

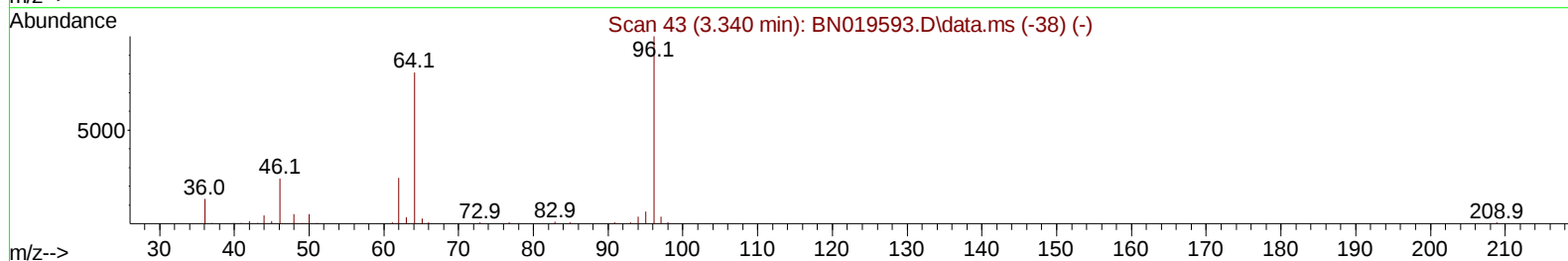
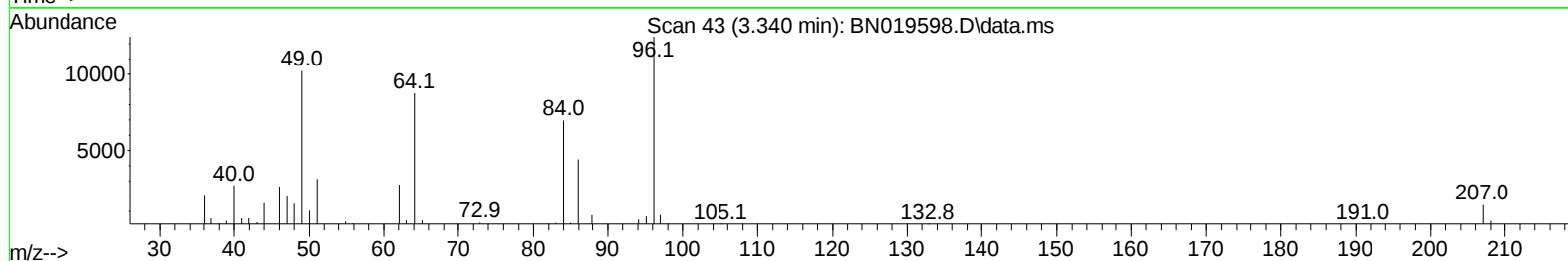
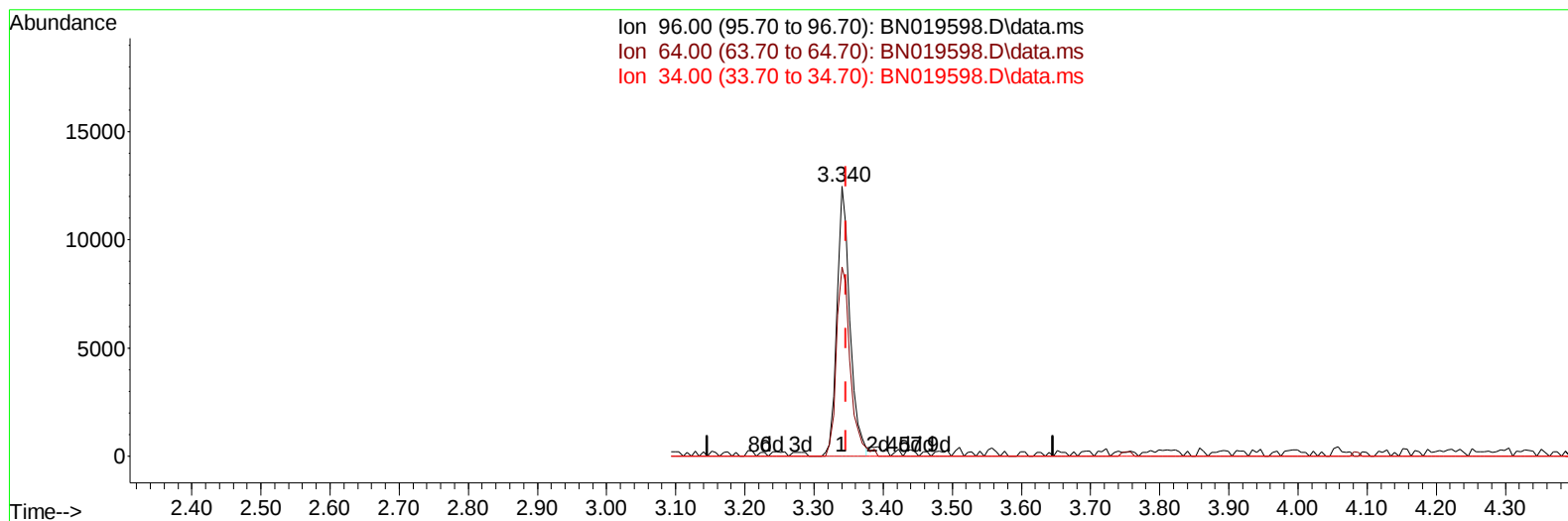
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN043022\
 Data File : BN019598.D
 Acq On : 29 Apr 2022 19:58
 Operator : CG/JU
 Sample : N2612-12
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BH0B6

Manual IntegrationsAPPROVED

Quant Time: Apr 30 03:06:32 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: BN019598.D\data.ms

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

3.340min (-0.006) 3.75 ng/uL

response 16277

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	80.20	70.15
34.00	0.00	0.00
0.00	0.00	0.00

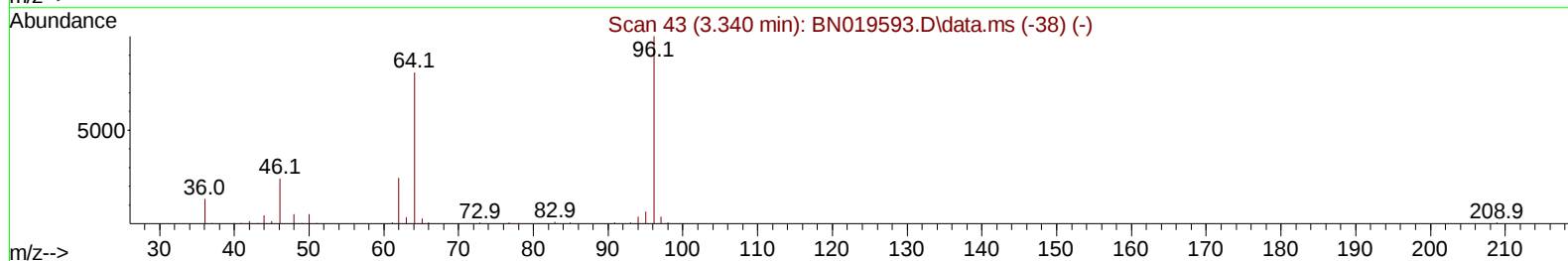
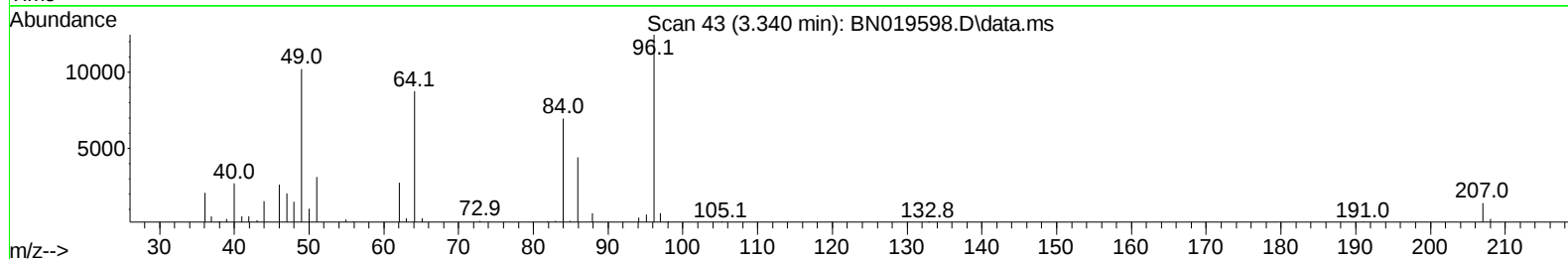
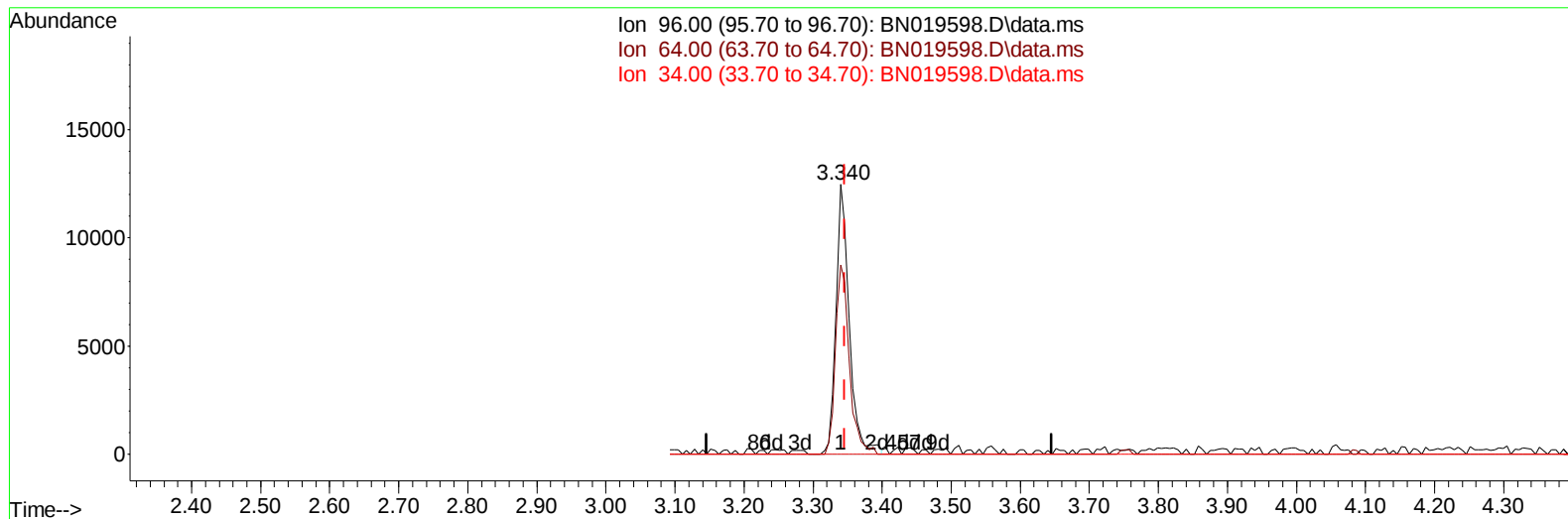
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TIC: BN019598.D\data.ms

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

3.340min (-0.006) 3.87 ng/uL m

response 16793

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	80.20	70.15
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	163622	20.000	ng/ul	0.00
20) Naphthalene-d8	10.663	136	690875	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.492	164	410646	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.233	188	826231	20.000	ng/ul	0.00
79) Chrysene-d12	21.416	240	766342	20.000	ng/ul	0.00
88) Perylene-d12	23.768	264	774044	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	16793m	3.873	ng/uL	0.00
4) Pyridine-d5	3.752	84	86609	7.888	ng/ul	0.00
7) Phenol-d5	7.034	99	76821	5.418	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.205	67	207184	25.191	ng/ul	0.00
11) 2-Chlorophenol-d4	7.405	132	222805	19.810	ng/ul	0.00
15) 4-Methylphenol-d8	8.569	113	144169	13.273	ng/ul	-0.01
21) Nitrobenzene-d5	9.022	128	136738	27.145	ng/ul	0.00
24) 2-Nitrophenol-d4	9.746	143	131932	26.577	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.281	165	228040	22.948	ng/ul	0.00
31) 4-Chloroaniline-d4	10.793	131	359861	23.856	ng/ul	0.00
46) Dimethylphthalate-d6	13.904	166	847427	30.045	ng/ul	0.00
49) Acenaphthylene-d8	14.187	160	956029	25.069	ng/ul	0.00
54) 4-Nitrophenol-d4	14.669	143	31683	5.776	ng/ul	-0.01
60) Fluorene-d10	15.481	176	725095	29.122	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.592	200	101276	24.010	ng/ul	0.00
73) Anthracene-d10	17.328	188	1236831	31.787	ng/ul	0.00
81) Pyrene-d10	19.622	212	1412608	32.061	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.615	264	1342178	33.920	ng/ul	0.00

Target Compounds Qvalue

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

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