

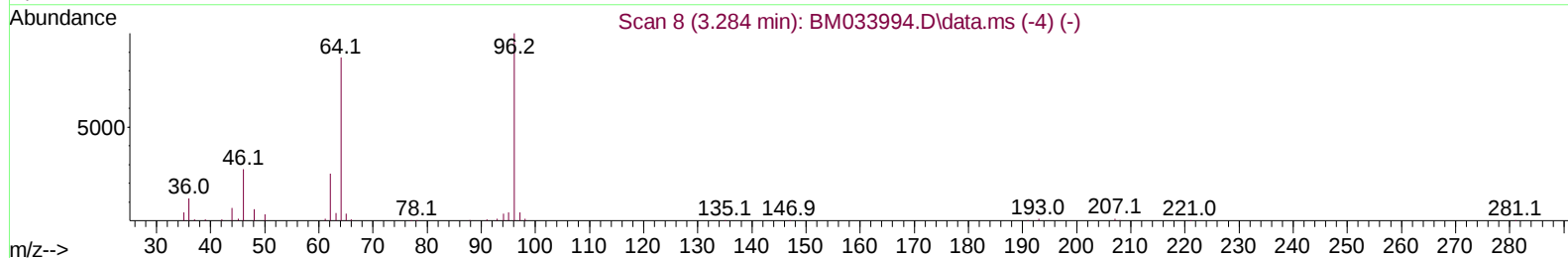
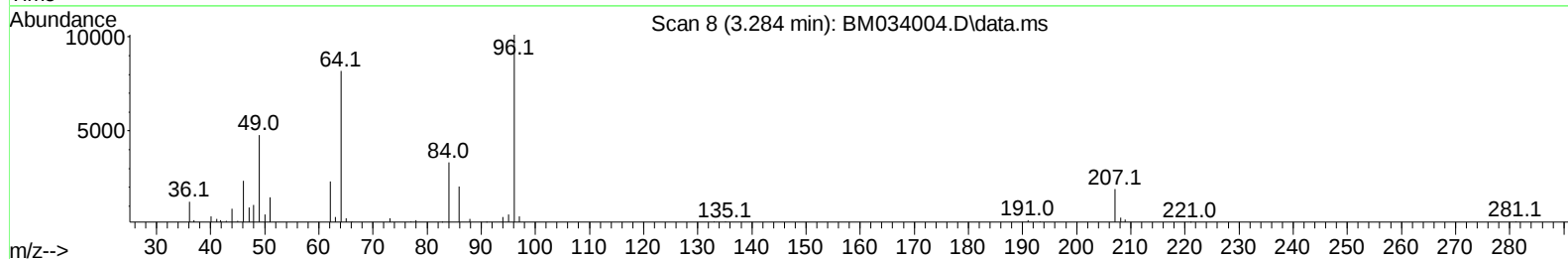
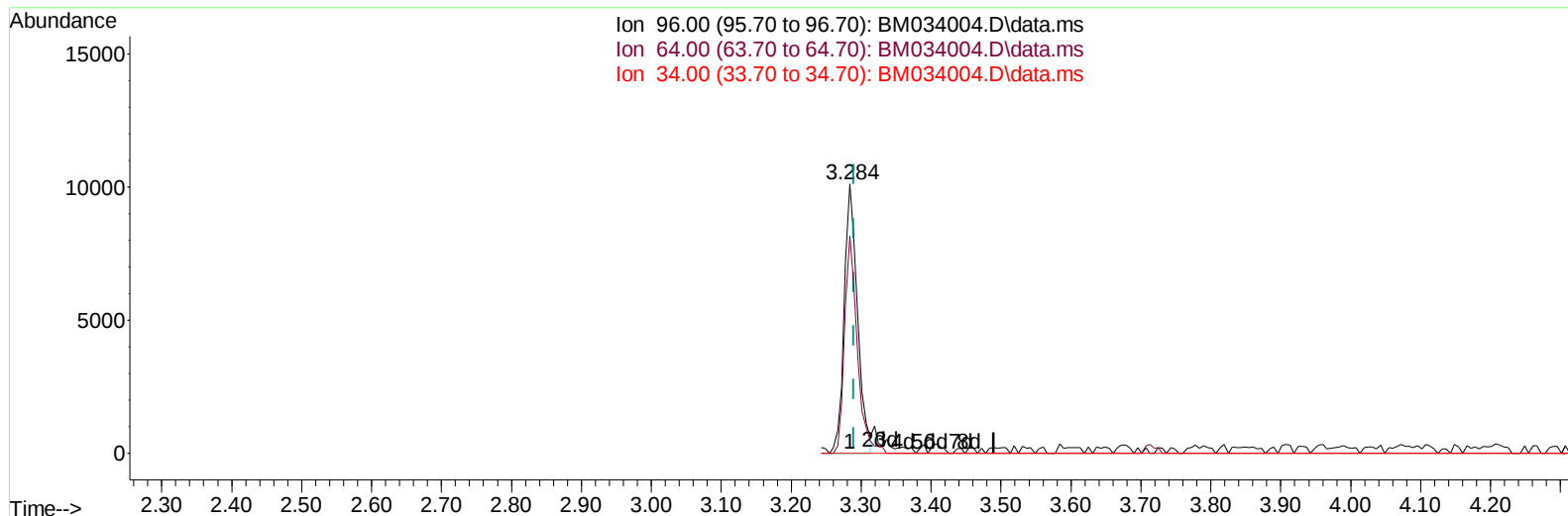
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020322\
Data File : BM034004.D
Acq On : 03 Feb 2022 16:24
Operator : CG/JU
Sample : N1365-03
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BGLY9

Manual IntegrationsAPPROVED

Quant Time: Feb 04 00:07:28 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: BM034004.D\data.ms

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

3.284min (-0.006) 5.07 ng/uL

response 13360

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.20	80.85
34.00	0.00	0.00
0.00	0.00	0.00

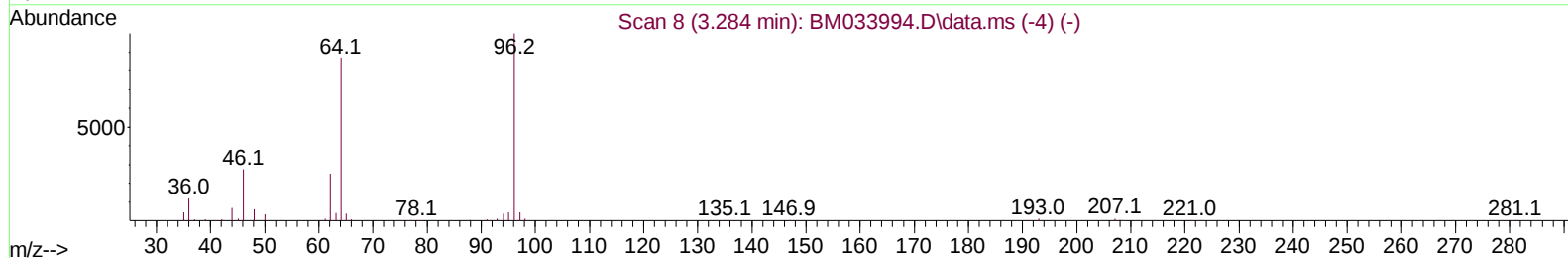
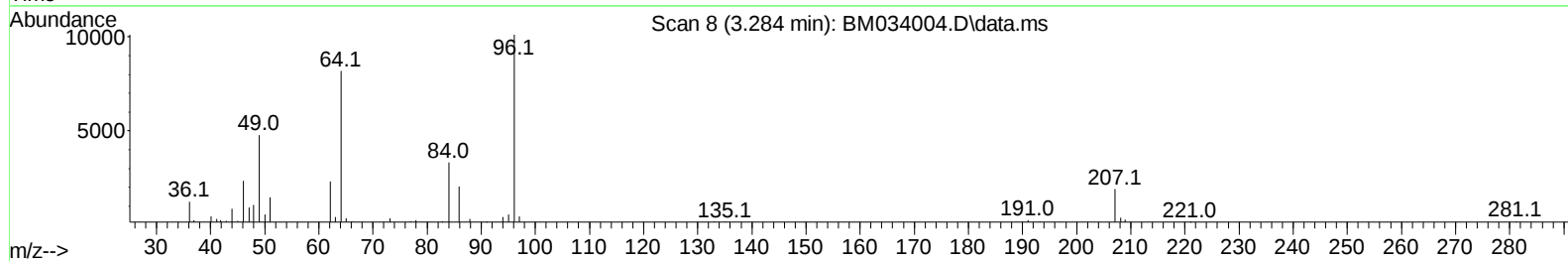
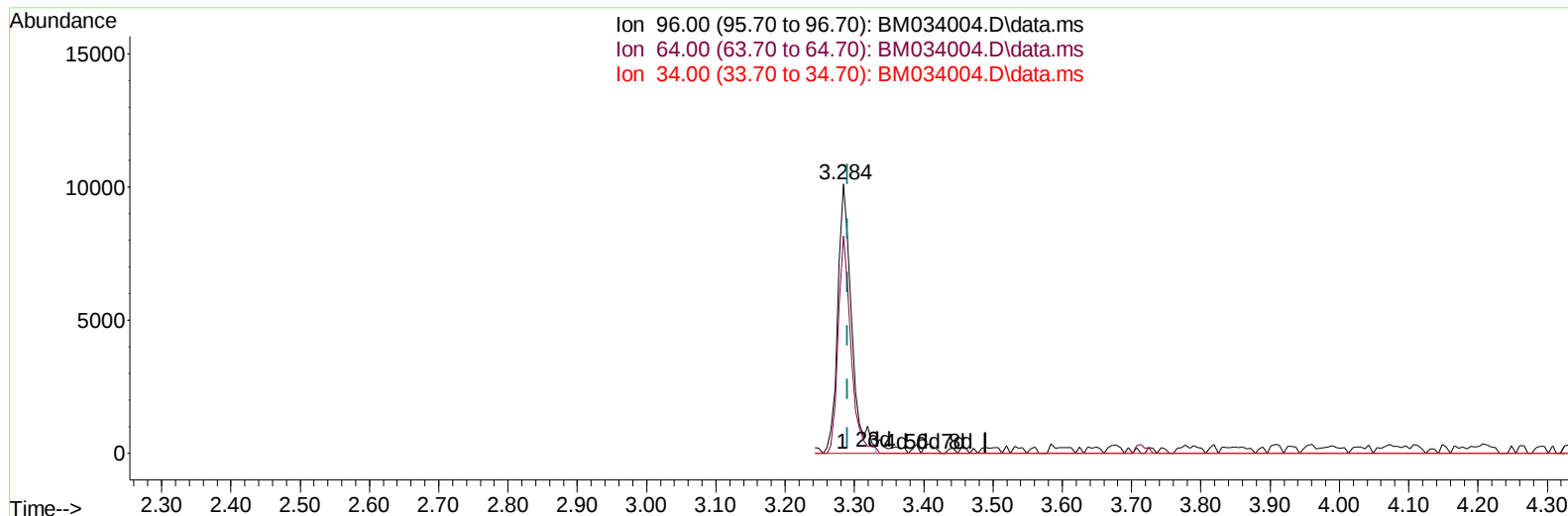
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TIC: BM034004.D\data.ms

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

3.284min (-0.006) 5.29 ng/uL m

response 13945

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.20	80.85
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	103880	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.730	136	446634	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.560	164	289939	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.295	188	585037	20.000	ng/ul	-0.01
79) Chrysene-d12	21.459	240	482407	20.000	ng/ul	0.00
88) Perylene-d12	23.847	264	447560	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.284	96	13945m	5.294	ng/uL	0.00
4) Pyridine-d5	3.713	84	135569	18.440	ng/ul	0.00
7) Phenol-d5	7.078	99	238592	27.416	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.242	67	153812	28.990	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.448	132	195923	28.100	ng/ul	-0.01
15) 4-Methylphenol-d8	8.631	113	198812	28.830	ng/ul	-0.01
21) Nitrobenzene-d5	9.083	128	100883	28.150	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.807	143	108703	28.088	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.348	165	187083	26.713	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.866	131	276430	28.583	ng/ul	0.00
46) Dimethylphthalate-d6	13.965	166	667324	30.800	ng/ul	-0.01
49) Acenaphthylene-d8	14.254	160	767960	28.934	ng/ul	0.00
54) 4-Nitrophenol-d4	14.748	143	73103	19.749	ng/ul	0.00
60) Fluorene-d10	15.548	176	543863	30.062	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.659	200	65472	17.826	ng/ul	-0.01
73) Anthracene-d10	17.395	188	835777	29.831	ng/ul	0.00
81) Pyrene-d10	19.677	212	886405	31.123	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.694	264	697941	29.690	ng/ul	-0.01
Target Compounds						
					Qvalue	
8) Phenol	7.101	94	18610	2.085	ng/ul#	90

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

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