

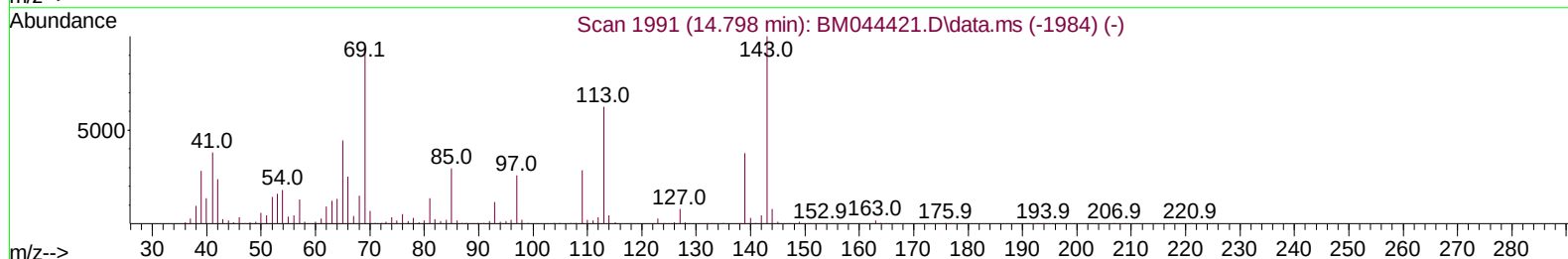
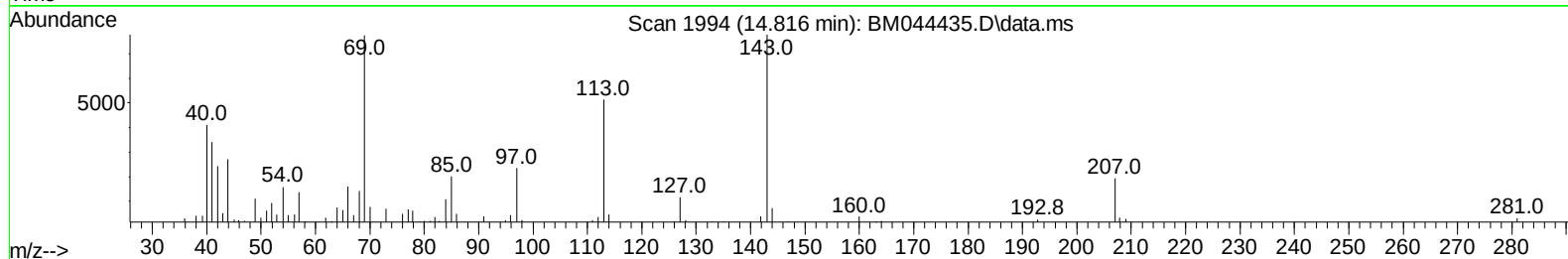
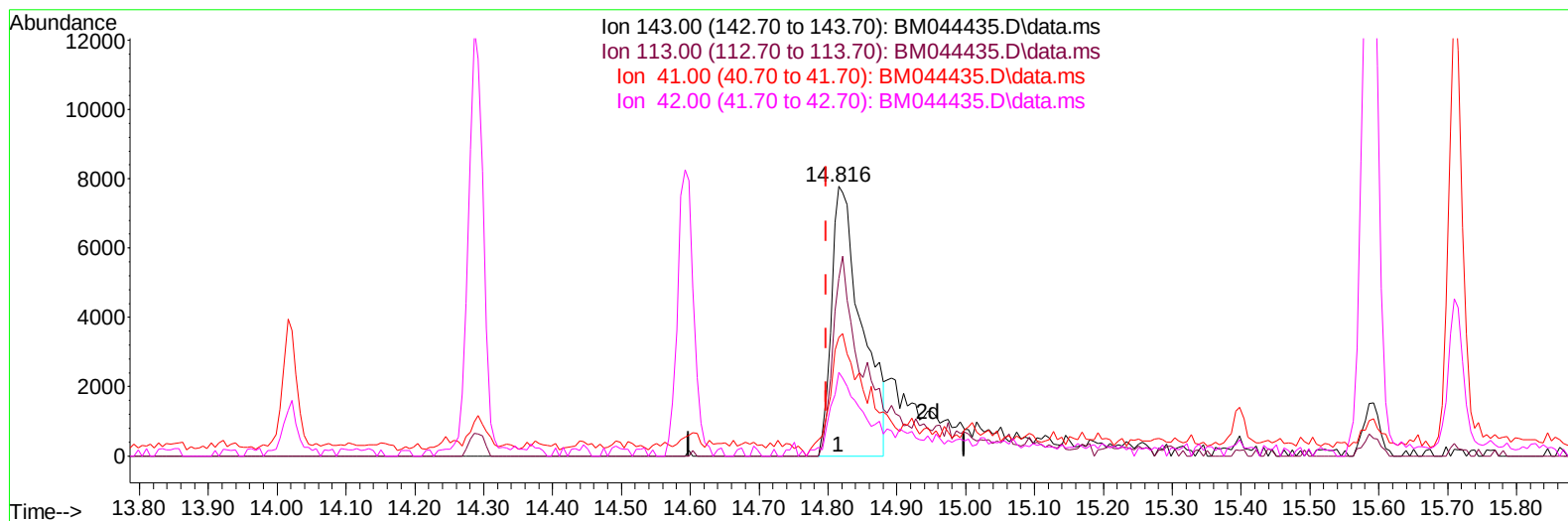
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM022824\
Data File : BM044435.D
Acq On : 29 Feb 2024 12:24
Operator : MA/JU
Sample : P1535-14
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0AC0

Manual IntegrationsAPPROVED

Quant Time: Feb 29 13:21:13 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM022324.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Feb 29 03:37:20 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 03/01/2024
Supervised By :mohammad ahmed 03/02/2024



TIC: BM044435.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.816min (+ 0.018) 3.26 ng/ul

response 23557

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	59.00	65.99
41.00	38.20	43.96
42.00	25.30	31.08#

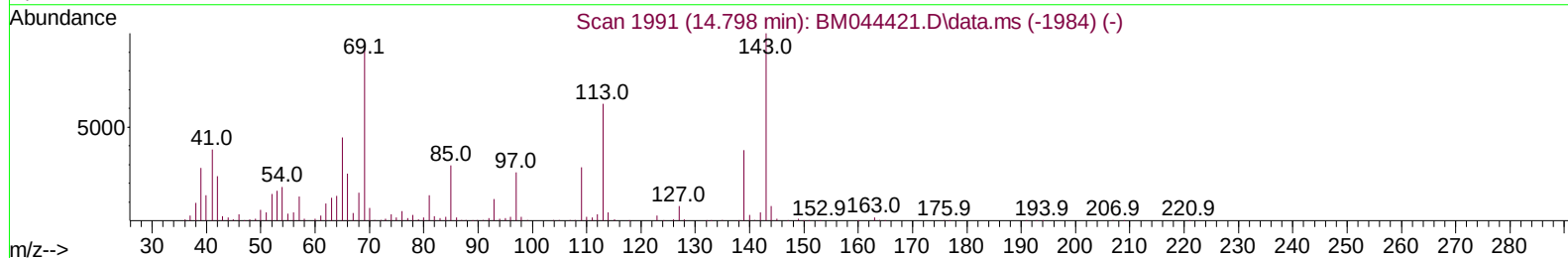
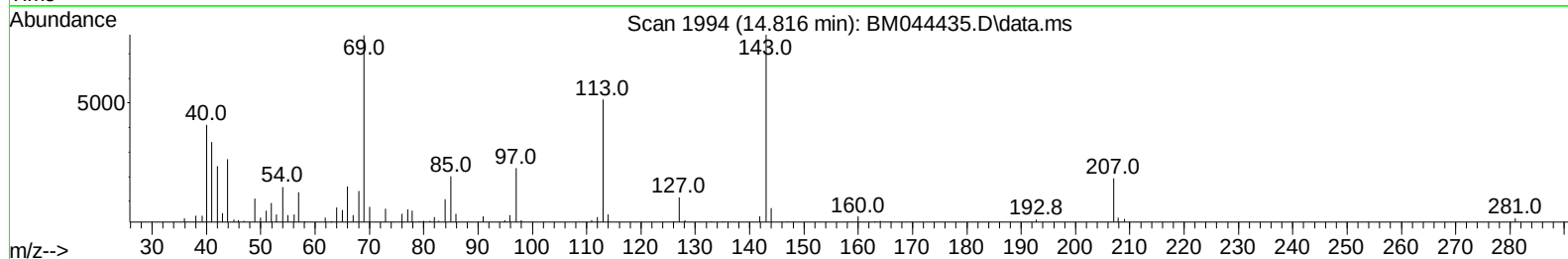
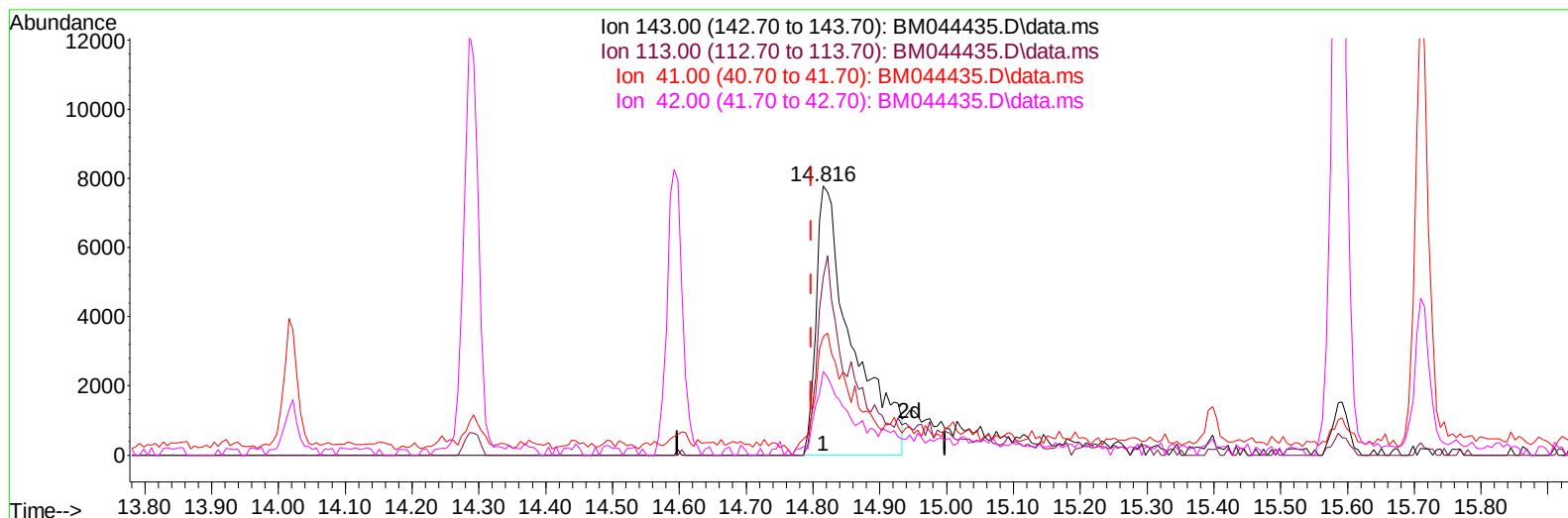
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TIC: BM044435.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.816min (+ 0.018) 4.01 ng/ul m

response 28962

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	59.00	65.99
41.00	38.20	43.96
42.00	25.30	31.08#

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 03/01/2024
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Quant Time: Feb 29 13:22:22 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.957	152	180460	20.000	ng/ul	0.00
20) Naphthalene-d8	10.763	136	809370	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.592	164	471096	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.345	188	970795	20.000	ng/ul	0.00
79) Chrysene-d12	21.539	240	798369	20.000	ng/ul	0.00
88) Perylene-d12	23.962	264	877947	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	20297	3.549	ng/uL	0.00
4) Pyridine-d5	3.805	84	64198	4.110	ng/ul	0.02
7) Phenol-d5	7.122	99	119828	6.932	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.298	67	366463	32.735	ng/ul	0.00
11) 2-Chlorophenol-d4	7.481	132	358474	27.532	ng/ul	0.00
15) 4-Methylphenol-d8	8.669	113	225358	17.229	ng/ul	0.00
21) Nitrobenzene-d5	9.128	128	228547	34.729	ng/ul	0.00
24) 2-Nitrophenol-d4	9.851	143	236553	35.233	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.381	165	386953	32.672	ng/ul	0.00
31) 4-Chloroaniline-d4	10.916	131	233357	11.732	ng/ul	0.01
46) Dimethylphthalate-d6	14.022	166	1481111	40.765	ng/ul	0.00
49) Acenaphthylene-d8	14.286	160	1667636	39.893	ng/ul	0.00
54) 4-Nitrophenol-d4	14.816	143	28962m	4.006	ng/ul	0.02
60) Fluorene-d10	15.586	176	1280586	41.642	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.716	200	191925	31.862	ng/ul	0.00
73) Anthracene-d10	17.445	188	2042869	44.133	ng/ul	0.00
81) Pyrene-d10	19.739	212	2323008	48.365	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.803	264	2071638	46.050	ng/ul	0.00

Target Compounds Qvalue

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

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