

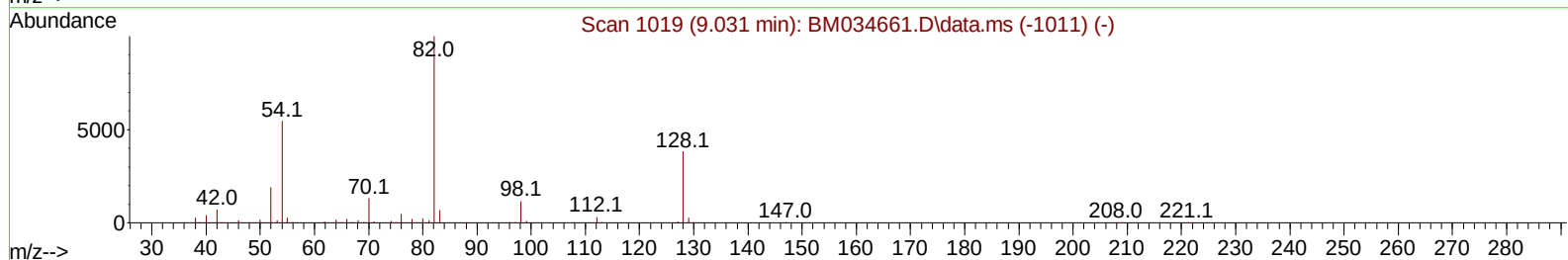
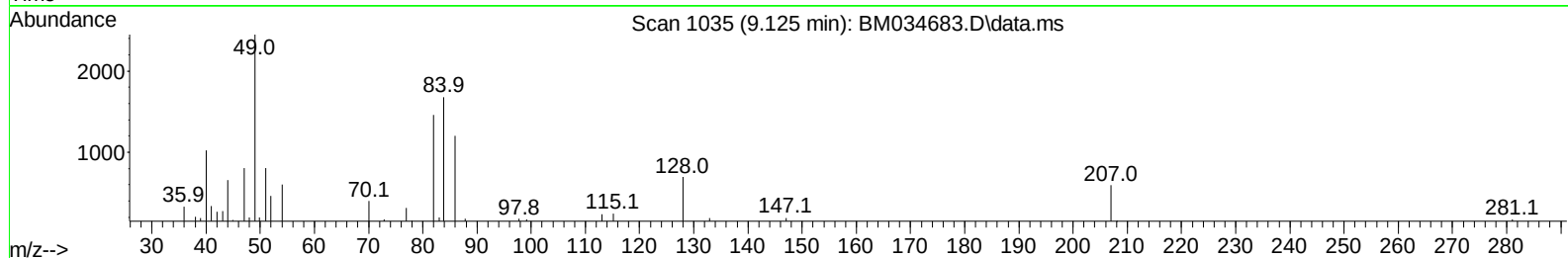
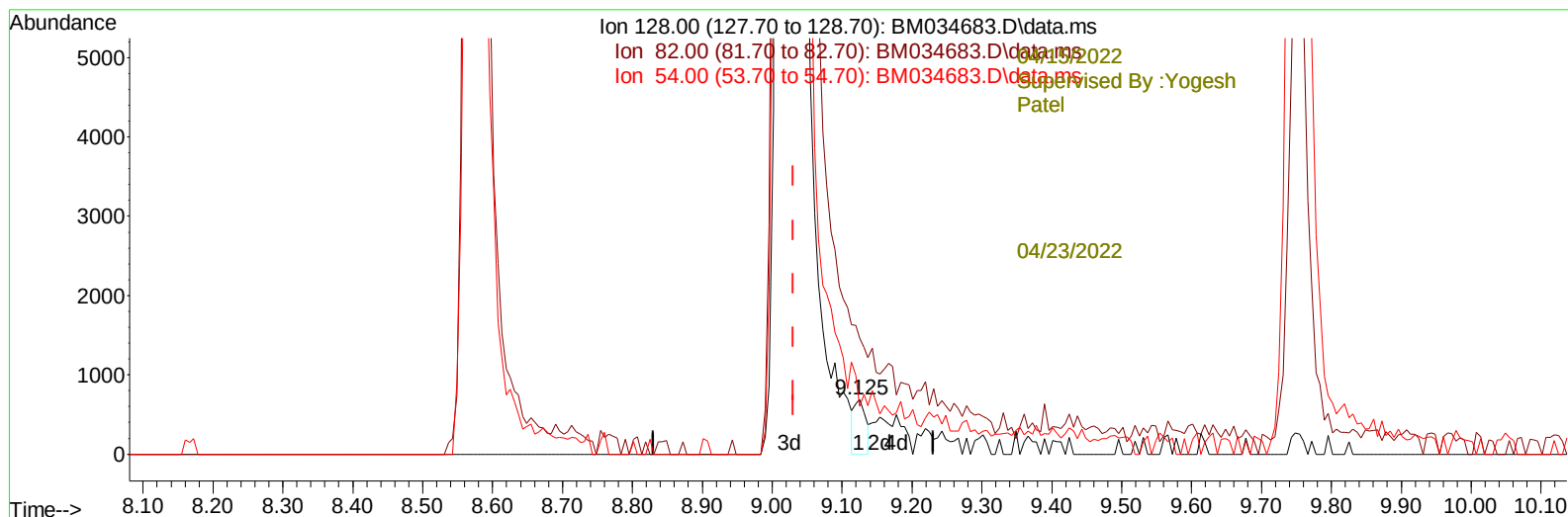
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041422\
 Data File : BM034683.D
 Acq On : 15 Apr 2022 00:39
 Operator : CG/JU
 Sample : N2316-22
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ETNS0

Manual IntegrationsAPPROVED

Quant Time: Apr 15 01:22:40 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 14 23:42:00 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/15/2022
 Supervised By :Yogesh Patel 04/23/2022



TIC: BM034683.D\data.ms

(21) Nitrobenzene-d5 (S)

9.125min (+ 0.094) 0.22 ng/u1

response 785

Ion	Exp%	Act%
128.00	100.00	100.00
82.00	228.00	211.27
54.00	100.50	87.43
0.00	0.00	0.00

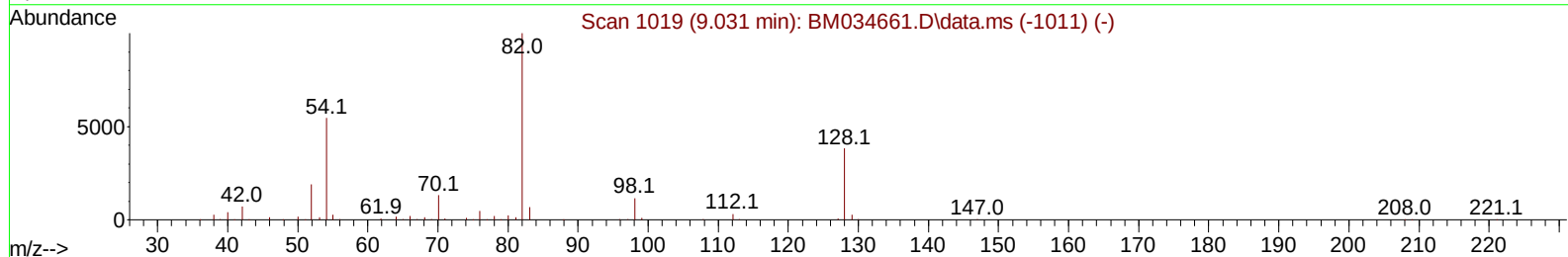
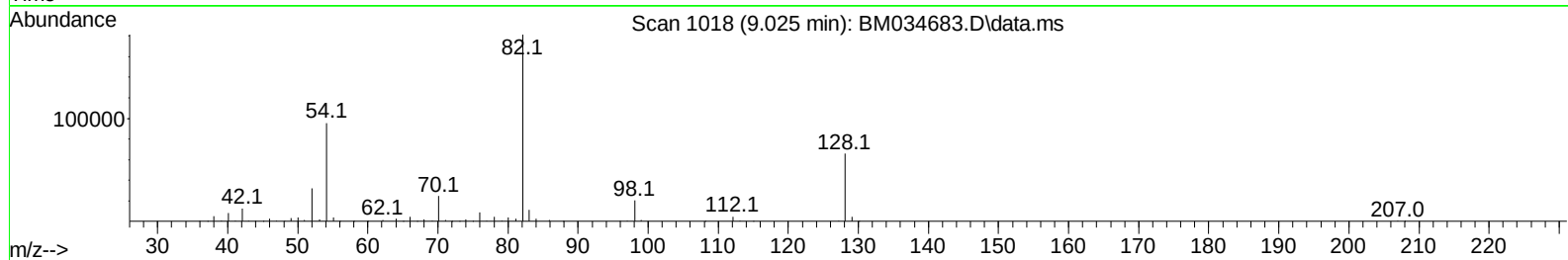
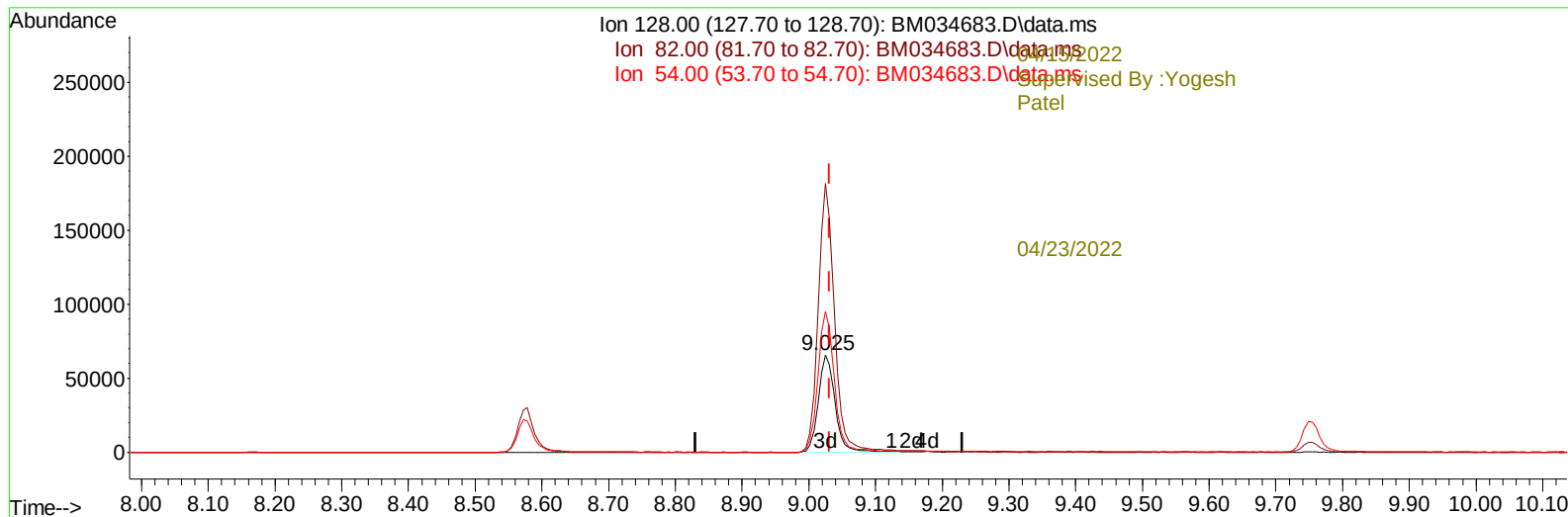
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(21) Nitrobenzene-d5 (S)

9.025min (-0.006) 32.15 ng/ul m

response 114093

Ion	Exp%	Act%
128.00	100.00	100.00
82.00	228.00	276.08#
54.00	100.50	145.19#
0.00	0.00	0.00

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 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)
-----04/15/2022							
Supervised By :Yogesh Patel							
Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.866	152		117415	20.000	ng/ul	0.00
20) Naphthalene-d8	10.672	136		482926	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.513	164		288613	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.254	188		567722	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.430	240		535654	20.000	ng/ul	# 0.00
88) Perylene-d12	23.801	264		605962	20.000	ng/ul	-0.01
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.243	96		16391	4.978	ng/uL	0.00
4) Pyridine-d5	3.667	84		203211	25.573	ng/ul	0.00
7) Phenol-d5	7.025	99		281793	28.835	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.190	67		208994	32.093	ng/ul	0.00
11) 2-Chlorophenol-d4	7.390	132		239930	31.759	ng/ul	0.00
15) 4-Methylphenol-d8	8.578	113		239953	31.433	ng/ul	0.00
21) Nitrobenzene-d5	9.025	128		114093m	32.146	ng/ul	0.00
24) 2-Nitrophenol-d4	9.754	143		119200	32.258	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.296	165		207250	29.786	ng/ul	0.00
31) 4-Chloroaniline-d4	10.807	131		280913	29.263	ng/ul	0.00
46) Dimethylphthalate-d6	13.925	166		726379	34.831	ng/ul	0.00
49) Acenaphthylene-d8	14.201	160		859316	32.059	ng/ul	0.00
54) 4-Nitrophenol-d4	14.713	143		55456	17.440	ng/ul	0.00
60) Fluorene-d10	15.501	176		623891	34.619	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.619	200		97213	28.049	ng/ul	0.00
73) Anthracene-d10	17.354	188		916243	33.851	ng/ul	0.00
81) Pyrene-d10	19.642	212		1135769	38.849	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.648	264		1063315	34.884	ng/ul	-0.01

Target Compounds Qvalue

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

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