

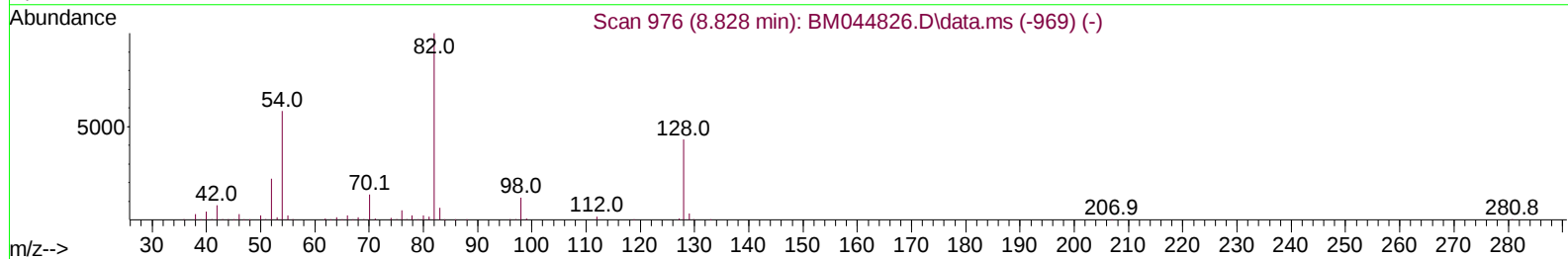
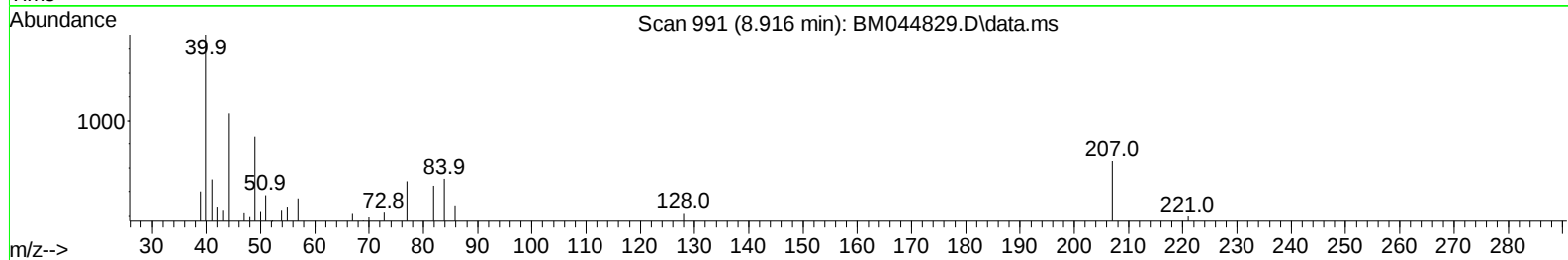
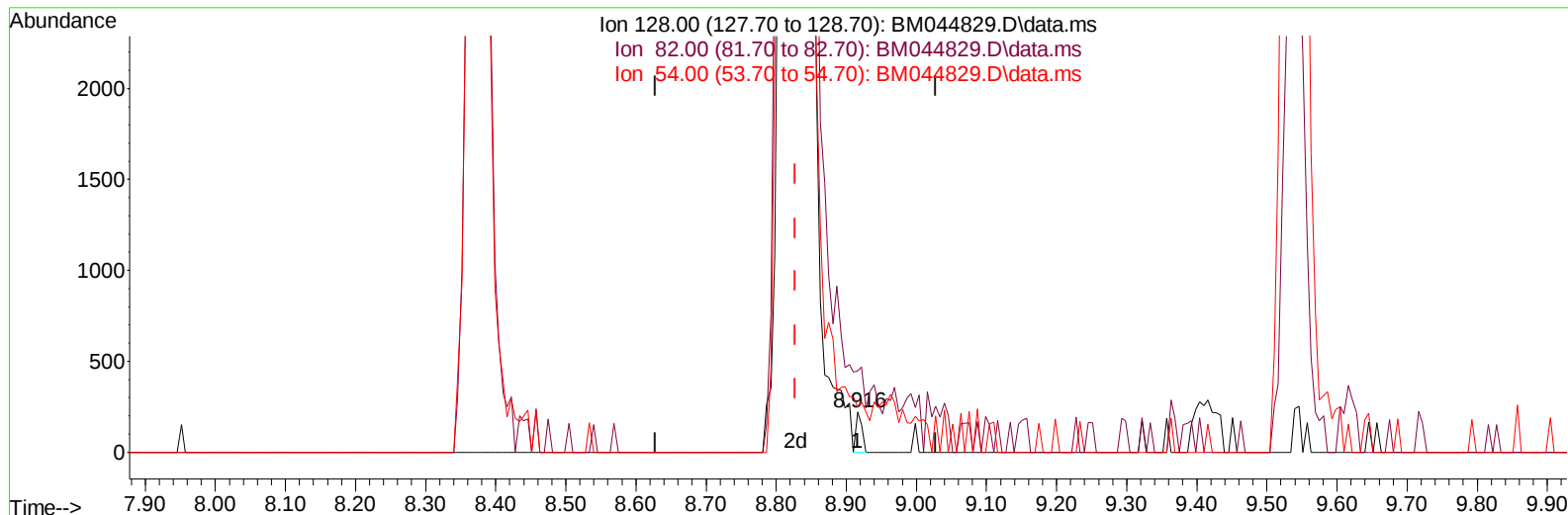
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM033024\
Data File : BM044829.D
Acq On : 30 Mar 2024 13:15
Operator : MA/JU
Sample : P1887-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
E0795

Manual IntegrationsAPPROVED

Quant Time: Mar 31 13:58:11 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

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



TIC: BM044829.D\data.ms

(21) Nitrobenzene-d5 (S)

8.916min (+ 0.088) 0.04 ng/u1

response 131

Ion	Exp%	Act%
128.00	100.00	100.00
82.00	224.70	202.71
54.00	110.50	112.22
0.00	0.00	0.00

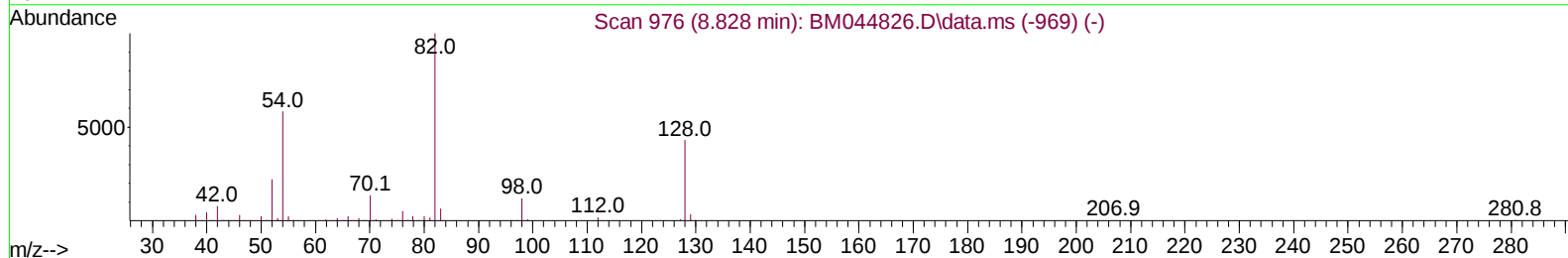
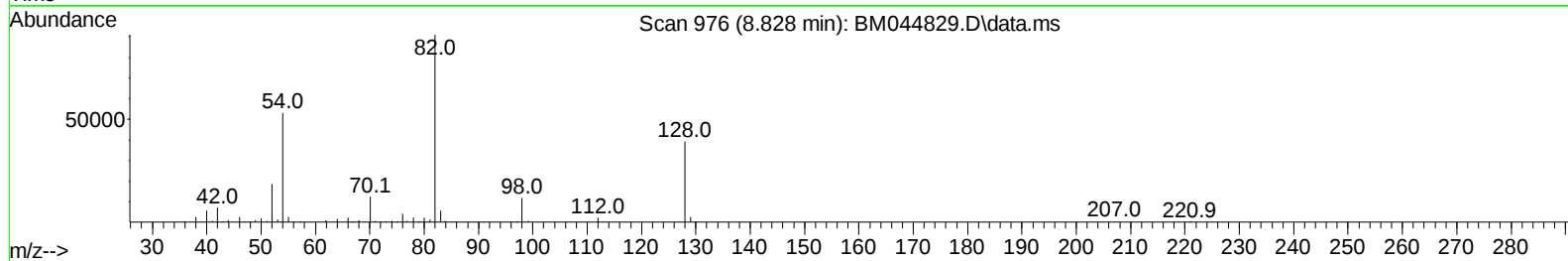
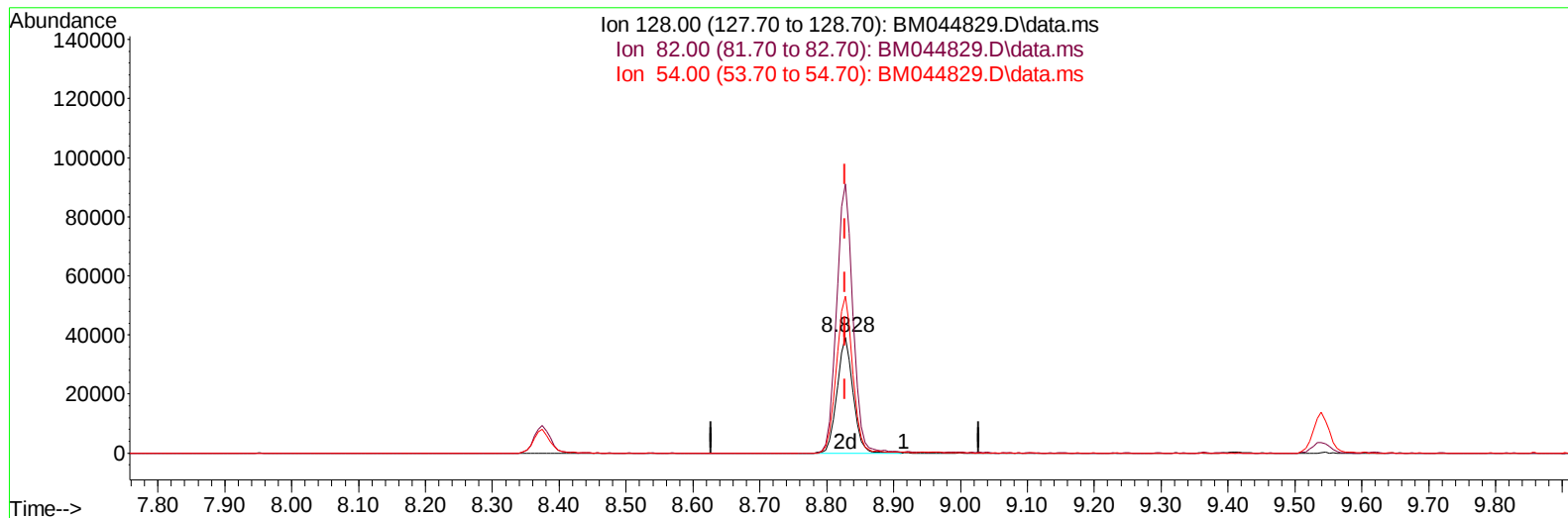
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TIC: BM044829.D\data.ms

(21) Nitrobenzene-d5 (S)

8.828min (+ 0.000) 20.52 ng/ul m

response 65141

Ion	Exp%	Act%
128.00	100.00	100.00
82.00	224.70	232.13
54.00	110.50	135.62#
0.00	0.00	0.00

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 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 E0795

Manual IntegrationsAPPROVED

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

Quant Time: Mar 31 15:12:17 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.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.669	152		94433	20.000	ng/ul	0.00
20) Naphthalene-d8	10.445	136		377046	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.316	164		214986	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.068	188		421640	20.000	ng/ul	0.00
79) Chrysene-d12	21.268	240		382665	20.000	ng/ul	0.00
88) Perylene-d12	23.503	264		430903	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.234	96		8044	2.889	ng/uL	0.00
4) Pyridine-d5	3.652	84		18999	2.318	ng/ul	0.02
7) Phenol-d5	6.846	99		153703	15.831	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.016	67		107424	17.913	ng/ul	0.00
11) 2-Chlorophenol-d4	7.199	132		129271	18.119	ng/ul	0.00
15) 4-Methylphenol-d8	8.375	113		75307	9.877	ng/ul	0.00
21) Nitrobenzene-d5	8.828	128		65141m	20.517	ng/ul	0.00
24) 2-Nitrophenol-d4	9.540	143		71743	20.961	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.075	165		107592	18.470	ng/ul	0.00
31) 4-Chloroaniline-d4	10.616	131		3135	0.324	ng/ul	0.02
46) Dimethylphthalate-d6	13.733	166		334522	19.836	ng/ul	0.00
49) Acenaphthylene-d8	14.004	160		391910	20.320	ng/ul	0.00
54) 4-Nitrophenol-d4	14.545	143		43729	12.504	ng/ul	0.01
60) Fluorene-d10	15.316	176		294974	20.814	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.451	200		32707	11.950	ng/ul	0.00
73) Anthracene-d10	17.168	188		434394	21.554	ng/ul	0.00
81) Pyrene-d10	19.468	212		506026	21.560	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.362	264		497934	21.985	ng/ul	0.00
Target Compounds							
						Qvalue	
16) Acetophenone	8.498	105		13947	1.179	ng/ul#	85

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

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