

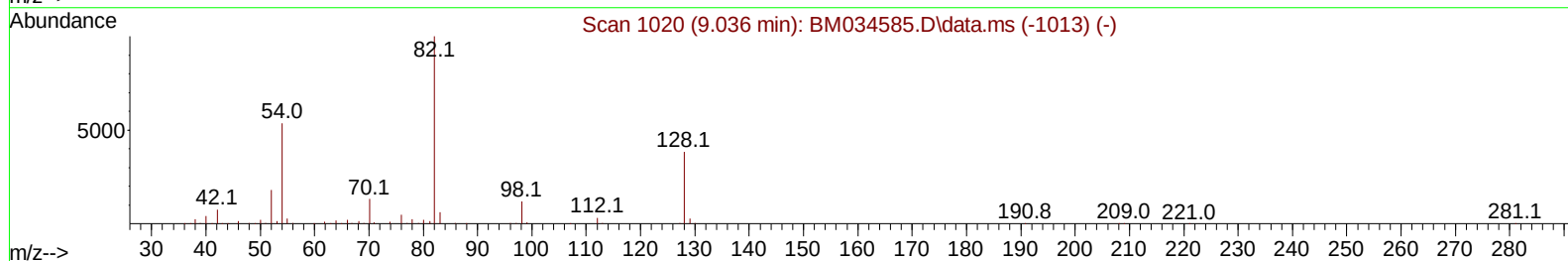
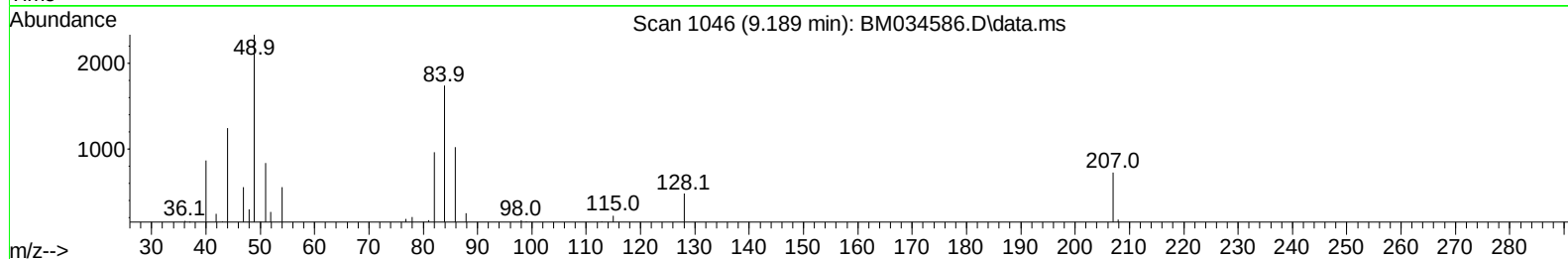
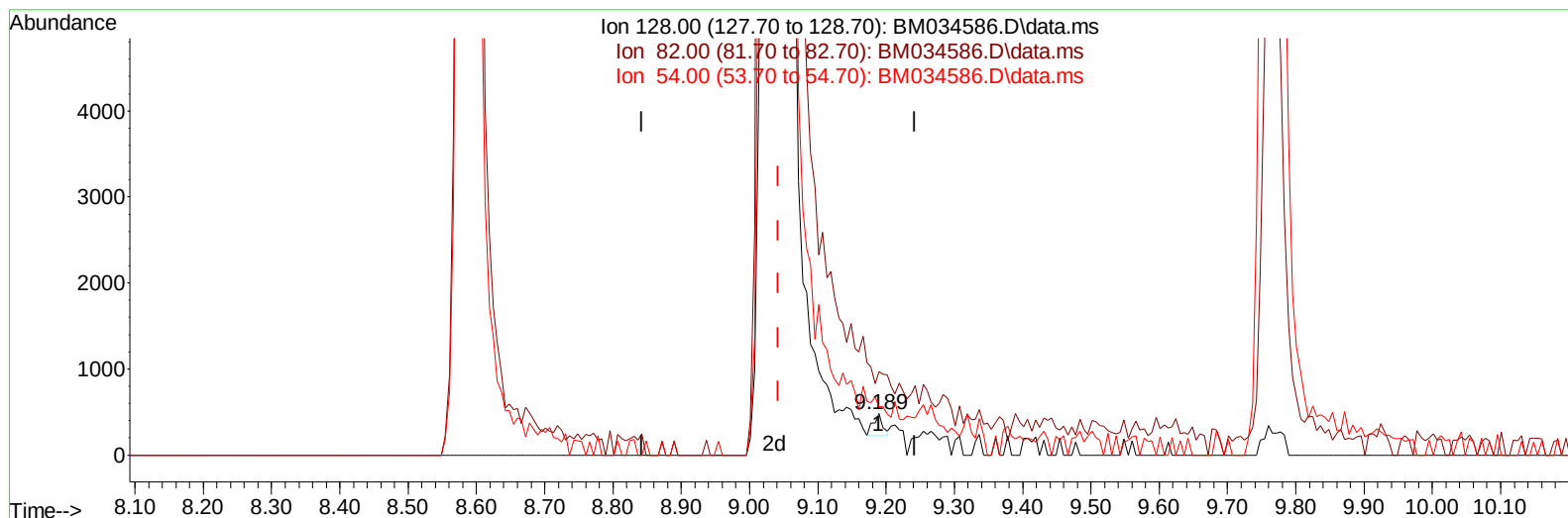
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040922\
 Data File : BM034586.D
 Acq On : 09 Apr 2022 07:16
 Operator : CG/JU
 Sample : PB143976BL
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK976

Manual IntegrationsAPPROVED

Quant Time: Apr 11 00:57:45 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Apr 08 14:30:08 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/11/2022
 Supervised By :Yogesh Patel 04/12/2022



TIC: BM034586.D\data.ms

(21) Nitrobenzene-d5 (S)

9.189min (+ 0.147) 0.07 ng/u1

response 247

Ion	Exp%	Act%
128.00	100.00	100.00
82.00	228.00	200.41
54.00	100.50	116.18
0.00	0.00	0.00

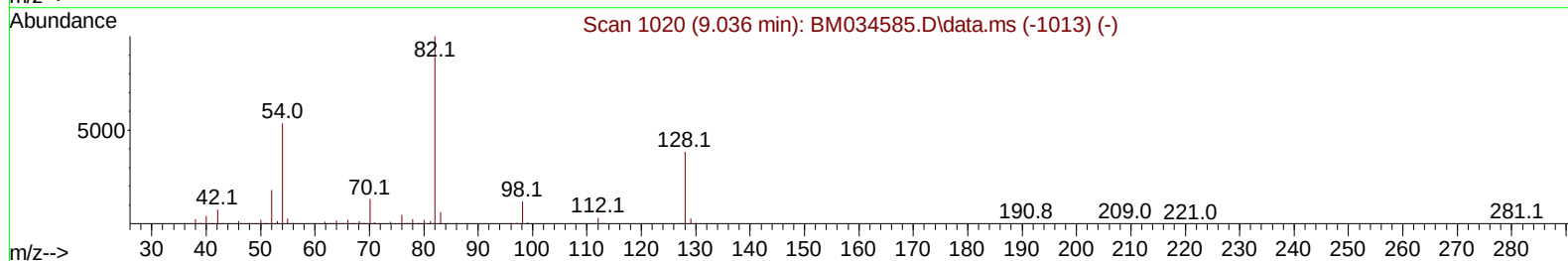
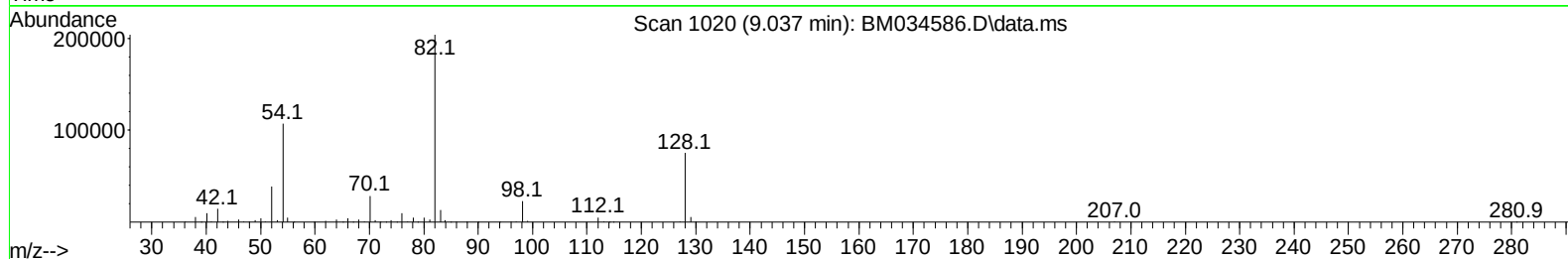
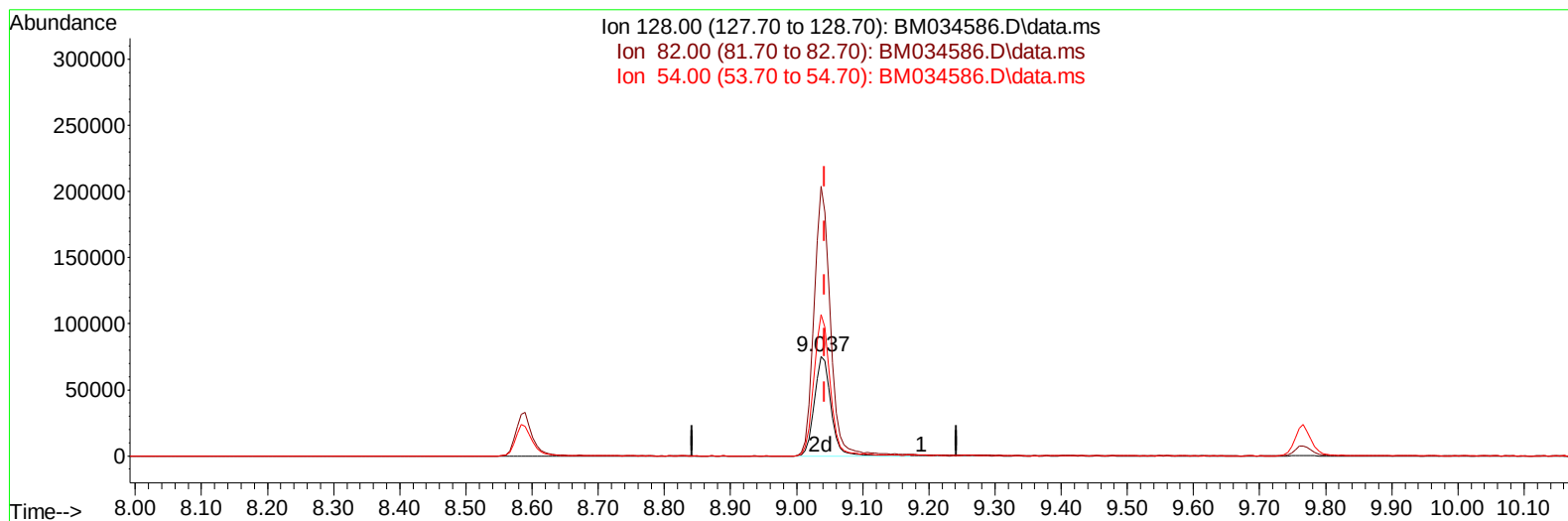
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TIC: BM034586.D\data.ms

(21) Nitrobenzene-d5 (S)

9.037min (-0.006) 38.15 ng/ul m

response 131825

Ion	Exp%	Act%
128.00	100.00	100.00
82.00	228.00	270.40
54.00	100.50	142.11#
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	7.878	152	119782	20.000	ng/ul	0.00
20) Naphthalene-d8	10.684	136	470177	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.524	164	261441	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.265	188	492761	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.442	240	454333	20.000	ng/ul	# 0.00
88) Perylene-d12	23.824	264	526225	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.255	96	24564	7.313	ng/uL	0.00
4) Pyridine-d5	3.678	84	292321	36.060	ng/ul	0.00
7) Phenol-d5	7.037	99	329571	33.057	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.201	67	242849	36.555	ng/ul	0.00
11) 2-Chlorophenol-d4	7.401	132	288631	37.451	ng/ul	0.00
15) 4-Methylphenol-d8	8.590	113	267279	34.321	ng/ul	0.00
21) Nitrobenzene-d5	9.037	128	131825m	38.149	ng/ul	0.00
24) 2-Nitrophenol-d4	9.766	143	135098	37.552	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.307	165	227143	33.530	ng/ul	0.00
31) 4-Chloroaniline-d4	10.825	131	314296	33.629	ng/ul	0.00
46) Dimethylphthalate-d6	13.936	166	727028	38.485	ng/ul	0.00
49) Acenaphthylene-d8	14.219	160	872166	35.920	ng/ul	0.00
54) 4-Nitrophenol-d4	14.719	143	53499	18.573	ng/ul	0.00
60) Fluorene-d10	15.513	176	618036	37.858	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.630	200	95941	31.893	ng/ul	0.00
73) Anthracene-d10	17.365	188	912014	38.820	ng/ul	0.00
81) Pyrene-d10	19.654	212	1092803	44.069	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.671	264	988533	37.345	ng/ul	0.00

Target Compounds Qvalue

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

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