

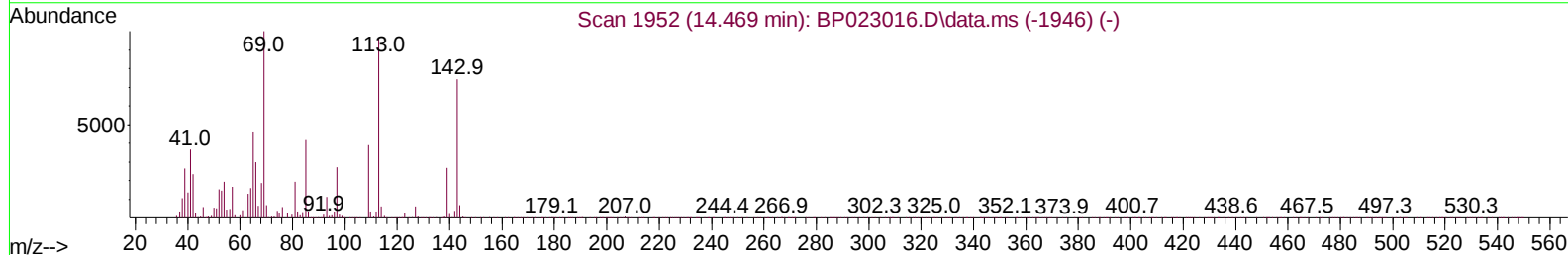
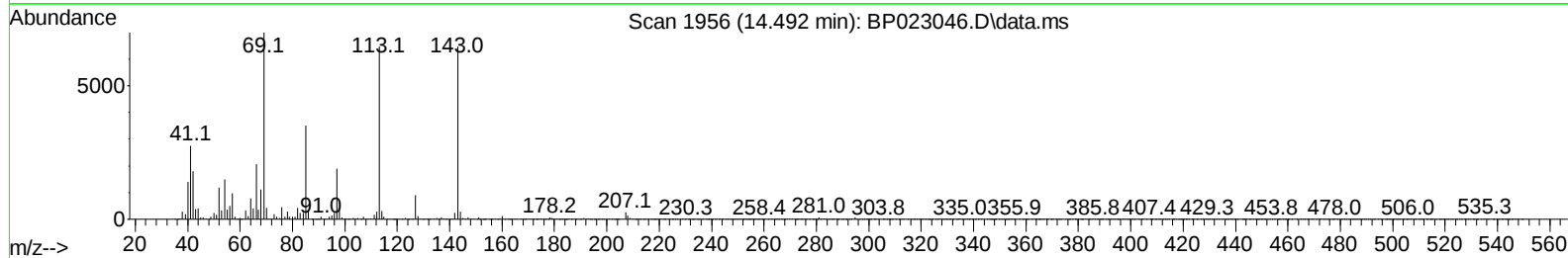
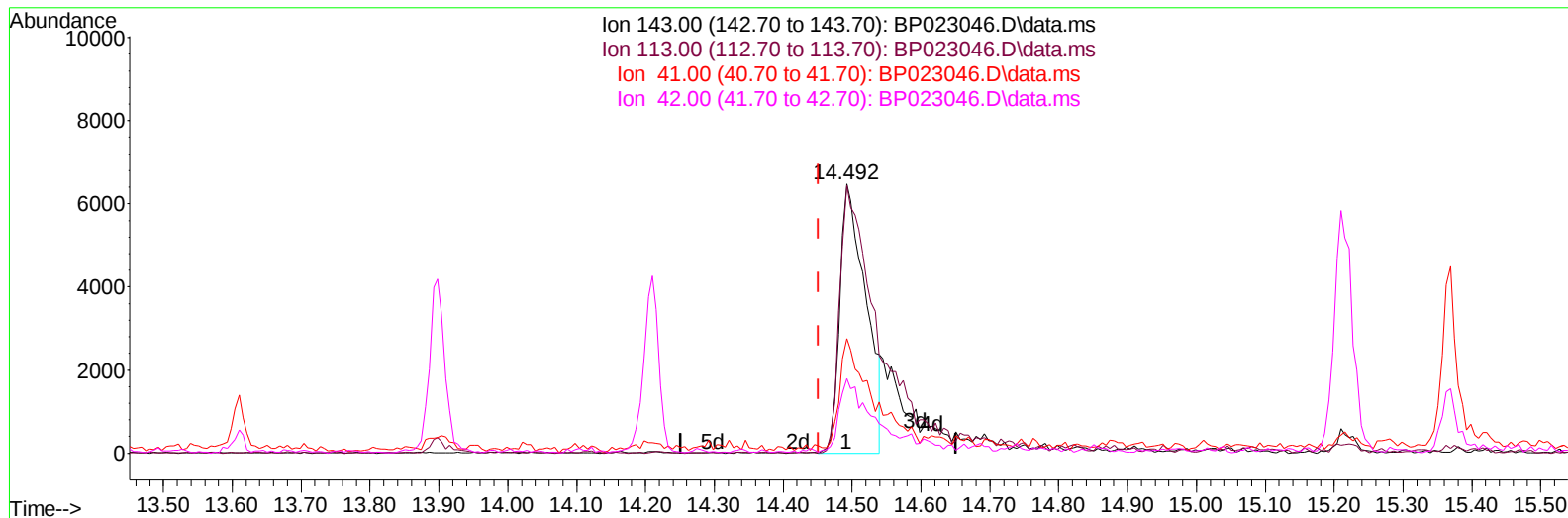
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111224\
Data File : BP023046.D
Acq On : 12 Nov 2024 14:33
Operator : RC/JU
Sample : P4741-07
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E27L4

Manual IntegrationsAPPROVED

Quant Time: Nov 13 00:02:02 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Nov 12 23:59:19 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/14/2024
Supervised By :mohammad ahmed 11/14/2024



TIC: BP023046.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.492min (+ 0.041) 7.99 ng/u1

response 16984

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	99.68
41.00	39.20	42.52
42.00	25.90	27.68

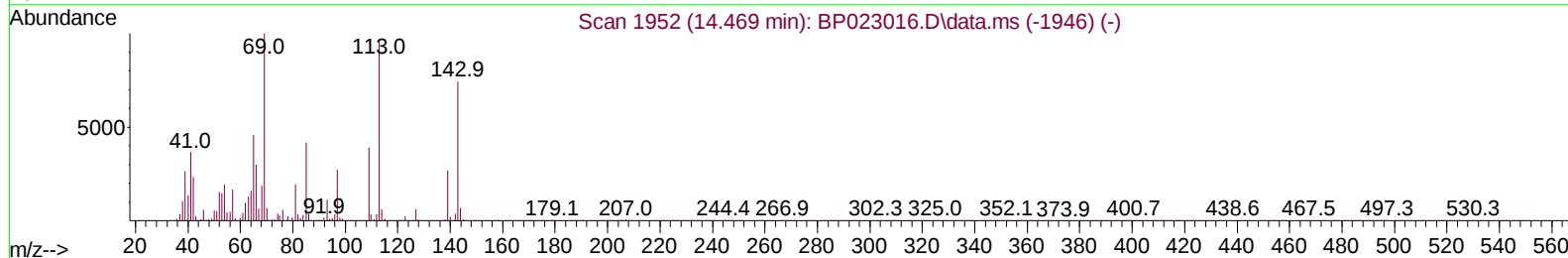
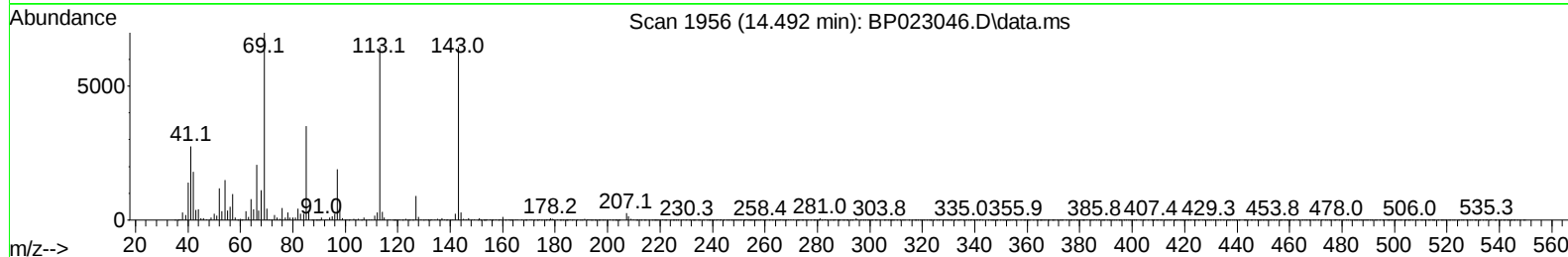
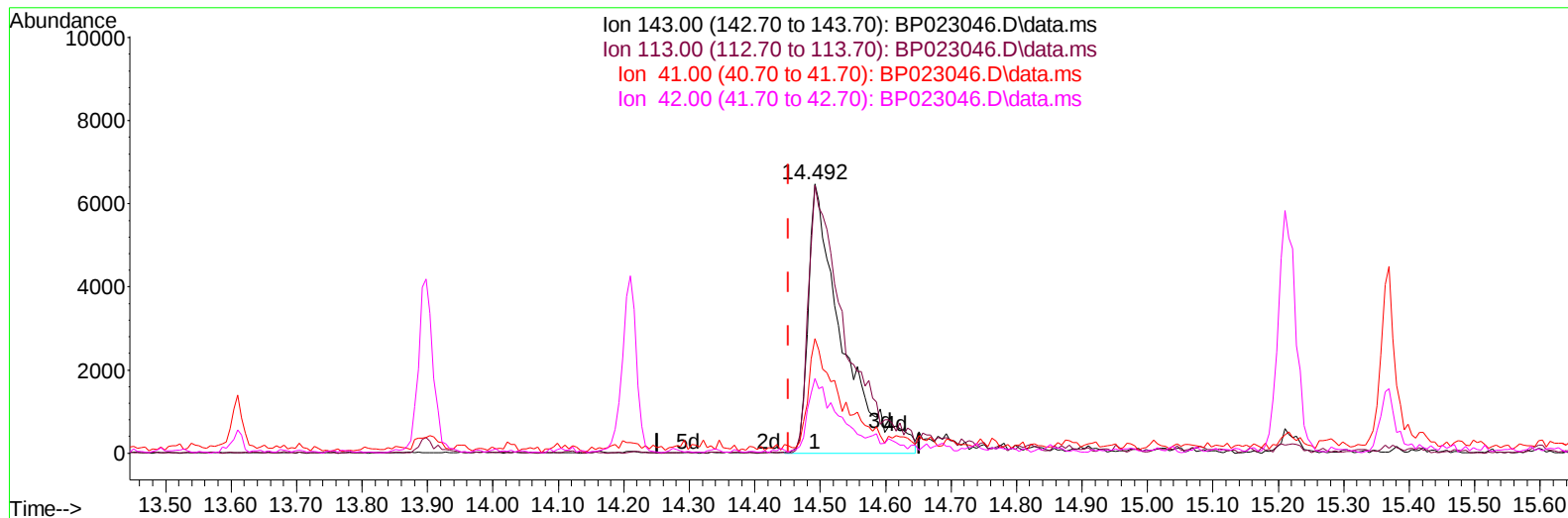
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TIC: BP023046.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.492min (+ 0.041) 10.93 ng/ul m

response 23234

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	99.68
41.00	39.20	42.52
42.00	25.90	27.68

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/14/2024
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Quant Time: Nov 13 01:45:52 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 12 23:59:19 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.587	152	61982	20.000	ng/ul	0.00
20) Naphthalene-d8	10.339	136	240640	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.210	164	196188	20.000	ng/ul	0.01
64) Phenanthrene-d10	17.010	188	502118	20.000	ng/ul	0.00
79) Chrysene-d12	21.439	240	600413	20.000	ng/ul	-0.01
88) Perylene-d12	24.650	264	725660	20.000	ng/ul	-0.04
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.140	96	7322	5.678	ng/uL	0.00
4) Pyridine-d5	3.557	84	50003	13.052	ng/ul	0.00
7) Phenol-d5	6.793	99	72038	15.720	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.934	67	91315	33.917	ng/ul	0.00
11) 2-Chlorophenol-d4	7.134	132	109762	32.705	ng/ul	0.00
15) 4-Methylphenol-d8	8.298	113	103788	27.523	ng/ul	0.00
21) Nitrobenzene-d5	8.728	128	57473	34.307	ng/ul	0.00
24) 2-Nitrophenol-d4	9.439	143	69558	34.827	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.987	165	142405	34.047	ng/ul	0.00
31) 4-Chloroaniline-d4	10.486	131	106353	21.314	ng/ul	0.00
46) Dimethylphthalate-d6	13.610	166	524304	37.080	ng/ul	0.00
49) Acenaphthylene-d8	13.898	160	495917	33.613	ng/ul	0.00
54) 4-Nitrophenol-d4	14.492	143	23234m	10.934	ng/ul	0.04
60) Fluorene-d10	15.216	176	442448	35.690	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.369	200	92547	30.768	ng/ul	0.00
73) Anthracene-d10	17.116	188	804088	37.821	ng/ul	0.00
81) Pyrene-d10	19.492	212	1050882	37.999	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.439	264	1361571	39.390	ng/ul	-0.02

Target Compounds Qvalue

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

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