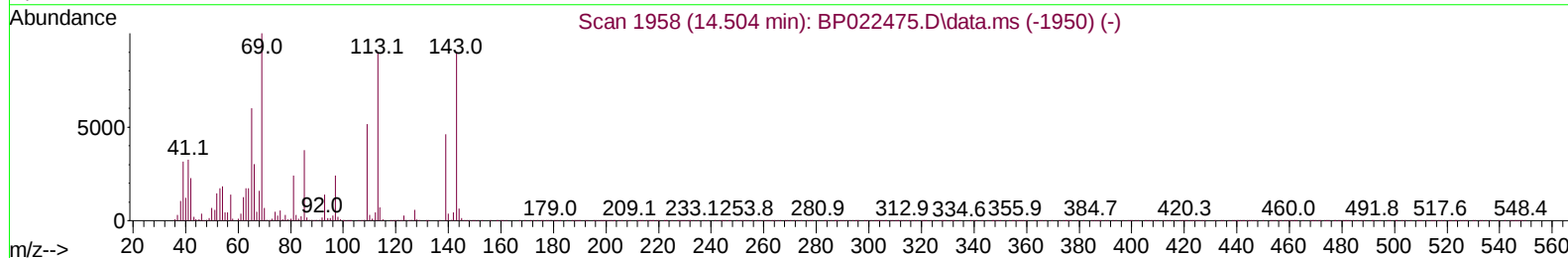
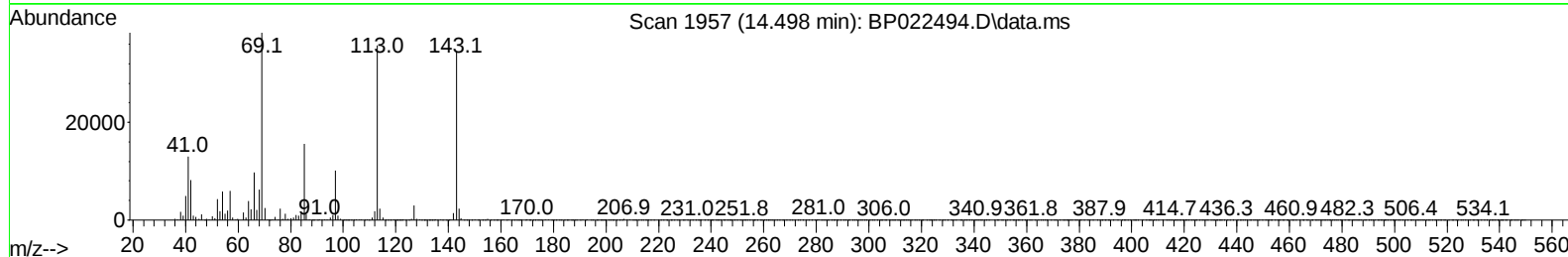
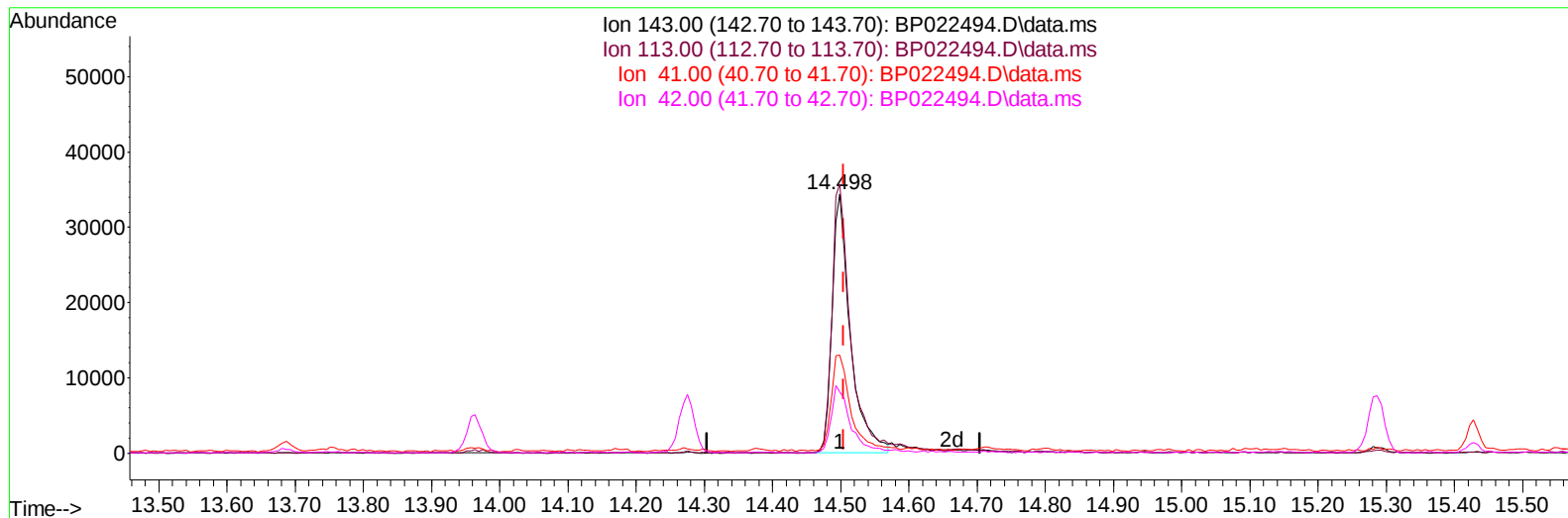


Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101924\
Data File : BP022494.D
Acq On : 20 Oct 2024 03:48
Operator : RC/JU
Sample : P4361-07
Misc :
ALS Vial : 22 Sample Multiplier: 1

Quant Time: Oct 20 08:49:55 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



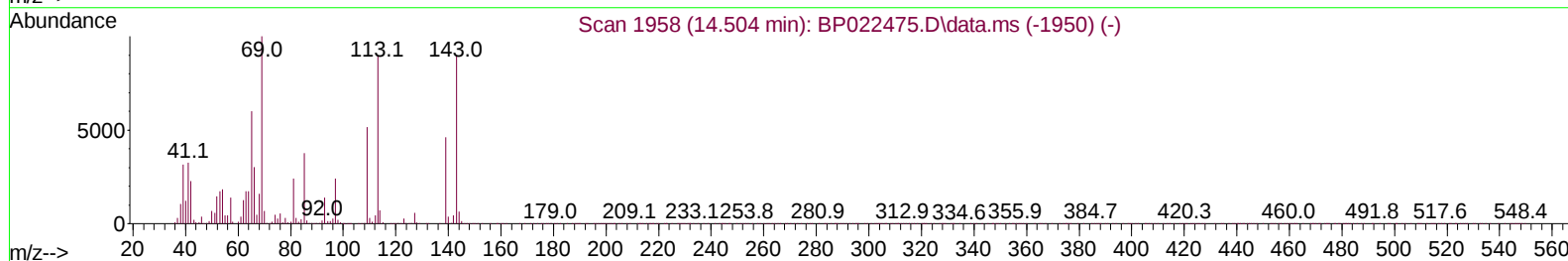
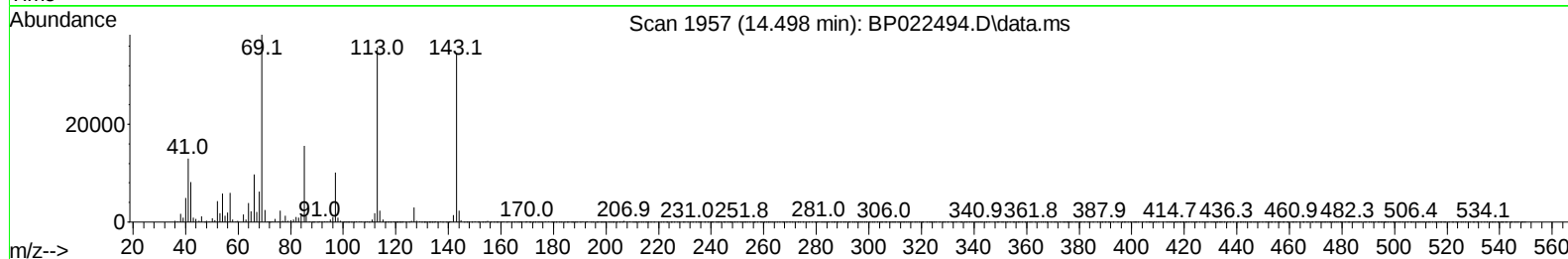
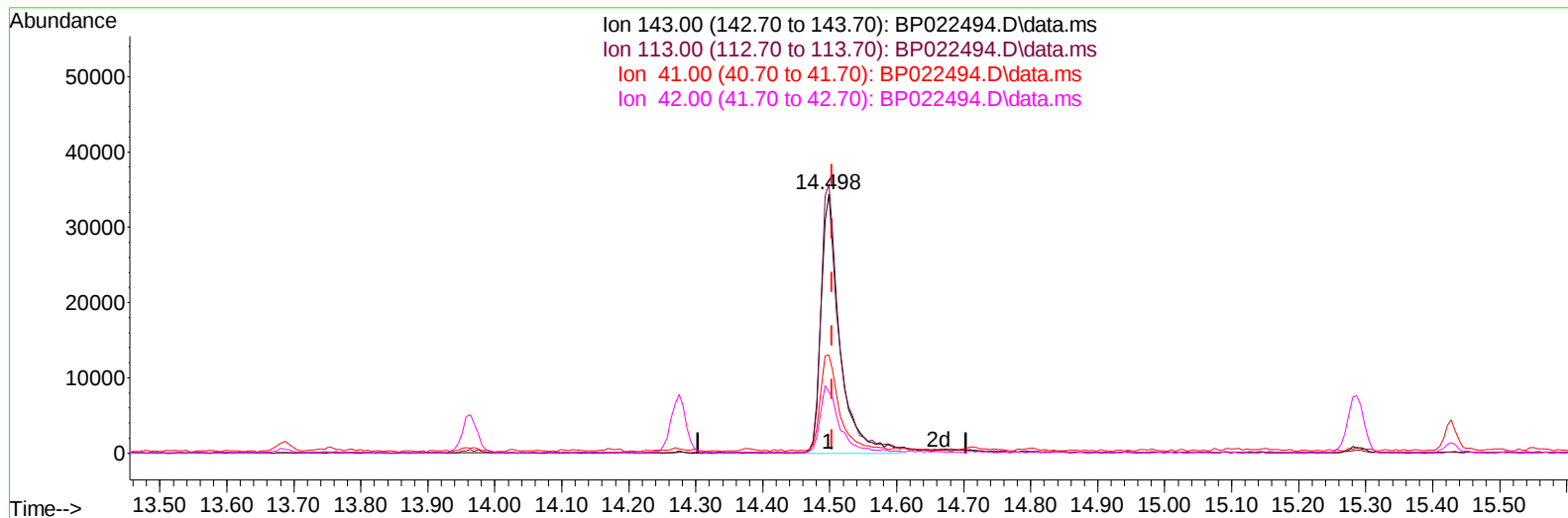
TIC: BP022494.D\data.ms

(54) 4-Nitrophenol-d4 (S)**14.498min (-0.005) 11.24 ng/ul****response 64297**

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	103.79
41.00	35.90	37.84
42.00	23.70	23.65

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101924\
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Misc :
ALS Vial : 22 Sample Multiplier: 1

Quant Time: Oct 20 08:49:55 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



TIC: BP022494.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (-0.005) 11.62 ng/ul m

response 66459

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	103.79
41.00	35.90	37.84
42.00	23.70	23.65

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101924\
 Data File : BP022494.D
 Acq On : 20 Oct 2024 03:48
 Operator : RC/JU
 Sample : P4361-07
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Quant Time: Oct 21 03:26:29 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	7.658	152	119167	20.000	ng/ul	0.00
20) Naphthalene-d8	10.416	136	503604	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.275	164	429172	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.087	188	1041307	20.000	ng/ul	0.00
79) Chrysene-d12	21.498	240	930260	20.000	ng/ul	0.00
88) Perylene-d12	24.757	264	915072	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.193	96	8263	2.695	ng/uL	0.00
4) Pyridine-d5	3.605	84	115023	13.663	ng/ul	0.00
7) Phenol-d5	6.852	99	169175	16.865	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.005	67	101408	17.200	ng/ul	0.00
11) 2-Chlorophenol-d4	7.205	132	123720	17.382	ng/ul	0.00
15) 4-Methylphenol-d8	8.363	113	154617	19.314	ng/ul	0.00
21) Nitrobenzene-d5	8.799	128	66946	17.289	ng/ul	0.00
24) 2-Nitrophenol-d4	9.510	143	82514	18.406	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.052	165	177242	18.602	ng/ul	0.00
31) 4-Chloroaniline-d4	10.557	131	200458	17.805	ng/ul	0.00
46) Dimethylphthalate-d6	13.687	166	676798	19.848	ng/ul	0.00
49) Acenaphthylene-d8	13.963	160	682888	18.856	ng/ul	0.00
54) 4-Nitrophenol-d4	14.498	143	66459m	11.618	ng/ul	0.00
60) Fluorene-d10	15.287	176	585346	19.791	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.428	200	75598	11.039	ng/ul	0.01
73) Anthracene-d10	17.175	188	983123	20.070	ng/ul	0.00
81) Pyrene-d10	19.557	212	1181851	24.024	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.533	264	1015070	20.771	ng/ul	0.00
Target Compounds						
					Qvalue	
72) Phenanthrene	17.122	178	78914	1.416	ng/ul	98
80) Fluoranthene	19.210	202	170313	3.012	ng/ul	99
82) Pyrene	19.586	202	141971	2.342	ng/ul	98
85) Benzo(a)anthracene	21.480	228	71118	1.091	ng/ul	98
87) Chrysene	21.545	228	79368	1.313	ng/ul	96
90) Benzo(b)fluoranthene	23.727	252	107275	1.862	ng/ul	98
93) Benzo(a)pyrene	24.604	252	63725	1.202	ng/ul	98

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

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