

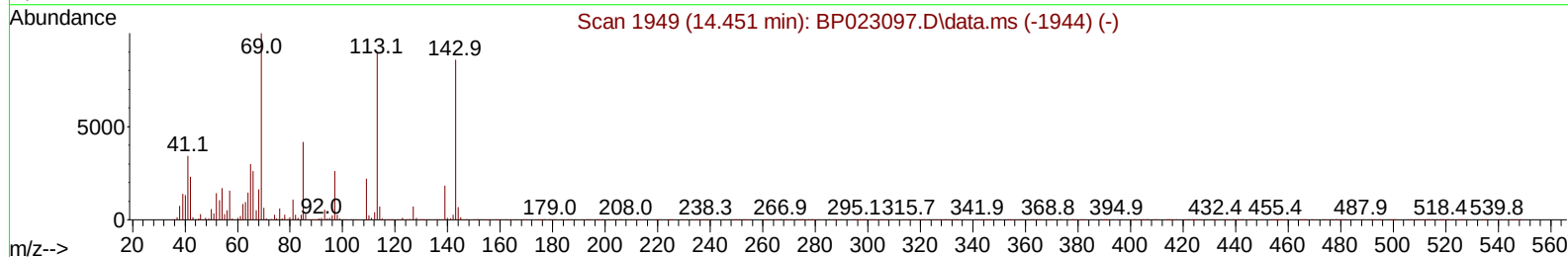
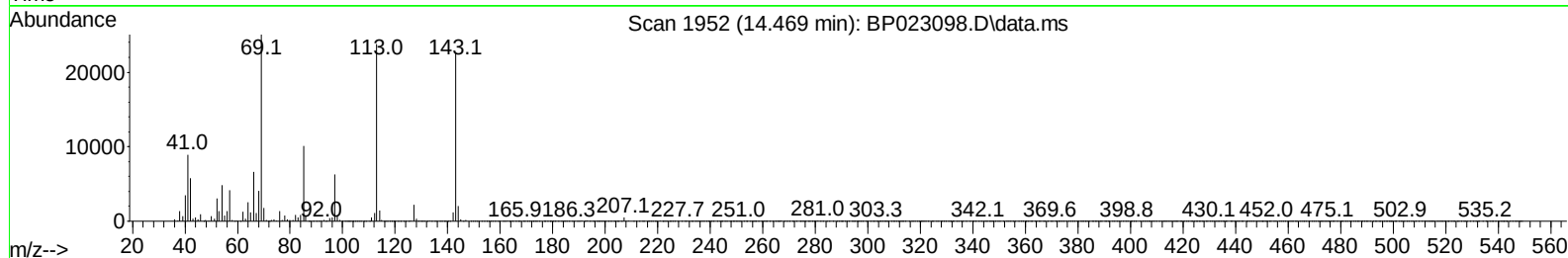
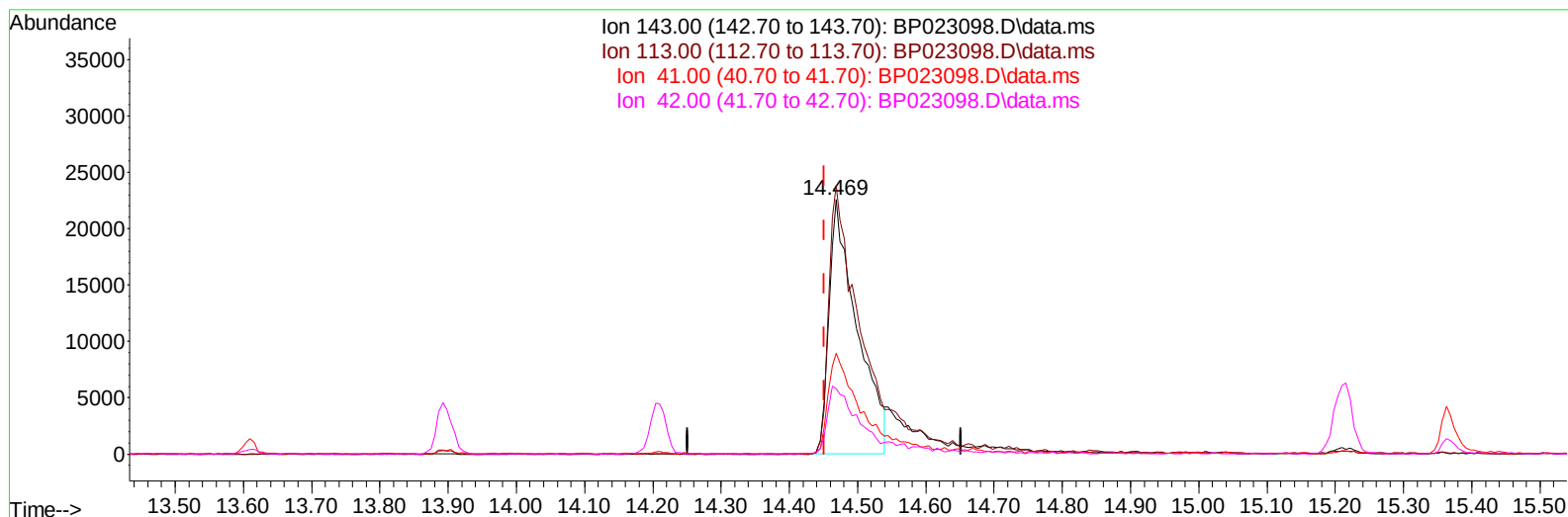
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111324\
Data File : BP023098.D
Acq On : 14 Nov 2024 02:58
Operator : RC/JU
Sample : PB164918BL
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK918

Manual IntegrationsAPPROVED

Quant Time: Nov 14 03:24:57 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 13 04:49:01 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/15/2024
Supervised By :mohammad ahmed 11/18/2024



TIC: BP023098.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.469min (+ 0.017) 23.73 ng/ul

response 64255

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	105.42
41.00	39.20	39.58
42.00	25.90	25.71

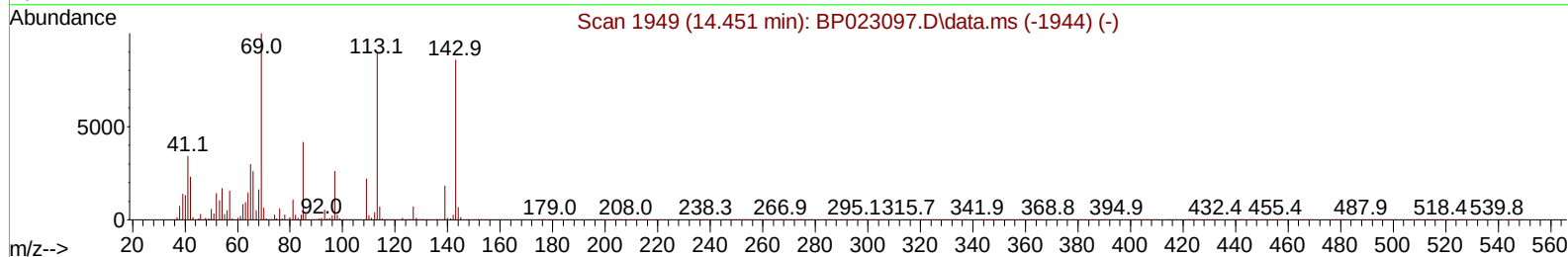
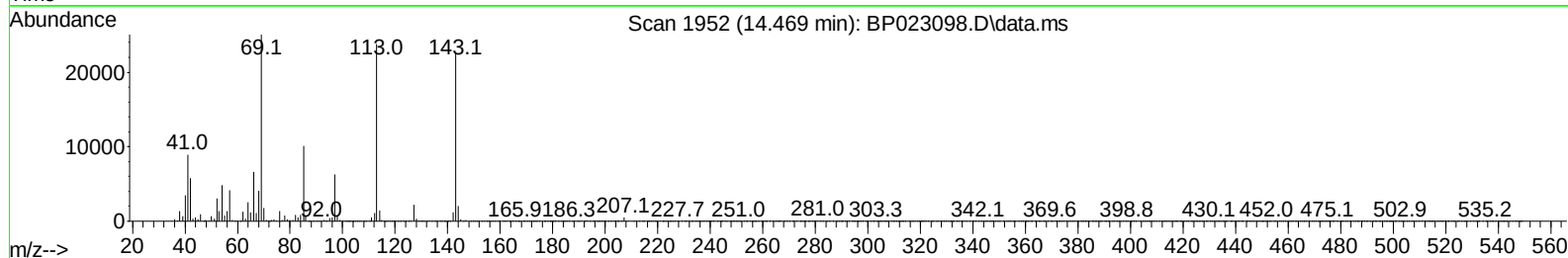
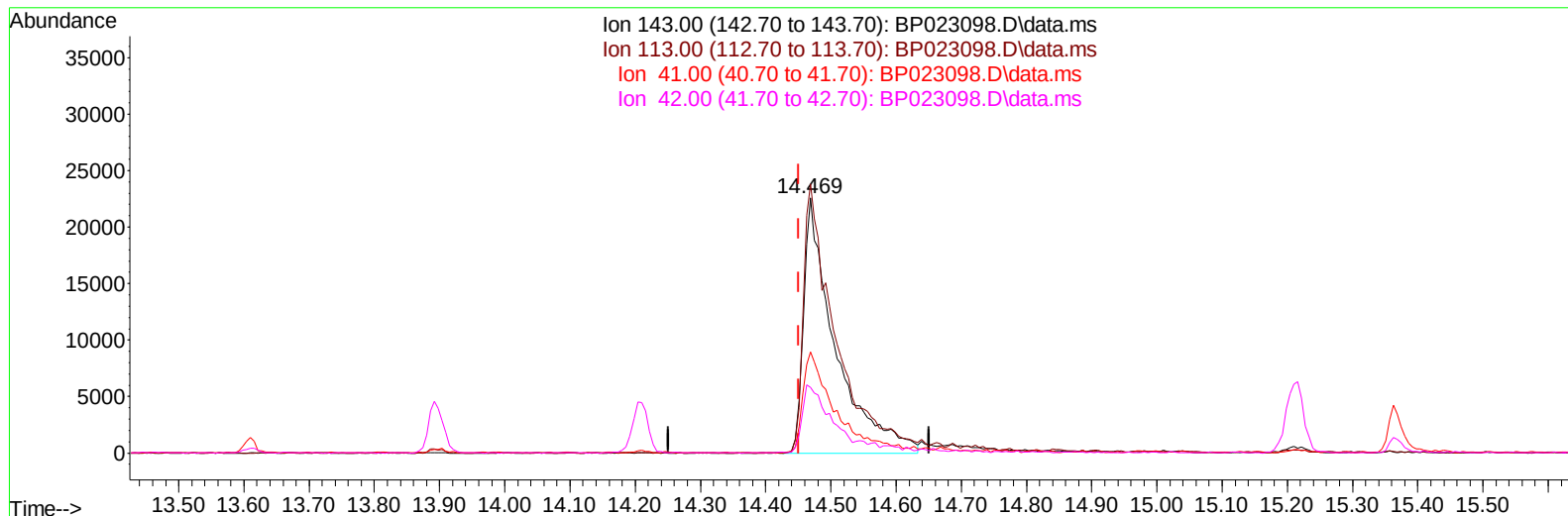
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TIC: BP023098.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.469min (+ 0.017) 28.12 ng/ul m

response 76138

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	105.42
41.00	39.20	39.58
42.00	25.90	25.71

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/15/2024
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Quant Time: Nov 14 03:26:07 2024
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 QLast Update : Wed Nov 13 04:49:01 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.581	152	82774	20.000	ng/ul	0.00
20) Naphthalene-d8	10.333	136	316422	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.210	164	249960	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.004	188	621630	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.445	240	699402	20.000	ng/ul	0.00
88) Perylene-d12	24.656	264	803953	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.140	96	11402	6.621	ng/uL	0.00
4) Pyridine-d5	3.551	84	143321	28.013	ng/ul	0.00
7) Phenol-d5	6.787	99	181281	29.623	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.928	67	110317	30.682	ng/ul	0.00
11) 2-Chlorophenol-d4	7.128	132	139262	31.072	ng/ul	0.00
15) 4-Methylphenol-d8	8.292	113	148440	29.476	ng/ul	0.00
21) Nitrobenzene-d5	8.722	128	72349	32.843	ng/ul	0.00
24) 2-Nitrophenol-d4	9.434	143	84030	31.997	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.981	165	167953	30.539	ng/ul	0.01
31) 4-Chloroaniline-d4	10.475	131	189483	28.879	ng/ul	0.00
46) Dimethylphthalate-d6	13.610	166	565696	31.401	ng/ul	0.00
49) Acenaphthylene-d8	13.898	160	611319	32.521	ng/ul	0.00
54) 4-Nitrophenol-d4	14.469	143	76138m	28.122	ng/ul	0.02
60) Fluorene-d10	15.210	176	510420	32.315	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.368	200	97349	26.142	ng/ul	0.00
73) Anthracene-d10	17.110	188	884864	33.619	ng/ul	0.00
81) Pyrene-d10	19.486	212	1149195	35.673	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.444	264	1297488	33.881	ng/ul	0.00

Target Compounds Qvalue

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

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