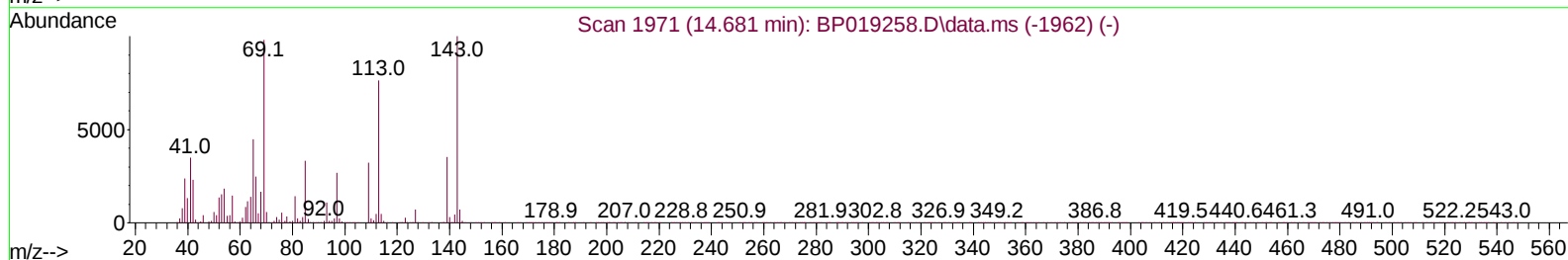
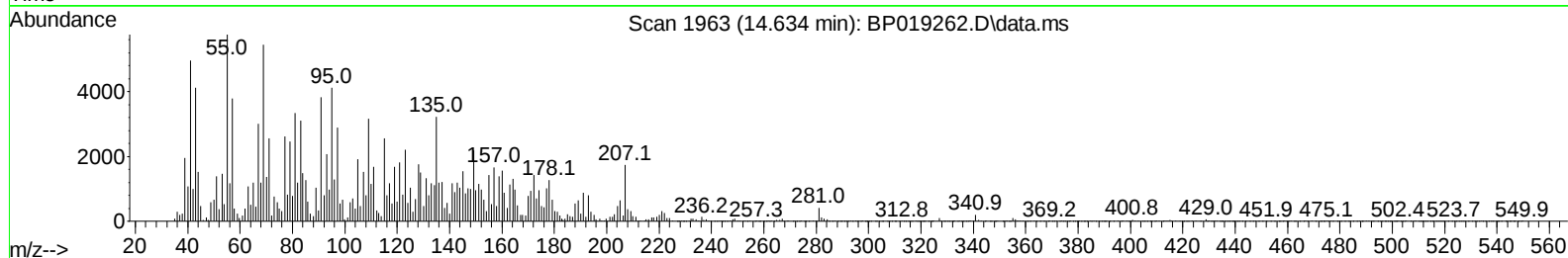
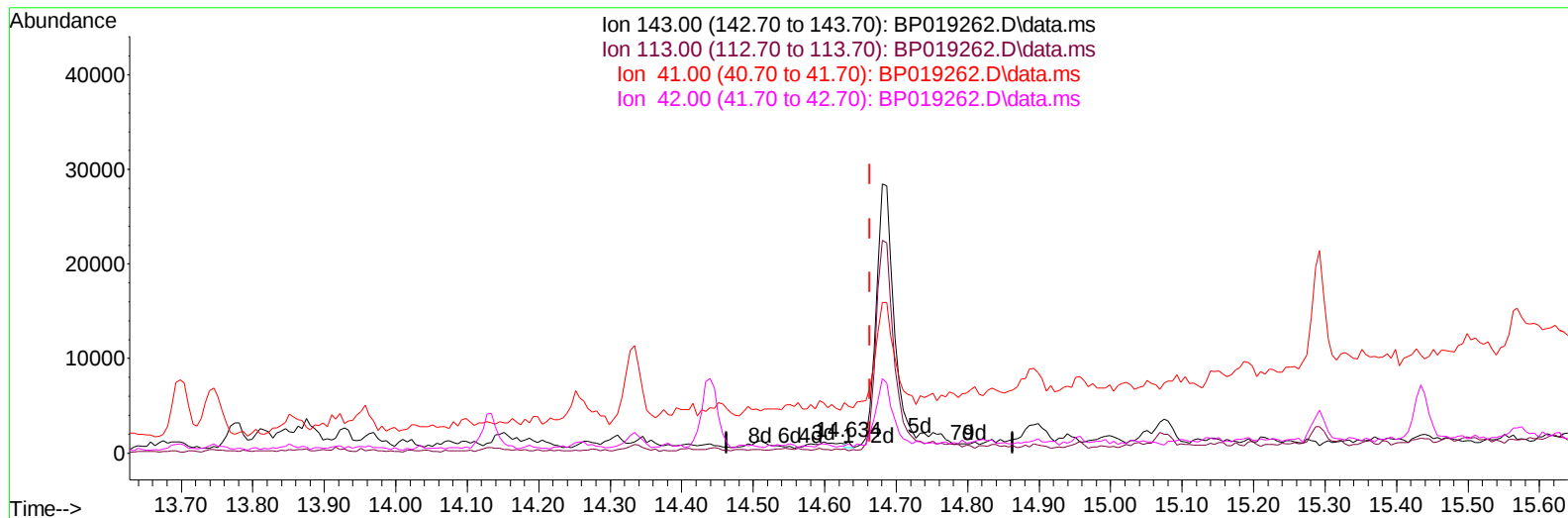


Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019262.D
Acq On : 04 Jan 2024 14:16
Operator : MA/JU
Sample : 05951-09
Misc :
ALS Vial : 6 Sample Multiplier: 1

Quant Time: Jan 04 16:58:07 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: BP019262.D\data.ms

(54) 4-Nitrophenol-d4 (S)

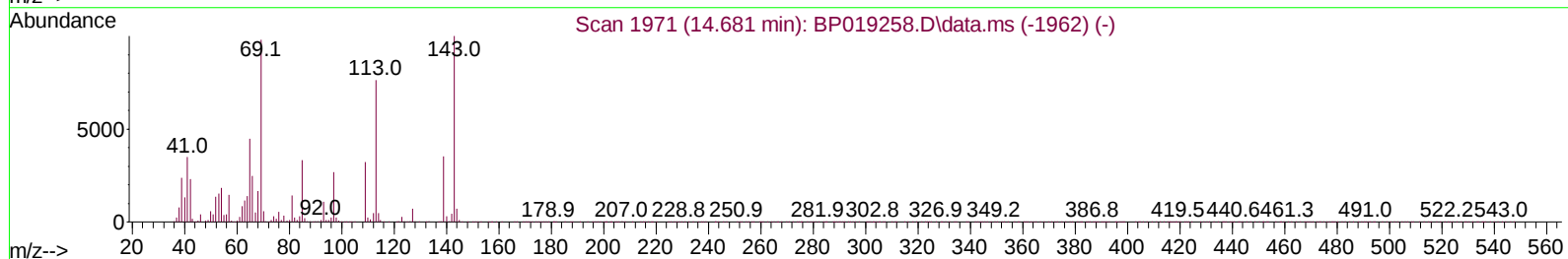
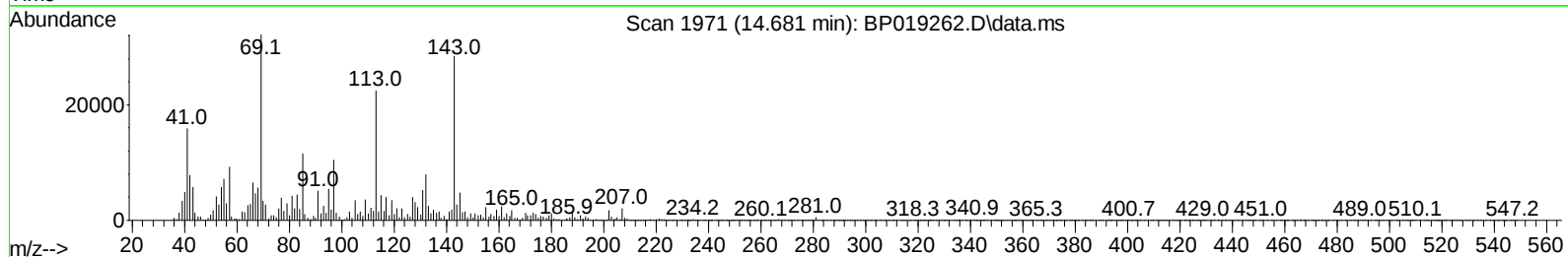
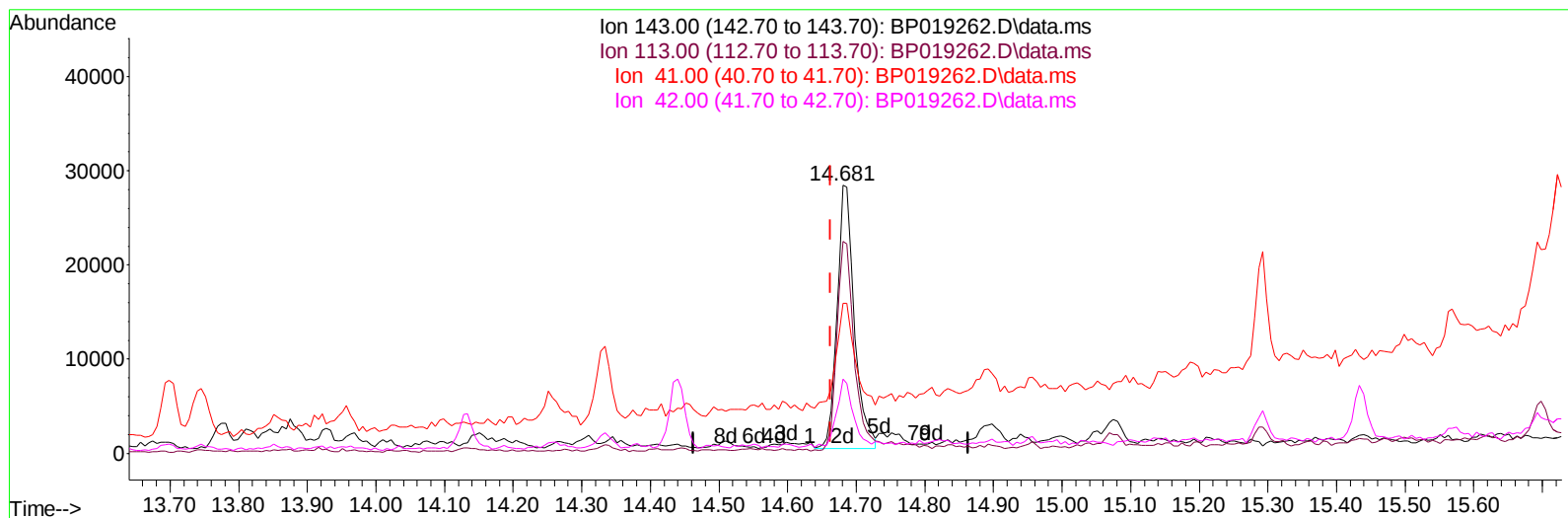
14.634min (-0.030) 0.04 ng/ul

response 210

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	73.50	21.93#
41.00	33.40	416.72#
42.00	23.00	84.20#

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
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Sample : 05951-09
Misc :
ALS Vial : 6 Sample Multiplier: 1

Quant Time: Jan 04 16:58:07 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: BP019262.D\data.ms

(54) 4-Nitrophenol-d4 (S)**14.681min (+ 0.017) 8.33 ng/ul m****response 49014**

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	73.50	79.16
41.00	33.40	55.92#
42.00	23.00	27.73#

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
 Data File : BP019262.D
 Acq On : 04 Jan 2024 14:16
 Operator : MA/JU
 Sample : 05951-09
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Quant Time: Jan 04 16:58:49 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	7.793	152	190564	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.593	136	666889	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.439	164	394315	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.204	188	831198	20.000	ng/ul	0.00
79) Chrysene-d12	21.304	240	995581	20.000	ng/ul	0.00
88) Perylene-d12	23.680	264	1240602	20.000	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	10358	1.856	ng/uL	0.00
4) Pyridine-d5	3.675	84	11429	0.816	ng/ul	0.02
7) Phenol-d5	6.987	99	178375	10.438	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.134	67	111340	11.262	ng/ul	0.00
11) 2-Chlorophenol-d4	7.334	132	147552	10.710	ng/ul	0.00
15) 4-Methylphenol-d8	8.522	113	136399	10.252	ng/ul	0.00
21) Nitrobenzene-d5	8.963	128	67793	11.760	ng/ul	0.00
24) 2-Nitrophenol-d4	9.681	143	76489	12.277	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.228	165	142247	11.307	ng/ul	0.00
31) 4-Chloroaniline-d4	10.751	131	21800	1.409	ng/ul	0.00
46) Dimethylphthalate-d6	13.857	166	418708	12.022	ng/ul	-0.01
49) Acenaphthylene-d8	14.134	160	482787	12.775	ng/ul	0.00
54) 4-Nitrophenol-d4	14.681	143	49014m	8.334	ng/ul	0.02
60) Fluorene-d10	15.434	176	381522	12.680	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.569	200	54926	10.080	ng/ul	0.00
73) Anthracene-d10	17.298	188	576605	13.283	ng/ul	0.00
81) Pyrene-d10	19.551	212	768514	12.666	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.521	264	900856	12.633	ng/ul	0.00
Target Compounds						
					Qvalue	
58) 2,3,4,6-Tetrachlorophenol	15.075	232	163571	18.198	ng/ul#	96
71) Pentachlorophenol	16.869	266	5482268	739.588	ng/ul	97
74) Anthracene	17.333	178	179683	3.583	ng/ul#	92
82) Pyrene	19.574	202	86058	1.180	ng/ul#	93
86) Bis(2-ethylhexyl)phtha...	21.216	149	64456	1.680	ng/ul#	83
87) Chrysene	21.321	228	89736	1.223	ng/ul#	78

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

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