

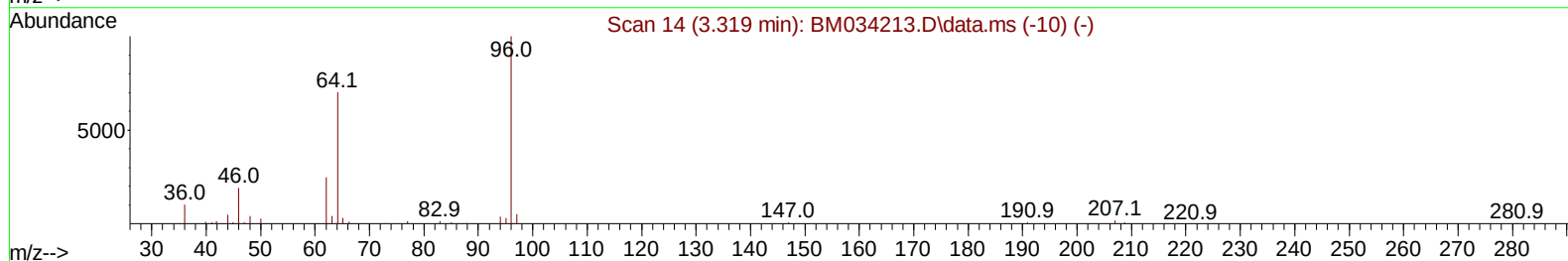
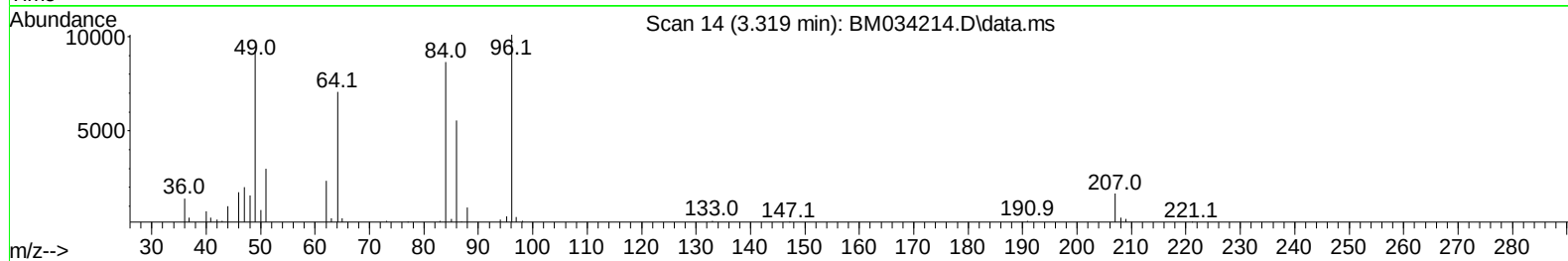
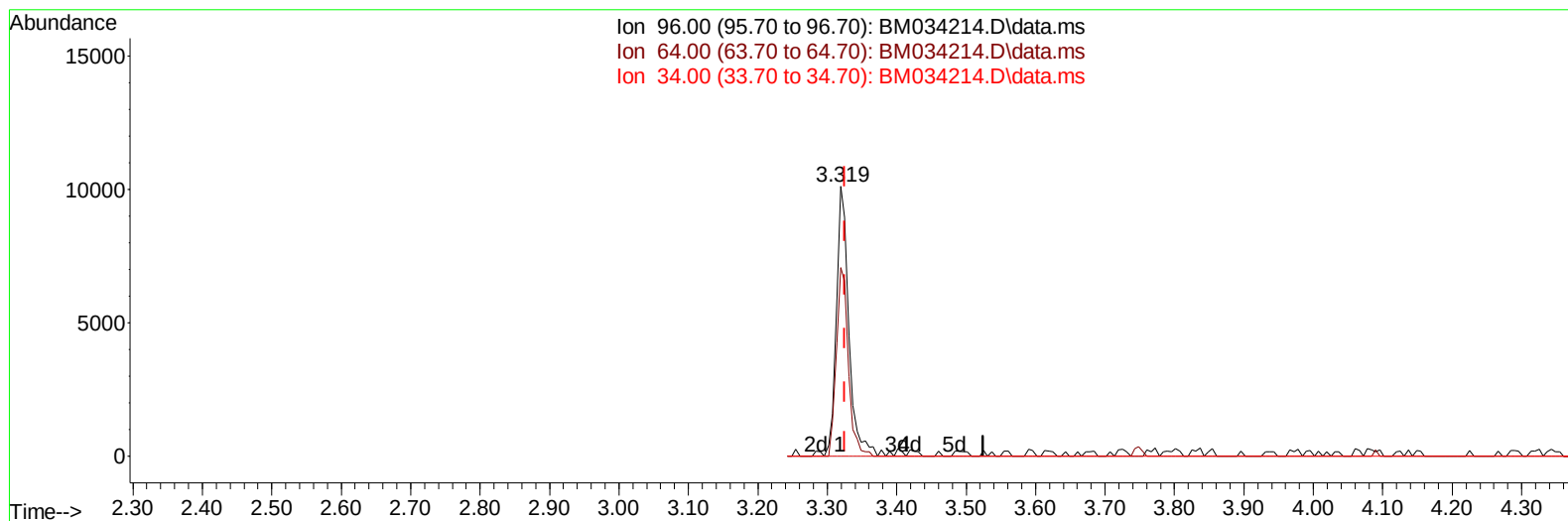
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM031122\
Data File : BM034214.D
Acq On : 11 Mar 2022 10:09
Operator : CG/JU
Sample : PB143172BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SBLK172

Manual IntegrationsAPPROVED

Quant Time: Mar 11 23:38:54 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: BM034214.D\data.ms

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

3.319min (-0.006) 4.65 ng/uL

response 12706

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	68.90	69.84
34.00	0.00	0.00
0.00	0.00	0.00

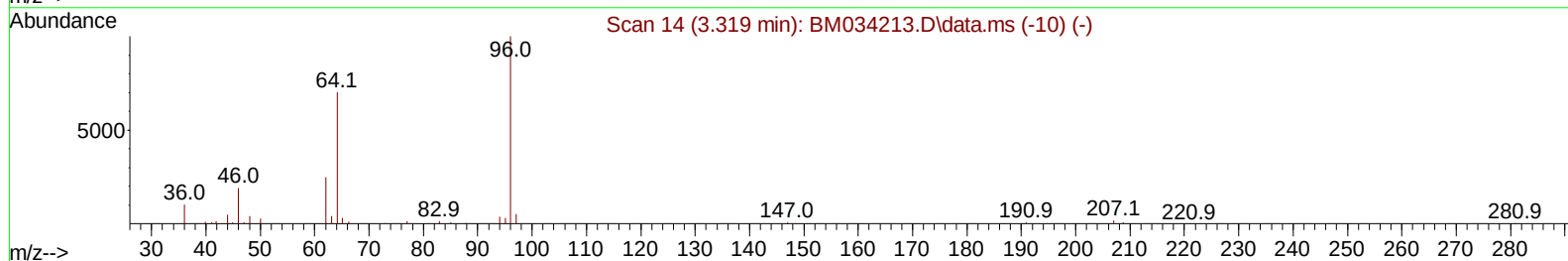
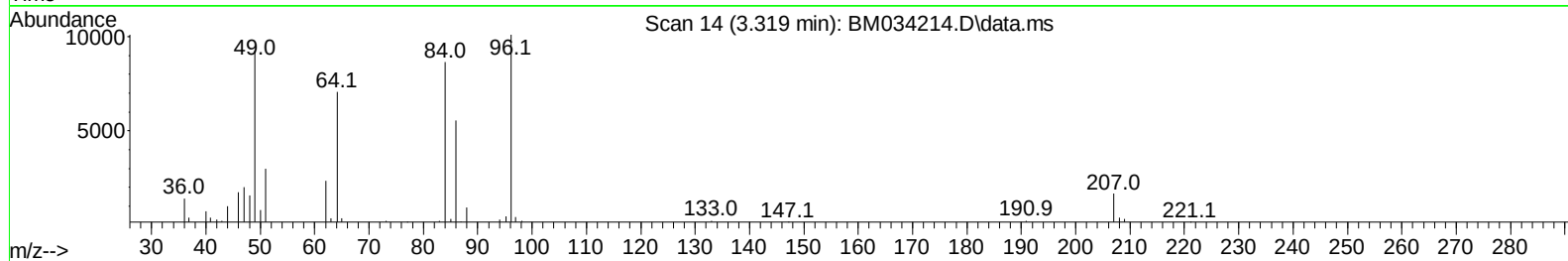
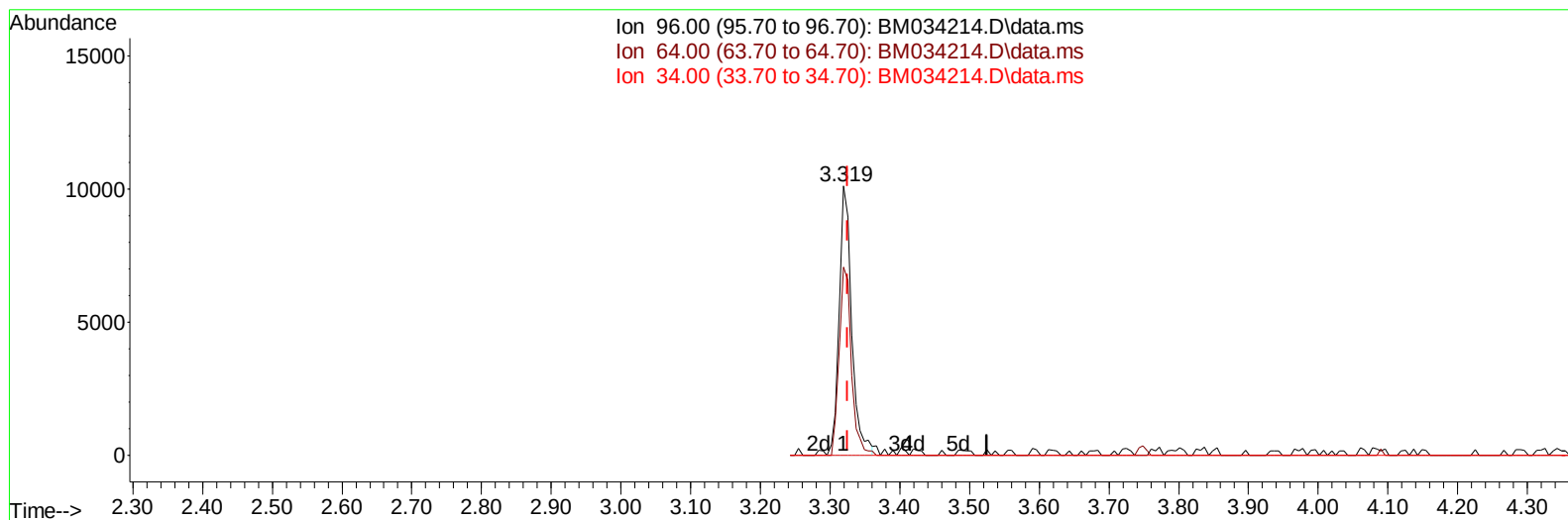
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TIC: BM034214.D\data.ms

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

3.319min (-0.006) 4.60 ng/uL m

response 12588

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	68.90	69.84
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	7.948	152	119362	20.000	ng/ul	0.00
20) Naphthalene-d8	10.766	136	473569	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.589	164	316724	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.330	188	626523	20.000	ng/ul	0.00
79) Chrysene-d12	21.500	240	544726	20.000	ng/ul	0.00
88) Perylene-d12	23.918	264	552880	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.319	96	12588m	4.603	ng/uL	0.00
4) Pyridine-d5	3.743	84	148229	22.768	ng/ul	0.00
7) Phenol-d5	7.101	99	158346	19.824	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.272	67	107378	22.914	ng/ul	0.00
11) 2-Chlorophenol-d4	7.478	132	155441	22.517	ng/ul	0.00
15) 4-Methylphenol-d8	8.654	113	136057	19.916	ng/ul	0.00
21) Nitrobenzene-d5	9.107	128	74262	21.862	ng/ul	0.00
24) 2-Nitrophenol-d4	9.842	143	79561	21.122	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.378	165	153380	20.312	ng/ul	0.00
31) 4-Chloroaniline-d4	10.895	131	186130	19.820	ng/ul	0.00
46) Dimethylphthalate-d6	13.995	166	545042	23.651	ng/ul	0.00
49) Acenaphthylene-d8	14.283	160	647808	22.132	ng/ul	0.00
54) 4-Nitrophenol-d4	14.772	143	46714	13.014	ng/ul	0.00
60) Fluorene-d10	15.577	176	465655	23.083	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.689	200	67890	17.127	ng/ul	0.00
73) Anthracene-d10	17.430	188	691132	23.355	ng/ul	0.00
81) Pyrene-d10	19.712	212	816370	27.529	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.759	264	685529	25.133	ng/ul	0.00

Target Compounds Qvalue

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

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