

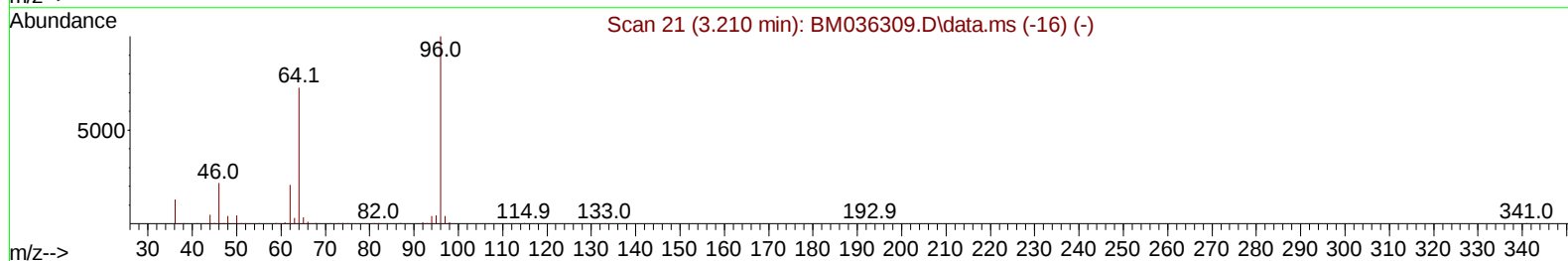
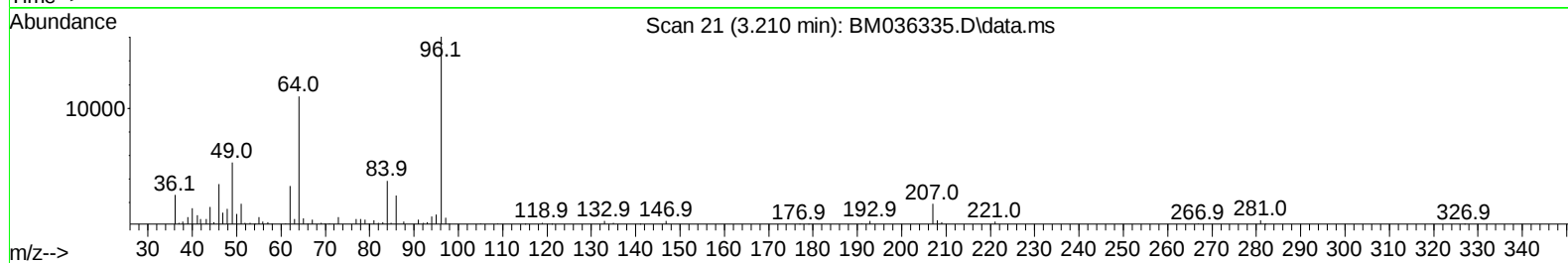
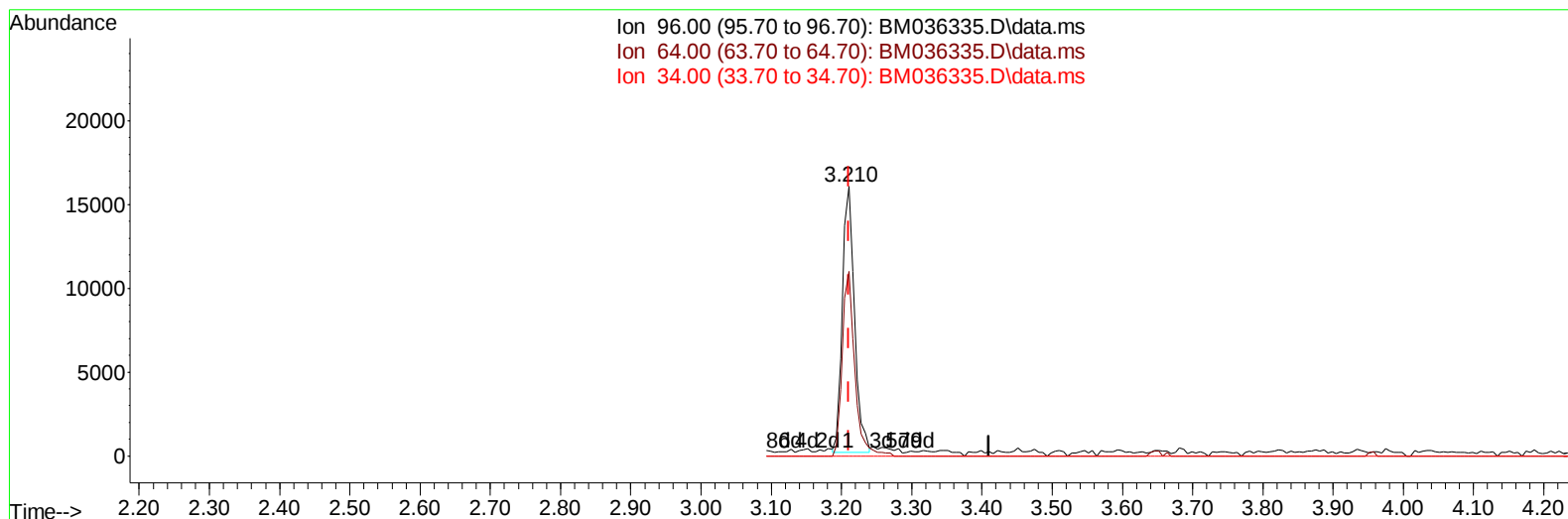
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081722\
 Data File : BM036335.D
 Acq On : 18 Aug 2022 03:01
 Operator : CG/JU
 Sample : N4176-04
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EX7P9

Manual IntegrationsAPPROVED

Quant Time: Aug 18 04:02:06 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM081622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Aug 16 15:00:33 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 08/18/2022
 Supervised By :Sohil Jodhani 08/19/2022



TIC: BM036335.D\data.ms

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

3.210min (+ 0.000) 2.67 ng/uL

response 19084

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	75.40	68.51
34.00	0.00	0.00
0.00	0.00	0.00

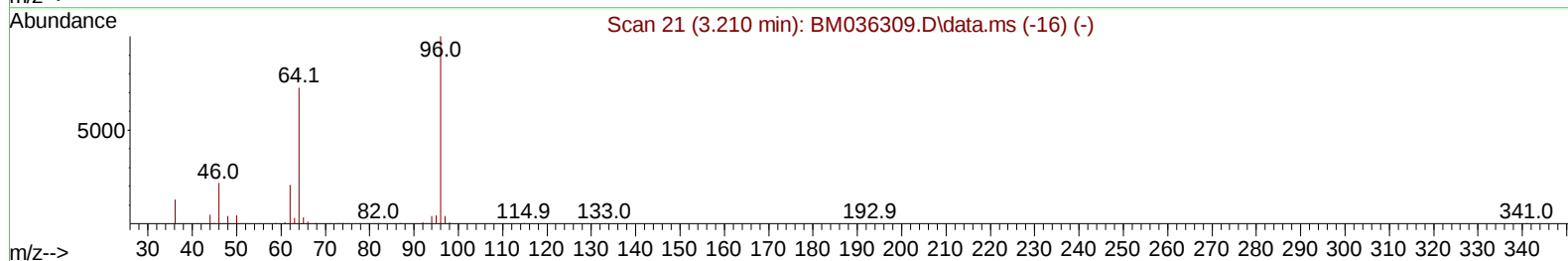
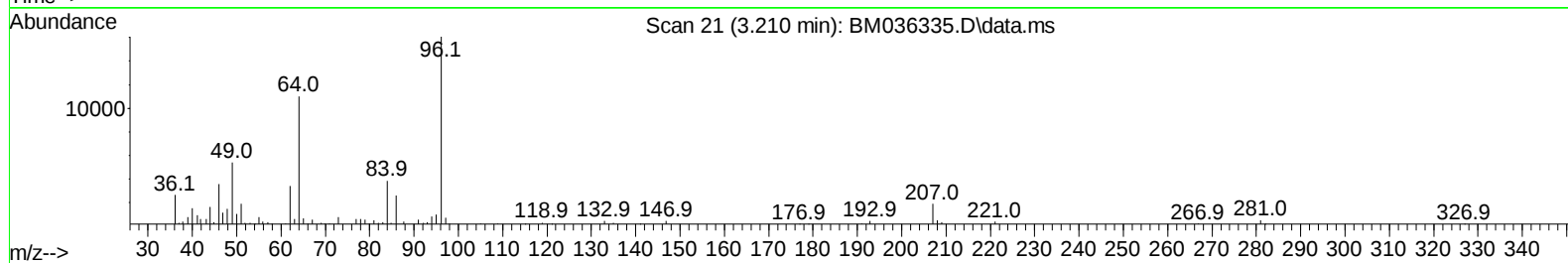
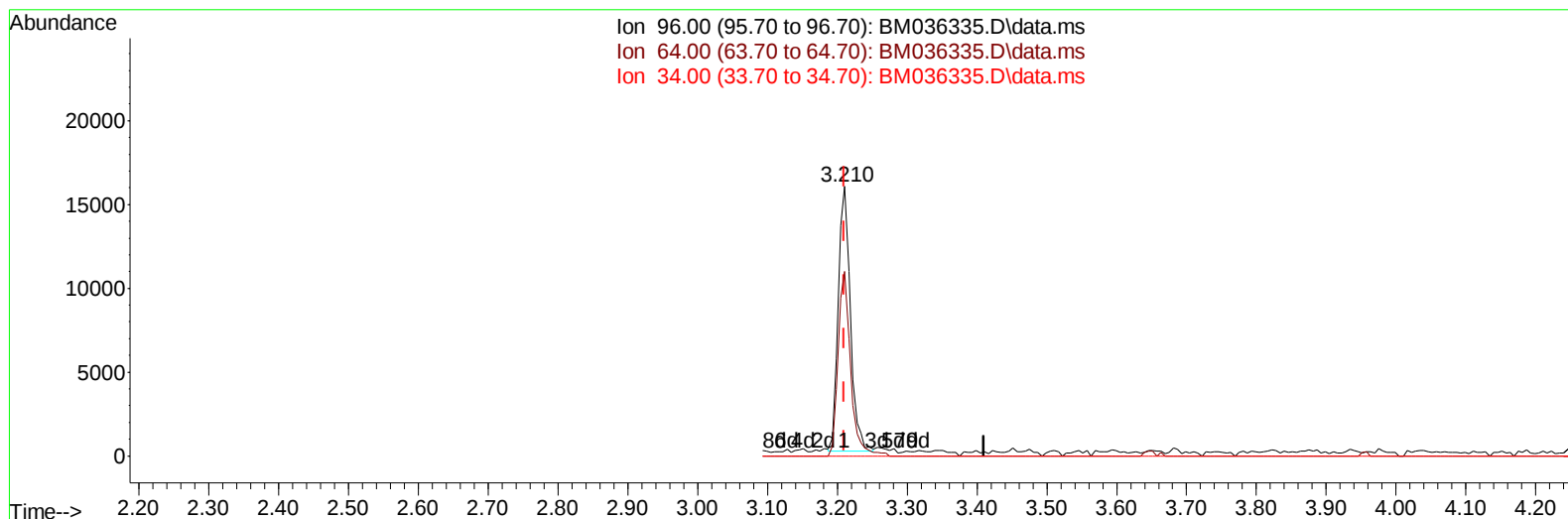
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TIC: BM036335.D\data.ms

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

3.210min (+ 0.000) 2.66 ng/uL m

response 19013

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	75.40	68.51
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	7.798	152	272106	20.000	ng/ul	0.00
20) Naphthalene-d8	10.598	136	1179723	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.439	164	646134	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.186	188	1195936	20.000	ng/ul	0.00
79) Chrysene-d12	21.368	240	1059382	20.000	ng/ul	0.00
88) Perylene-d12	23.697	264	1141852	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.210	96	19013m	2.663	ng/uL	0.00
4) Pyridine-d5	3.640	84	115270	5.854	ng/ul	0.00
7) Phenol-d5	6.981	99	356569	15.352	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.134	67	234754	16.057	ng/ul	0.00
11) 2-Chlorophenol-d4	7.334	132	304790	16.702	ng/ul	0.00
15) 4-Methylphenol-d8	8.522	113	283381	16.340	ng/ul	0.00
21) Nitrobenzene-d5	8.969	128	151862	16.978	ng/ul	0.00
24) 2-Nitrophenol-d4	9.687	143	155748	17.167	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.228	165	262502	16.193	ng/ul	0.00
31) 4-Chloroaniline-d4	10.751	131	370590	14.746	ng/ul	0.00
46) Dimethylphthalate-d6	13.863	166	749543	18.336	ng/ul	0.00
49) Acenaphthylene-d8	14.133	160	909545	17.840	ng/ul	0.00
54) 4-Nitrophenol-d4	14.680	143	94153	12.156	ng/ul	0.00
60) Fluorene-d10	15.433	176	661129	17.898	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.568	200	57083	9.284	ng/ul	0.00
73) Anthracene-d10	17.286	188	1001385	18.831	ng/ul	0.00
81) Pyrene-d10	19.580	212	1149496	20.917	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.550	264	1139091	20.347	ng/ul	0.00

Target Compounds Qvalue

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

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