

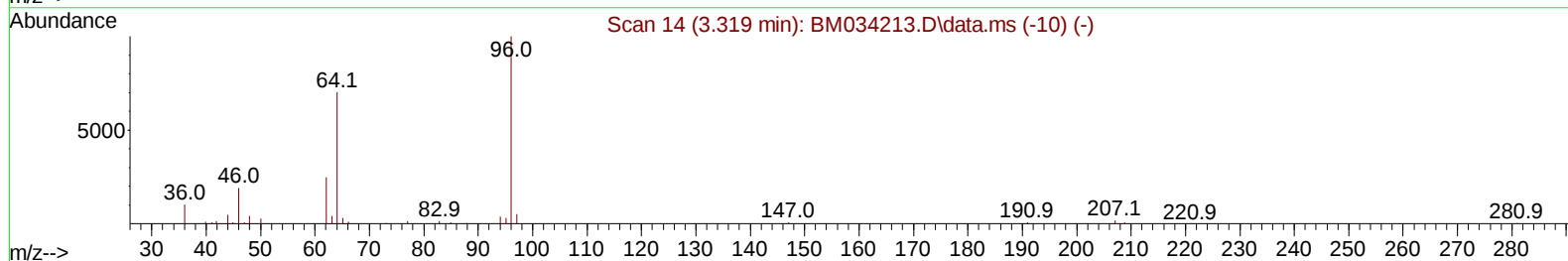
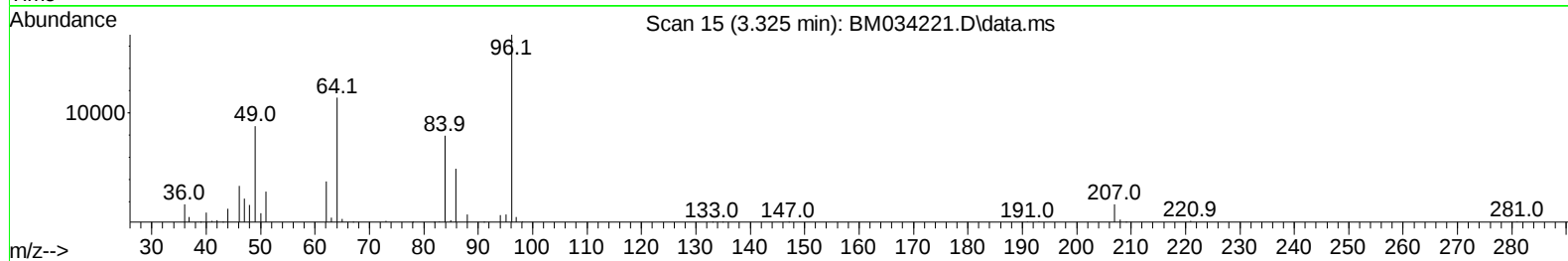
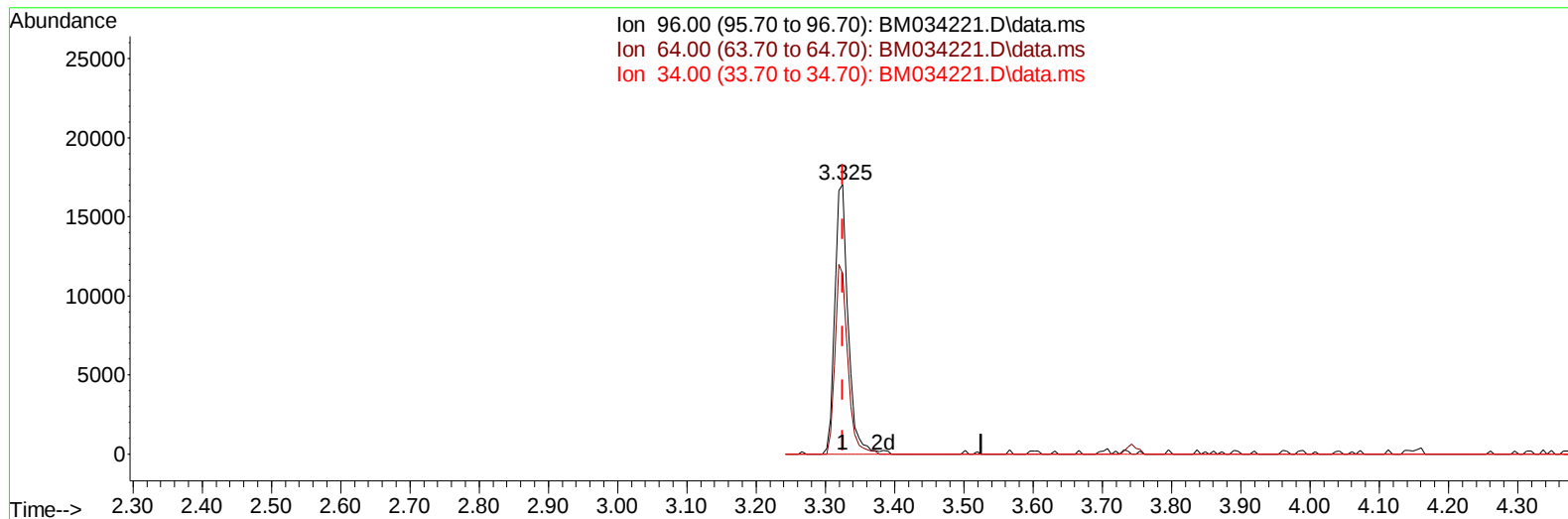
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM031122\
 Data File : BM034221.D
 Acq On : 11 Mar 2022 14:22
 Operator : CG/JU
 Sample : N1816-02
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBSC9

Manual IntegrationsAPPROVED

Quant Time: Mar 11 23:40:24 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM031022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Mar 11 02:00:05 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 03/14/2022
 Supervised By :mohammad ahmed 03/14/2022



TIC: BM034221.D\data.ms

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

3.325min (+ 0.000) 8.05 ng/uL

response 22632

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	68.90	66.74
34.00	0.00	0.00
0.00	0.00	0.00

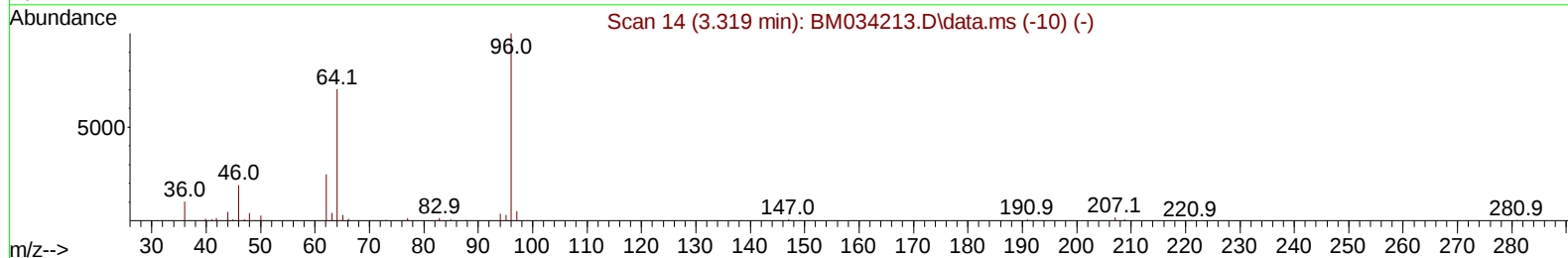
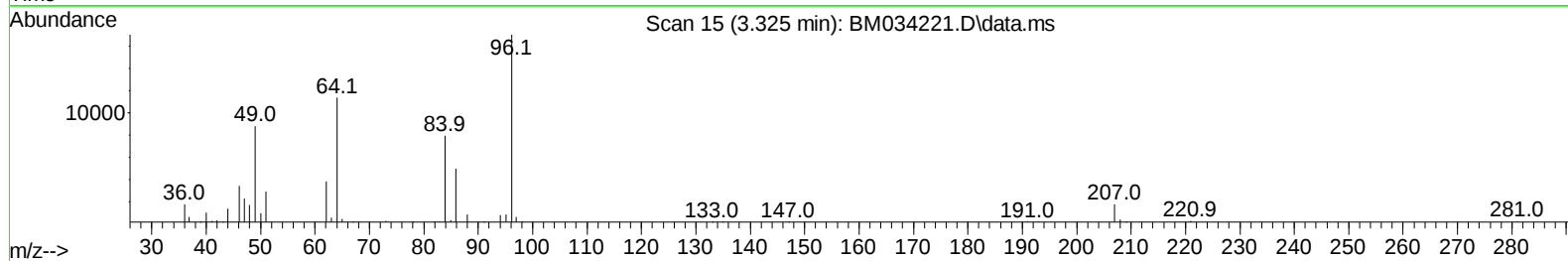
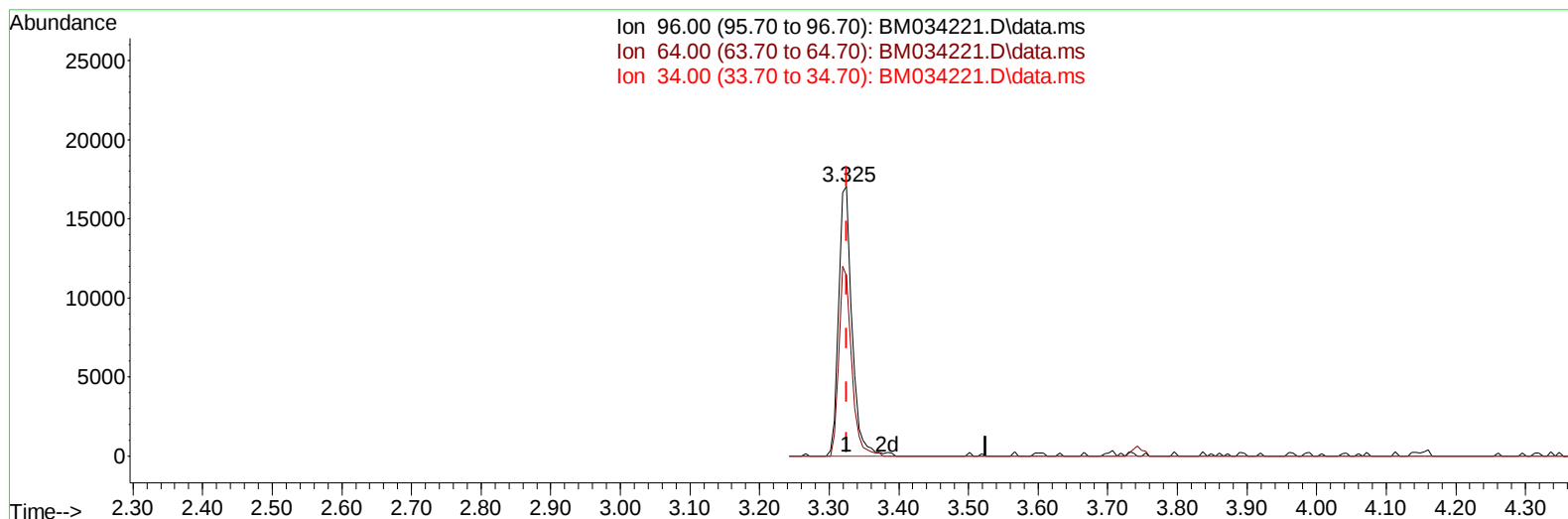
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TIC: BM034221.D\data.ms

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

3.325min (+ 0.000) 8.00 ng/uL m

response 22495

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	68.90	66.74
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	7.954	152	122731	20.000	ng/ul	0.00
20) Naphthalene-d8	10.766	136	482606	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.595	164	305412	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.330	188	597260	20.000	ng/ul	0.00
79) Chrysene-d12	21.500	240	553809	20.000	ng/ul	0.00
88) Perylene-d12	23.918	264	615207	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.325	96	22495m	8.001	ng/uL	0.00
4) Pyridine-d5	3.743	84	267443	39.952	ng/ul	0.00
7) Phenol-d5	7.101	99	320625	39.038	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.272	67	199737	41.454	ng/ul	0.00
11) 2-Chlorophenol-d4	7.478	132	298615	42.070	ng/ul	0.00
15) 4-Methylphenol-d8	8.654	113	275824	39.266	ng/ul	0.00
21) Nitrobenzene-d5	9.113	128	144187	41.652	ng/ul	0.00
24) 2-Nitrophenol-d4	9.842	143	161473	42.066	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.378	165	298959	38.849	ng/ul	0.00
31) 4-Chloroaniline-d4	10.895	131	365070	38.147	ng/ul	0.00
46) Dimethylphthalate-d6	14.001	166	921653	41.474	ng/ul	0.00
49) Acenaphthylene-d8	14.283	160	1156277	40.967	ng/ul	0.00
54) 4-Nitrophenol-d4	14.771	143	99254	28.674	ng/ul	0.00
60) Fluorene-d10	15.583	176	797931	41.019	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.689	200	121446	32.138	ng/ul	0.00
73) Anthracene-d10	17.430	188	1176584	41.708	ng/ul	0.00
81) Pyrene-d10	19.718	212	1350150	44.781	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.765	264	1279246	42.148	ng/ul	0.00

Target Compounds Qvalue

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

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