

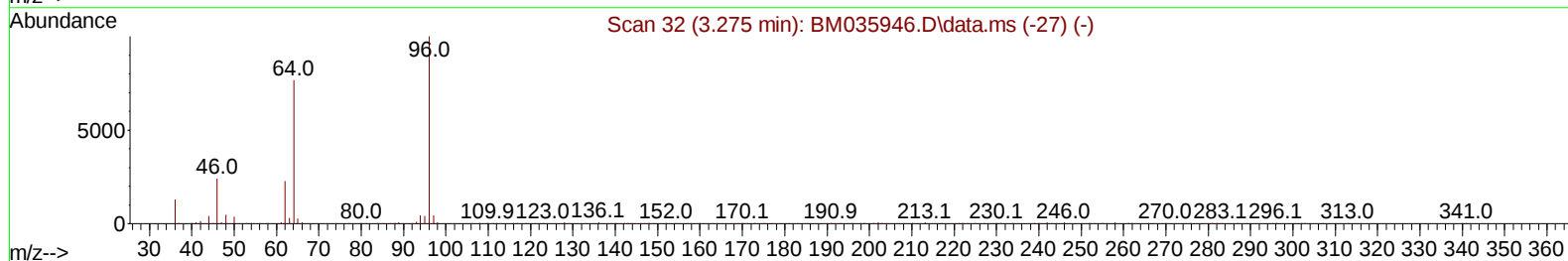
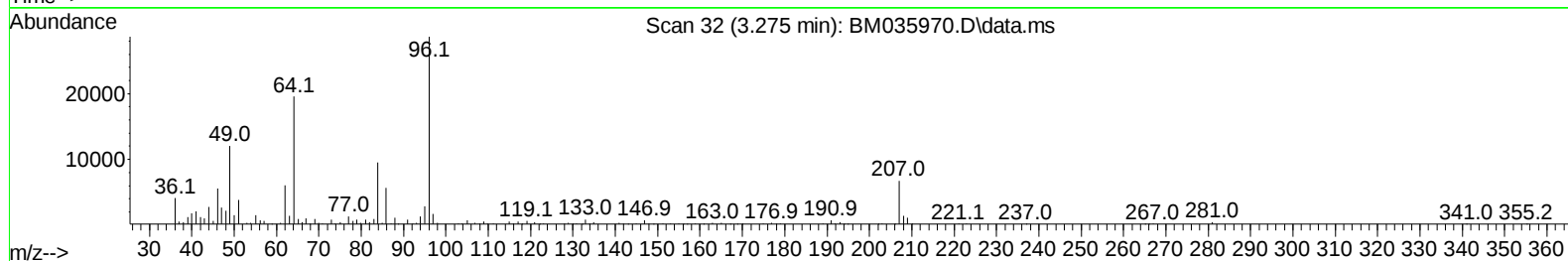
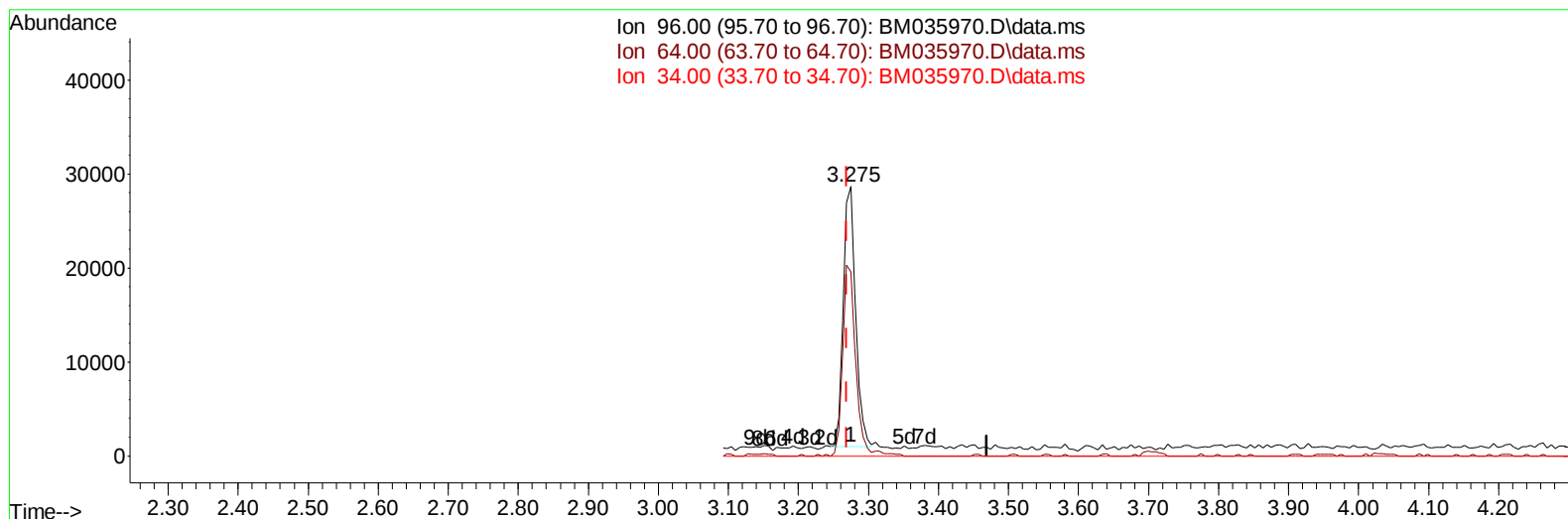
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072722\
 Data File : BM035970.D
 Acq On : 28 Jul 2022 03:31
 Operator : CG/JU
 Sample : N3741-07
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 H0B15

Manual IntegrationsAPPROVED

Quant Time: Jul 28 04:07:56 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM072522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jul 27 23:55:50 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 07/28/2022
 Supervised By :mohammad ahmed 07/29/2022



TIC: BM035970.D\data.ms

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

3.275min (+ 0.006) 4.70 ng/uL

response 34013

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	75.40	68.21
34.00	0.00	0.00
0.00	0.00	0.00

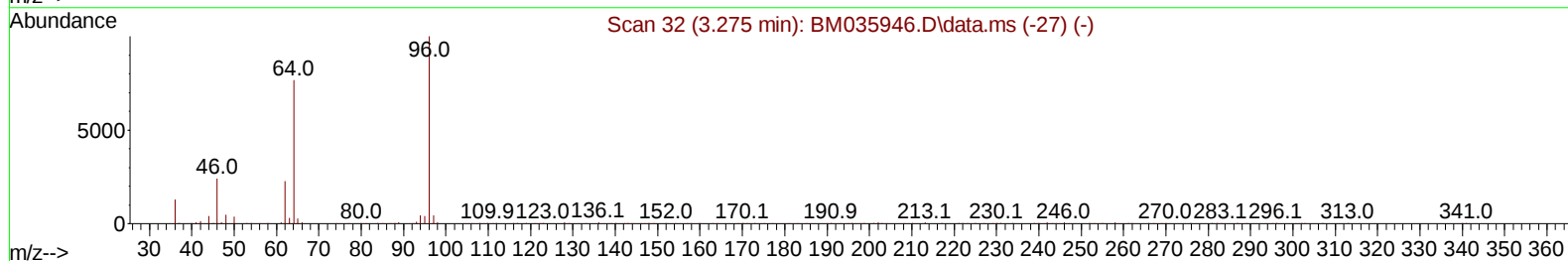
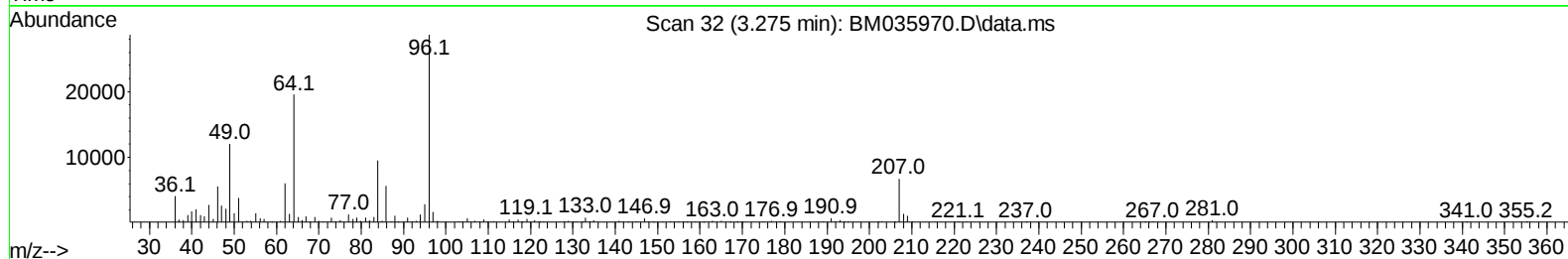
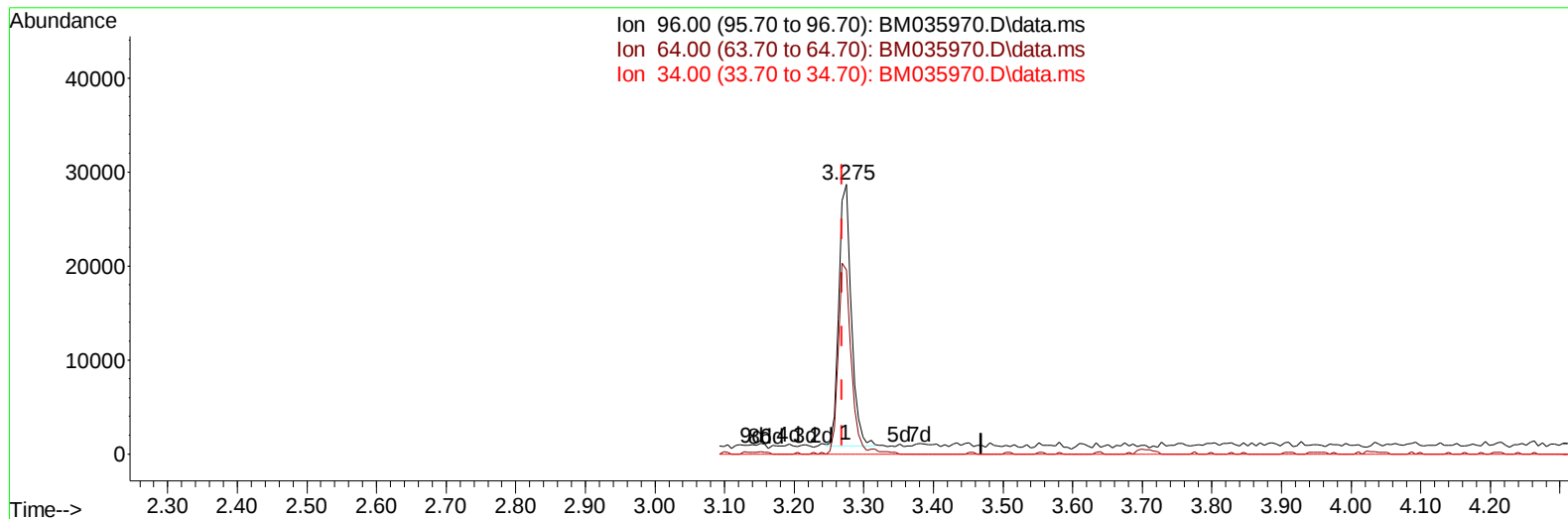
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TIC: BM035970.D\data.ms

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

3.275min (+ 0.006) 4.81 ng/uL m

response 34809

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	75.40	68.21
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	7.881	152	272215	20.000	ng/ul	0.00
20) Naphthalene-d8	10.686	136	1143061	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.521	164	680781	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.256	188	1406713	20.000	ng/ul	0.00
79) Chrysene-d12	21.433	240	1329801	20.000	ng/ul	0.00
88) Perylene-d12	23.803	264	1370169	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	34809m	4.812	ng/uL	0.00
4) Pyridine-d5	3.698	84	240307	12.126	ng/ul	0.00
7) Phenol-d5	7.045	99	625501	27.401	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.216	67	386299	25.384	ng/ul	0.00
11) 2-Chlorophenol-d4	7.410	132	547402	28.889	ng/ul	0.00
15) 4-Methylphenol-d8	8.592	113	469405	25.382	ng/ul	0.00
21) Nitrobenzene-d5	9.051	128	268038	29.592	ng/ul	0.00
24) 2-Nitrophenol-d4	9.769	143	278424	31.551	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.310	165	459423	28.785	ng/ul	0.00
31) 4-Chloroaniline-d4	10.827	131	701772	25.973	ng/ul	0.00
46) Dimethylphthalate-d6	13.933	166	1468464	31.666	ng/ul	0.00
49) Acenaphthylene-d8	14.215	160	1836245	31.035	ng/ul	0.00
54) 4-Nitrophenol-d4	14.721	143	271852	28.548	ng/ul	0.00
60) Fluorene-d10	15.509	176	1319804	31.495	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.633	200	175352	23.190	ng/ul	0.00
73) Anthracene-d10	17.356	188	2253377	35.431	ng/ul	0.00
81) Pyrene-d10	19.644	212	2986826	40.527	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.656	264	2855548	41.051	ng/ul	0.00

Target Compounds Qvalue

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

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