

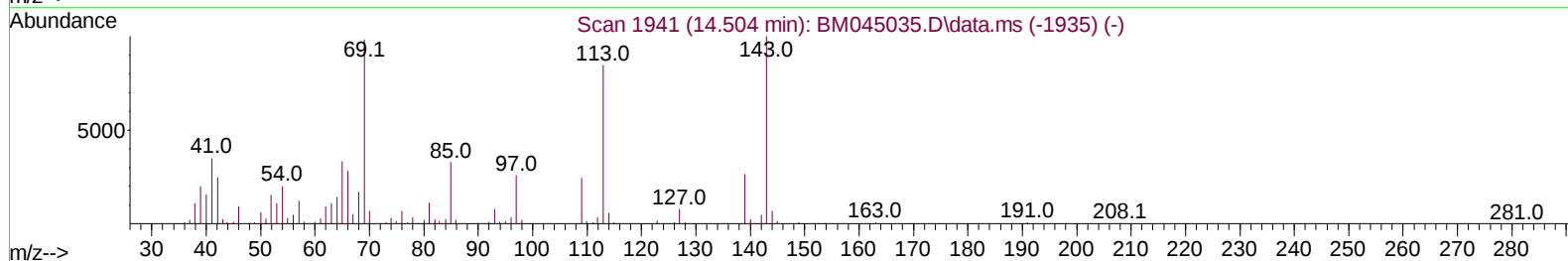
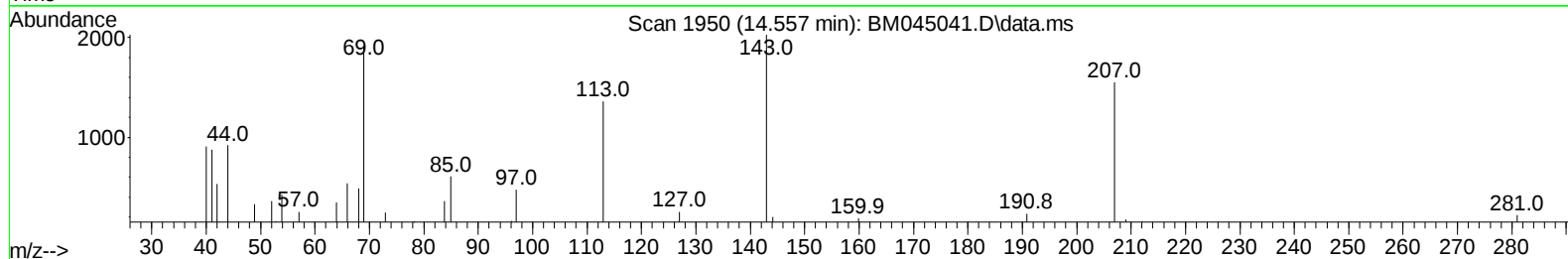
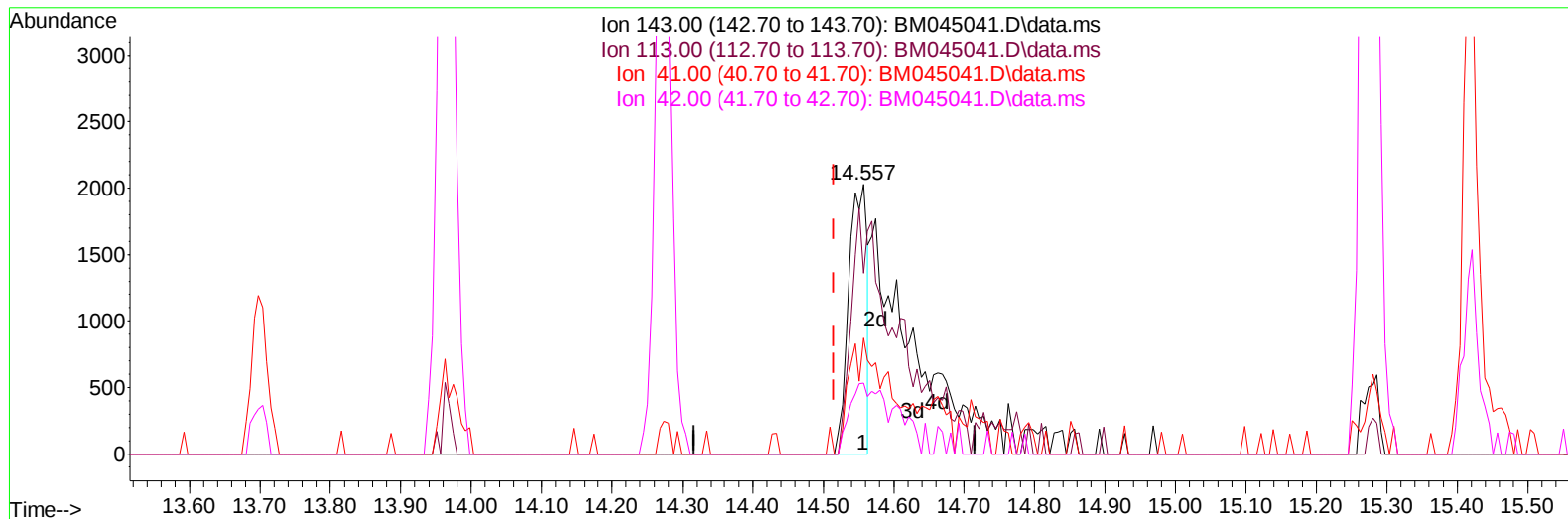
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040724\
Data File : BM045041.D
Acq On : 08 Apr 2024 14:42
Operator : MA/JU
Sample : P1997-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BH7G9

Manual IntegrationsAPPROVED

Quant Time: Apr 08 15:13:08 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040524.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Apr 08 05:51:38 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 04/09/2024
Supervised By :mohammad ahmed 04/09/2024



TIC: BM045041.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.557min (+ 0.041) 1.44 ng/u1

response 3694

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	78.80	67.09
41.00	36.50	43.07
42.00	23.30	26.25

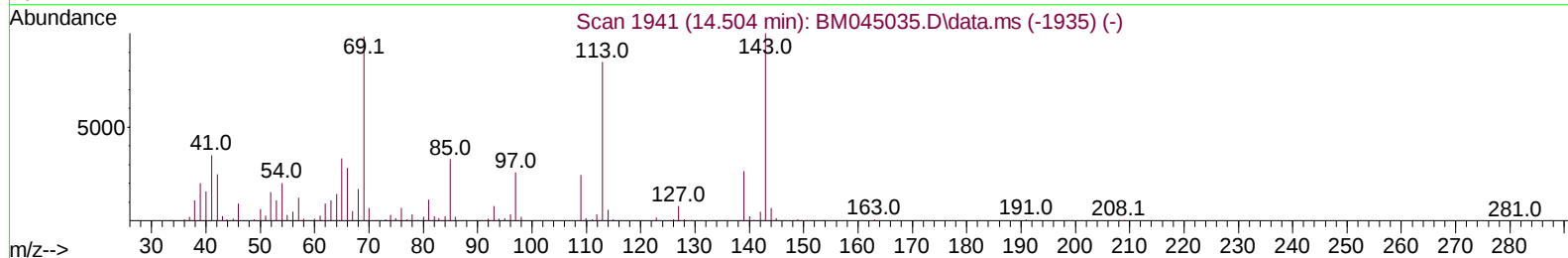
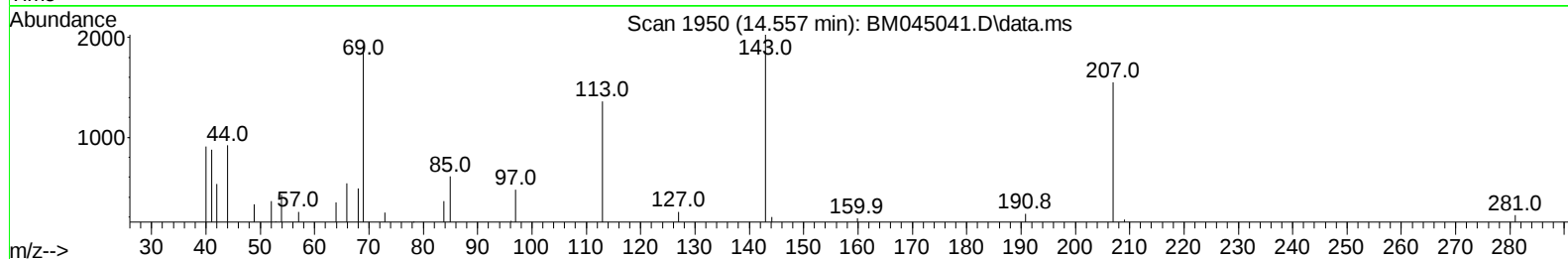
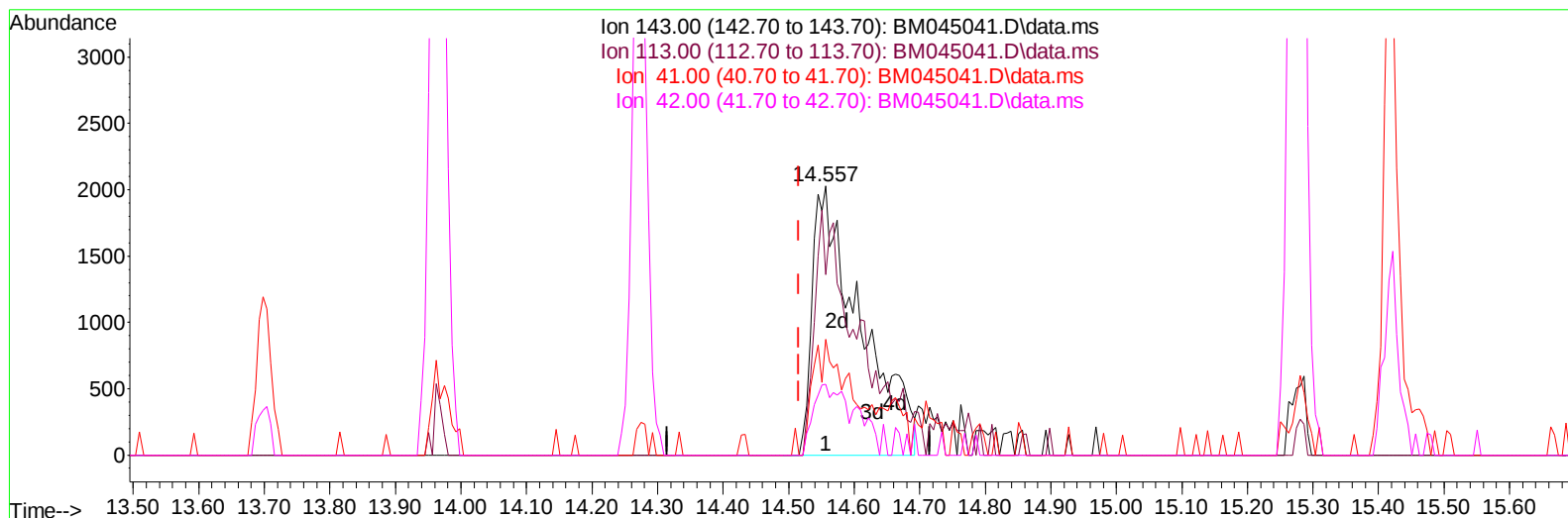
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TIC: BM045041.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.557min (+ 0.041) 3.99 ng/ul m

response 10280

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	78.80	67.09
41.00	36.50	43.07
42.00	23.30	26.25

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.622	152	64262	20.000	ng/ul	0.00
20) Naphthalene-d8	10.398	136	280814	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.274	164	177094	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.033	188	384770	20.000	ng/ul	0.00
79) Chrysene-d12	21.245	240	376767	20.000	ng/ul	0.00
88) Perylene-d12	23.462	264	444967	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	7328	4.258	ng/uL	0.00
4) Pyridine-d5	3.616	84	44414	9.491	ng/ul	0.01
7) Phenol-d5	6.810	99	39552	7.260	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.981	67	141893	43.717	ng/ul	0.00
11) 2-Chlorophenol-d4	7.157	132	143028	33.511	ng/ul	0.00
15) 4-Methylphenol-d8	8.334	113	82994	19.463	ng/ul	0.00
21) Nitrobenzene-d5	8.787	128	92307	42.794	ng/ul	0.00
24) 2-Nitrophenol-d4	9.498	143	96310	42.302	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.028	165	152176	36.456	ng/ul	0.00
31) 4-Chloroaniline-d4	10.557	131	220554	33.187	ng/ul	0.00
46) Dimethylphthalate-d6	13.704	166	536544	43.513	ng/ul	0.00
49) Acenaphthylene-d8	13.969	160	638684	43.845	ng/ul	0.00
54) 4-Nitrophenol-d4	14.557	143	10280m	3.995	ng/ul	0.04
60) Fluorene-d10	15.274	176	490408	44.611	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.421	200	71843	31.491	ng/ul	0.00
73) Anthracene-d10	17.133	188	808126	46.688	ng/ul	0.00
81) Pyrene-d10	19.439	212	961396	51.822	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.327	264	1051927	48.461	ng/ul	0.00

Target Compounds Qvalue

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

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