

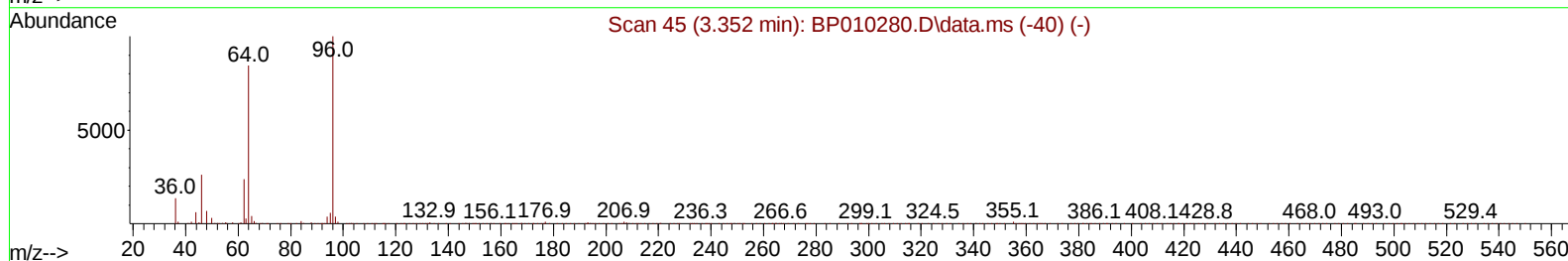
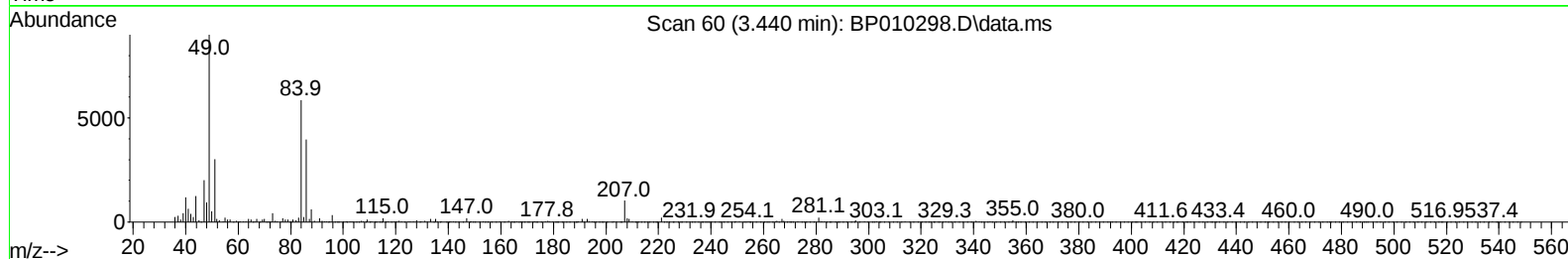
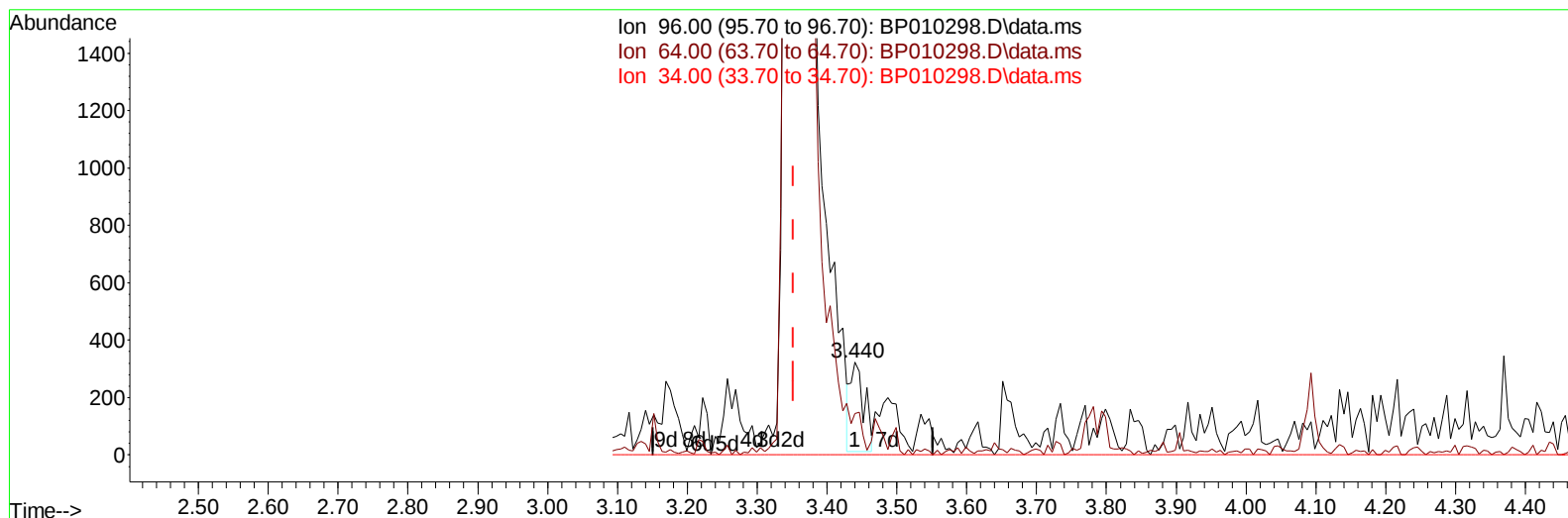
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051122\
 Data File : BP010298.D
 Acq On : 11 May 2022 22:47
 Operator : CG/JU
 Sample : N2707-08
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EXL89

Manual IntegrationsAPPROVED

Quant Time: May 12 01:16:06 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu May 12 01:10:15 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/12/2022
 Supervised By :Yogesh Patel 05/17/2022



TIC: BP010298.D\data.ms

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

3.440min (+ 0.088) 0.10 ng/uL

response 423

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	44.89
34.00	0.00	0.00
0.00	0.00	0.00

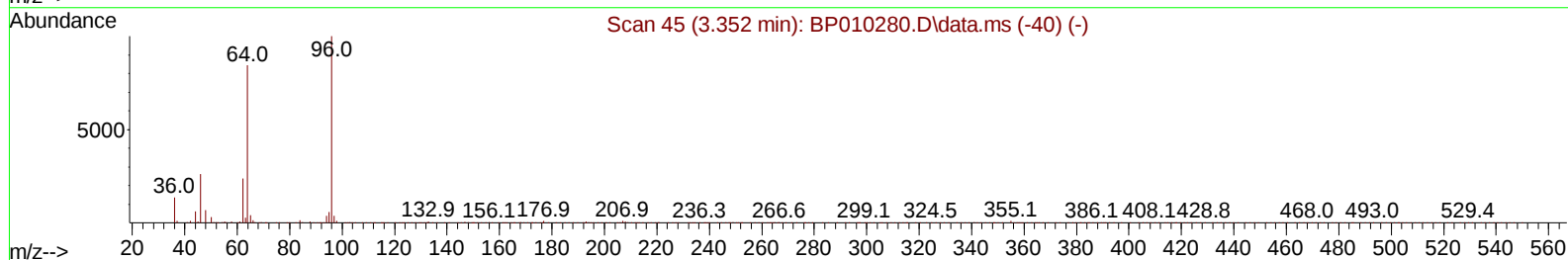
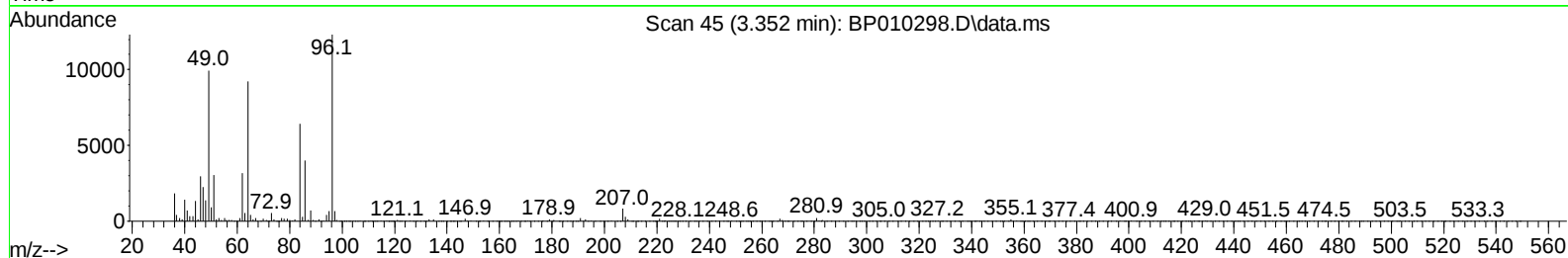
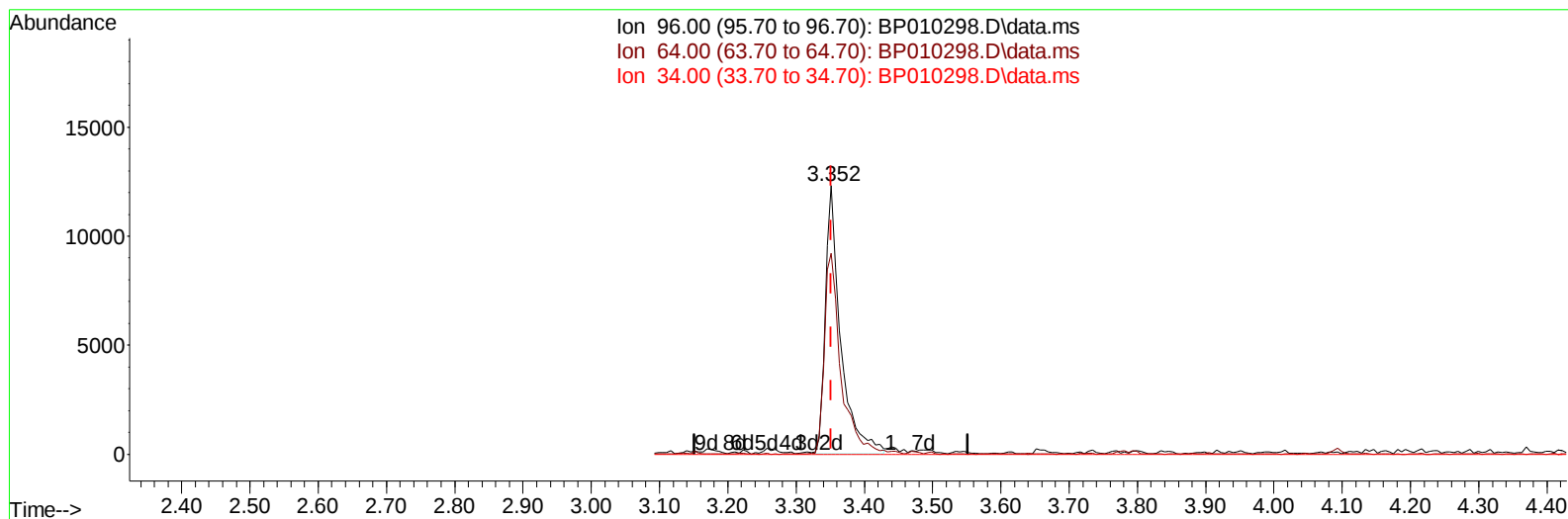
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TIC: BP010298.D\data.ms

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

3.352min (-0.000) 4.70 ng/uL m

response 19072

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	74.94#
34.00	0.00	0.00
0.00	0.00	0.00

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Instrument :
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.904	152	172376	20.000	ng/ul	0.00
20) Naphthalene-d8	10.710	136	764342	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.539	164	550577	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.286	188	1196069	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.368	240	1173618	20.000	ng/ul	# 0.00
88) Perylene-d12	23.804	264	1118209	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	19072m	4.700	ng/uL	0.00
4) Pyridine-d5	3.775	84	76469	6.772	ng/ul	0.00
7) Phenol-d5	7.069	99	76883	4.802	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.234	67	286152	28.892	ng/ul	0.00
11) 2-Chlorophenol-d4	7.440	132	250655	21.258	ng/ul	0.00
15) 4-Methylphenol-d8	8.616	113	163572	12.427	ng/ul	0.00
21) Nitrobenzene-d5	9.063	128	178803	29.175	ng/ul	0.00
24) 2-Nitrophenol-d4	9.787	143	186528	26.694	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.328	165	295604	23.242	ng/ul	0.00
31) 4-Chloroaniline-d4	10.845	131	444601	23.679	ng/ul	0.00
46) Dimethylphthalate-d6	13.945	166	1285285	29.303	ng/ul	0.00
49) Acenaphthylene-d8	14.228	160	1428235	27.311	ng/ul	0.00
54) 4-Nitrophenol-d4	14.739	143	27619	3.283	ng/ul	0.00
60) Fluorene-d10	15.533	176	1122743	30.324	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.651	200	211814	25.861	ng/ul	0.00
73) Anthracene-d10	17.386	188	1885107	32.185	ng/ul	0.00
81) Pyrene-d10	19.621	212	2288649	35.289	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.645	264	2046062	34.310	ng/ul	-0.01

Target Compounds Qvalue

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

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