

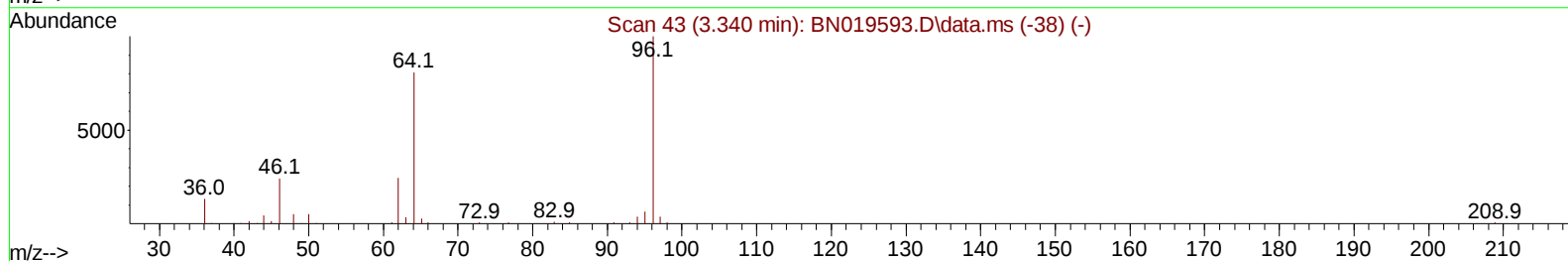
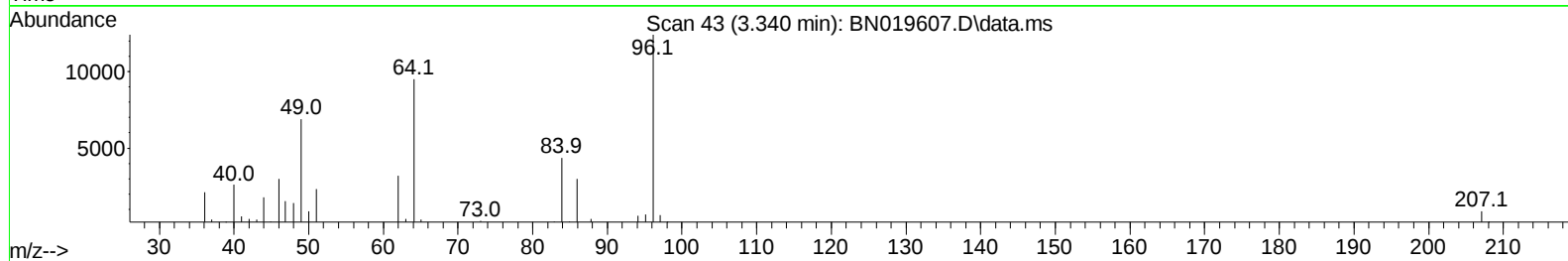
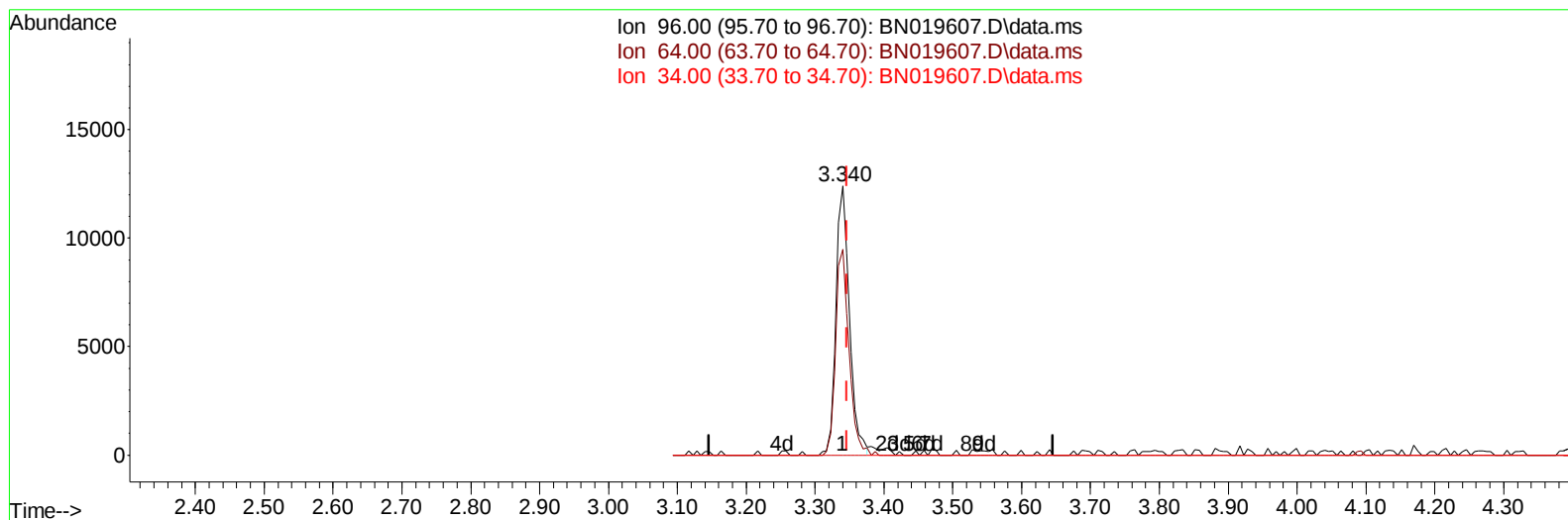
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN043022\
Data File : BN019607.D
Acq On : 30 Apr 2022 01:19
Operator : CG/JU
Sample : N2612-11
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BH0B5

Manual IntegrationsAPPROVED

Quant Time: Apr 30 03:09:30 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: BN019607.D\data.ms

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

3.340min (-0.006) 4.17 ng/uL

response 16758

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	80.20	76.51
34.00	0.00	0.00
0.00	0.00	0.00

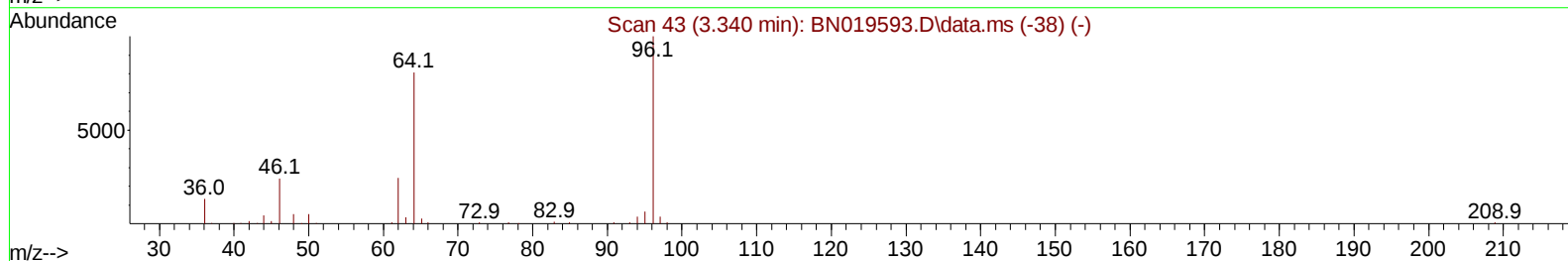
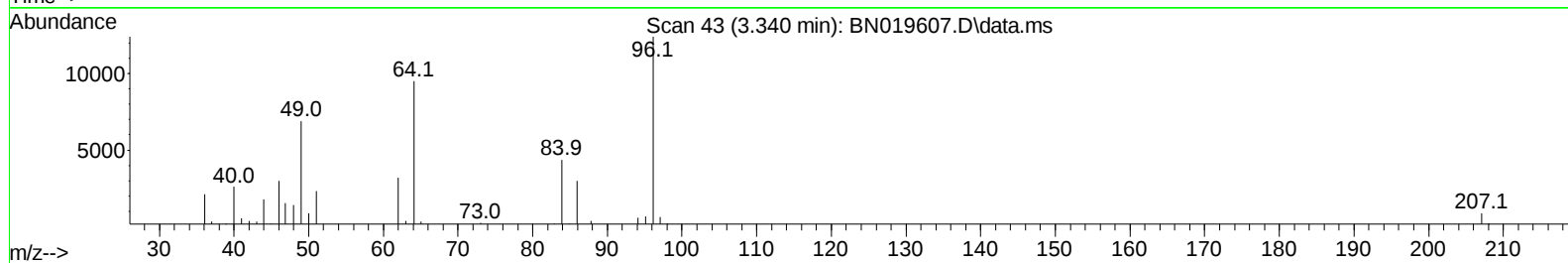
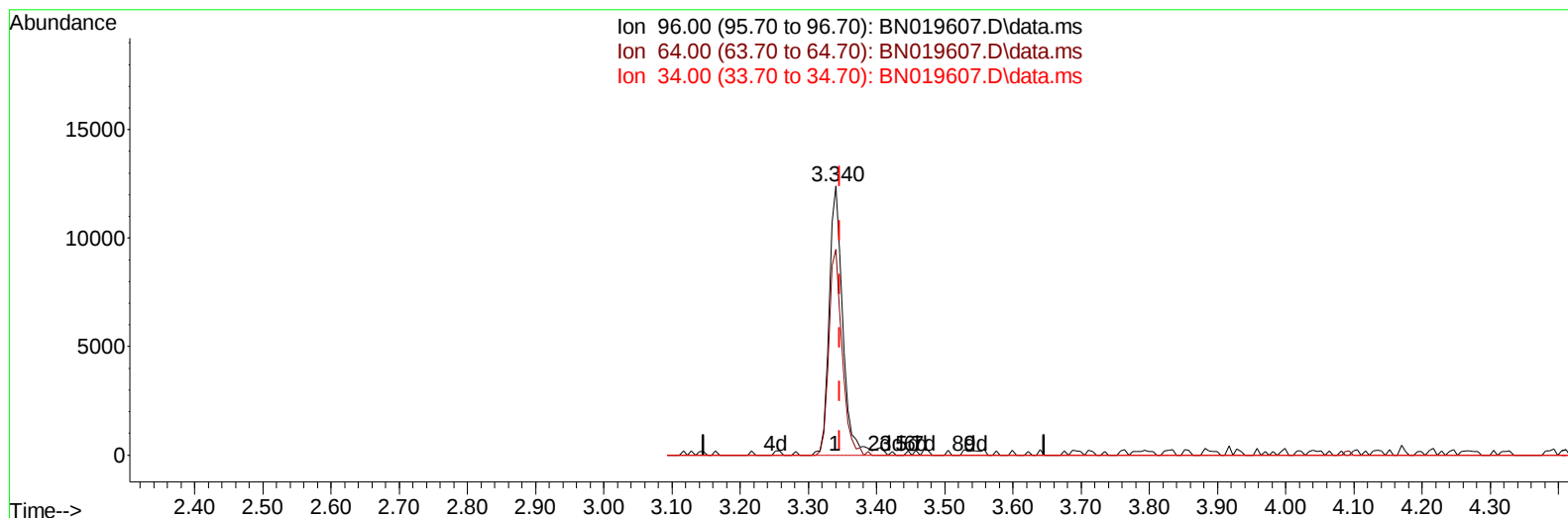
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TIC: BN019607.D\data.ms

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

3.340min (-0.006) 4.32 ng/uL m

response 17357

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	80.20	76.51
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	7.875	152	151544	20.000	ng/ul	0.00
20) Naphthalene-d8	10.663	136	626900	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.492	164	359608	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.228	188	710120	20.000	ng/ul	-0.01
79) Chrysene-d12	21.410	240	672981	20.000	ng/ul	-0.01
88) Perylene-d12	23.763	264	696080	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	17357m	4.322	ng/uL	0.00
4) Pyridine-d5	3.752	84	36632	3.602	ng/ul	0.00
7) Phenol-d5	7.034	99	54854	4.177	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.199	67	175893	23.091	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.405	132	177929	17.081	ng/ul	0.00
15) 4-Methylphenol-d8	8.569	113	105216	10.458	ng/ul	-0.01
21) Nitrobenzene-d5	9.022	128	111325	24.356	ng/ul	0.00
24) 2-Nitrophenol-d4	9.746	143	104312	23.157	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.281	165	172060	19.082	ng/ul	0.00
31) 4-Chloroaniline-d4	10.793	131	236567	17.283	ng/ul	0.00
46) Dimethylphthalate-d6	13.898	166	620107	25.106	ng/ul	-0.01
49) Acenaphthylene-d8	14.187	160	738758	22.121	ng/ul	0.00
54) 4-Nitrophenol-d4	14.669	143	17783	3.702	ng/ul	-0.01
60) Fluorene-d10	15.481	176	539580	24.747	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.592	200	61960	17.091	ng/ul	0.00
73) Anthracene-d10	17.328	188	918159	27.455	ng/ul	0.00
81) Pyrene-d10	19.622	212	1038168	26.832	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.610	264	983293	27.634	ng/ul	-0.01

Target Compounds Qvalue

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

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