

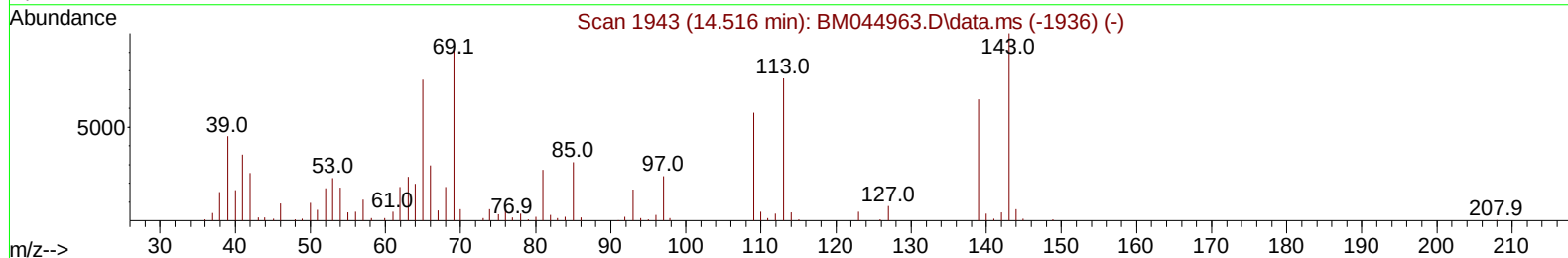
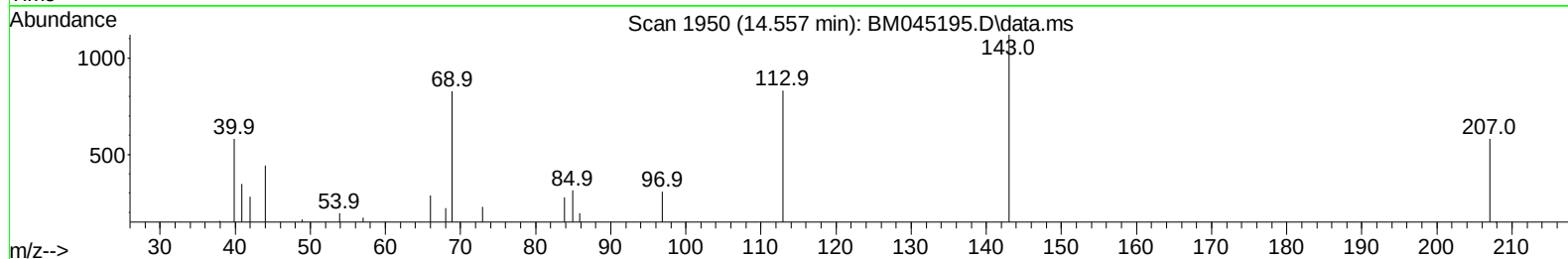
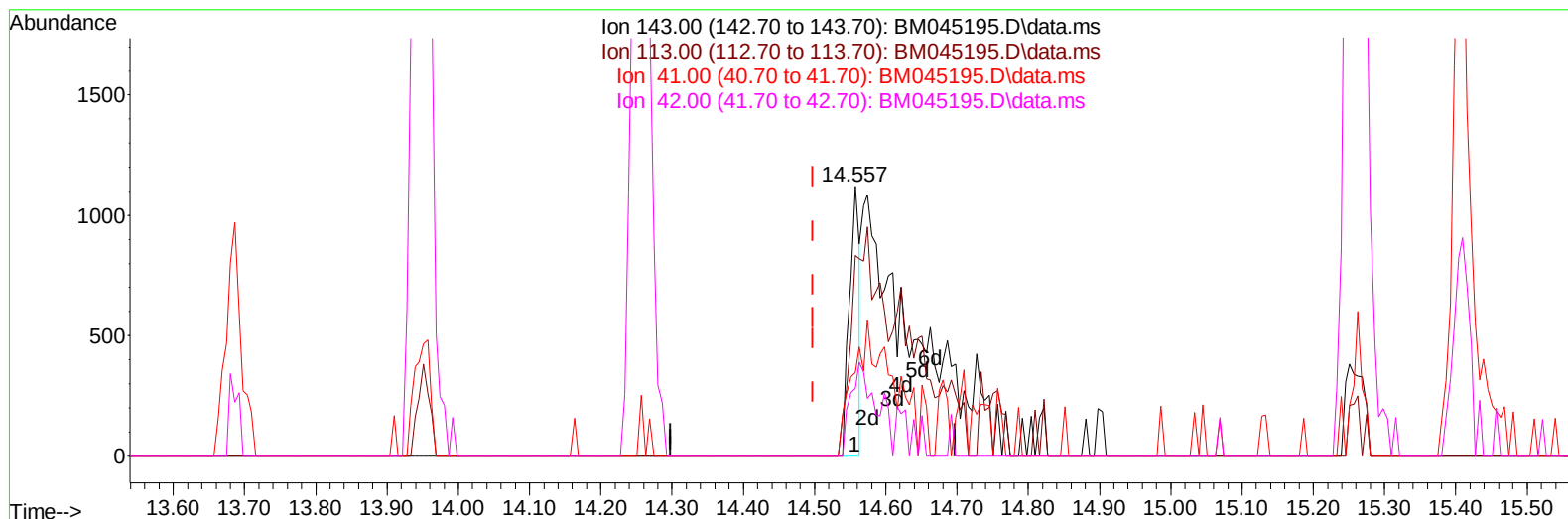
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041124\
Data File : BM045195.D
Acq On : 13 Apr 2024 06:54
Operator : MA/JU
Sample : P2014-06
Misc :
ALS Vial : 59 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
E0MD3

Manual IntegrationsAPPROVED

Quant Time: Apr 15 02:38:48 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040524.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Apr 11 18:54:49 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 04/16/2024
Supervised By :mohammad ahmed 04/16/2024



TIC: BM045195.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.557min (+ 0.059) 0.51 ng/ul

response 1129

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	78.80	74.29
41.00	36.50	31.07
42.00	23.30	25.18

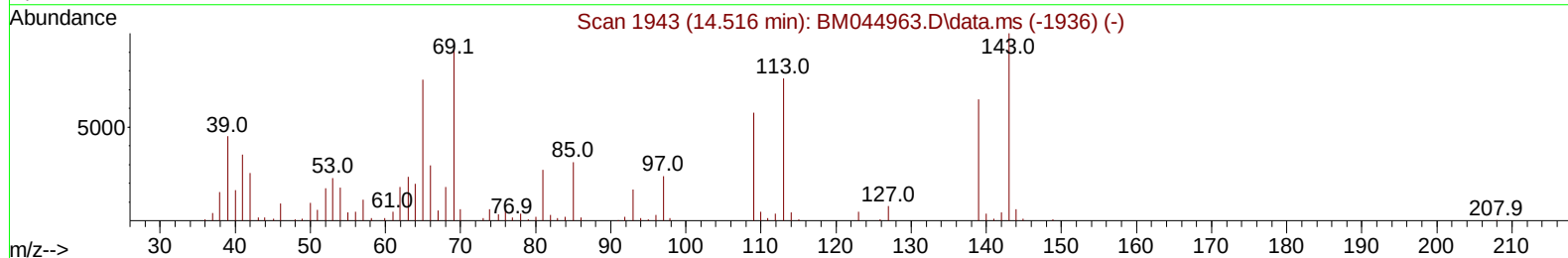
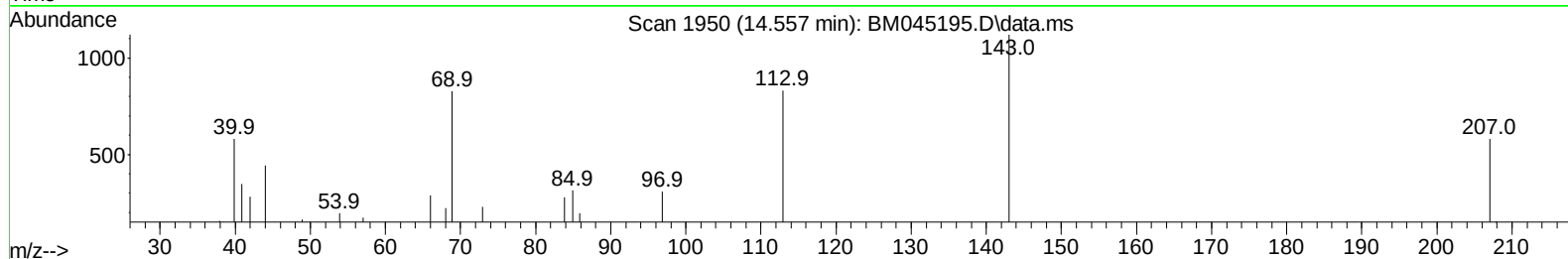
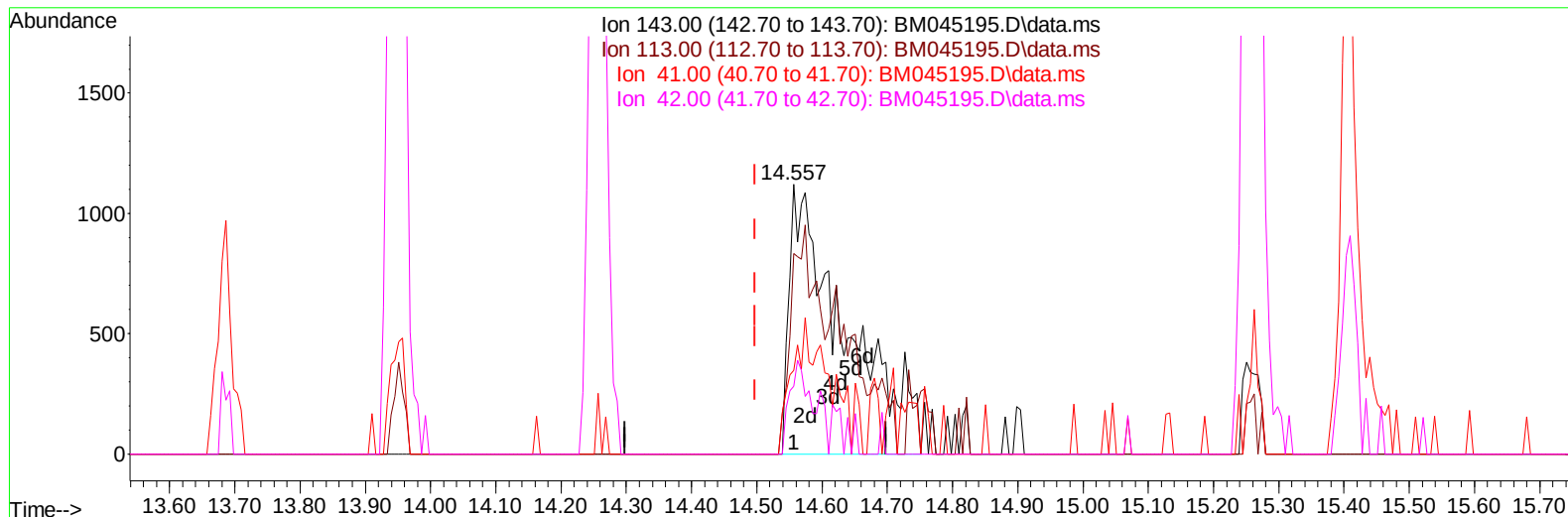
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(54) 4-Nitrophenol-d4 (S)

14.557min (+ 0.059) 2.99 ng/ul m

response 6589

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	78.80	74.29
41.00	36.50	31.07
42.00	23.30	25.18

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Reviewed By :Yogesh Patel 04/16/2024
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Quant Time: Apr 15 02:46:24 2024
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 QLast Update : Thu Apr 11 18:54:49 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.610	152	67603	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.381	136	268444	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.257	164	151574	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.015	188	264901	20.000	ng/ul	-0.01
79) Chrysene-d12	21.233	240	171262	20.000	ng/ul	0.00
88) Perylene-d12	23.444	264	203860	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	7227	3.992	ng/uL	0.00
4) Pyridine-d5	3.610	84	38988	7.919	ng/ul	0.00
7) Phenol-d5	6.798	99	34483	6.017	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.969	67	107932	31.610	ng/ul	0.00
11) 2-Chlorophenol-d4	7.146	132	112023	24.950	ng/ul	0.00
15) 4-Methylphenol-d8	8.322	113	62389	13.908	ng/ul	0.00
21) Nitrobenzene-d5	8.775	128	68557	33.248	ng/ul	0.00
24) 2-Nitrophenol-d4	9.481	143	67676	31.095	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.016	165	104578	26.208	ng/ul	0.00
31) 4-Chloroaniline-d4	10.545	131	145379	22.883	ng/ul	0.00
46) Dimethylphthalate-d6	13.686	166	363777	34.469	ng/ul	0.00
49) Acenaphthylene-d8	13.951	160	435433	34.924	ng/ul	0.00
54) 4-Nitrophenol-d4	14.557	143	6589m	2.992	ng/ul	0.06
60) Fluorene-d10	15.263	176	324698	34.510	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.410	200	41426	26.375	ng/ul	0.00
73) Anthracene-d10	17.115	188	487293	40.892	ng/ul	-0.01
81) Pyrene-d10	19.427	212	499238	59.201	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.303	264	412519	41.481	ng/ul	0.00

Target Compounds Qvalue

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

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