

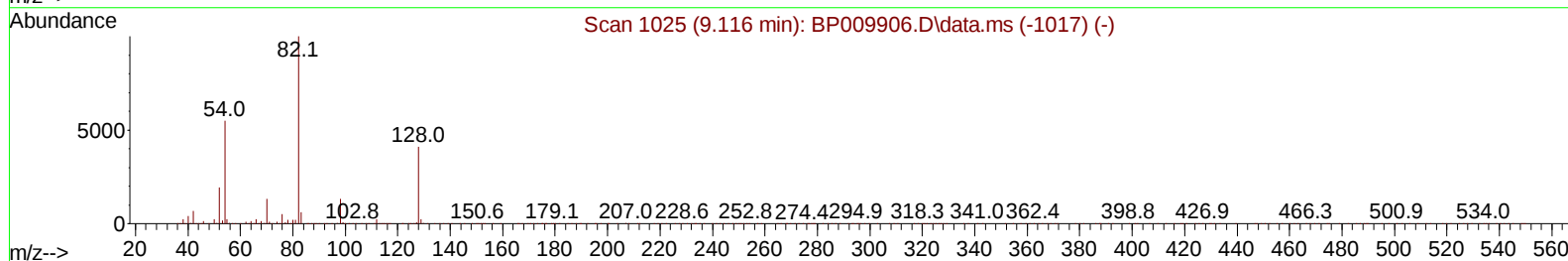
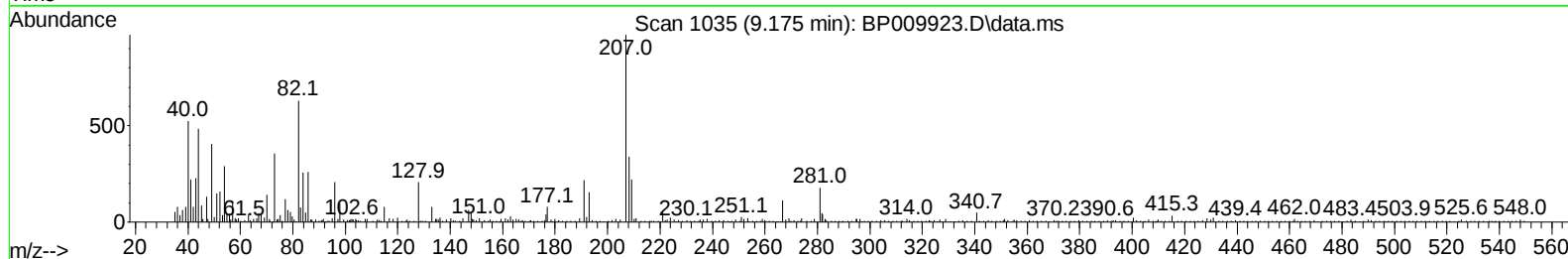
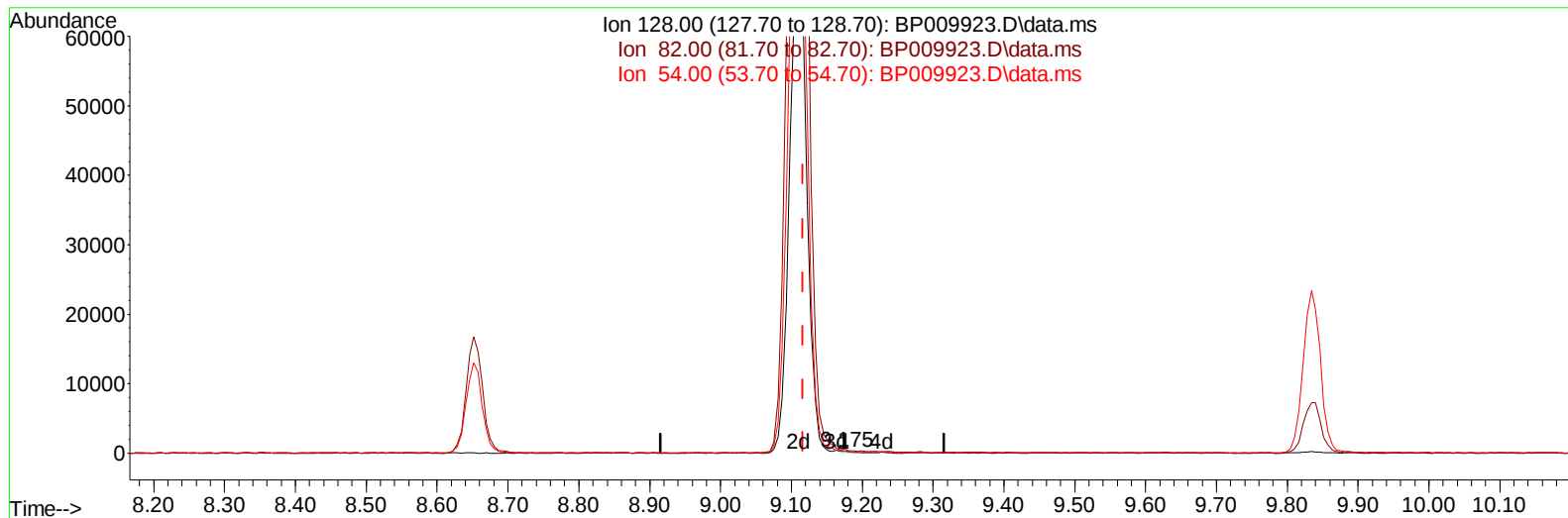
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041822\
 Data File : BP009923.D
 Acq On : 18 Apr 2022 19:29
 Operator : CG/JU
 Sample : N2421-01
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0J55

Manual IntegrationsAPPROVED

Quant Time: Apr 19 00:58:06 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Apr 18 00:59:30 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/19/2022
 Supervised By :Yogesh Patel 04/22/2022



TIC: BP009923.D\data.ms

(21) Nitrobenzene-d5 (S)

9.175min (+ 0.059) 0.04 ng/u1

response 167

Ion	Exp%	Act%
128.00	100.00	100.00
82.00	287.30	294.86
54.00	120.20	135.98
0.00	0.00	0.00

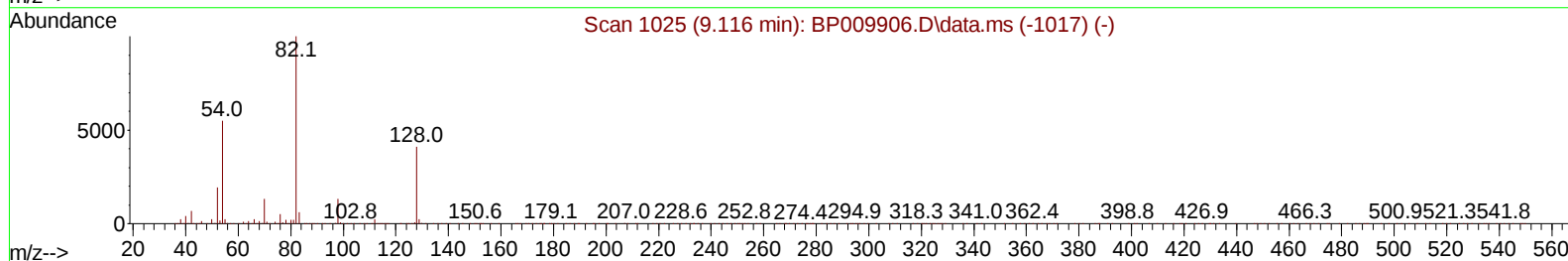
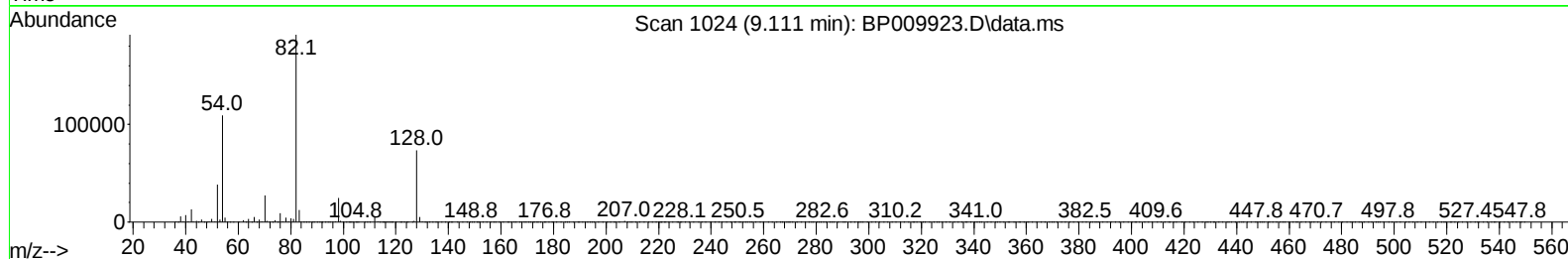
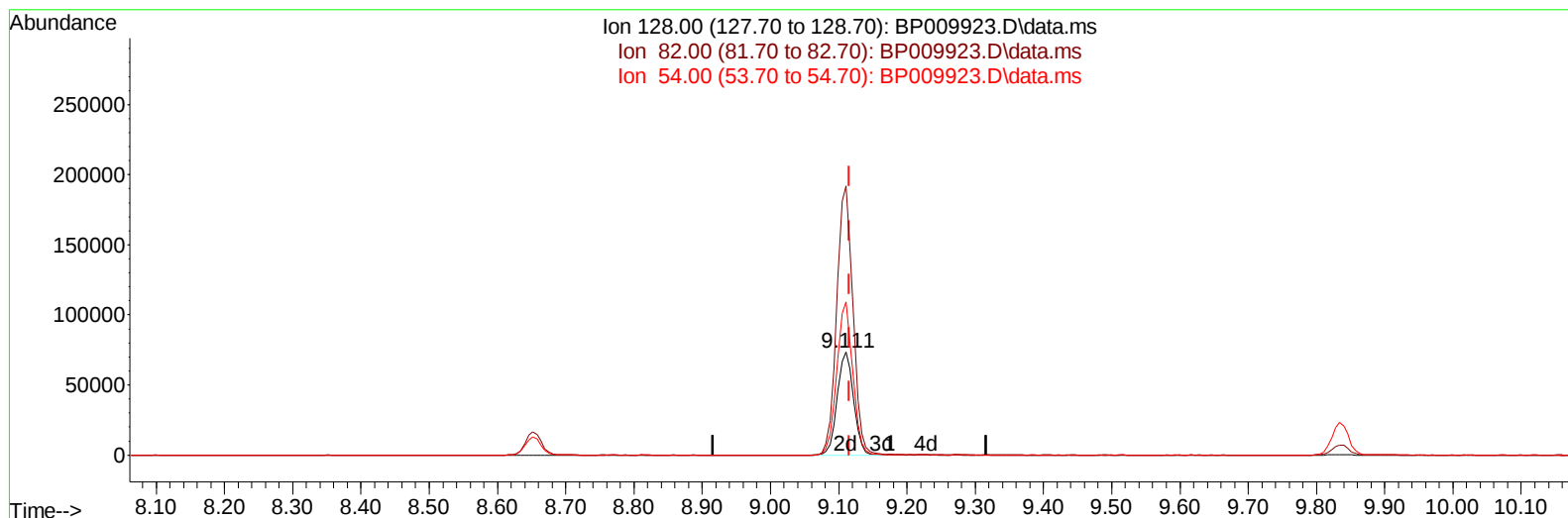
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TIC: BP009923.D\data.ms

(21) Nitrobenzene-d5 (S)

9.111min (-0.006) 32.10 ng/ul m

response 121672

Ion	Exp%	Act%
128.00	100.00	100.00
82.00	287.30	261.04
54.00	120.20	148.92#
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.952	152	113670	20.000	ng/ul	0.00
20) Naphthalene-d8	10.757	136	525687	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.581	164	388297	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.334	188	884309	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.416	240	885008	20.000	ng/ul	# 0.00
88) Perylene-d12	23.874	264	791595	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	11783	4.626	ng/uL	0.00
4) Pyridine-d5	3.799	84	65646	9.348	ng/ul	0.00
7) Phenol-d5	7.105	99	61598	6.172	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	183211	30.838	ng/ul	0.00
11) 2-Chlorophenol-d4	7.481	132	173445	23.041	ng/ul	0.00
15) 4-Methylphenol-d8	8.652	113	126057	15.131	ng/ul	-0.01
21) Nitrobenzene-d5	9.111	128	121672m	32.096	ng/ul	0.00
24) 2-Nitrophenol-d4	9.834	143	127274	30.623	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.375	165	224755	25.843	ng/ul	0.00
31) 4-Chloroaniline-d4	10.887	131	306860	24.773	ng/ul	0.00
46) Dimethylphthalate-d6	13.993	166	1006571	33.001	ng/ul	0.00
49) Acenaphthylene-d8	14.275	160	1062038	29.390	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.763	143	28855	5.216	ng/ul	0.00
60) Fluorene-d10	15.575	176	907704	35.376	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.687	200	158902	30.333	ng/ul	0.00
73) Anthracene-d10	17.434	188	1591334	37.585	ng/ul	0.00
81) Pyrene-d10	19.663	212	1934103	40.015	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.721	264	1602721	38.879	ng/ul	0.00

Target Compounds Qvalue

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

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