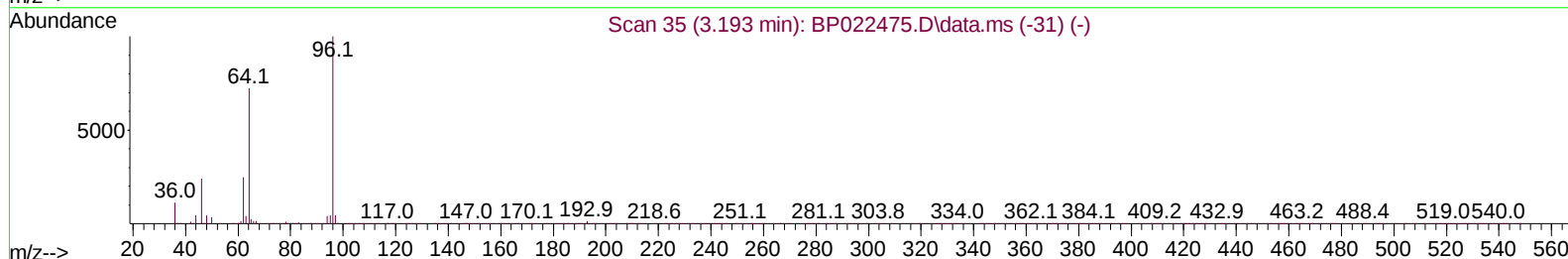
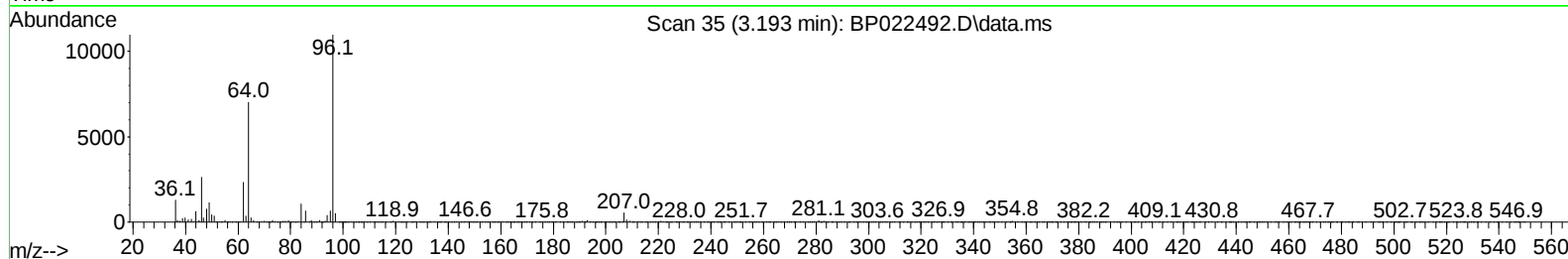
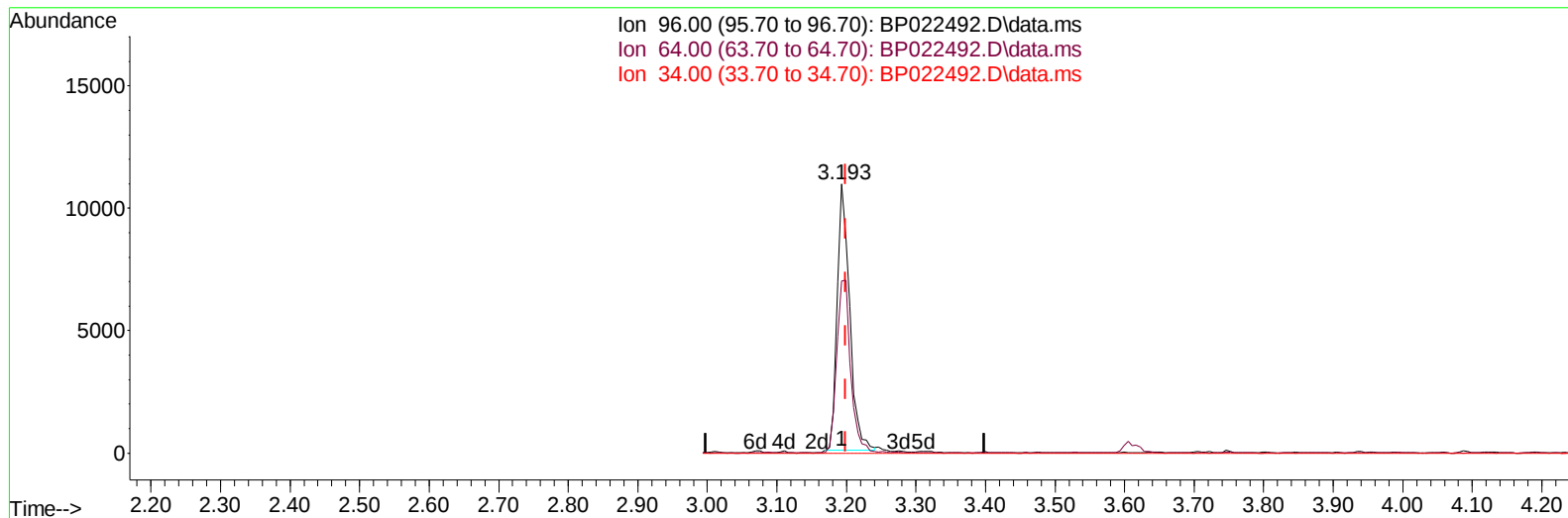


Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101924\
Data File : BP022492.D
Acq On : 20 Oct 2024 02:26
Operator : RC/JU
Sample : PB164183BL
Misc :
ALS Vial : 20 Sample Multiplier: 1

Quant Time: Oct 21 03:09:09 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Oct 18 11:08:24 2024
Response via : Initial Calibration



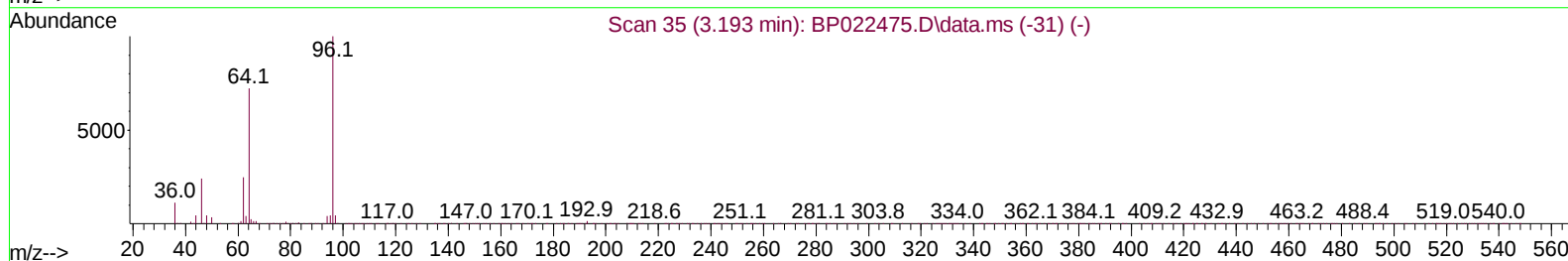
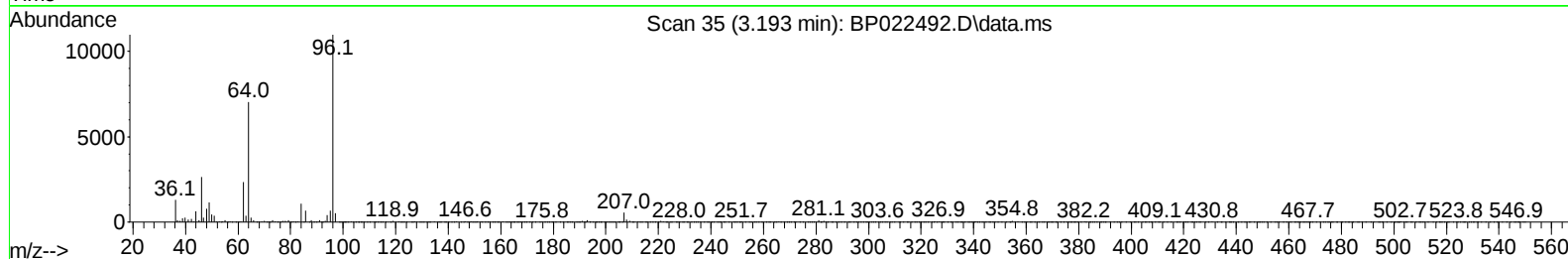
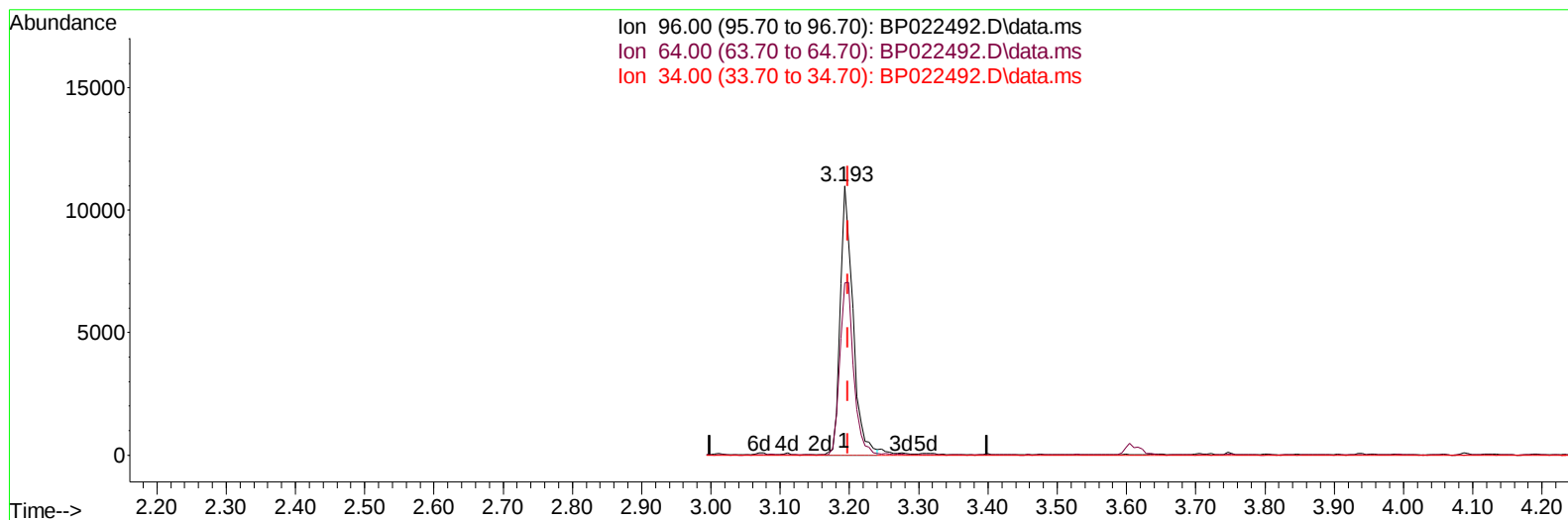
TIC: BP022492.D\data.ms

(3) 1,4-Dioxane-d8 (S)**3.193min (-0.006) 4.43 ng/uL****response 13594**

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	76.40	63.98
34.00	0.00	0.00
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101924\
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Sample : PB164183BL
Misc :
ALS Vial : 20 Sample Multiplier: 1

Quant Time: Oct 21 03:09:09 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Oct 18 11:08:24 2024
Response via : Initial Calibration



TIC: BP022492.D\data.ms

(3) 1,4-Dioxane-d8 (S)**3.193min (-0.006) 4.61 ng/uL m****response 14135**

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	76.40	63.98
34.00	0.00	0.00
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101924\
 Data File : BP022492.D
 Acq On : 20 Oct 2024 02:26
 Operator : RC/JU
 Sample : PB164183BL
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Quant Time: Oct 21 03:09:55 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 18 11:08:24 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.663	152	119204	20.000	ng/ul	0.00
20) Naphthalene-d8	10.416	136	485101	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.275	164	410503	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.069	188	942311	20.000	ng/ul	-0.01
79) Chrysene-d12	21.504	240	763328	20.000	ng/ul	0.00
88) Perylene-d12	24.751	264	885399	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.193	96	14135m	4.608	ng/uL	0.00
4) Pyridine-d5	3.605	84	192076	22.809	ng/ul	0.00
7) Phenol-d5	6.852	99	230514	22.973	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.005	67	145617	24.691	ng/ul	0.00
11) 2-Chlorophenol-d4	7.205	132	176901	24.846	ng/ul	0.00
15) 4-Methylphenol-d8	8.363	113	195856	24.458	ng/ul	0.00
21) Nitrobenzene-d5	8.805	128	93912	25.178	ng/ul	0.00
24) 2-Nitrophenol-d4	9.516	143	114478	26.510	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.052	165	227245	24.760	ng/ul	0.00
31) 4-Chloroaniline-d4	10.552	131	273287	25.199	ng/ul	0.00
46) Dimethylphthalate-d6	13.681	166	860790	26.392	ng/ul	0.00
49) Acenaphthylene-d8	13.963	160	880616	25.421	ng/ul	0.00
54) 4-Nitrophenol-d4	14.498	143	100194	18.312	ng/ul	0.00
60) Fluorene-d10	15.275	176	725637	25.650	ng/ul	-0.02
65) 4,6-Dinitro-2-methylph...	15.422	200	131407	21.203	ng/ul	0.00
73) Anthracene-d10	17.181	188	1138777	25.689	ng/ul	0.01
81) Pyrene-d10	19.551	212	1237953	30.668	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.521	264	1245387	26.338	ng/ul	0.00

Target Compounds	Qvalue

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

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