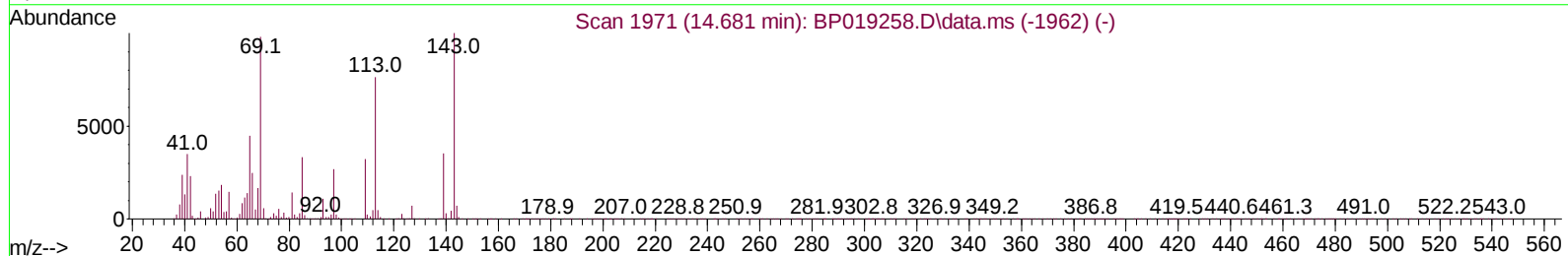
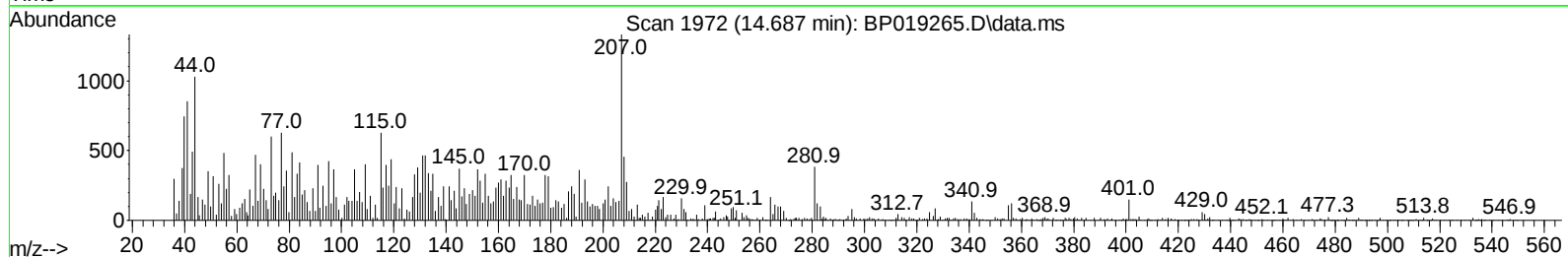
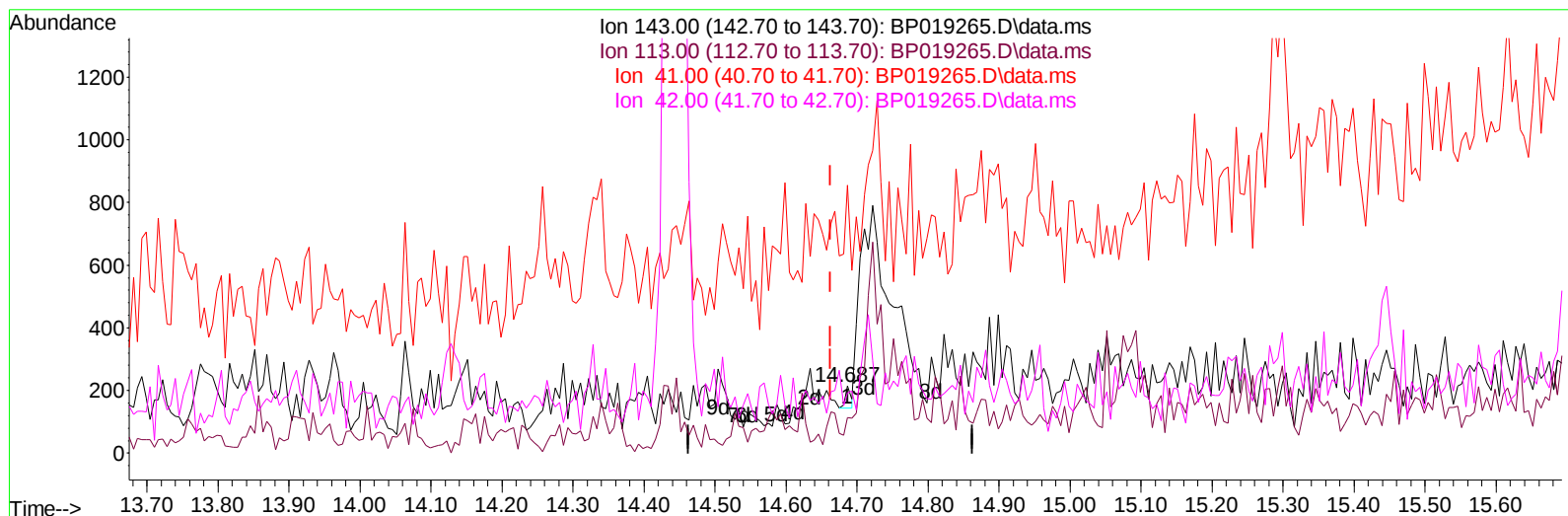


Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019265.D
Acq On : 04 Jan 2024 15:58
Operator : MA/JU
Sample : 05951-09DL 20X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Quant Time: Jan 04 16:39:44 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Dec 29 03:48:11 2023
Response via : Initial Calibration



TIC: BP019265.D\data.ms

(54) 4-Nitrophenol-d4 (S)

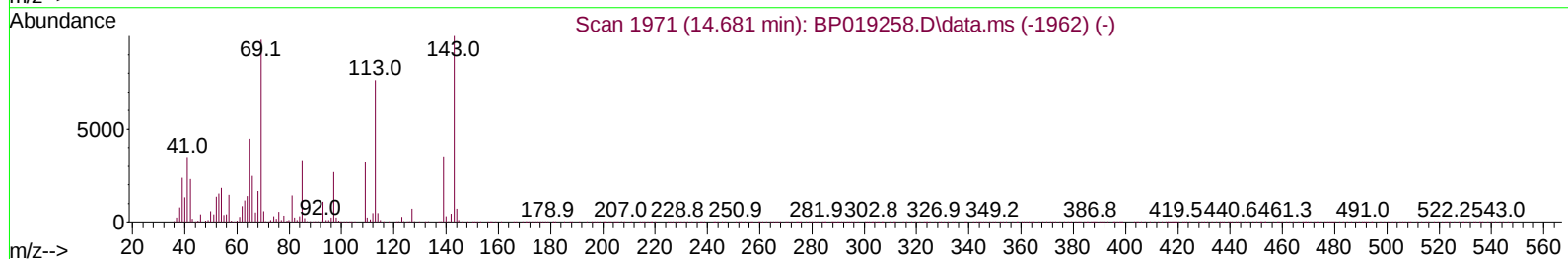
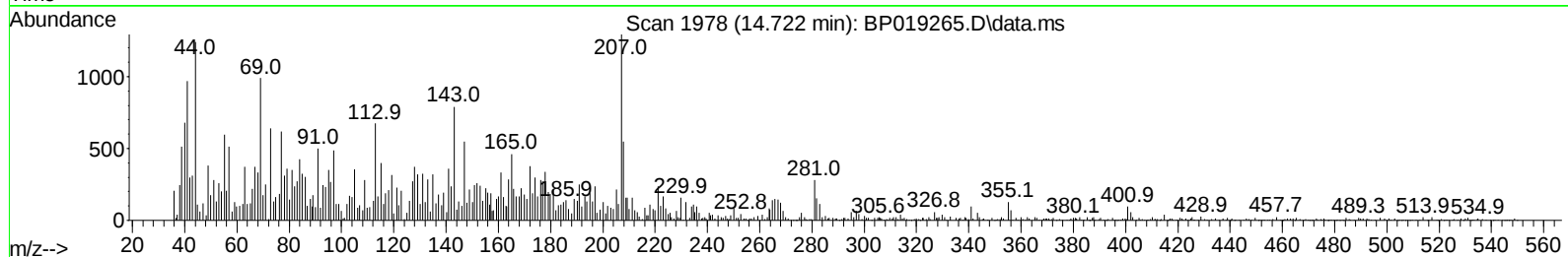
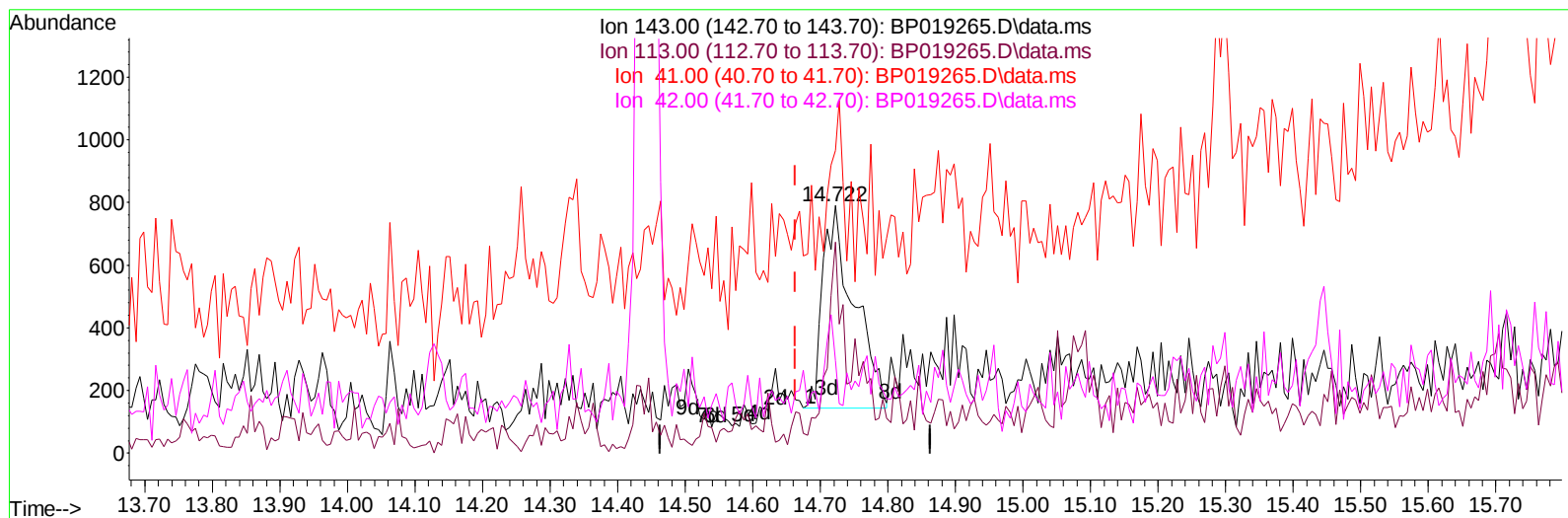
14.687min (+ 0.023) 0.01 ng/u1

response 38

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	73.50	62.62
41.00	33.40	399.53#
42.00	23.00	89.72#

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019265.D
Acq On : 04 Jan 2024 15:58
Operator : MA/JU
Sample : 05951-09DL 20X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Quant Time: Jan 04 16:39:44 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Dec 29 03:48:11 2023
Response via : Initial Calibration



TIC: BP019265.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.722min (+ 0.059) 0.48 ng/ul m

response 2081

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	73.50	85.19
41.00	33.40	122.41#
42.00	23.00	37.47#

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
 Data File : BP019265.D
 Acq On : 04 Jan 2024 15:58
 Operator : MA/JU
 Sample : 05951-09DL 20X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Quant Time: Jan 05 01:22:33 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 29 03:48:11 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.799	152	136013	20.000	ng/ul	0.00
20) Naphthalene-d8	10.593	136	485769	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.445	164	291174	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.198	188	665785	20.000	ng/ul	0.00
79) Chrysene-d12	21.304	240	727873	20.000	ng/ul	0.00
88) Perylene-d12	23.674	264	913189	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	592	0.149	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	6.993	99	7968	0.653	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.140	67	5434	0.770	ng/ul	0.00
11) 2-Chlorophenol-d4	7.334	132	6778	0.689	ng/ul	0.00
15) 4-Methylphenol-d8	8.534	113	5453	0.574	ng/ul	0.01
21) Nitrobenzene-d5	8.969	128	2902	0.691	ng/ul	0.00
24) 2-Nitrophenol-d4	9.687	143	3228	0.711	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.240	165	5731	0.625	ng/ul	0.01
31) 4-Chloroaniline-d4	0.000	131	0d	0.000	ng/ul	
46) Dimethylphthalate-d6	13.863	166	19919	0.774	ng/ul	0.00
49) Acenaphthylene-d8	14.134	160	24006	0.860	ng/ul	0.00
54) 4-Nitrophenol-d4	14.722	143	2081m	0.479	ng/ul	0.06
60) Fluorene-d10	15.439	176	19727	0.888	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.581	200	1888	0.433	ng/ul	0.00
73) Anthracene-d10	17.298	188	34079	0.980	ng/ul	0.00
81) Pyrene-d10	19.545	212	40908	0.922	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.516	264	45852	0.874	ng/ul	0.00
Target Compounds						
					Qvalue	
58) 2,3,4,6-Tetrachlorophenol	15.081	232	9283	1.399	ng/ul#	85
71) Pentachlorophenol	16.857	266	301669	50.808	ng/ul	98

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

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