

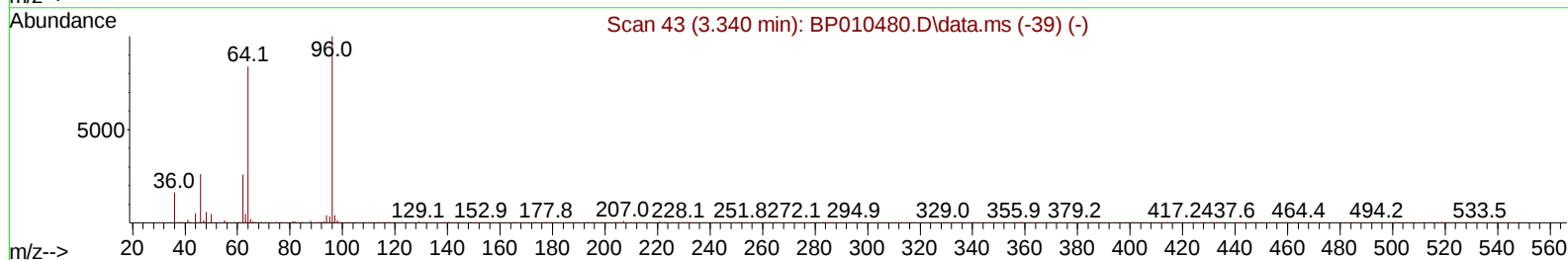
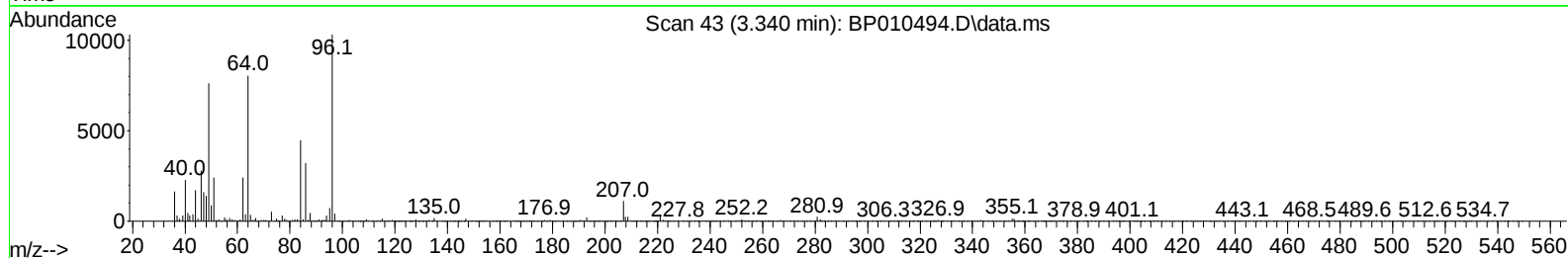
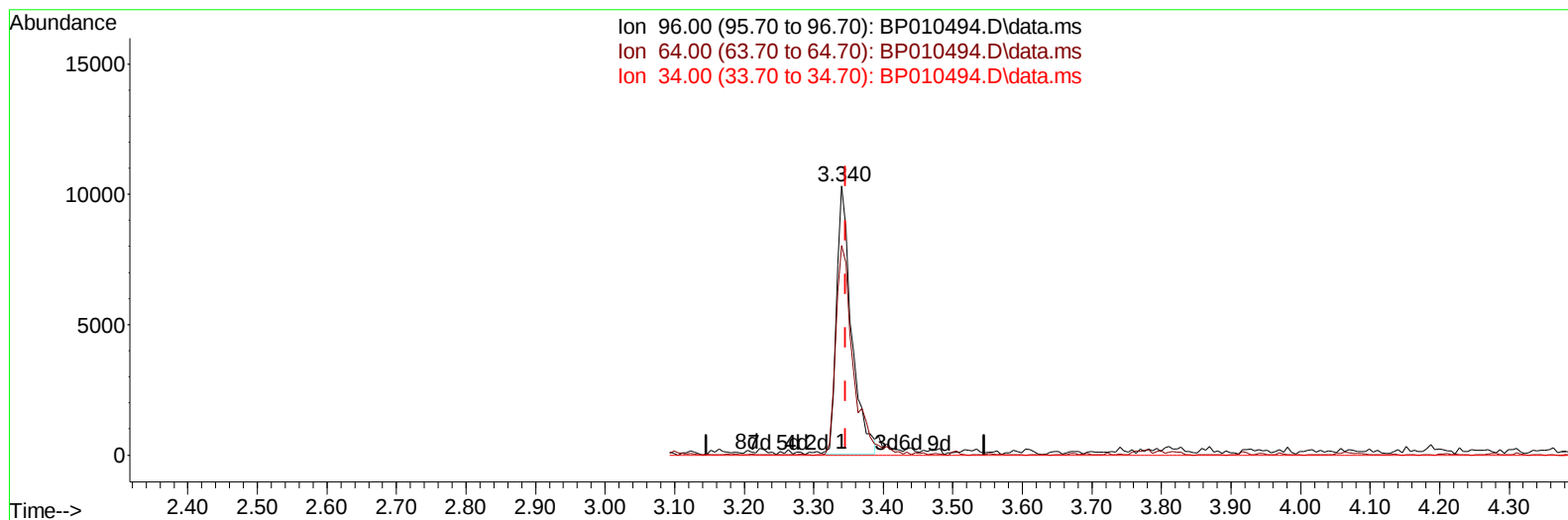
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052322\
 Data File : BP010494.D
 Acq On : 24 May 2022 01:06
 Operator : CG/JU
 Sample : N2950-04
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBTF0

Manual IntegrationsAPPROVED

Quant Time: May 24 02:39:31 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri May 13 15:04:31 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/24/2022
 Supervised By :Yogesh Patel 05/26/2022



TIC: BP010494.D\data.ms

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

3.340min (-0.006) 3.87 ng/uL

response 15400

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	78.04#
34.00	0.00	0.00
0.00	0.00	0.00

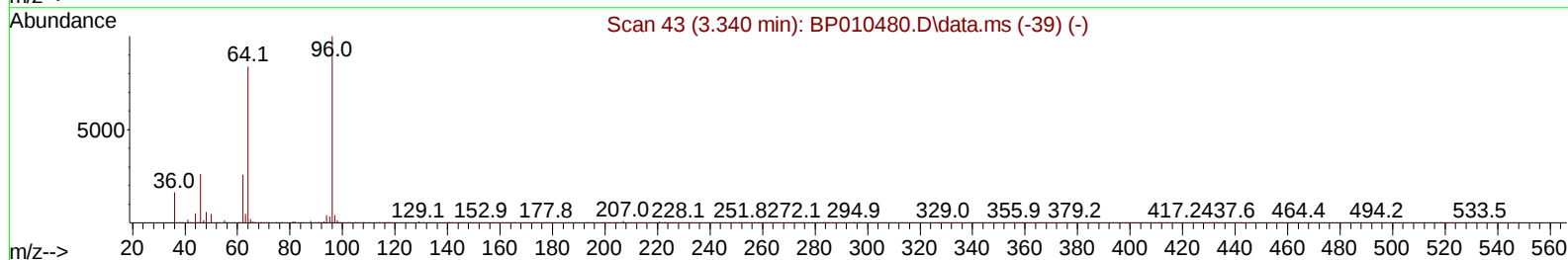
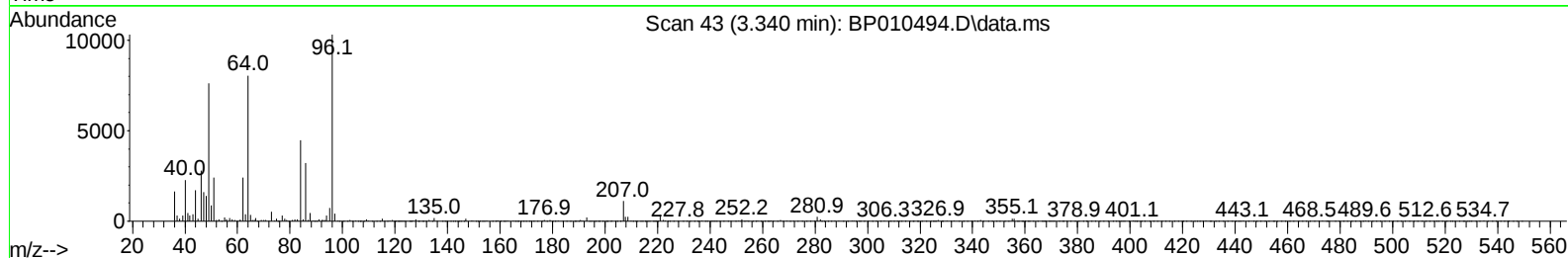
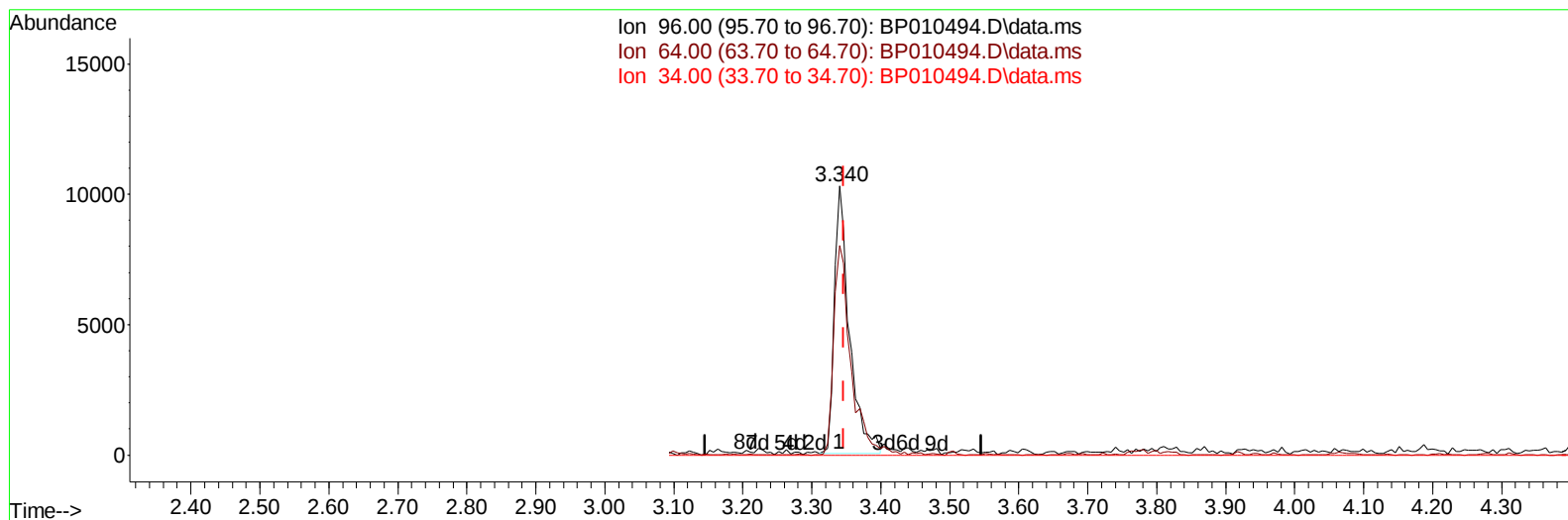
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TIC: BP010494.D\data.ms

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

3.340min (-0.006) 3.94 ng/uL m

response 15683

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	78.04#
34.00	0.00	0.00
0.00	0.00	0.00

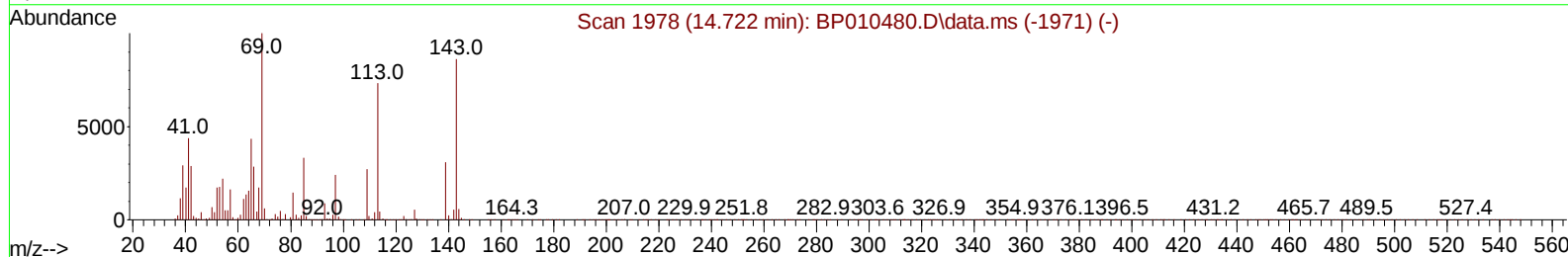
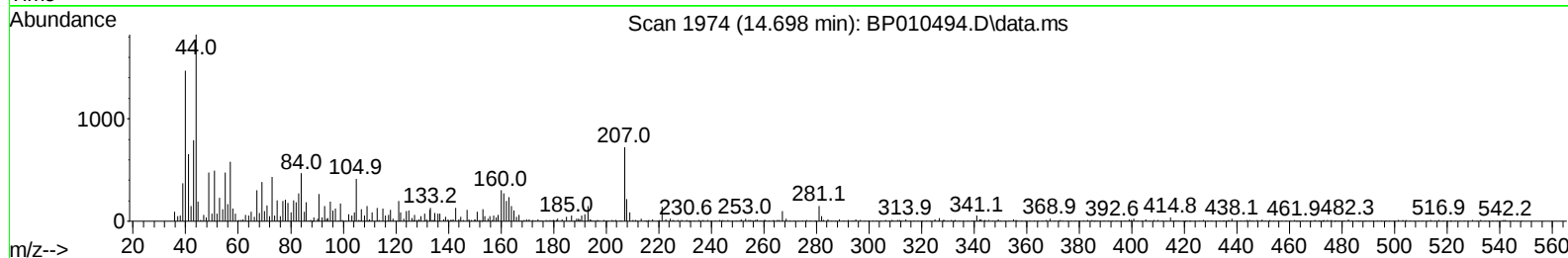
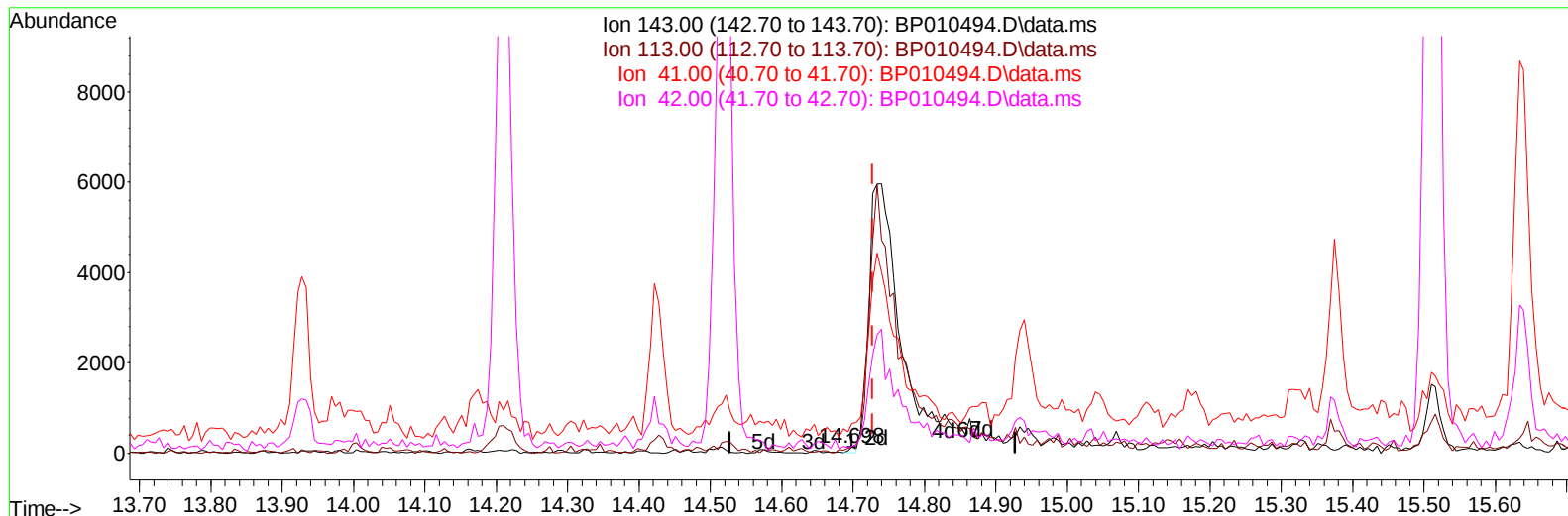
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TIC: BP010494.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.698min (-0.029) 0.01 ng/u1

response 83

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	109.30
41.00	23.20	506.20#
42.00	18.00	115.50#

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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	7.887	152	169782	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.687	136	772410	20.000	ng/ul	#-0.02
38) Acenaphthene-d10	14.522	164	517874	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.269	188	1081032	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.357	240	931293	20.000	ng/ul	# 0.00
88) Perylene-d12	23.774	264	741490	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	15683m	3.937	ng/uL	0.00
4) Pyridine-d5	3.764	84	69648	6.158	ng/ul	0.00
7) Phenol-d5	7.058	99	63507	4.457	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.216	67	263171	27.487	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.422	132	175279	16.075	ng/ul	-0.01
15) 4-Methylphenol-d8	8.599	113	131072	10.823	ng/ul	-0.01
21) Nitrobenzene-d5	9.046	128	165857	27.745	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.769	143	127017	20.049	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.310	165	224276	18.723	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.834	131	40513	2.275	ng/ul	0.00
46) Dimethylphthalate-d6	13.928	166	1176919	30.845	ng/ul	-0.01
49) Acenaphthylene-d8	14.210	160	1227320	27.887	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.734	143	19899m	2.930	ng/ul	0.00
60) Fluorene-d10	15.510	176	996976	30.034	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.634	200	110570	17.489	ng/ul	-0.01
73) Anthracene-d10	17.369	188	1614833	32.989	ng/ul	-0.01
81) Pyrene-d10	19.604	212	1919618	38.813	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.621	264	1274442	34.229	ng/ul	-0.01
Target Compounds						
					Qvalue	
47) Dimethylphthalate	13.975	163	478768	12.650	ng/ul#	92

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

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