

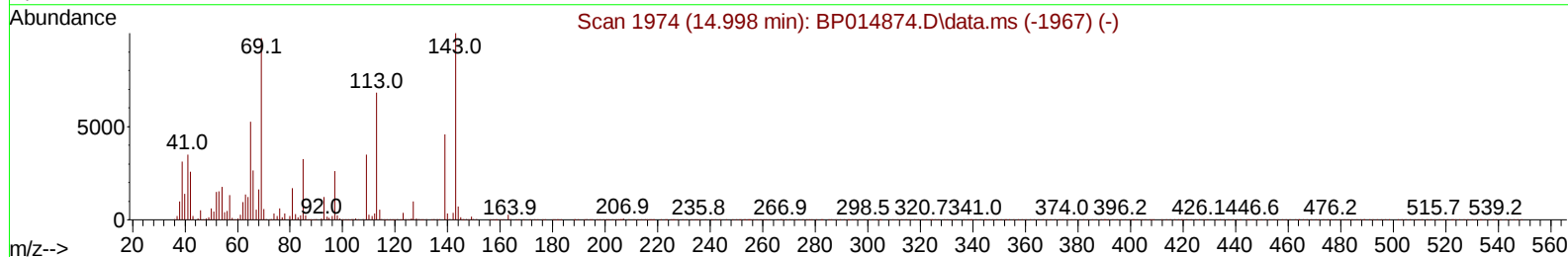
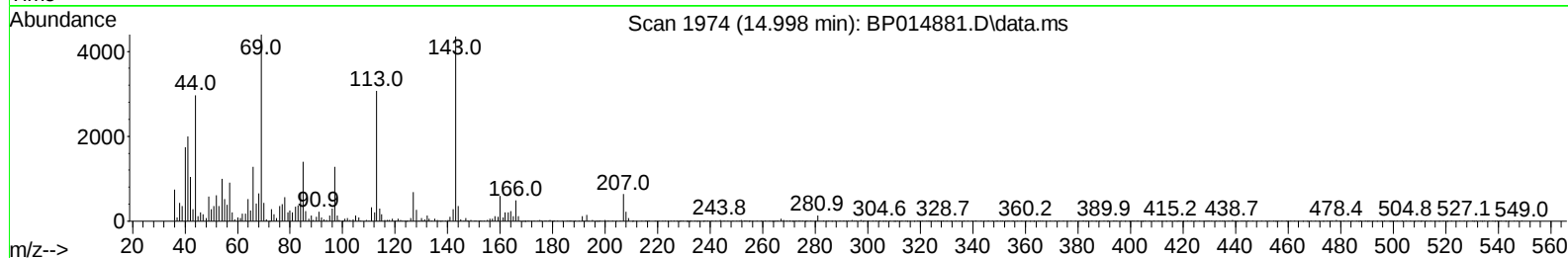
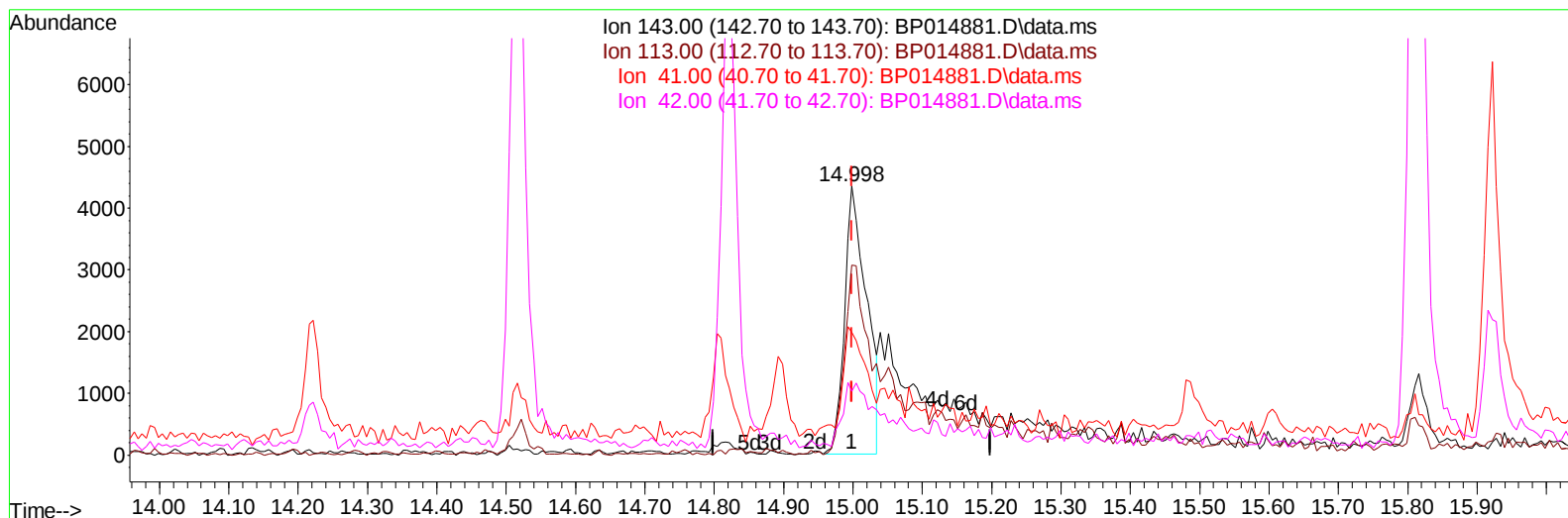
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042823\
Data File : BP014881.D
Acq On : 28 Apr 2023 18:49
Operator : CG/JU
Sample : 02488-15
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHE70

Manual IntegrationsAPPROVED

Quant Time: Apr 29 03:17:57 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042823.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Apr 29 00:40:33 2023
Response via : Initial Calibration

Reviewed By :Christian Giraldo 05/01/2023
Supervised By :Jagrut Upadhyay 05/01/2023



TIC: BP014881.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.998min (-0.000) 2.04 ng/u1

response 9508

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	68.20	70.51
41.00	36.10	45.67#
42.00	26.80	23.80

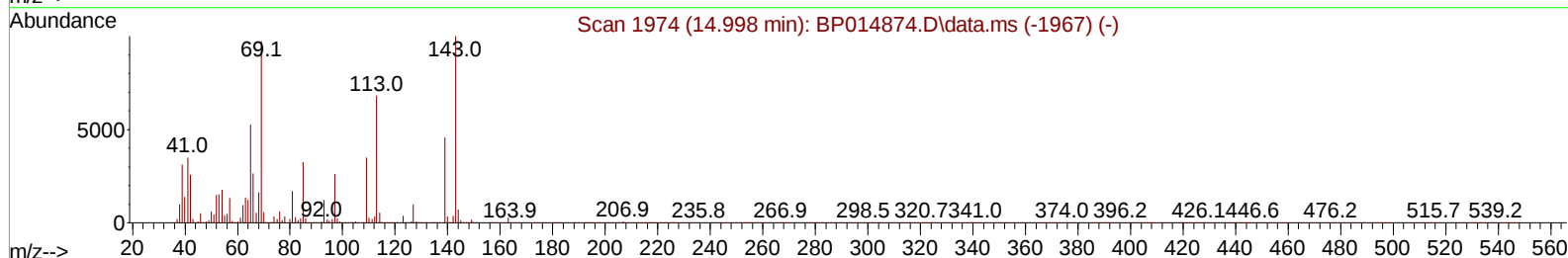
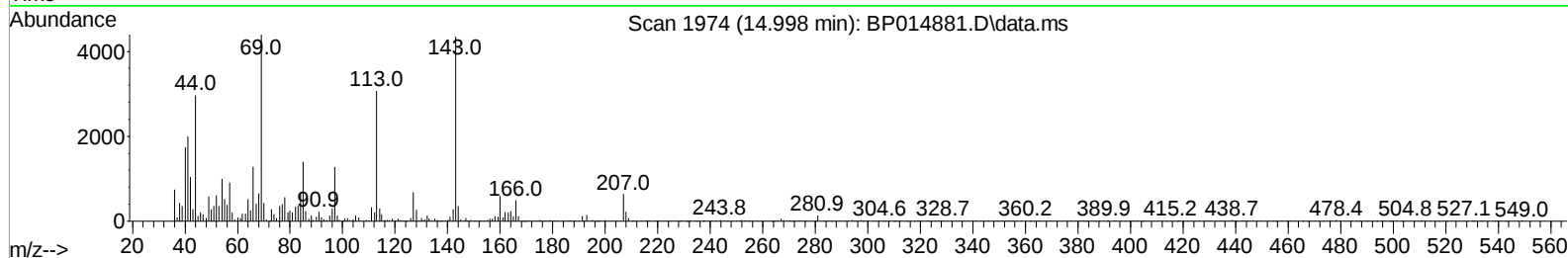
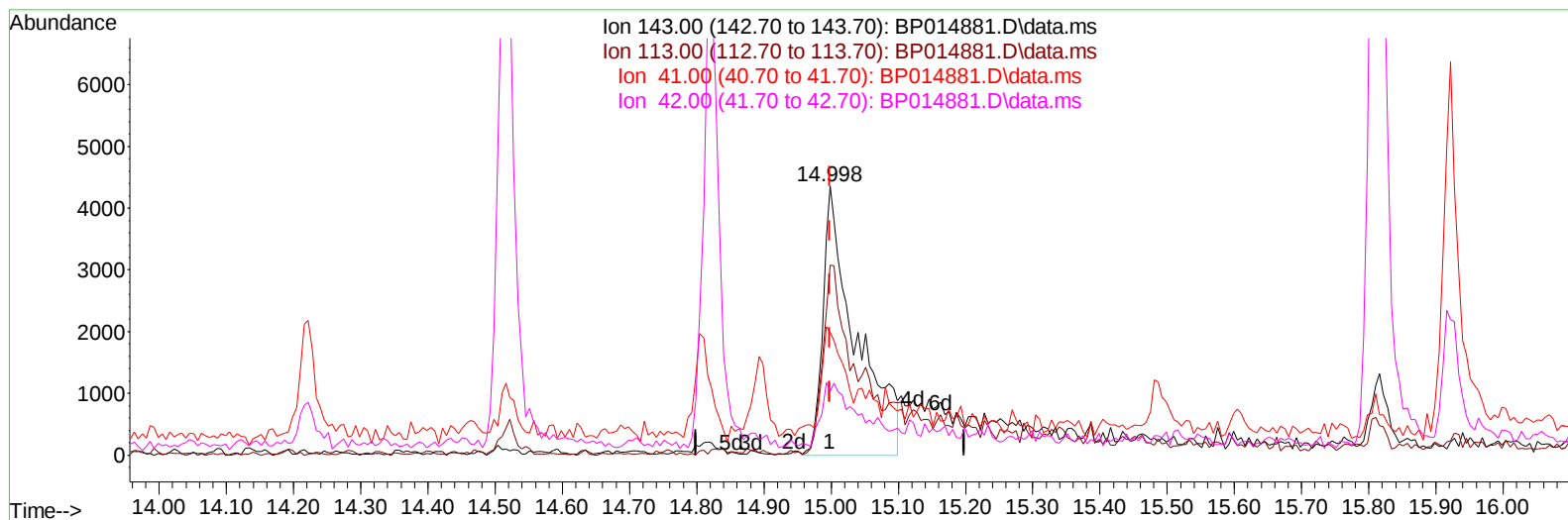
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TIC: BP014881.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.998min (-0.000) 3.18 ng/ul m

response 14788

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	68.20	70.51
41.00	36.10	45.67#
42.00	26.80	23.80

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Reviewed By :Christian Giraldo 05/01/2023
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Quant Time: Apr 29 03:25:39 2023
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 QLast Update : Sat Apr 29 00:40:33 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.175	152	175998	20.000	ng/ul	0.00
20) Naphthalene-d8	11.010	136	715297	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.822	164	436435	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.575	188	894550	20.000	ng/ul	0.00
79) Chrysene-d12	21.657	240	813689	20.000	ng/ul	0.00
88) Perylene-d12	24.263	264	935250	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.470	96	24906	5.100	ng/uL	0.00
4) Pyridine-d5	3.917	84	65548	5.584	ng/ul	0.01
7) Phenol-d5	7.316	99	106716	7.747	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.487	67	296319	34.298	ng/ul	0.00
11) 2-Chlorophenol-d4	7.699	132	287318	27.548	ng/ul	0.00
15) 4-Methylphenol-d8	8.881	113	183280	17.461	ng/ul	0.00
21) Nitrobenzene-d5	9.352	128	170795	33.255	ng/ul	0.00
24) 2-Nitrophenol-d4	10.081	143	144243	30.595	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.622	165	281407	27.791	ng/ul	0.00
31) 4-Chloroaniline-d4	11.146	131	269990	18.442	ng/ul	0.00
46) Dimethylphthalate-d6	14.222	166	1136604	36.339	ng/ul	0.00
49) Acenaphthylene-d8	14.516	160	1331213	35.631	ng/ul	0.00
54) 4-Nitrophenol-d4	14.998	143	14788m	3.177	ng/ul	0.00
60) Fluorene-d10	15.816	176	1057806	38.798	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.922	200	116705	30.843	ng/ul	0.00
73) Anthracene-d10	17.675	188	1766802	43.247	ng/ul	0.00
81) Pyrene-d10	19.892	212	2044316	45.235	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.092	264	2006041	43.427	ng/ul	0.00

Target Compounds Qvalue

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

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