

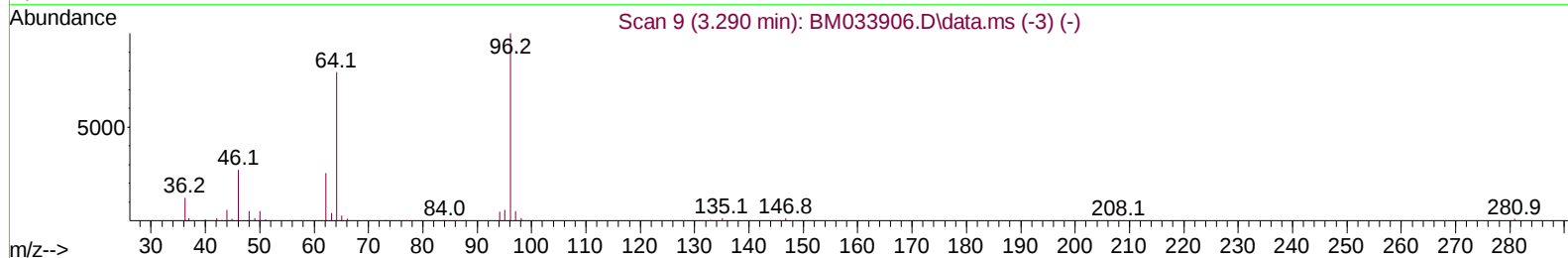
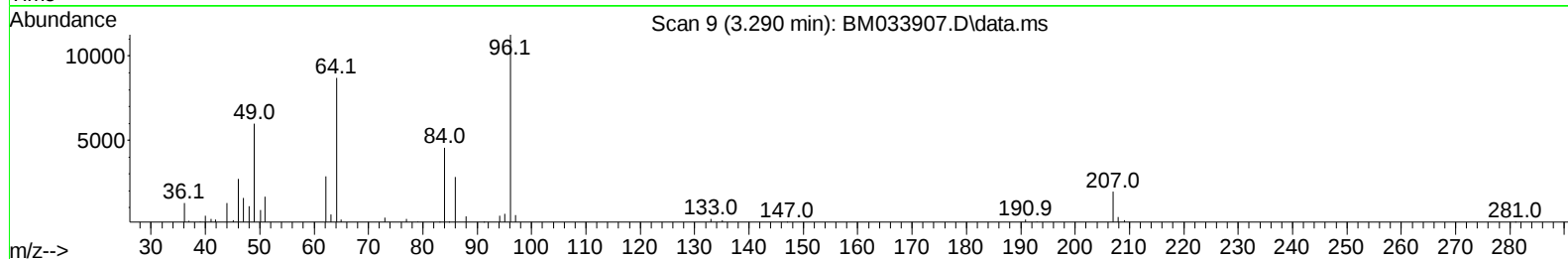
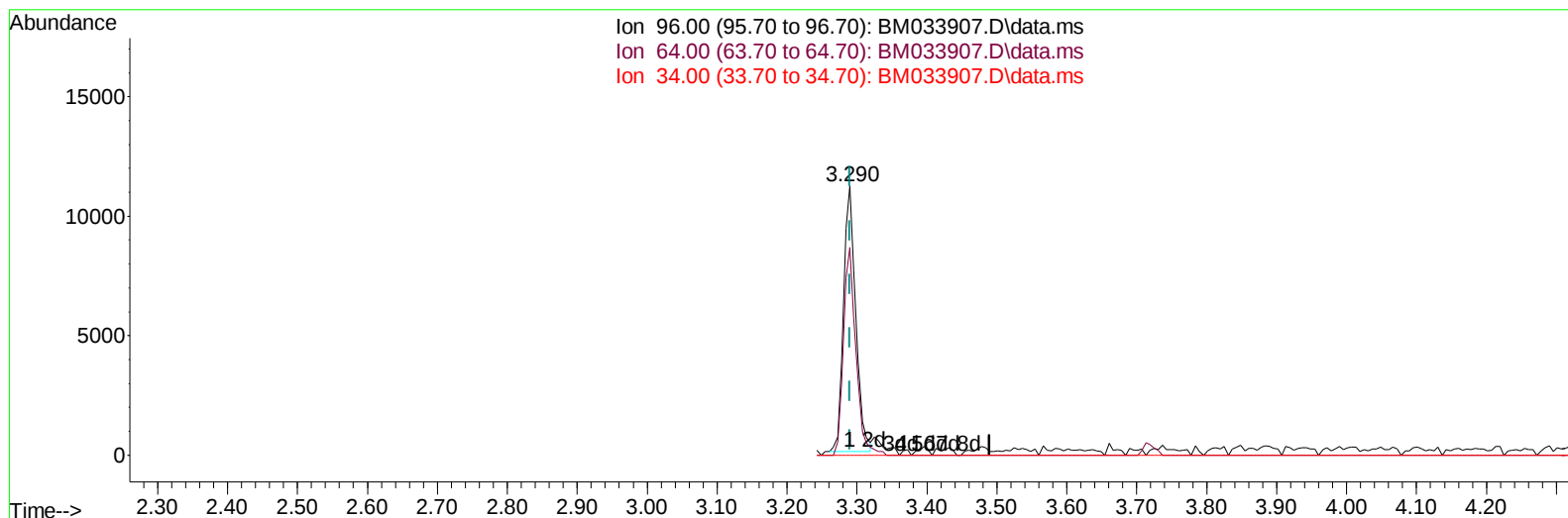
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012822\
 Data File : BM033907.D
 Acq On : 28 Jan 2022 17:08
 Operator : CG/JU
 Sample : PB142334BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK334

Manual IntegrationsAPPROVED

Quant Time: Jan 28 23:27:16 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM010622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jan 28 23:26:29 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 01/31/2022
 Supervised By :mohammad ahmed 02/01/2022



TIC: BM033907.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.290min (-0.000) 5.06 ng/uL

response 13597

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.20	77.11
34.00	0.00	0.00
0.00	0.00	0.00

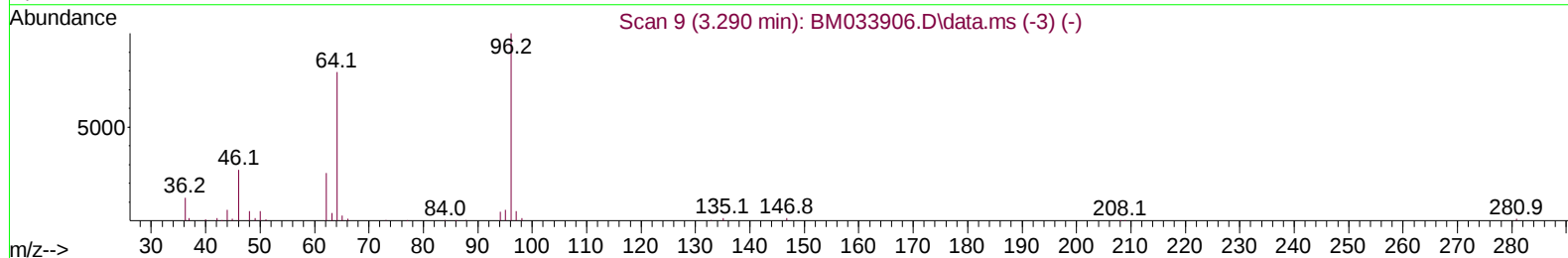
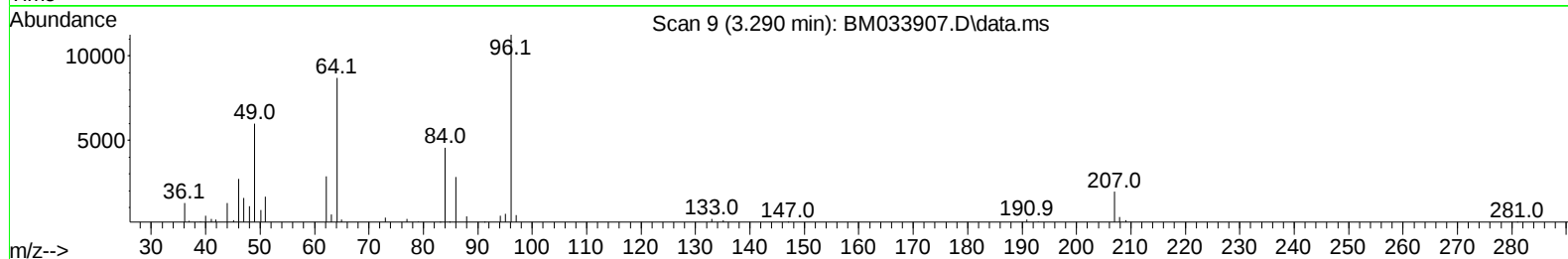
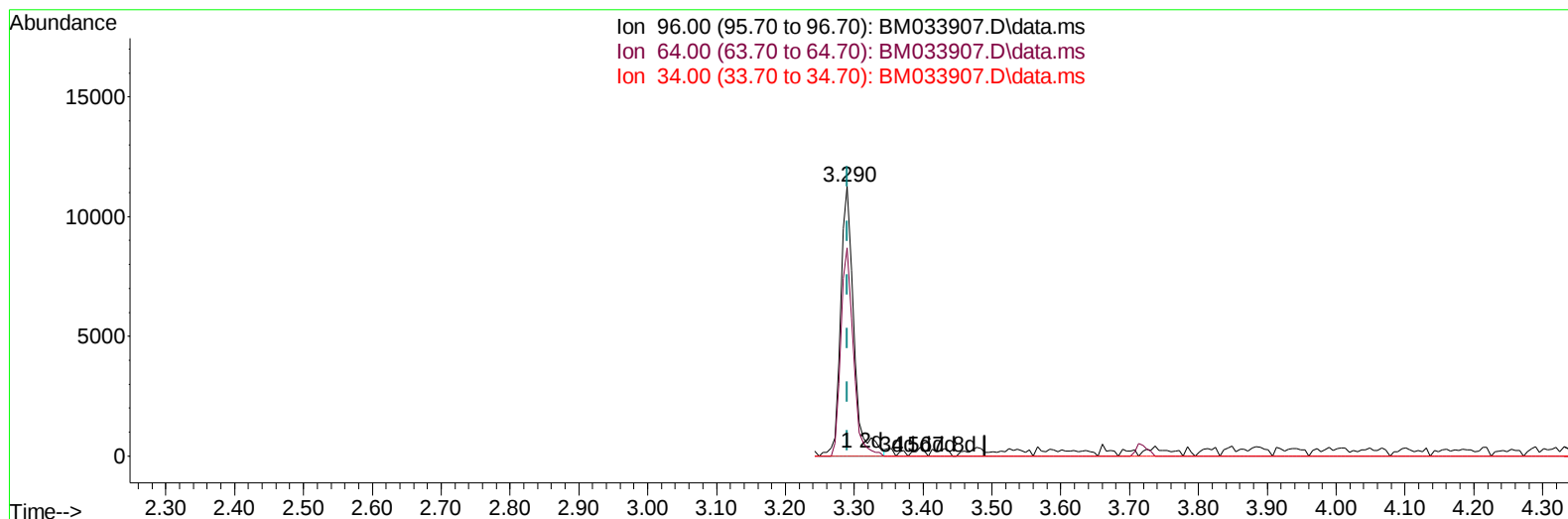
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TIC: BM033907.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.290min (-0.000) 5.55 ng/uL m

response 14933

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.20	77.11
34.00	0.00	0.00
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	7.931	152	101051	20.000	ng/ul	0.00
20) Naphthalene-d8	10.742	136	415163	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.571	164	260900	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.306	188	550325	20.000	ng/ul	0.00
79) Chrysene-d12	21.471	240	542905	20.000	ng/ul	0.00
88) Perylene-d12	23.865	264	542452	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.290	96	14933m	5.552	ng/uL	0.00
4) Pyridine-d5	3.719	84	182392	24.420	ng/ul	0.00
7) Phenol-d5	7.089	99	224330	24.717	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.254	67	144103	26.072	ng/ul	0.00
11) 2-Chlorophenol-d4	7.460	132	187785	27.062	ng/ul	0.00
15) 4-Methylphenol-d8	8.642	113	178664	23.780	ng/ul	0.00
21) Nitrobenzene-d5	9.095	128	88771	27.075	ng/ul	0.00
24) 2-Nitrophenol-d4	9.819	143	91668	26.381	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.360	165	158659	23.817	ng/ul	0.00
31) 4-Chloroaniline-d4	10.877	131	238411	24.077	ng/ul	0.00
46) Dimethylphthalate-d6	13.977	166	551375	26.823	ng/ul	0.00
49) Acenaphthylene-d8	14.265	160	636382	26.242	ng/ul	0.00
54) 4-Nitrophenol-d4	14.754	143	66558	16.871	ng/ul	0.00
60) Fluorene-d10	15.560	176	458050	26.859	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.671	200	64574	19.726	ng/ul	0.00
73) Anthracene-d10	17.401	188	767936	28.758	ng/ul	0.00
81) Pyrene-d10	19.689	212	891988	30.424	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.712	264	843500	29.885	ng/ul	0.00

Target Compounds Qvalue

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

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