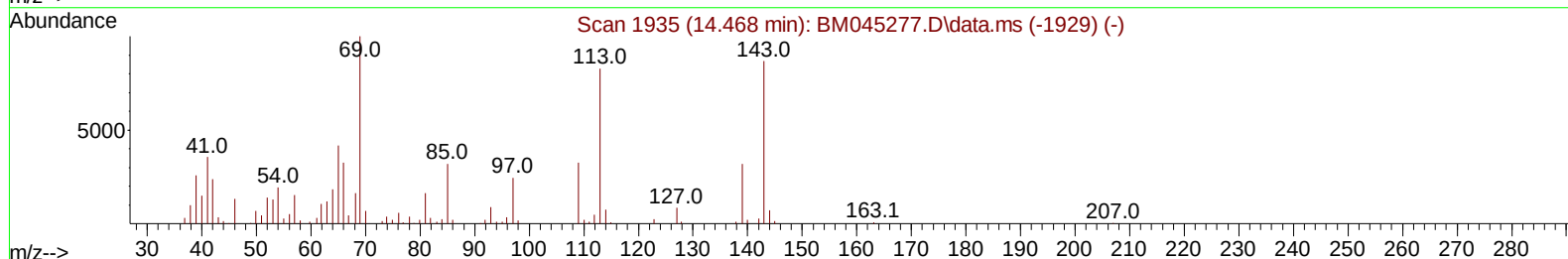
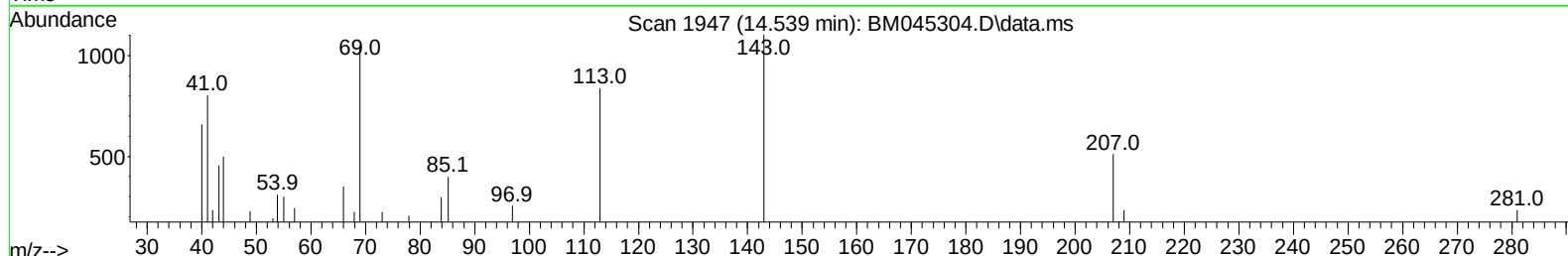
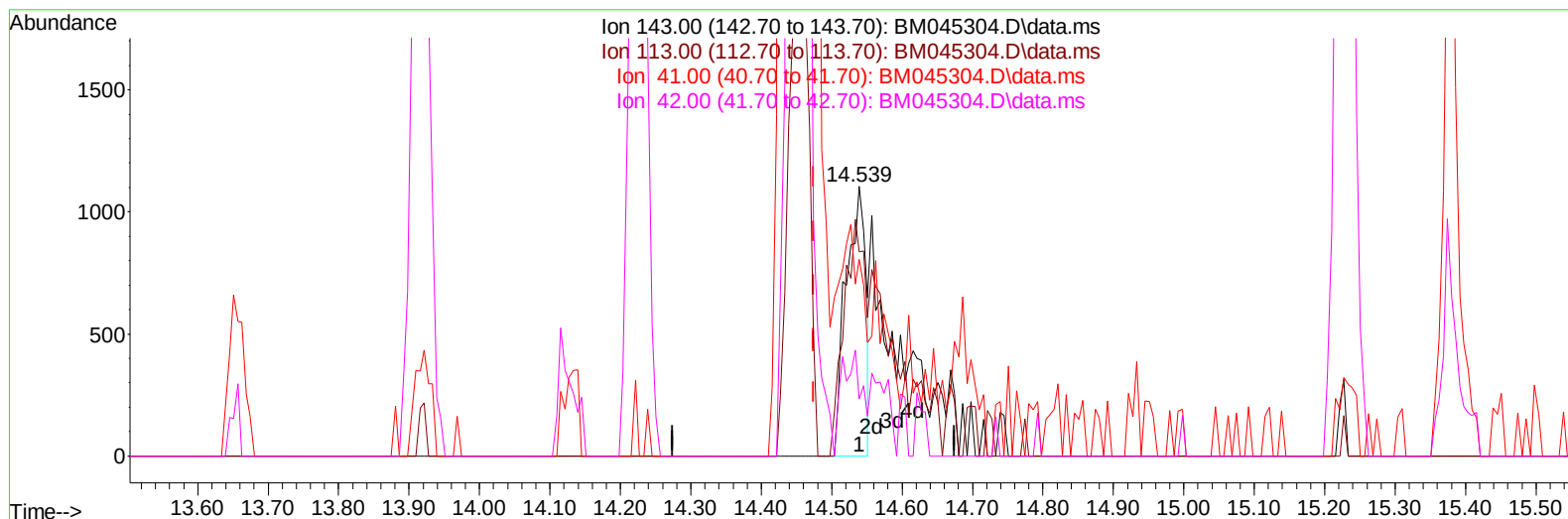
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041524\
 Data File : BM045304.D
 Acq On : 17 Apr 2024 05:50
 Operator : MA/JU
 Sample : P2099-08
 Misc :
 ALS Vial : 65 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0RJ5

Manual IntegrationsAPPROVED

Quant Time: Apr 17 06:47:17 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040524.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Apr 15 22:24:11 2024
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 04/17/2024
 Supervised By :mohammad ahmed 04/18/2024



TIC: BM045304.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.539min (+ 0.065) 1.20 ng/ul

response 2139

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	78.80	75.91
41.00	36.50	73.01#
42.00	23.30	21.20

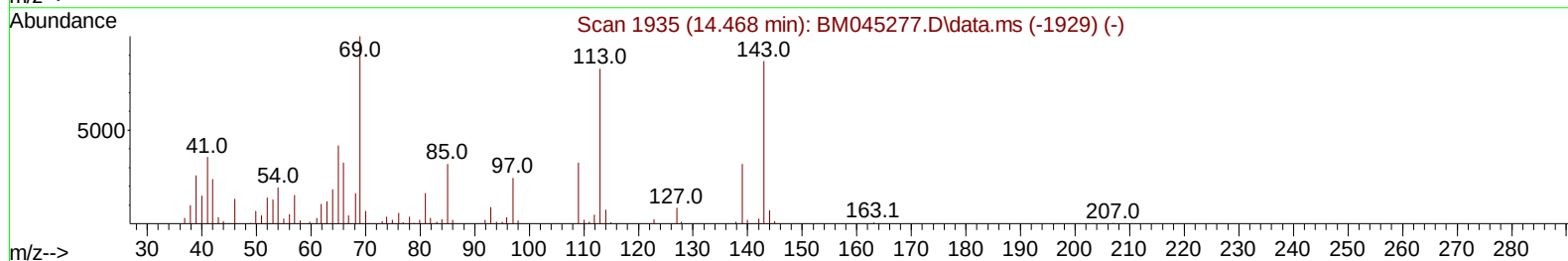
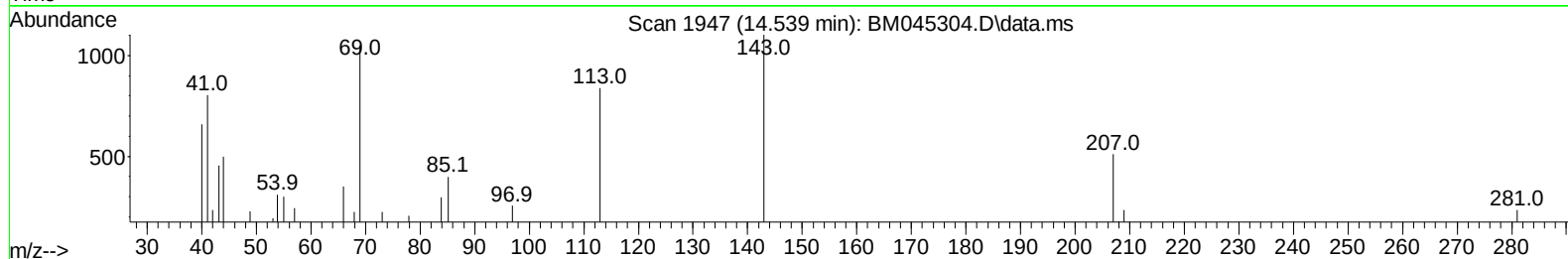
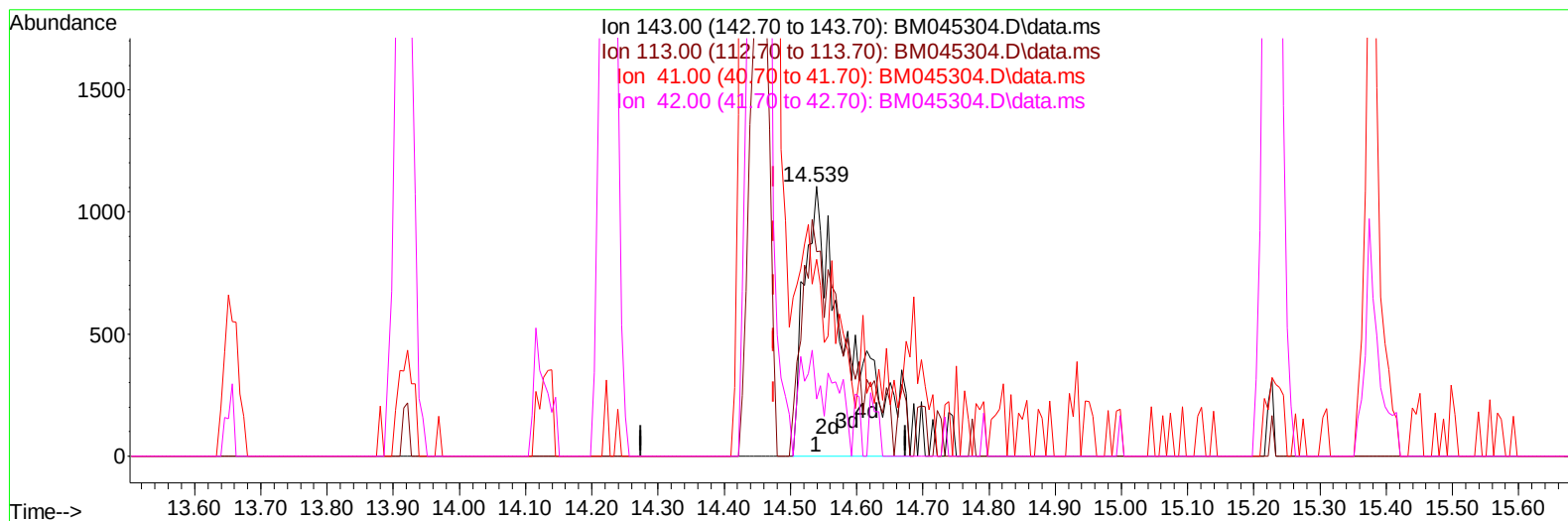
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TIC: BM045304.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.539min (+ 0.065) 2.84 ng/ul m

response 5059

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	78.80	75.91
41.00	36.50	73.01#
42.00	23.30	21.20

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 04/17/2024
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Quant Time: Apr 17 06:48:06 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.575	152	52824	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.345	136	204905	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.227	164	122530	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.986	188	258581	20.000	ng/ul	-0.01
79) Chrysene-d12	21.203	240	261112	20.000	ng/ul	0.00
88) Perylene-d12	23.397	264	350333	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.181	96	5656	3.998	ng/uL	0.00
4) Pyridine-d5	3.587	84	26282	6.832	ng/ul	0.00
7) Phenol-d5	6.769	99	17518	3.912	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.934	67	67638	25.351	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.110	132	67270	19.174	ng/ul	-0.01
15) 4-Methylphenol-d8	8.292	113	36643	10.454	ng/ul	0.00
21) Nitrobenzene-d5	8.739	128	41770	26.539	ng/ul	0.00
24) 2-Nitrophenol-d4	9.451	143	42224	25.417	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.981	165	67403	22.129	ng/ul	0.00
31) 4-Chloroaniline-d4	10.510	131	92976	19.173	ng/ul	-0.01
46) Dimethylphthalate-d6	13.657	166	253285	29.688	ng/ul	0.00
49) Acenaphthylene-d8	13.916	160	280761	27.857	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.539	143	5059m	2.841	ng/ul	0.06
60) Fluorene-d10	15.227	176	229668	30.196	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.380	200	32354	21.102	ng/ul	0.00
73) Anthracene-d10	17.086	188	381796	32.822	ng/ul	-0.01
81) Pyrene-d10	19.398	212	465711	36.222	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.262	264	570775	33.398	ng/ul	-0.01

Target Compounds Qvalue

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

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