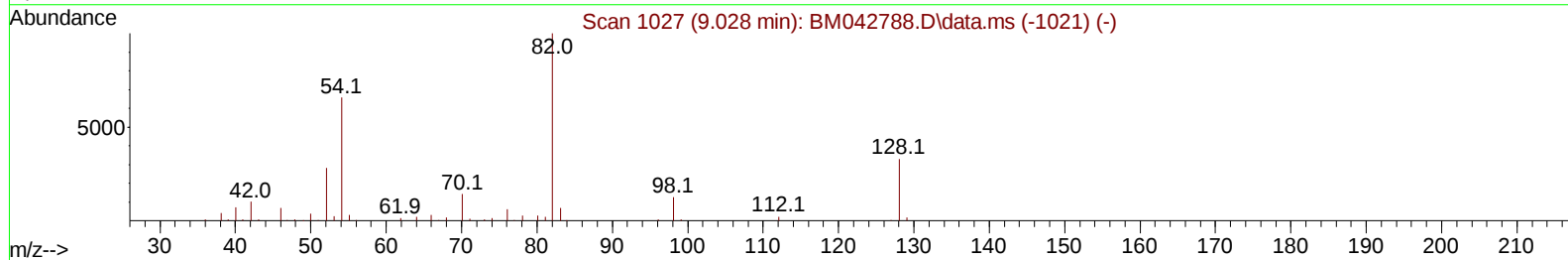
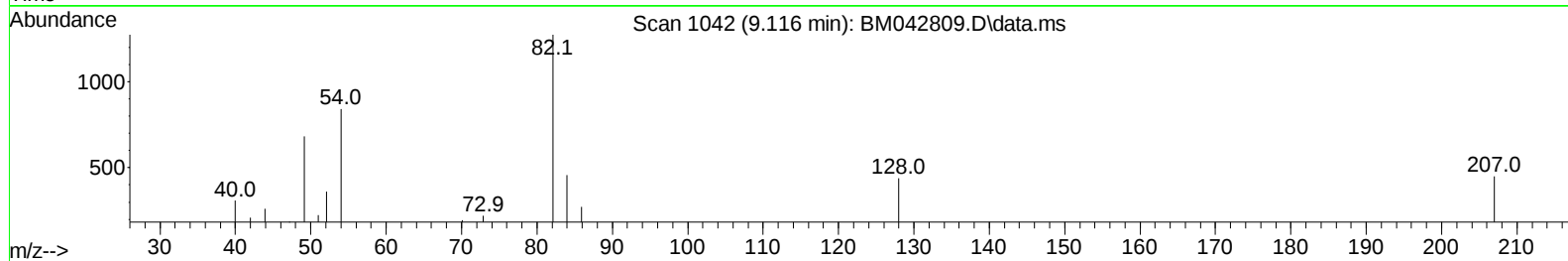
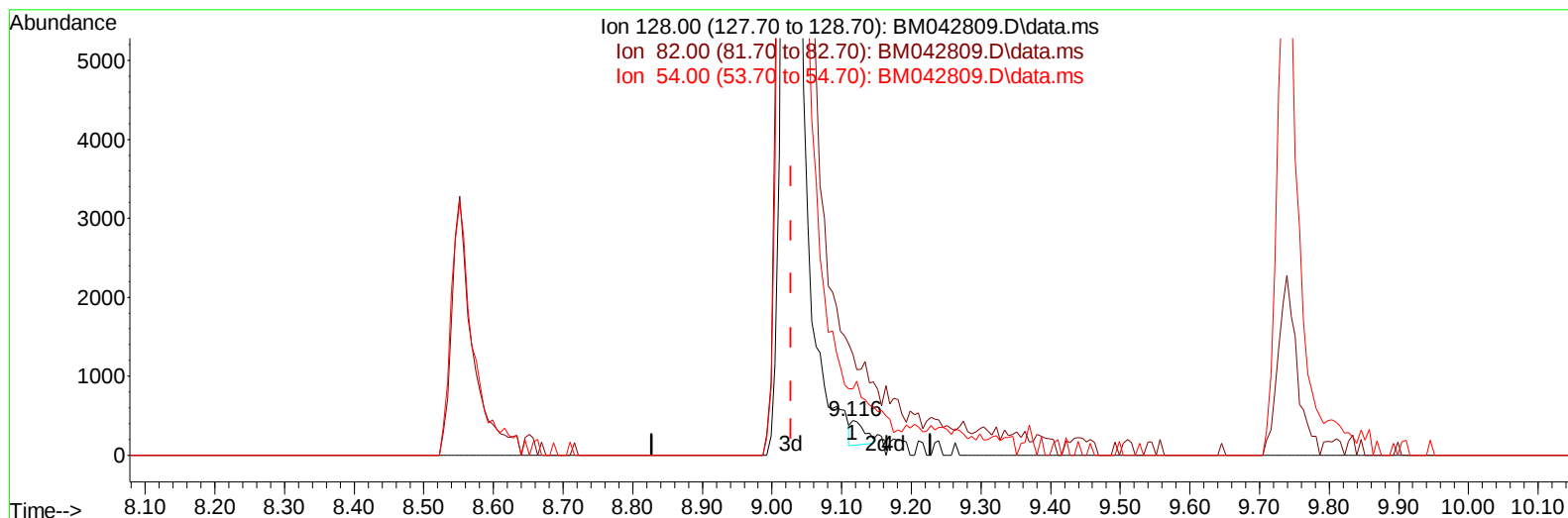


Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111323\
Data File : BM042809.D
Acq On : 16 Nov 2023 16:44
Operator : MA/JU
Sample : 05259-05
Misc :
ALS Vial : 49 Sample Multiplier: 1

Quant Time: Nov 16 23:54:35 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM111323.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 15 22:28:07 2023
Response via : Initial Calibration



TIC: BM042809.D\data.ms

(21) Nitrobenzene-d5 (S)

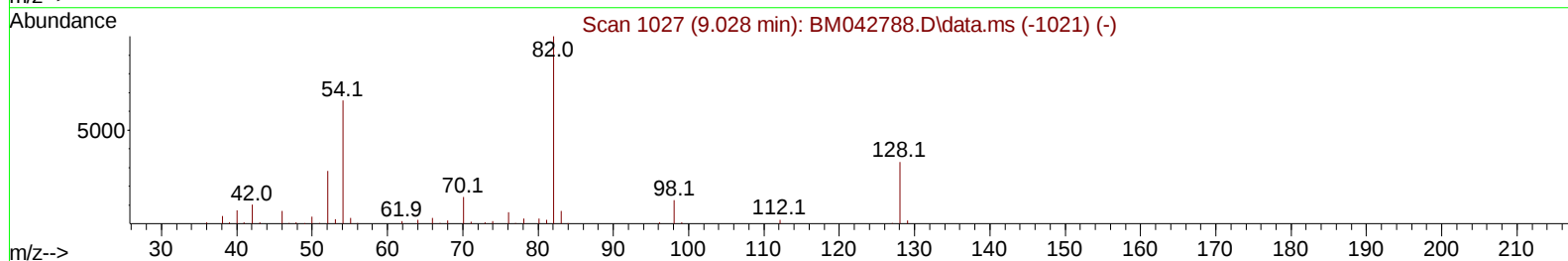
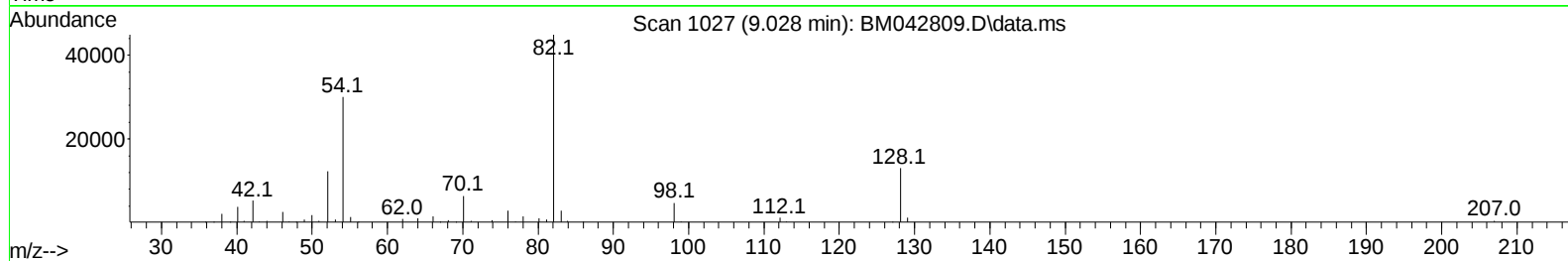
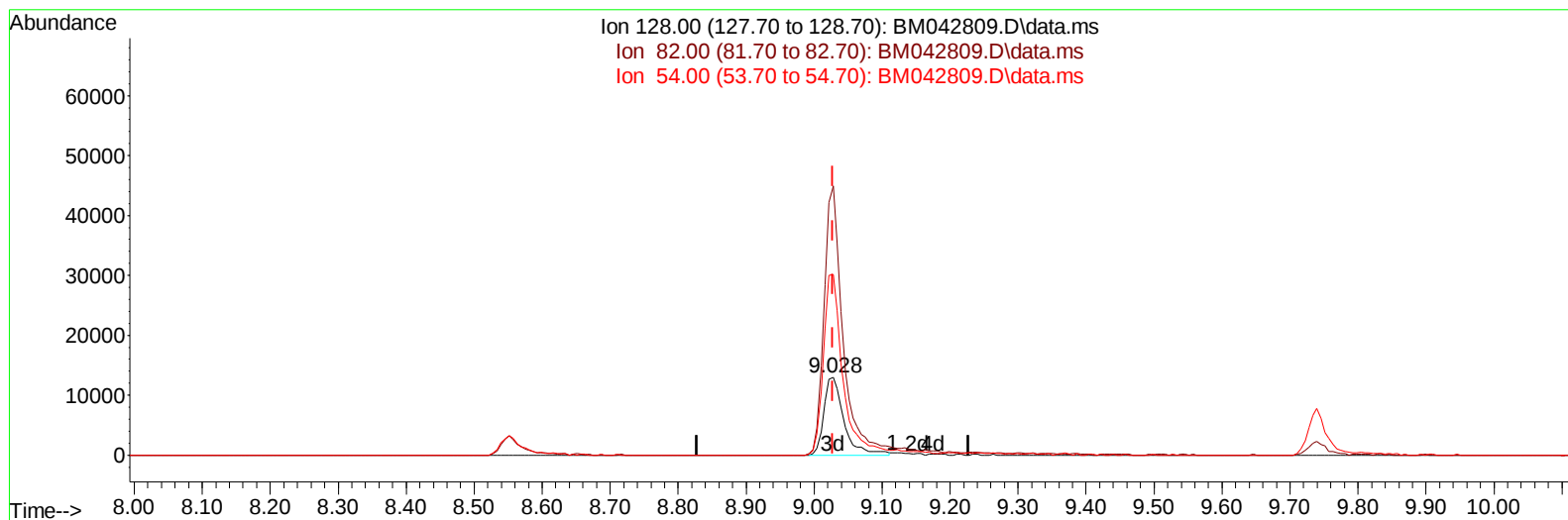
9.116min (+ 0.088) 0.53 ng/u1

response 418

Ion	Exp%	Act%
128.00	100.00	100.00
82.00	289.30	291.08
54.00	185.60	192.22
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111323\
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Acq On : 16 Nov 2023 16:44
Operator : MA/JU
Sample : 05259-05
Misc :
ALS Vial : 49 Sample Multiplier: 1

Quant Time: Nov 16 23:54:35 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM111323.MA.M
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TIC: BM042809.D\data.ms

(21) Nitrobenzene-d5 (S)

9.028min (-0.000) 33.67 ng/ul m

response 26504

Ion	Exp%	Act%
128.00	100.00	100.00
82.00	289.30	345.62
54.00	185.60	231.56#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111323\
 Data File : BM042809.D
 Acq On : 16 Nov 2023 16:44
 Operator : MA/JU
 Sample : 05259-05
 Misc :
 ALS Vial : 49 Sample Multiplier: 1

Quant Time: Nov 17 00:14:59 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM111323.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 15 22:28:07 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.828	152	22050	20.000	ng/ul	0.00
20) Naphthalene-d8	10.634	136	89777	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.480	164	55413	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.233	188	120798	20.000	ng/ul	0.00
79) Chrysene-d12	21.421	240	118601	20.000	ng/ul	# 0.00
88) Perylene-d12	23.780	264	146507	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.175	96	3445	4.902	ng/uL	0.00
4) Pyridine-d5	3.622	84	22621	12.941	ng/ul	0.00
7) Phenol-d5	6.999	99	12610	5.876	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.175	67	60906	40.390	ng/ul	0.00
11) 2-Chlorophenol-d4	7.346	132	43310	26.155	ng/ul	0.00
15) 4-Methylphenol-d8	8.551	113	24013	14.211	ng/ul	0.00
21) Nitrobenzene-d5	9.028	128	26504m	33.672	ng/ul	0.00
24) 2-Nitrophenol-d4	9.740	143	27540	29.909	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.263	165	45953	26.840	ng/ul	0.00
31) 4-Chloroaniline-d4	10.816	131	52201	24.013	ng/ul	0.00
46) Dimethylphthalate-d6	13.904	166	201338	38.391	ng/ul	0.00
49) Acenaphthylene-d8	14.180	160	209715	37.373	ng/ul	0.00
54) 4-Nitrophenol-d4	14.804	143	289	0.488	ng/ul	0.06
60) Fluorene-d10	15.475	176	174069	40.082	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.622	200	22031	24.077	ng/ul	0.00
73) Anthracene-d10	17.333	188	283176	41.829	ng/ul	0.00
81) Pyrene-d10	19.627	212	384250	49.476	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.621	264	397165	45.321	ng/ul	0.00

Target Compounds	Qvalue

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

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