

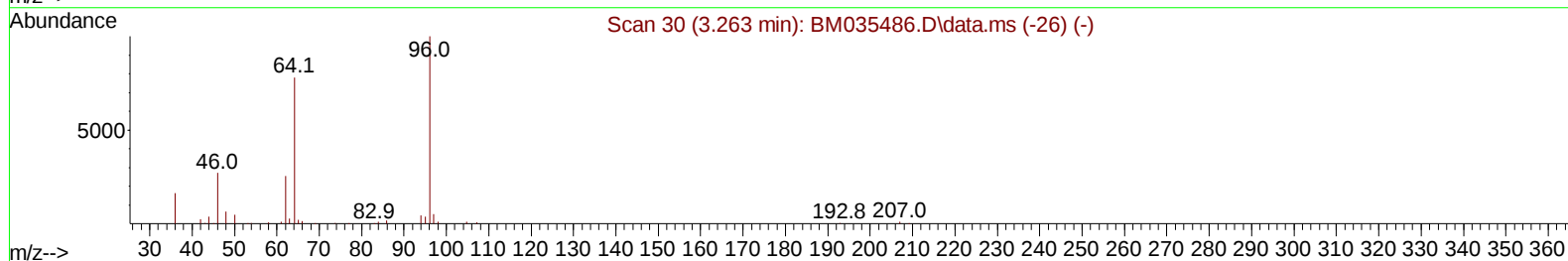
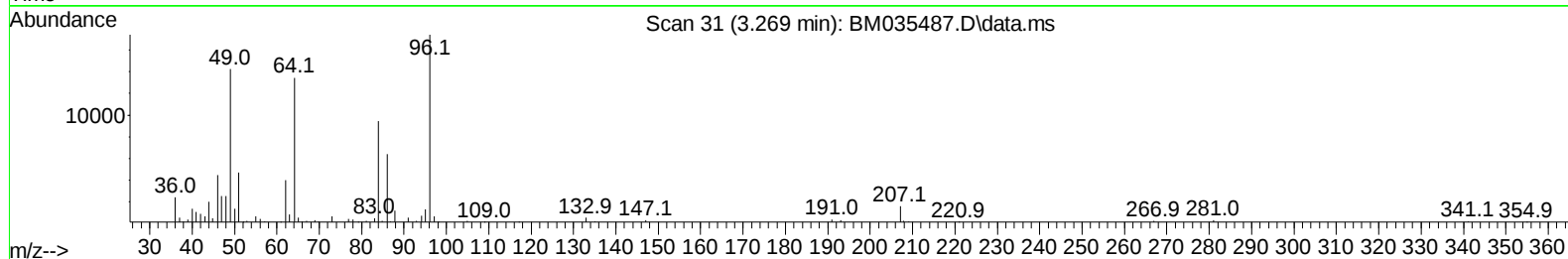
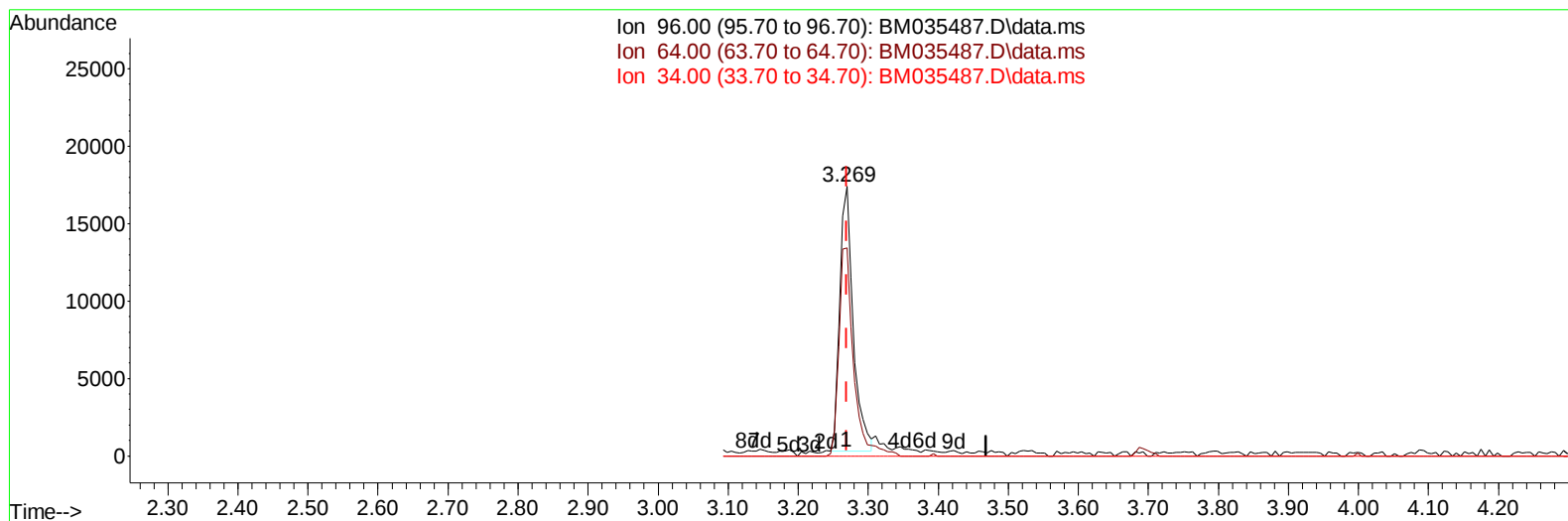
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061722\
 Data File : BM035487.D
 Acq On : 17 Jun 2022 10:15
 Operator : CG/JU
 Sample : PB145488BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK488

Manual IntegrationsAPPROVED

Quant Time: Jun 17 23:00:44 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM061622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jun 17 01:58:49 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/18/2022
 Supervised By :Yogesh Patel 06/22/2022



TIC: BM035487.D\data.ms

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

3.269min (+ 0.000) 4.50 ng/uL

response 23169

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	75.40	77.32
34.00	0.00	0.00
0.00	0.00	0.00

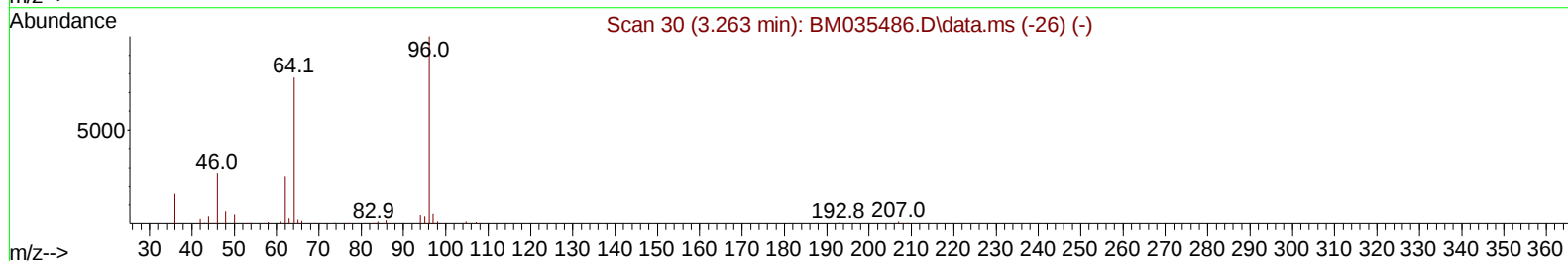
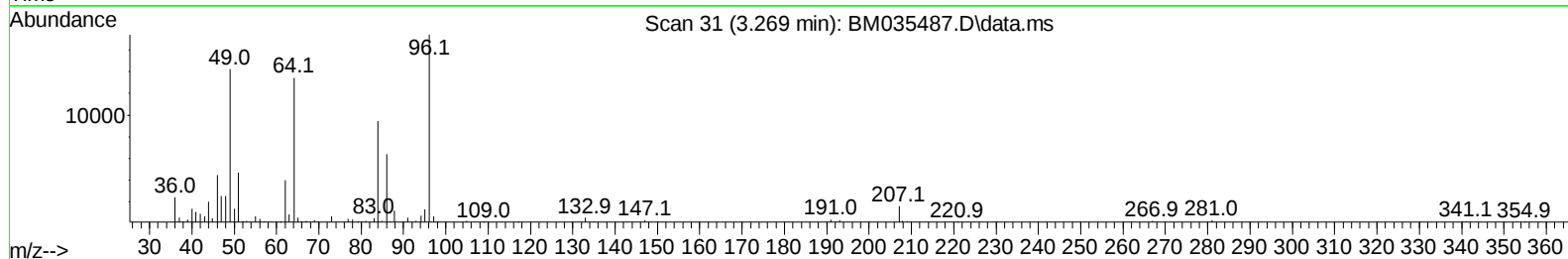
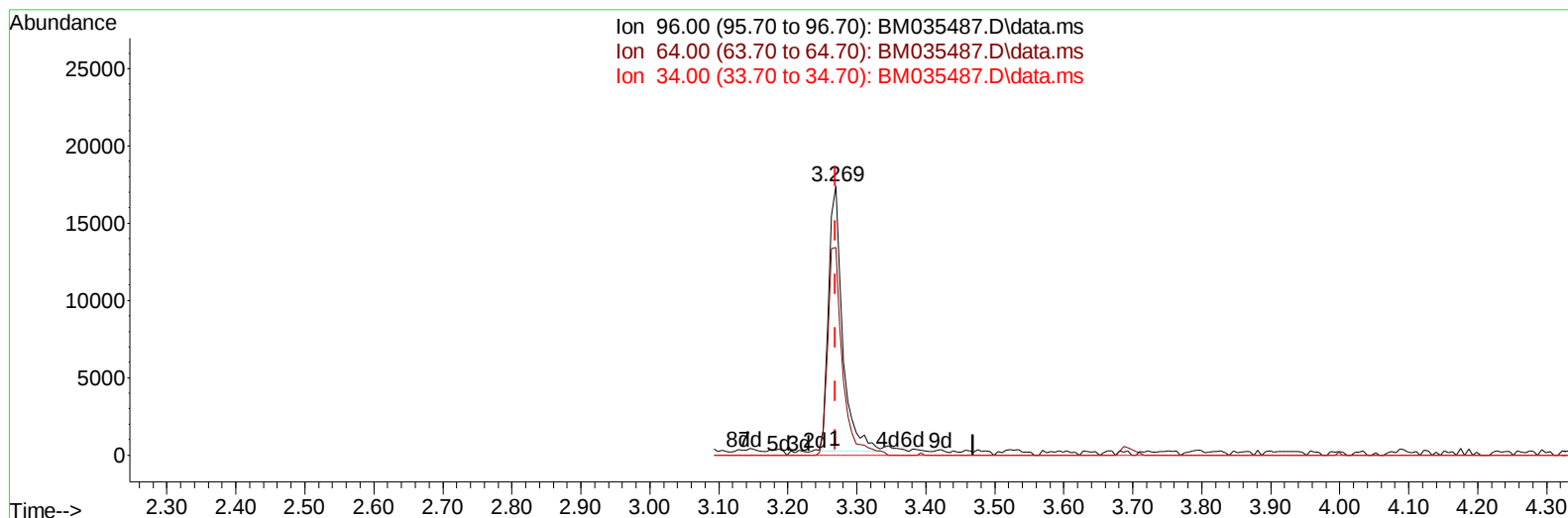
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TIC: BM035487.D\data.ms

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

3.269min (+ 0.000) 4.71 ng/uL m

response 24246

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	75.40	77.32
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	7.816	152	205667	20.000	ng/ul	0.00
20) Naphthalene-d8	10.610	136	874563	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.457	164	517275	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.204	188	1040565	20.000	ng/ul	0.00
79) Chrysene-d12	21.392	240	1063359	20.000	ng/ul	0.00
88) Perylene-d12	23.733	264	1158456	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	24246m	4.706	ng/uL	0.00
4) Pyridine-d5	3.687	84	287044	22.621	ng/ul	0.00
7) Phenol-d5	7.004	99	377839	23.971	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.146	67	247421	25.242	ng/ul	0.00
11) 2-Chlorophenol-d4	7.351	132	355625	26.048	ng/ul	0.00
15) 4-Methylphenol-d8	8.540	113	325688	24.683	ng/ul	0.00
21) Nitrobenzene-d5	8.975	128	159503	24.501	ng/ul	0.00
24) 2-Nitrophenol-d4	9.704	143	173225	24.676	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.257	165	232033	19.468	ng/ul	0.00
31) 4-Chloroaniline-d4	10.757	131	658503	35.114	ng/ul	0.00
46) Dimethylphthalate-d6	13.875	166	959313	27.264	ng/ul	0.00
49) Acenaphthylene-d8	14.151	160	1125104	25.803	ng/ul	0.00
54) 4-Nitrophenol-d4	14.733	143	95840	17.732	ng/ul	0.03
60) Fluorene-d10	15.451	176	833235	27.350	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.598	200	116153	18.723	ng/ul	0.00
73) Anthracene-d10	17.304	188	1342377	28.700	ng/ul	0.00
81) Pyrene-d10	19.598	212	1535831	29.378	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.580	264	1648048	29.495	ng/ul	0.00

Target Compounds Qvalue

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

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