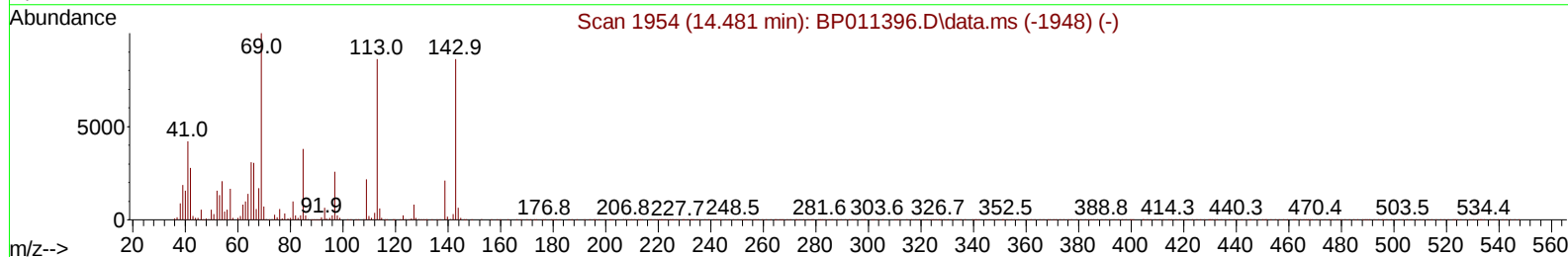
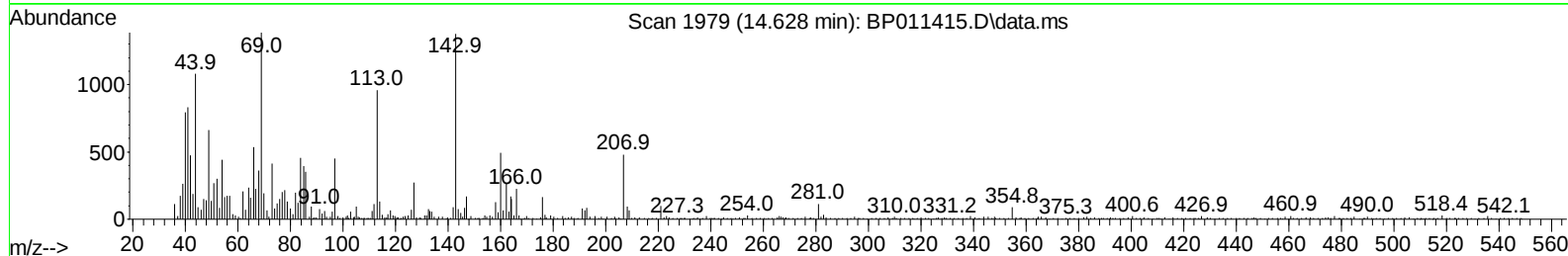
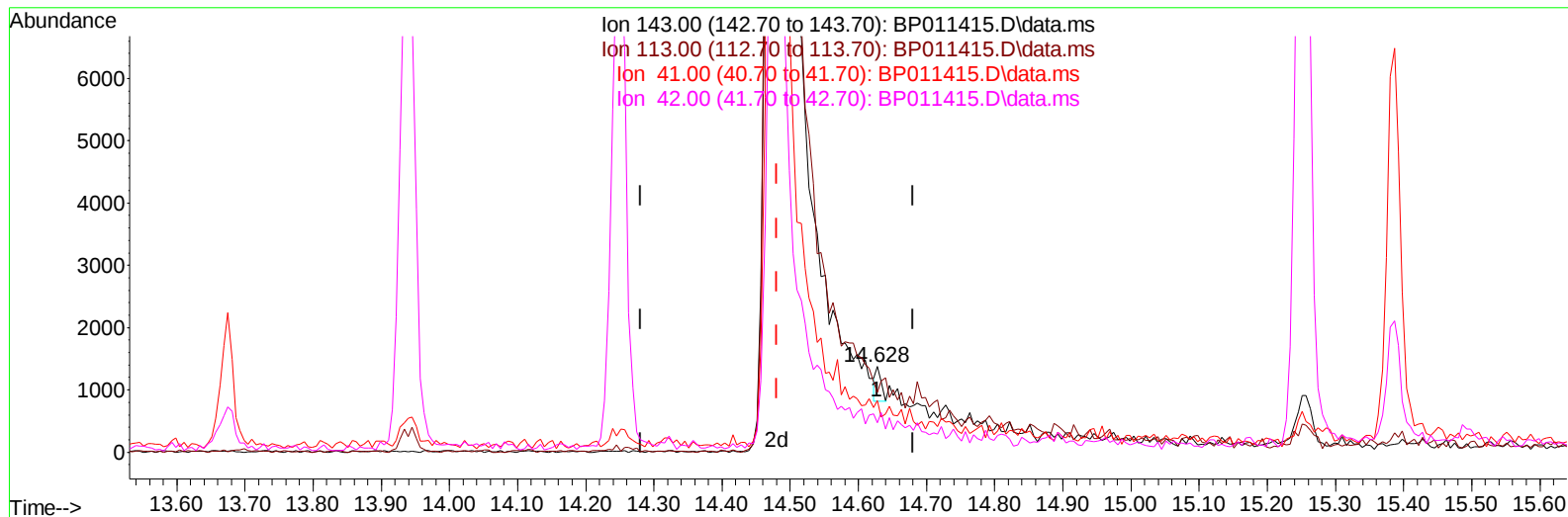


Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081122\
Data File : BP011415.D
Acq On : 12 Aug 2022 08:16
Operator : CG/JU
Sample : PB416917BL
Misc :
ALS Vial : 39 Sample Multiplier: 1

Quant Time: Aug 12 08:45:46 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Aug 11 23:05:53 2022
Response via : Initial Calibration



TIC: BP011415.D\data.ms

(54) 4-Nitrophenol-d4 (S)

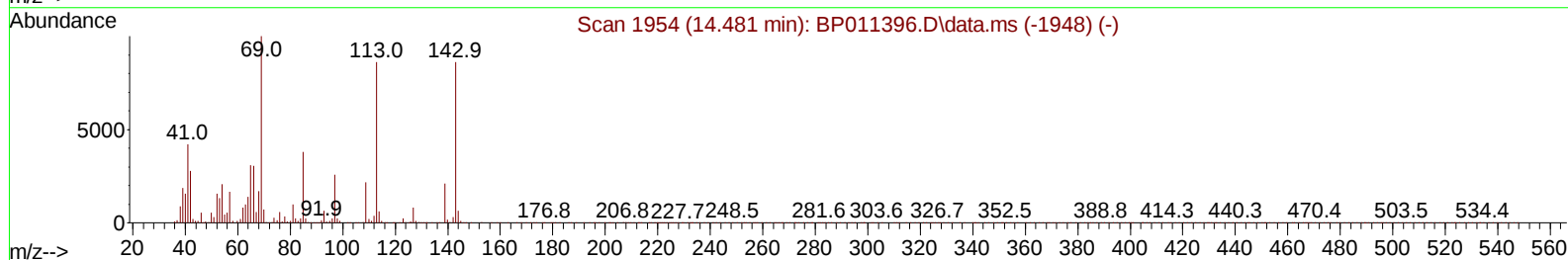
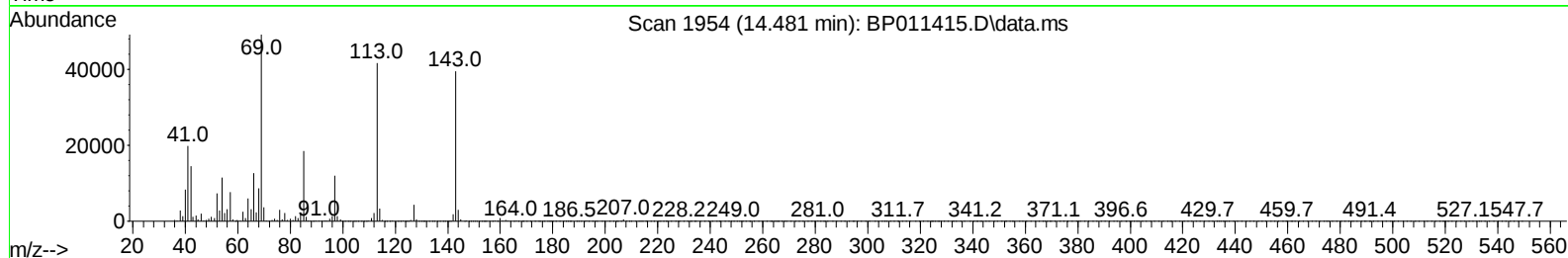
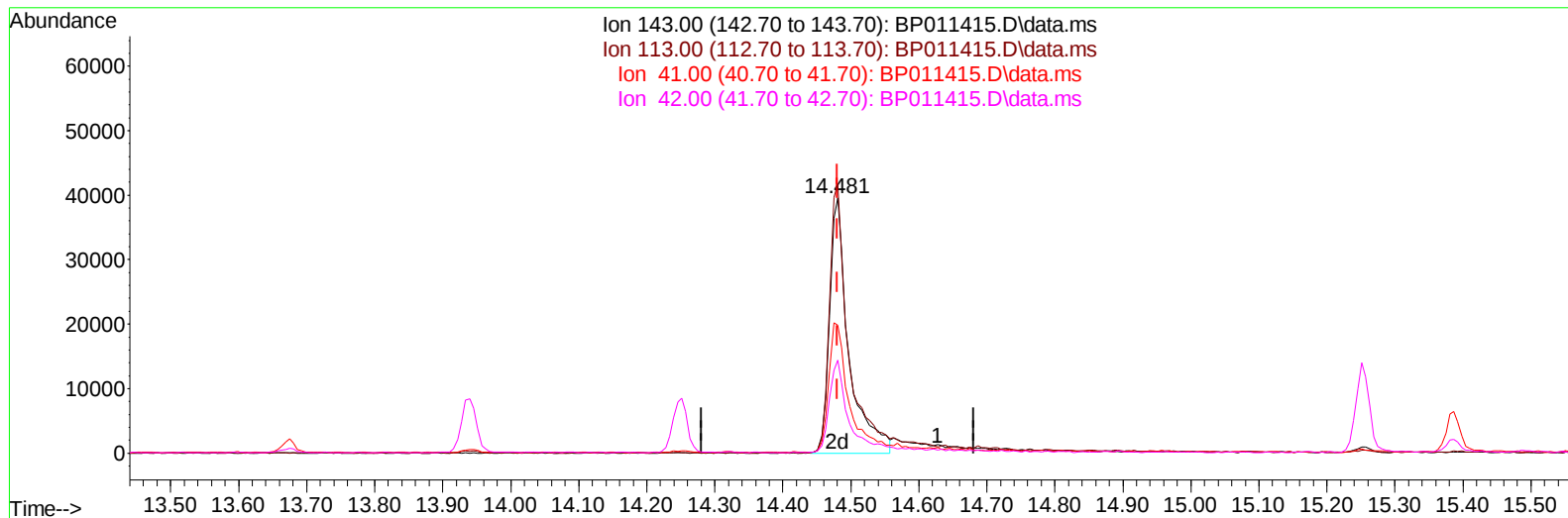
14.628min (+ 0.147) 0.10 ng/ul

response 307

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	69.62
41.00	54.40	60.33
42.00	35.50	34.45

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081122\
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Acq On : 12 Aug 2022 08:16
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Misc :
ALS Vial : 39 Sample Multiplier: 1

Quant Time: Aug 12 08:45:46 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION
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Response via : Initial Calibration



TIC: BP011415.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.481min (+ 0.000) 24.53 ng/ul m

response 78434

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	105.38#
41.00	54.40	49.97
42.00	35.50	36.52

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081122\
 Data File : BP011415.D
 Acq On : 12 Aug 2022 08:16
 Operator : CG/JU
 Sample : PB416917BL
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Quant Time: Aug 12 08:45:46 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Aug 11 23:05:53 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.575	152	101094	20.000	ng/ul	0.00
20) Naphthalene-d8	10.363	136	430663	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.251	164	252306	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.022	188	519975	20.000	ng/ul	0.00
79) Chrysene-d12	21.151	240	489101	20.000	ng/ul	0.00
88) Perylene-d12	23.457	264	502351	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.081	96	18235	5.530	ng/uL	0.00
4) Pyridine-d5	3.487	84	216750	26.057	ng/ul	0.00
7) Phenol-d5	6.758	99	245255	27.413	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.922	67	177787	28.026	ng/ul	0.00
11) 2-Chlorophenol-d4	7.105	132	208981	29.771	ng/ul	0.00
15) 4-Methylphenol-d8	8.293	113	201226	28.646	ng/ul	-0.01
21) Nitrobenzene-d5	8.740	128	98945	33.758	ng/ul	0.00
24) 2-Nitrophenol-d4	9.457	143	98824	36.827	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.993	165	158964	29.334	ng/ul	0.00
31) 4-Chloroaniline-d4	10.516	131	224994	23.748	ng/ul	0.00
46) Dimethylphthalate-d6	13.675	166	592216	33.829	ng/ul	0.00
49) Acenaphthylene-d8	13.940	160	656975	31.176	ng/ul	0.00
54) 4-Nitrophenol-d4	14.481	143	78434m	24.530	ng/ul	0.00
60) Fluorene-d10	15.251	176	466947	31.189	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.387	200	70375	29.138	ng/ul	0.00
73) Anthracene-d10	17.122	188	781634	33.444	ng/ul	0.00
81) Pyrene-d10	19.386	212	879018	32.621	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.310	264	839651	33.143	ng/ul	0.00

Target Compounds	Qvalue

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

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