

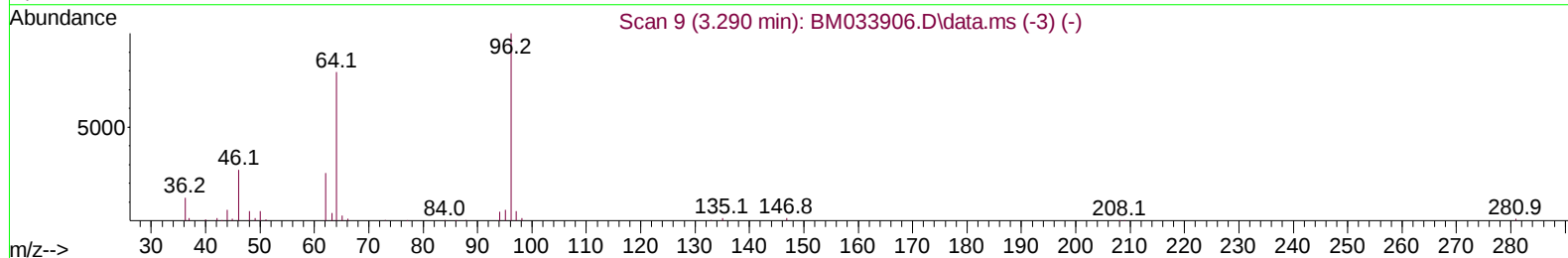
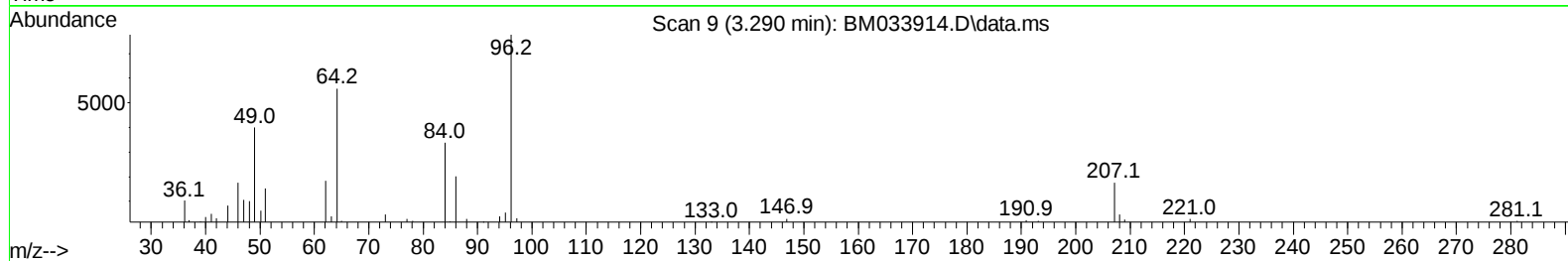
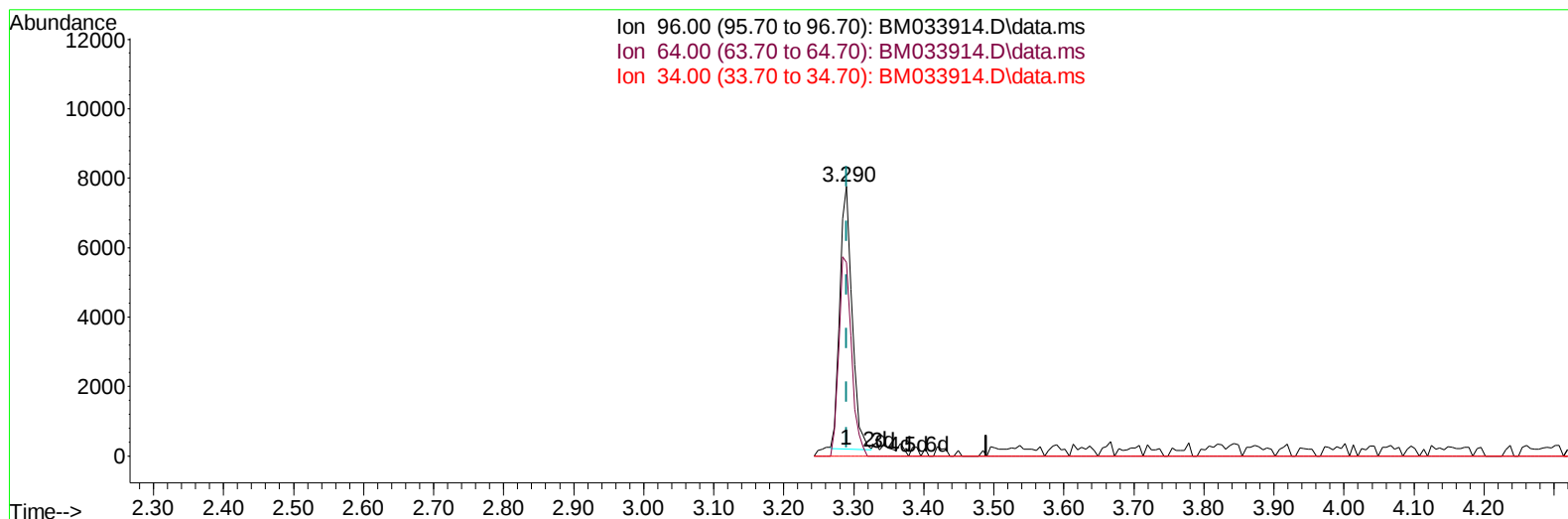
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012822\
 Data File : BM033914.D
 Acq On : 28 Jan 2022 21:17
 Operator : CG/JU
 Sample : N1302-11
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0F27

Manual IntegrationsAPPROVED

Quant Time: Jan 28 23:29:33 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: BM033914.D\data.ms

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

3.290min (+ 0.000) 4.10 ng/uL

response 9361

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.20	71.89
34.00	0.00	0.00
0.00	0.00	0.00

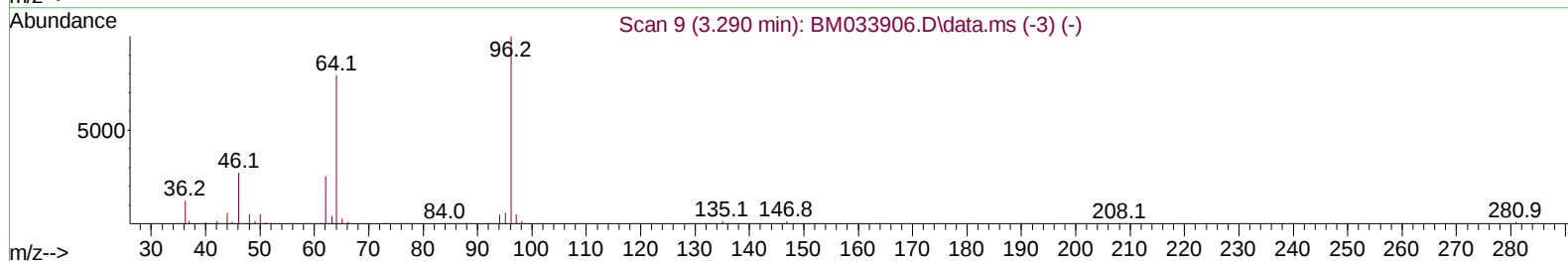
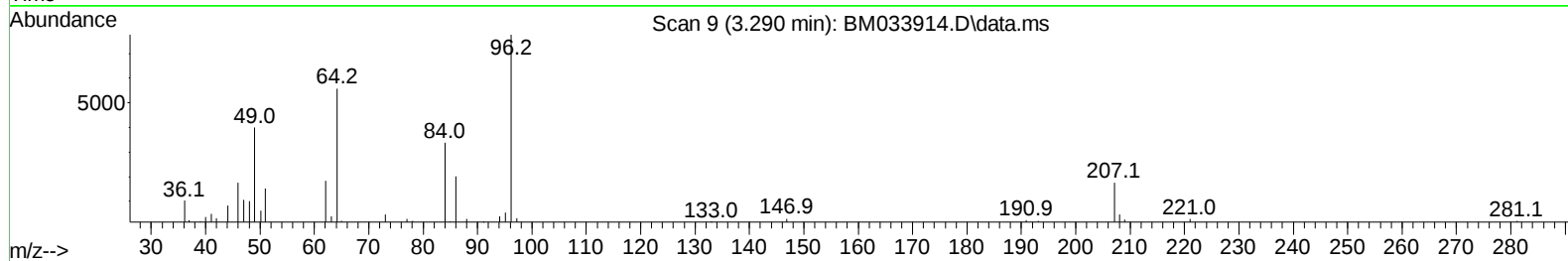
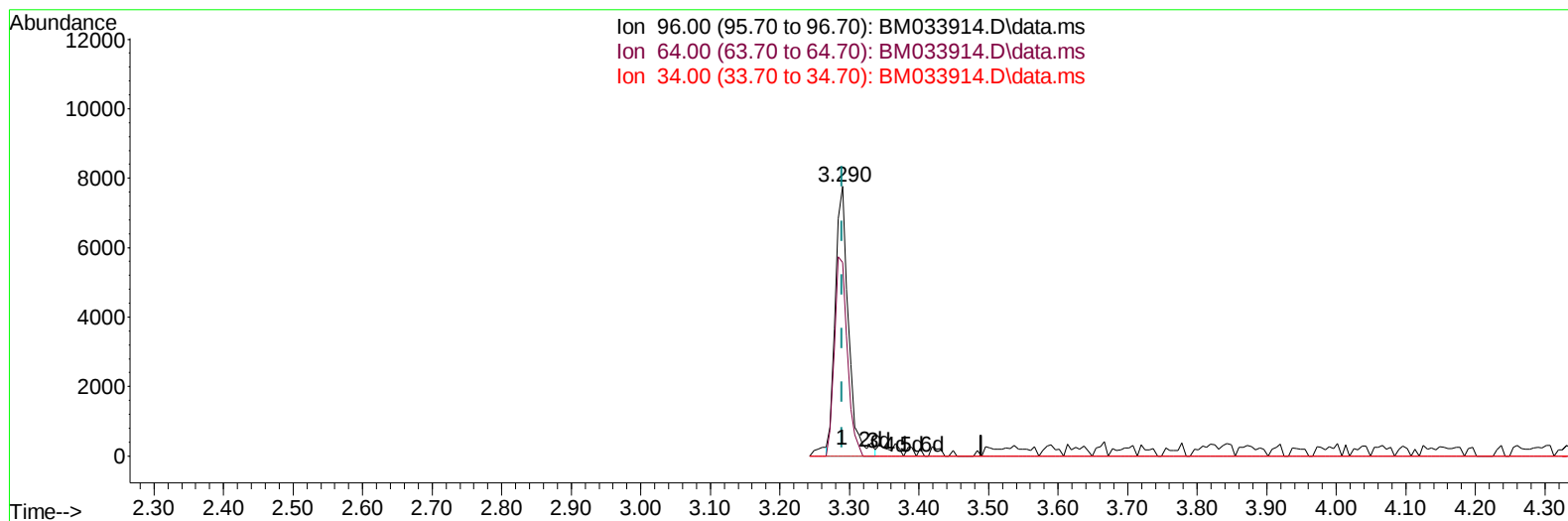
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TIC: BM033914.D\data.ms

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

3.290min (+ 0.000) 4.48 ng/uL m

response 10235

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.20	71.89
34.00	0.00	0.00
0.00	0.00	0.00

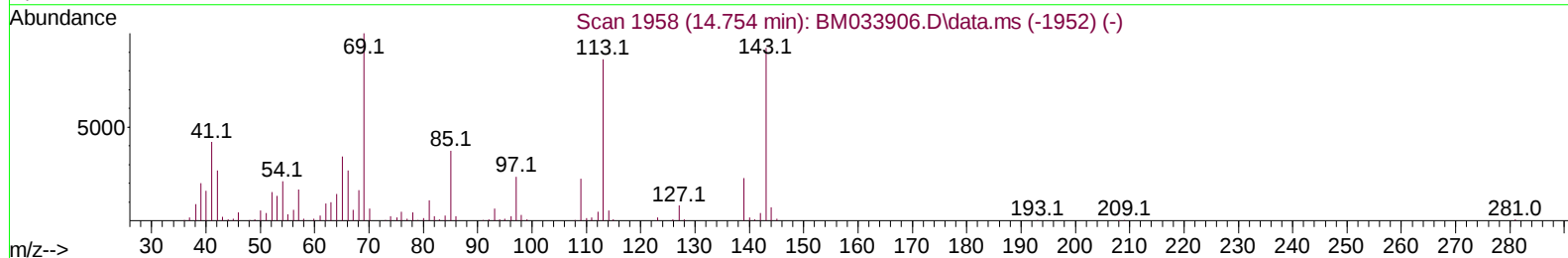
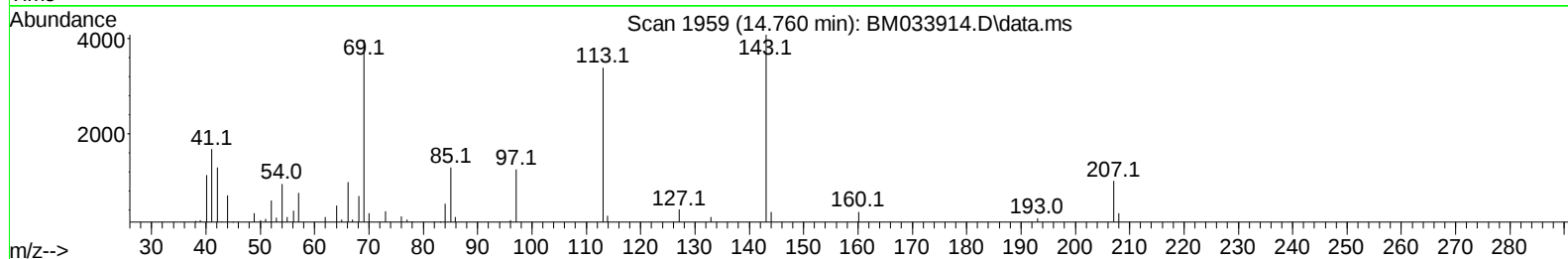
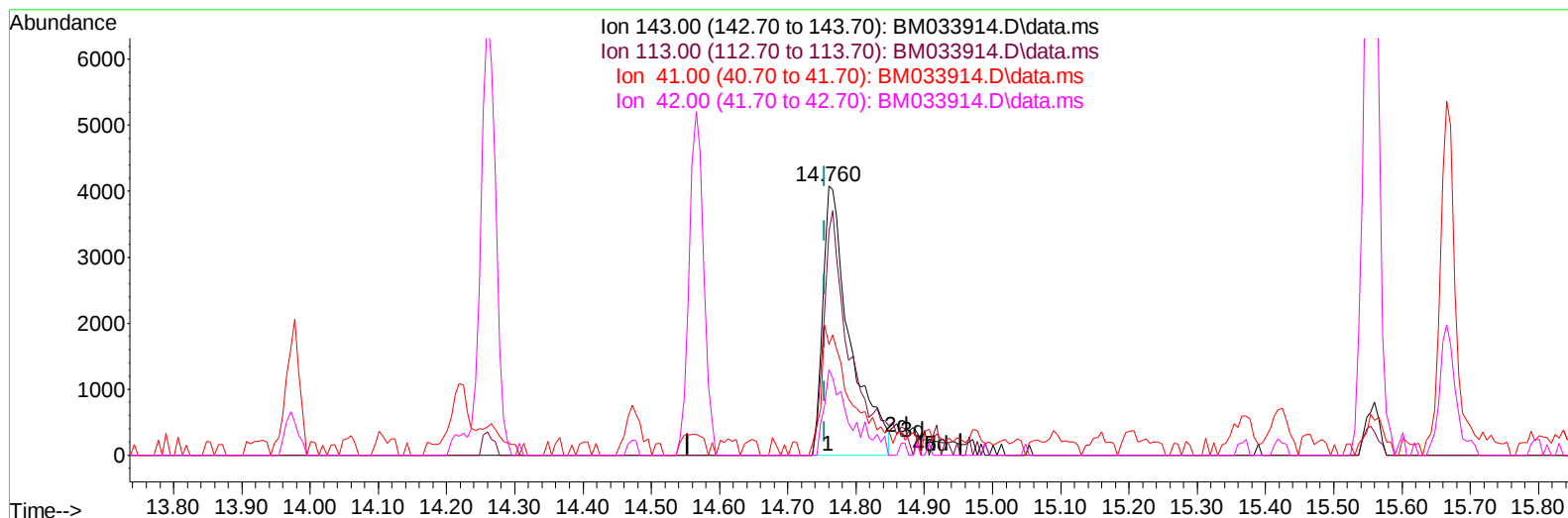
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TIC: BM033914.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.760min (+ 0.006) 3.42 ng/ul m

response 11196

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.00	83.06#
41.00	57.20	41.25#
42.00	39.50	31.73

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.931	152	85797	20.000	ng/ul	0.00
20) Naphthalene-d8	10.742	136	350350	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.566	164	216505	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.301	188	450882	20.000	ng/ul	0.00
79) Chrysene-d12	21.465	240	442895	20.000	ng/ul	0.00
88) Perylene-d12	23.859	264	452550	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.290	96	10235m	4.482	ng/uL	0.00
4) Pyridine-d5	3.719	84	59535	9.388	ng/ul	0.00
7) Phenol-d5	7.084	99	45458	5.899	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.254	67	166602	35.502	ng/ul	0.00
11) 2-Chlorophenol-d4	7.454	132	159915	27.143	ng/ul	0.00
15) 4-Methylphenol-d8	8.637	113	90946	14.257	ng/ul	0.00
21) Nitrobenzene-d5	9.095	128	103365	37.359	ng/ul	0.00
24) 2-Nitrophenol-d4	9.819	143	103625	35.339	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.360	165	169832	30.211	ng/ul	0.00
31) 4-Chloroaniline-d4	10.872	131	238751	28.571	ng/ul	0.00
46) Dimethylphthalate-d6	13.977	166	665797	39.031	ng/ul	0.00
49) Acenaphthylene-d8	14.260	160	769770	38.251	ng/ul	0.00
54) 4-Nitrophenol-d4	14.760	143	11196m	3.420	ng/ul	0.00
60) Fluorene-d10	15.554	176	563719	39.834	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.666	200	82876	30.901	ng/ul	0.00
73) Anthracene-d10	17.401	188	932683	42.630	ng/ul	0.00
81) Pyrene-d10	19.689	212	1081454	45.215	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.706	264	1051689	44.664	ng/ul	0.00

Target Compounds Qvalue

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

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