

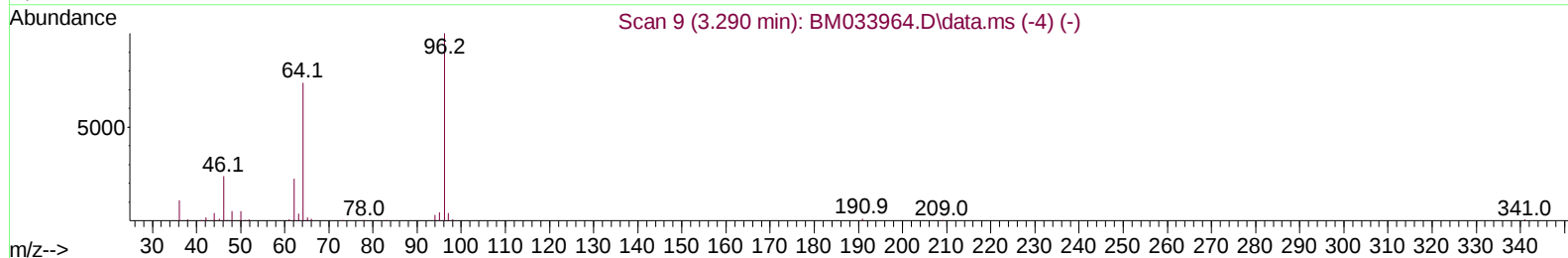
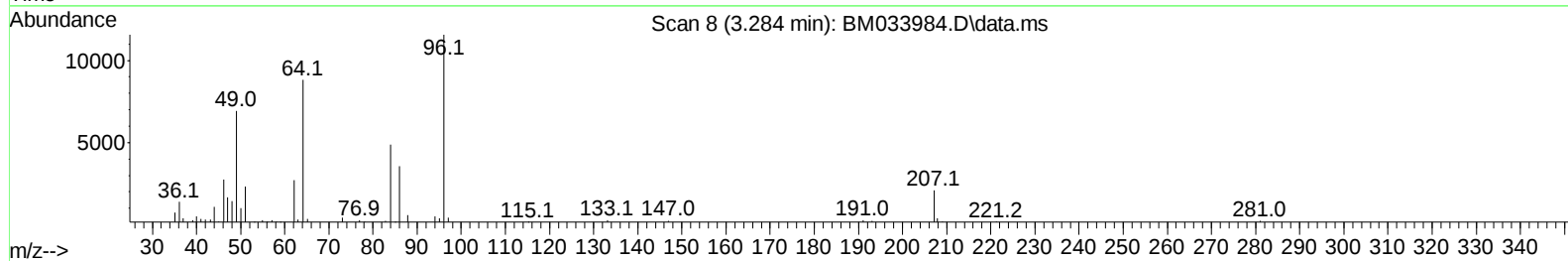
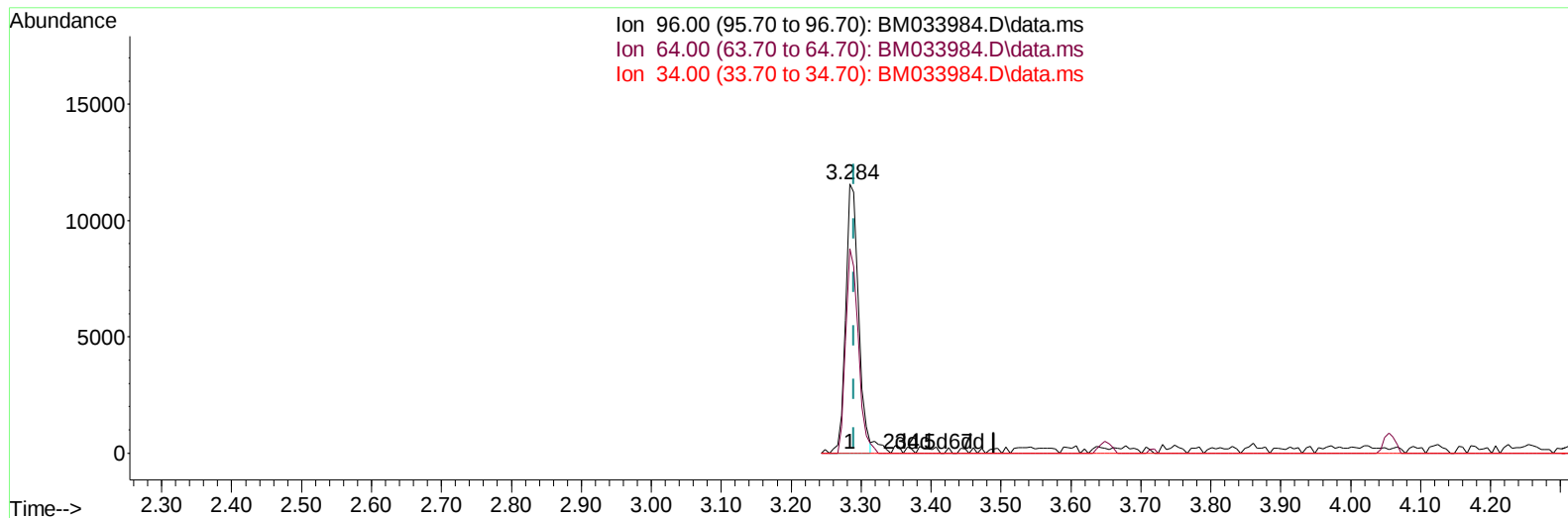
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
 Data File : BM033984.D
 Acq On : 02 Feb 2022 21:24
 Operator : CG/JU
 Sample : N1355-10
 Misc :
 ALS Vial : 55 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0VF4

Manual IntegrationsAPPROVED

Quant Time: Feb 03 00:15:51 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Feb 01 02:16:28 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 02/03/2022
 Supervised By :Yogesh Patel 02/04/2022



TIC: BM033984.D\data.ms

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

3.284min (-0.006) 4.84 ng/uL

response 14720

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.20	76.12
34.00	0.00	0.00
0.00	0.00	0.00

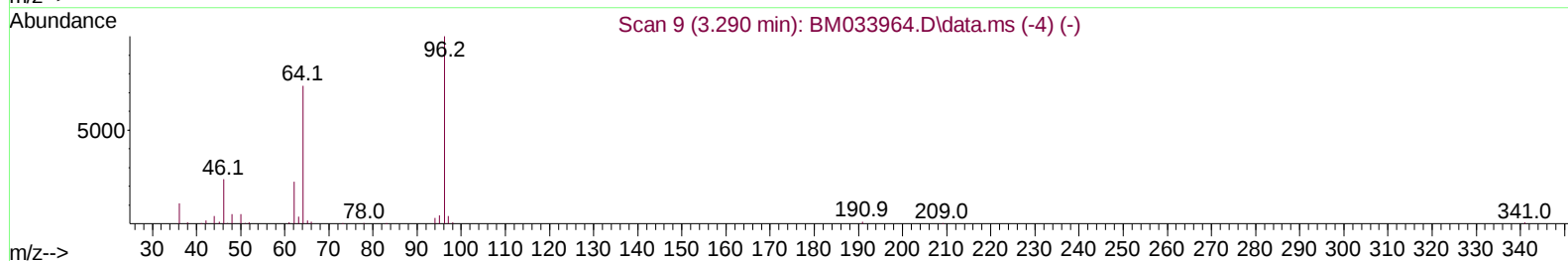
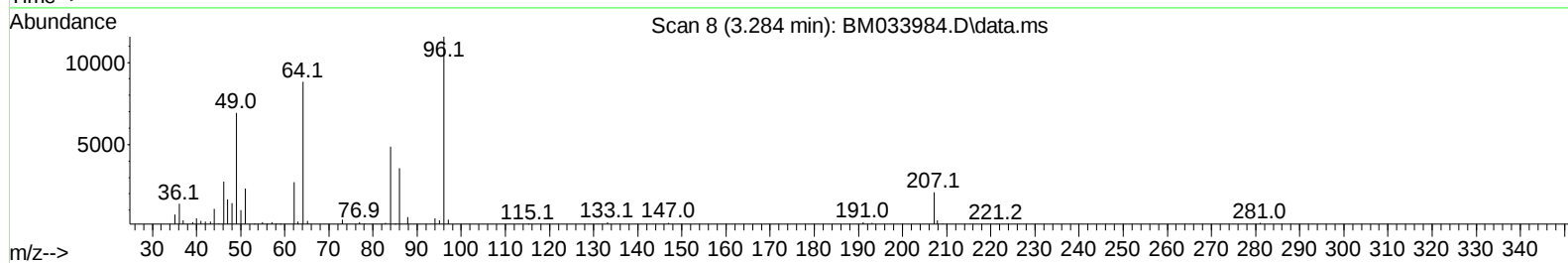
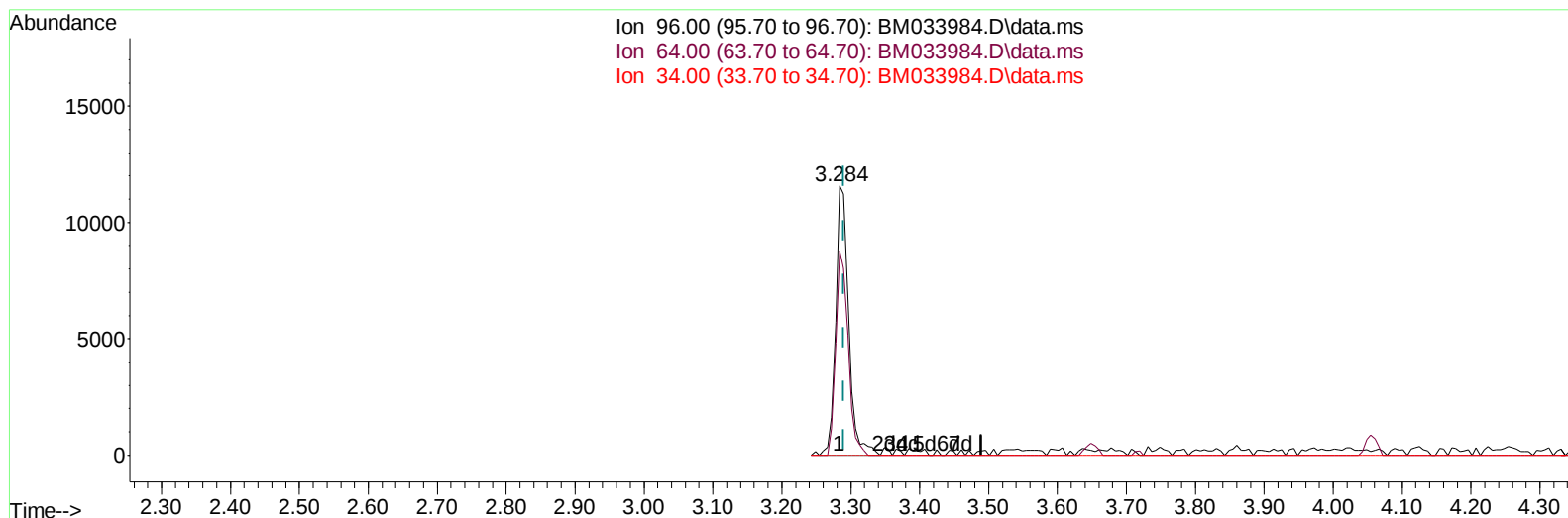
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TIC: BM033984.D\data.ms

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

3.284min (-0.006) 5.01 ng/uL m

response 15215

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.20	76.12
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	7.925	152	119847	20.000	ng/ul	0.00
20) Naphthalene-d8	10.736	136	502432	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.560	164	299790	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.295	188	550240	20.000	ng/ul	-0.01
79) Chrysene-d12	21.459	240	502892	20.000	ng/ul	0.00
88) Perylene-d12	23.847	264	537975	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.284	96	15215m	5.007	ng/uL	0.00
4) Pyridine-d5	3.719	84	69894	8.240	ng/ul	0.00
7) Phenol-d5	7.078	99	288617	28.746	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.248	67	177830	29.052	ng/ul	0.00
11) 2-Chlorophenol-d4	7.448	132	235846	29.319	ng/ul	-0.01
15) 4-Methylphenol-d8	8.631	113	234896	29.524	ng/ul	-0.01
21) Nitrobenzene-d5	9.083	128	115839	28.734	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.813	143	126428	29.040	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.354	165	218277	27.706	ng/ul	0.00
31) 4-Chloroaniline-d4	10.866	131	255257	23.463	ng/ul	0.00
46) Dimethylphthalate-d6	13.971	166	643631	28.731	ng/ul	0.00
49) Acenaphthylene-d8	14.254	160	796217	29.013	ng/ul	0.00
54) 4-Nitrophenol-d4	14.748	143	77716	20.305	ng/ul	0.00
60) Fluorene-d10	15.548	176	532328	28.458	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.659	200	76396	22.116	ng/ul	-0.01
73) Anthracene-d10	17.395	188	764324	29.005	ng/ul	0.00
81) Pyrene-d10	19.677	212	831776	28.016	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.694	264	805176	28.495	ng/ul	-0.01
Target Compounds						
					Qvalue	
8) Phenol	7.107	94	15923	1.546	ng/ul#	83

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

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