

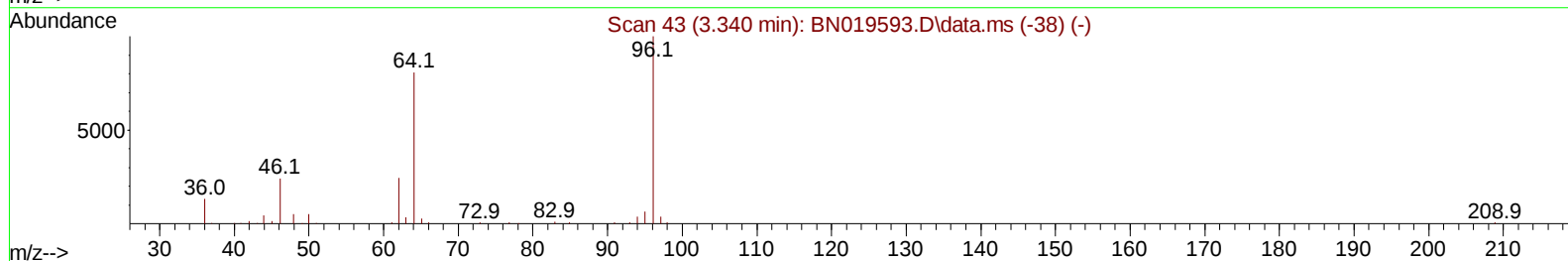
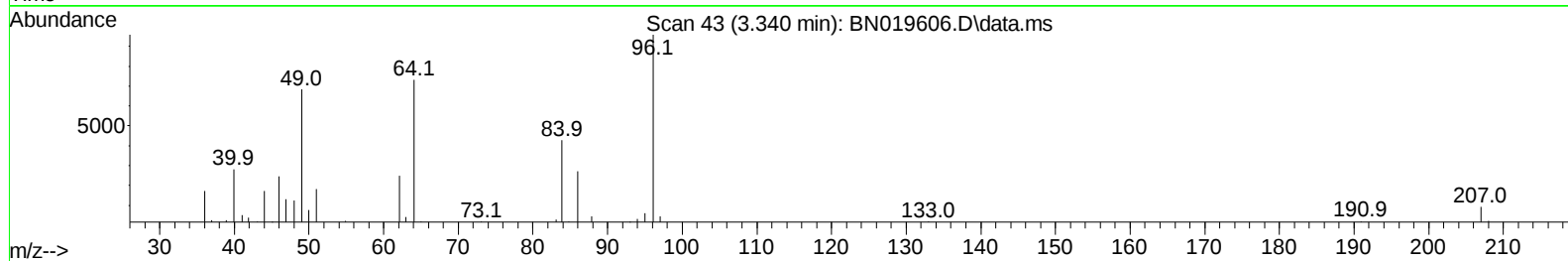
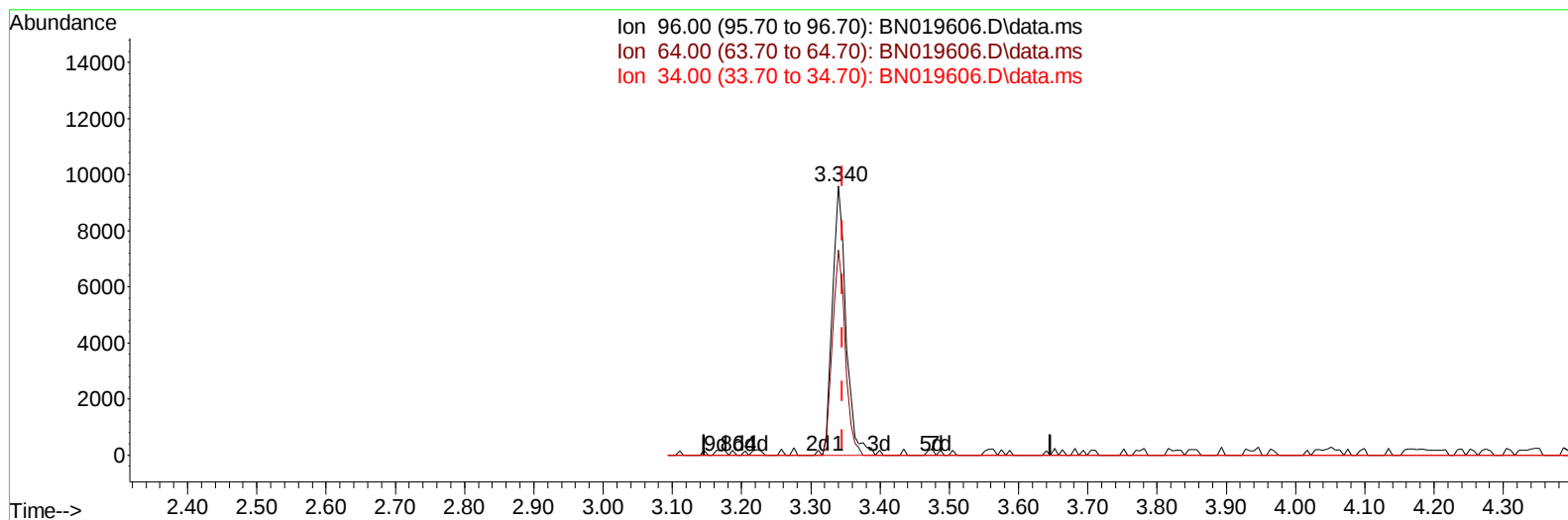
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN043022\
 Data File : BN019606.D
 Acq On : 30 Apr 2022 00:44
 Operator : CG/JU
 Sample : N2612-07
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BH0A5

Manual IntegrationsAPPROVED

Quant Time: Apr 30 03:09:04 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN042722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Apr 27 13:41:44 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/02/2022
 Supervised By :Yogesh Patel 05/05/2022



TIC: BN019606.D\data.ms

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

3.340min (-0.006) 3.12 ng/uL

response 12559

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	80.20	76.22
34.00	0.00	0.00
0.00	0.00	0.00

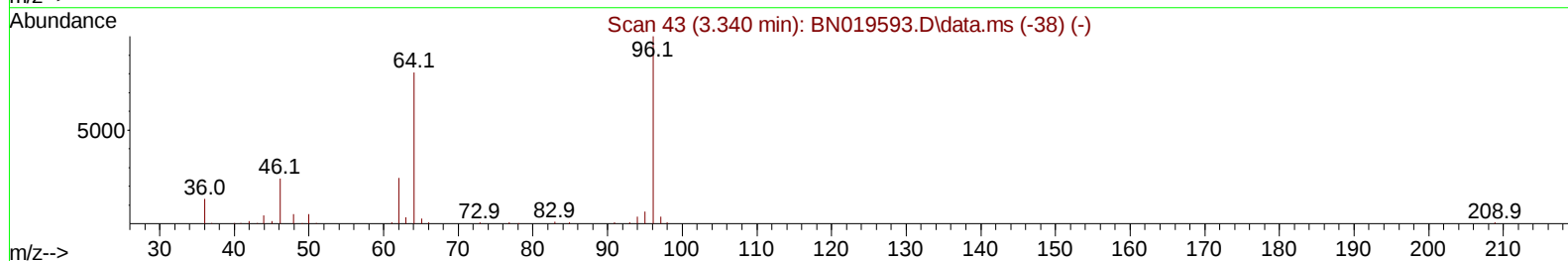
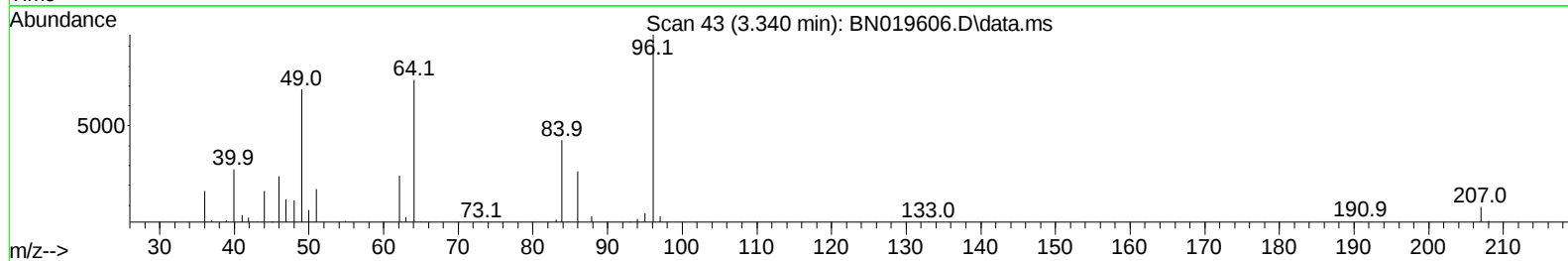
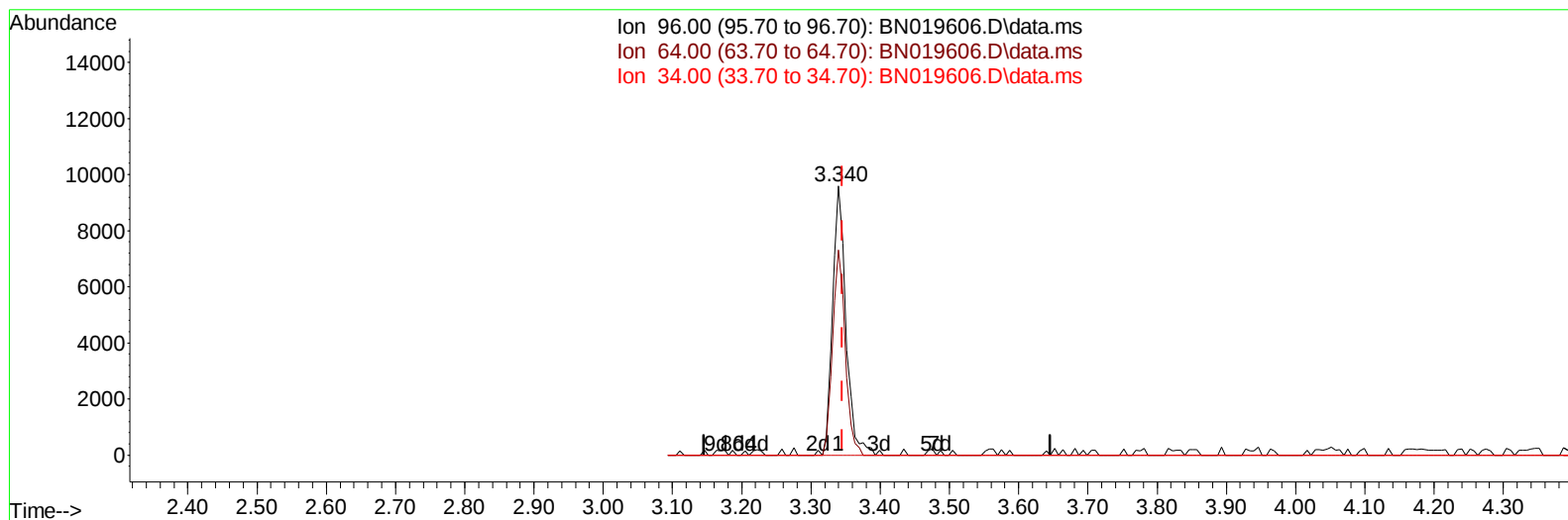
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TIC: BN019606.D\data.ms

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

3.340min (-0.006) 3.21 ng/uL m

response 12906

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	80.20	76.22
34.00	0.00	0.00
0.00	0.00	0.00

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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.869	152		151905	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.663	136		630977	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.492	164		365527	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.228	188		729370	20.000	ng/ul	-0.01
79) Chrysene-d12	21.410	240		689833	20.000	ng/ul	-0.01
88) Perylene-d12	23.762	264		680526	20.000	ng/ul	-0.02
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.340	96		12906m	3.206	ng/uL	0.00
4) Pyridine-d5	3.746	84		82053	8.050	ng/ul	0.00
7) Phenol-d5	7.034	99		12115	0.920	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.199	67		106017	13.885	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.405	132		27796	2.662	ng/ul	0.00
15) 4-Methylphenol-d8	8.569	113		18952	1.879	ng/ul	-0.01
21) Nitrobenzene-d5	9.022	128		66792	14.518	ng/ul	0.00
24) 2-Nitrophenol-d4	9.746	143		13226	2.917	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.281	165		24665	2.718	ng/ul	0.00
31) 4-Chloroaniline-d4	10.793	131		247945	17.997	ng/ul	0.00
46) Dimethylphthalate-d6	13.904	166		381556	15.198	ng/ul	0.00
49) Acenaphthylene-d8	14.187	160		442597	13.038	ng/ul	0.00
54) 4-Nitrophenol-d4	14.675	143		2921	0.598	ng/ul	0.00
60) Fluorene-d10	15.481	176		330335	14.905	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.592	200		4675	1.256	ng/ul	0.00
73) Anthracene-d10	17.328	188		556743	16.209	ng/ul	0.00
81) Pyrene-d10	19.616	212		655473	16.527	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.610	264		598977	17.218	ng/ul	-0.01

Target Compounds Qvalue

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

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