

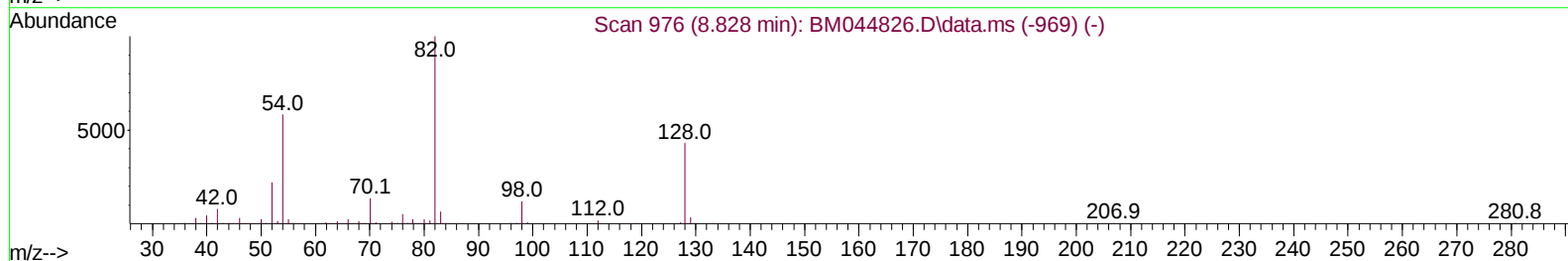
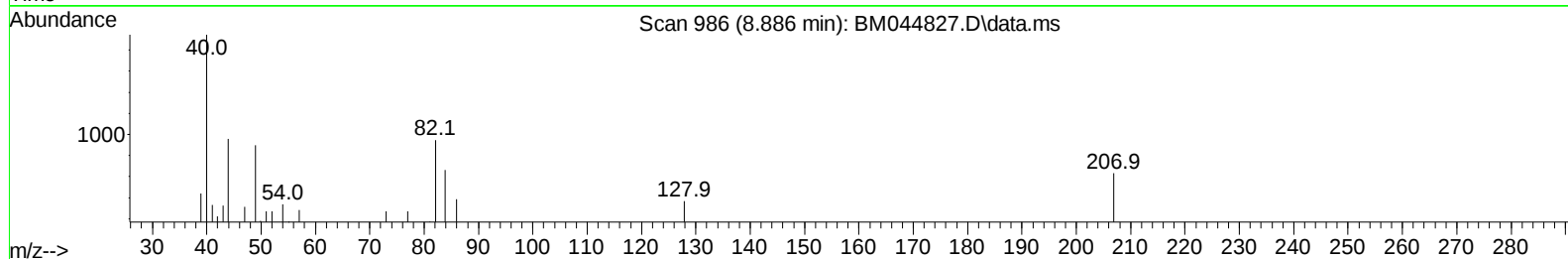
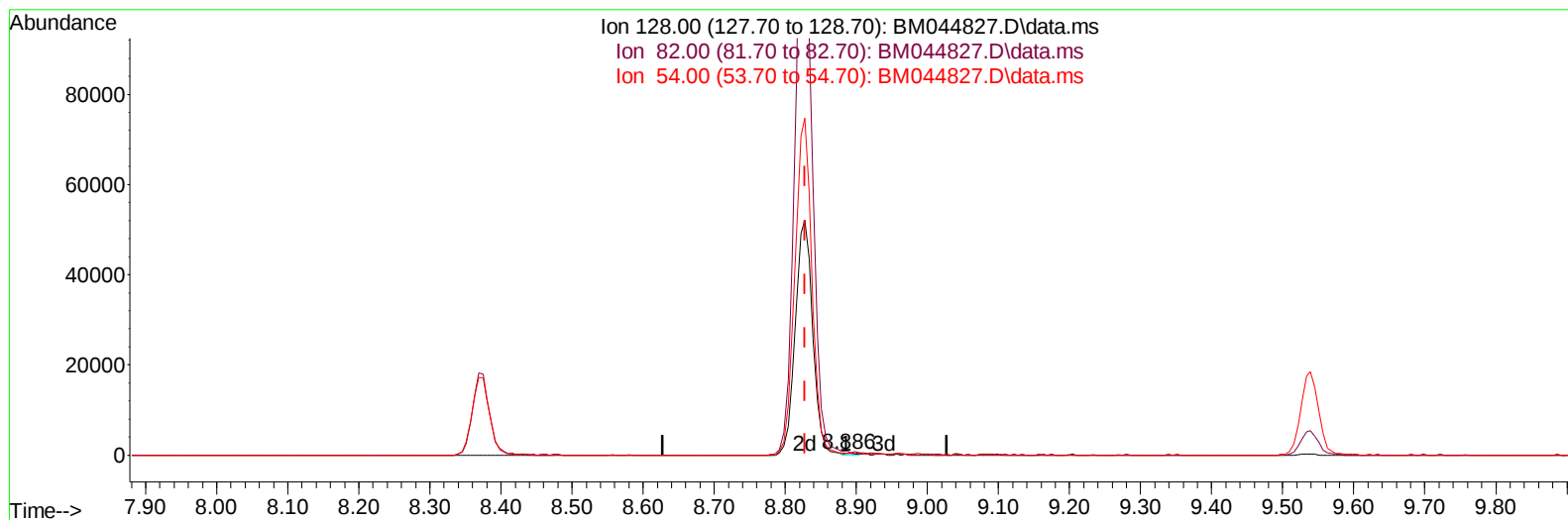
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM033024\
Data File : BM044827.D
Acq On : 30 Mar 2024 12:02
Operator : MA/JU
Sample : PB159819BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SBLK819

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 04/01/2024
Supervised By :mohammad ahmed 04/02/2024

Quant Time: Mar 31 13:57:55 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030524.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Sun Mar 31 13:54:26 2024
Response via : Initial Calibration



TIC: BM044827.D\data.ms

(21) Nitrobenzene-d5 (S)

8.886min (+ 0.059) 0.12 ng/u1

response 418

Ion	Exp%	Act%
128.00	100.00	100.00
82.00	224.70	258.58
54.00	110.50	92.64
0.00	0.00	0.00

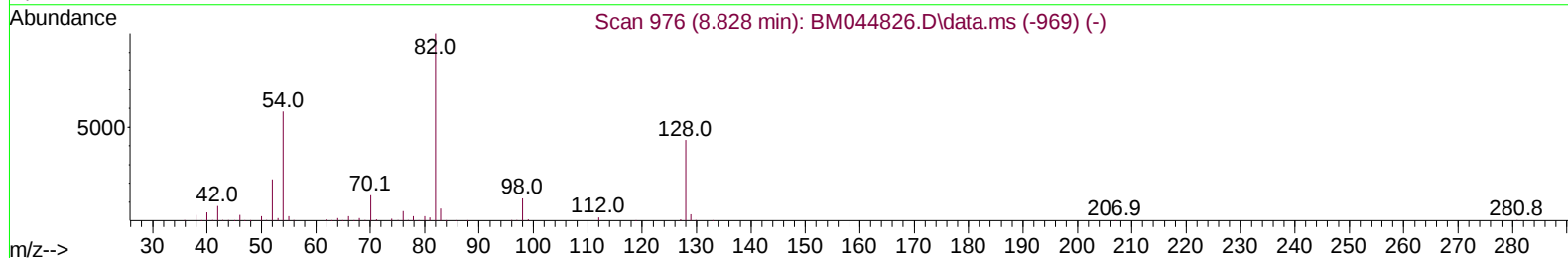
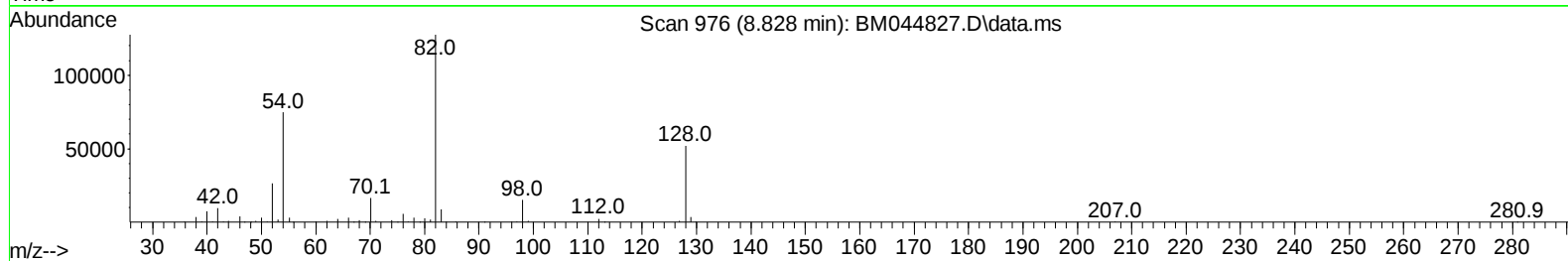
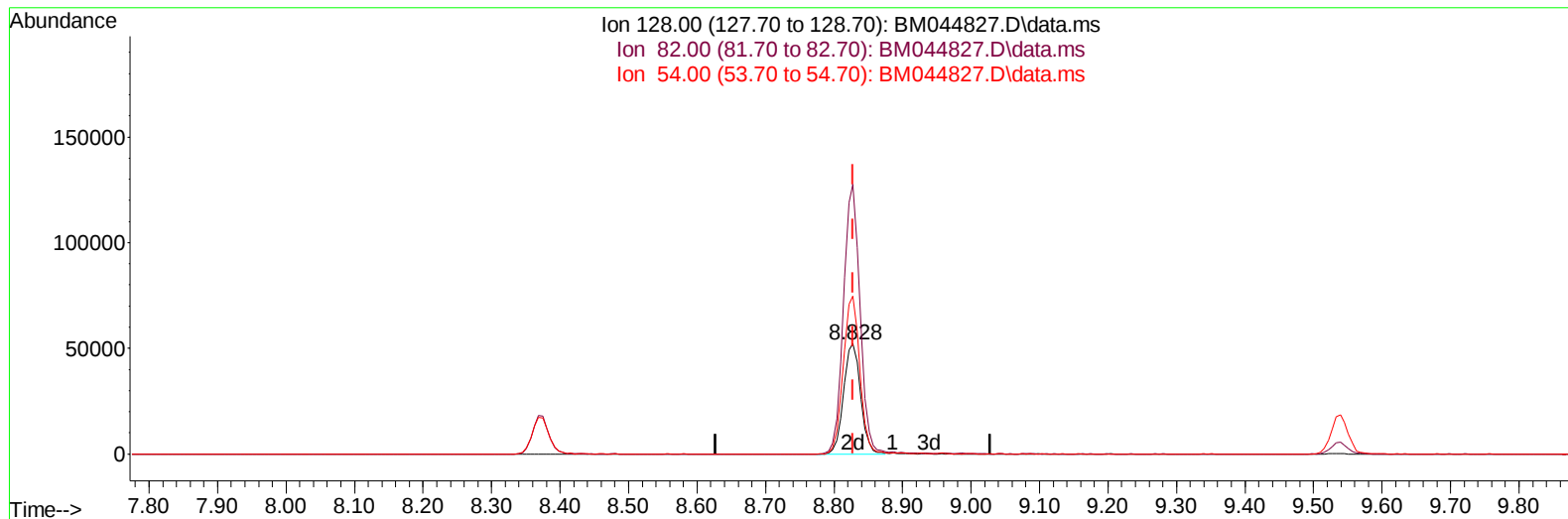
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TIC: BM044827.D\data.ms

(21) Nitrobenzene-d5 (S)

8.828min (-0.000) 25.90 ng/ul m

response 88707

Ion	Exp%	Act%
128.00	100.00	100.00
82.00	224.70	244.54
54.00	110.50	143.36#
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 04/01/2024
 Supervised By :mohammad ahmed 04/02/2024

Quant Time: Mar 31 15:00:33 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030524.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sun Mar 31 13:54:26 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.663	152	97896	20.000	ng/ul	0.00
20) Naphthalene-d8	10.445	136	406697	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.310	164	243341	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.068	188	482289	20.000	ng/ul	0.00
79) Chrysene-d12	21.268	240	444690	20.000	ng/ul	0.00
88) Perylene-d12	23.497	264	497944	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	13074	4.530	ng/uL	0.00
4) Pyridine-d5	3.634	84	125400	14.761	ng/ul	0.00
7) Phenol-d5	6.845	99	205928	20.460	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.016	67	146766	23.608	ng/ul	0.00
11) 2-Chlorophenol-d4	7.198	132	178384	24.118	ng/ul	0.00
15) 4-Methylphenol-d8	8.369	113	146668	18.555	ng/ul	0.00
21) Nitrobenzene-d5	8.828	128	88707m	25.902	ng/ul	0.00
24) 2-Nitrophenol-d4	9.539	143	98666	26.725	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.069	165	156261	24.869	ng/ul	0.00
31) 4-Chloroaniline-d4	10.604	131	25393	2.433	ng/ul	0.01
46) Dimethylphthalate-d6	13.733	166	490873	25.716	ng/ul	0.00
49) Acenaphthylene-d8	14.004	160	515337	23.606	ng/ul	0.00
54) 4-Nitrophenol-d4	14.539	143	92072	23.259	ng/ul	0.00
60) Fluorene-d10	15.309	176	402077	25.065	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.445	200	79889	25.518	ng/ul	0.00
73) Anthracene-d10	17.168	188	620946	26.935	ng/ul	0.00
81) Pyrene-d10	19.468	212	728416	26.707	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.362	264	627407	23.972	ng/ul	0.00

Target Compounds Qvalue

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

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