

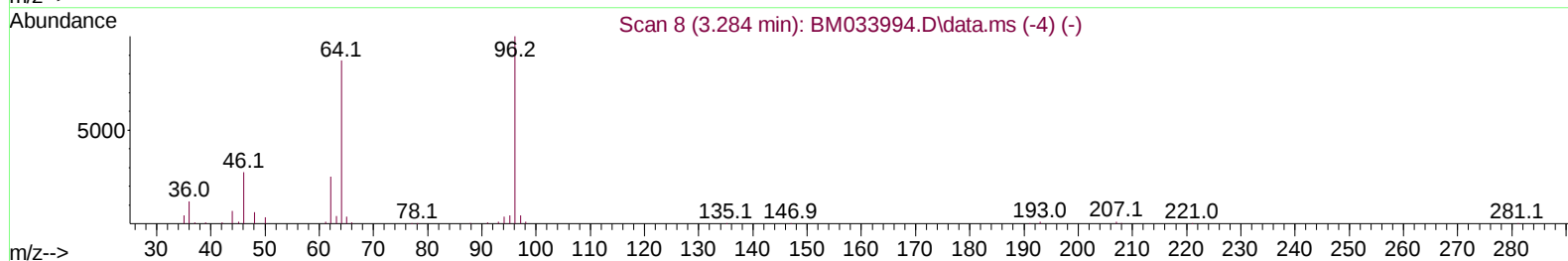
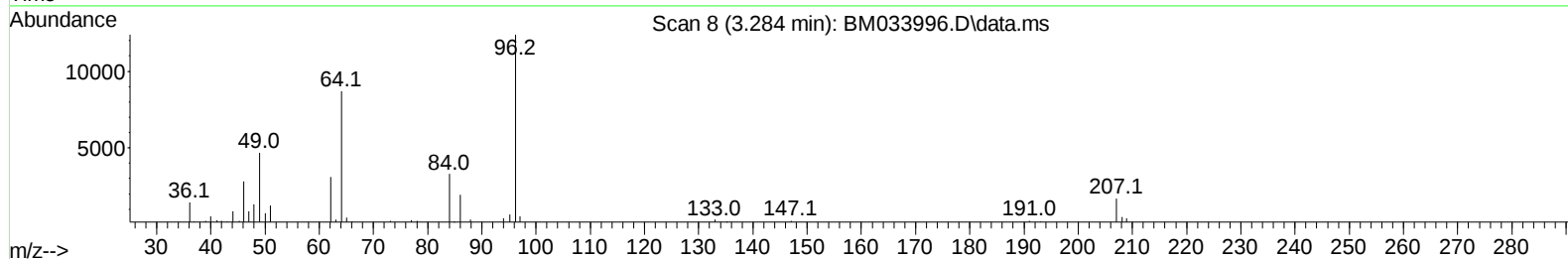
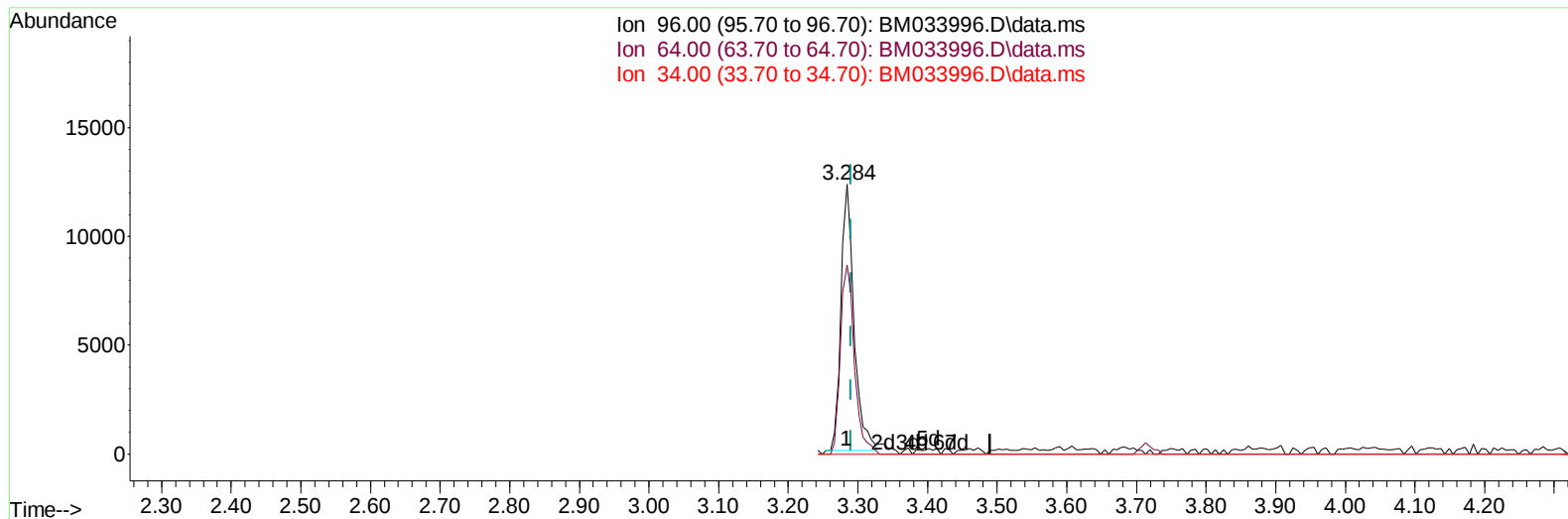
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020322\
Data File : BM033996.D
Acq On : 03 Feb 2022 11:37
Operator : CG/JU
Sample : PB142431BL
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SBLK431

Manual IntegrationsAPPROVED

Quant Time: Feb 04 00:04:57 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Feb 01 02:16:28 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 02/04/2022
Supervised By :mohammad ahmed 02/04/2022



TIC: BM033996.D\data.ms

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

3.284min (-0.006) 6.26 ng/uL

response 16040

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.20	70.11
34.00	0.00	0.00
0.00	0.00	0.00

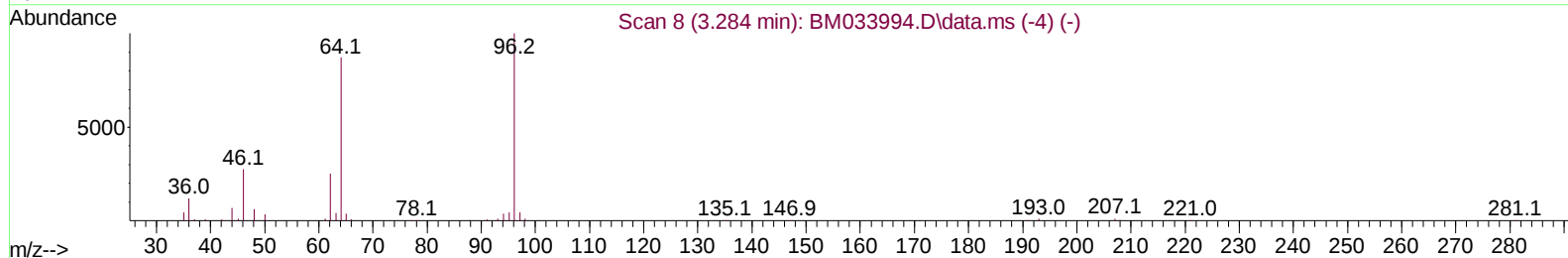
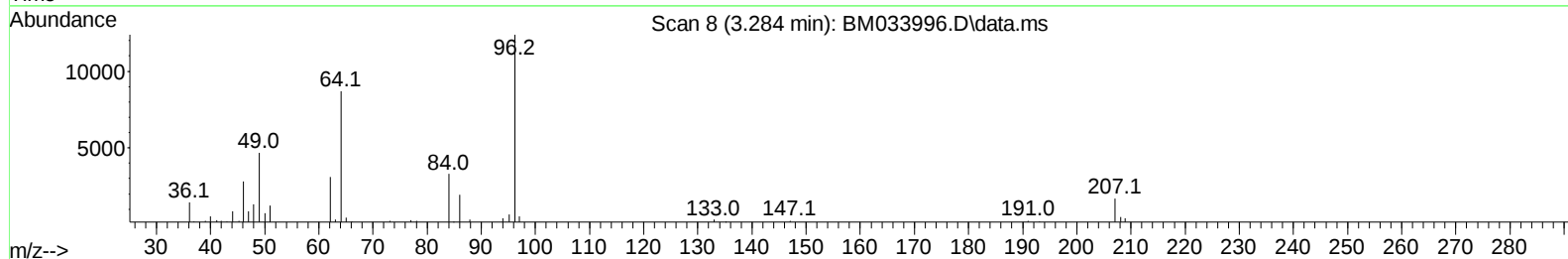
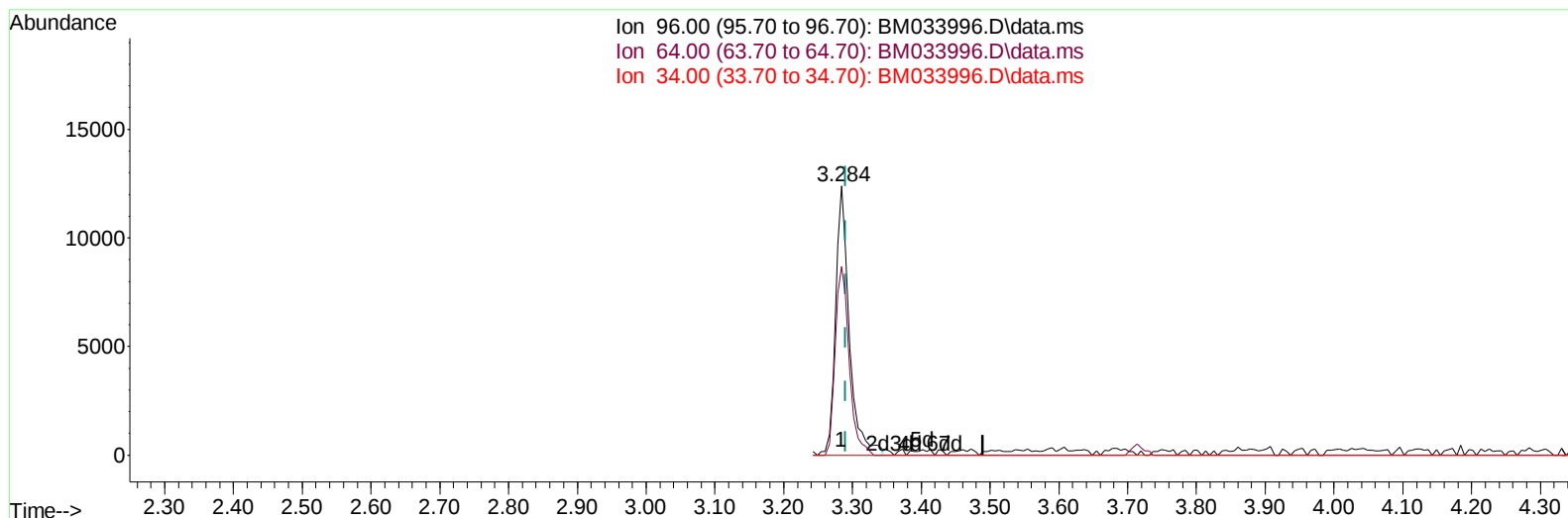
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TIC: BM033996.D\data.ms

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

3.284min (-0.006) 6.70 ng/uL m

response 17157

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.20	70.11
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	7.919	152	100986	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.731	136	435556	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.560	164	294057	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.295	188	614359	20.000	ng/ul	-0.01
79) Chrysene-d12	21.459	240	515072	20.000	ng/ul	0.00
88) Perylene-d12	23.847	264	465626	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.284	96	17157m	6.700	ng/uL	0.00
4) Pyridine-d5	3.713	84	219570	30.722	ng/ul	0.00
7) Phenol-d5	7.078	99	279644	33.054	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.243	67	178456	34.599	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.448	132	228527	33.715	ng/ul	-0.01
15) 4-Methylphenol-d8	8.631	113	238815	35.623	ng/ul	-0.01
21) Nitrobenzene-d5	9.084	128	118946	34.034	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.813	143	130957	34.699	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.348	165	226131	33.110	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.866	131	342751	36.342	ng/ul	0.00
46) Dimethylphthalate-d6	13.972	166	872515	39.707	ng/ul	0.00
49) Acenaphthylene-d8	14.254	160	968422	35.976	ng/ul	0.00
54) 4-Nitrophenol-d4	14.748	143	116600	31.059	ng/ul	0.00
60) Fluorene-d10	15.548	176	695006	37.879	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.666	200	113710	29.482	ng/ul	0.00
73) Anthracene-d10	17.395	188	1131040	38.442	ng/ul	0.00
81) Pyrene-d10	19.683	212	1241407	40.824	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.694	264	941596	38.501	ng/ul	-0.01

Target Compounds Qvalue

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

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