

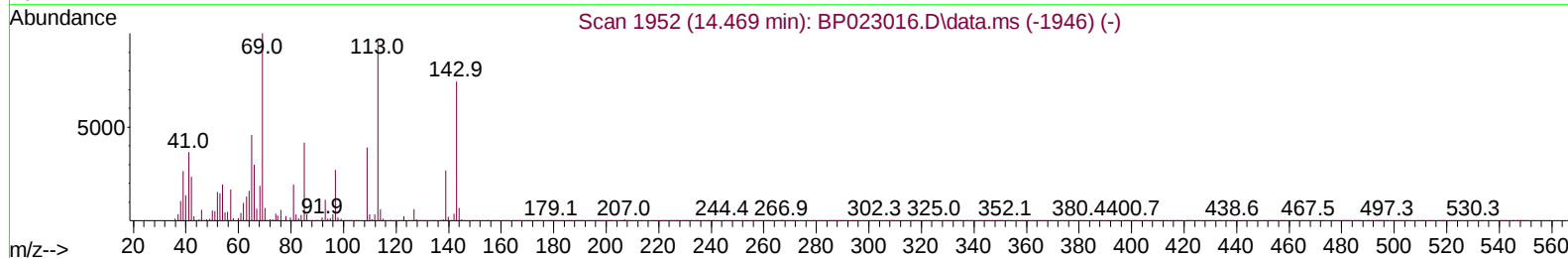
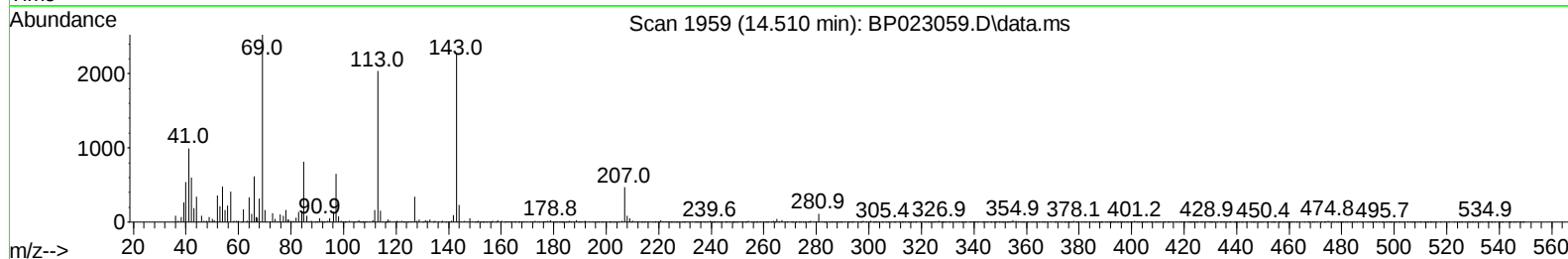
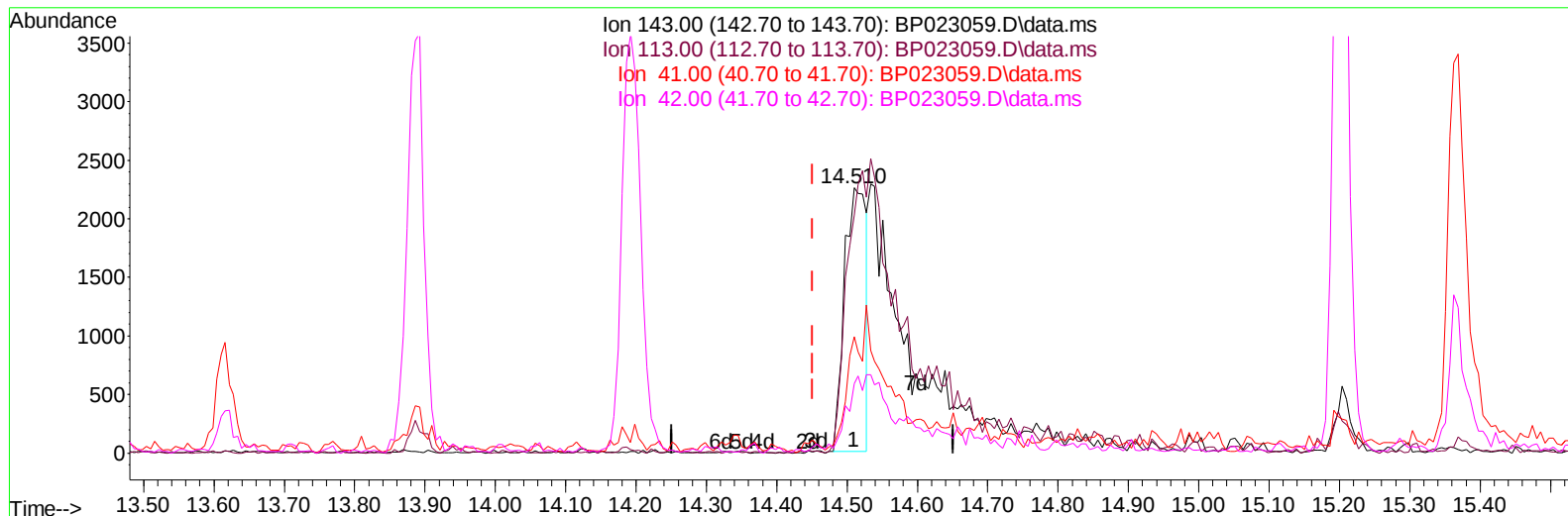
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111224\
Data File : BP023059.D
Acq On : 12 Nov 2024 23:20
Operator : RC/JU
Sample : P4687-08
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0CC5

Manual IntegrationsAPPROVED

Quant Time: Nov 13 03:16:36 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: BP023059.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.510min (+ 0.058) 2.09 ng/u1

response 4808

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	89.76
41.00	39.20	43.89
42.00	25.90	26.27

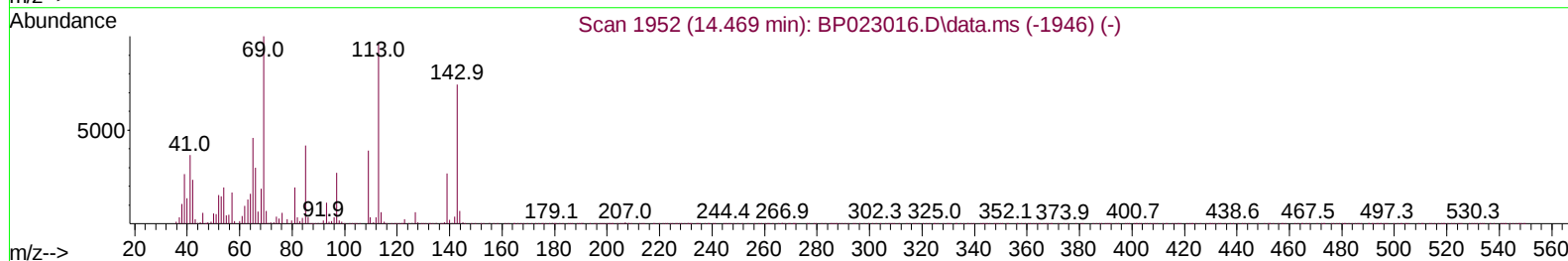
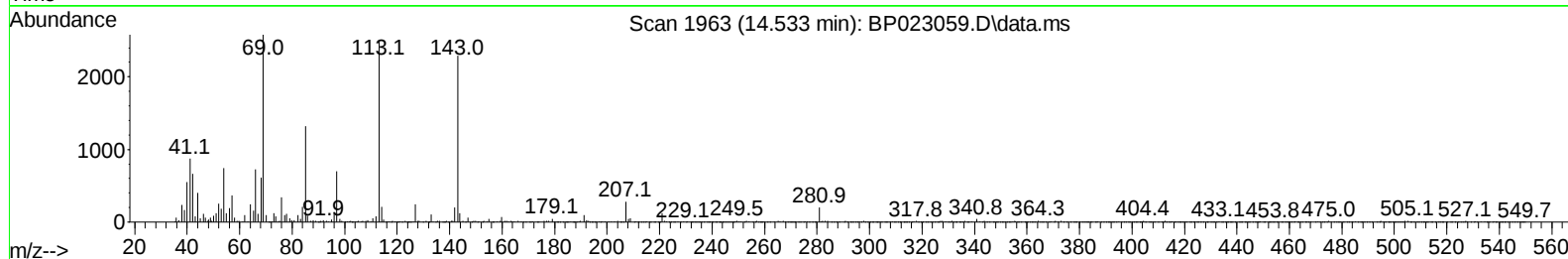
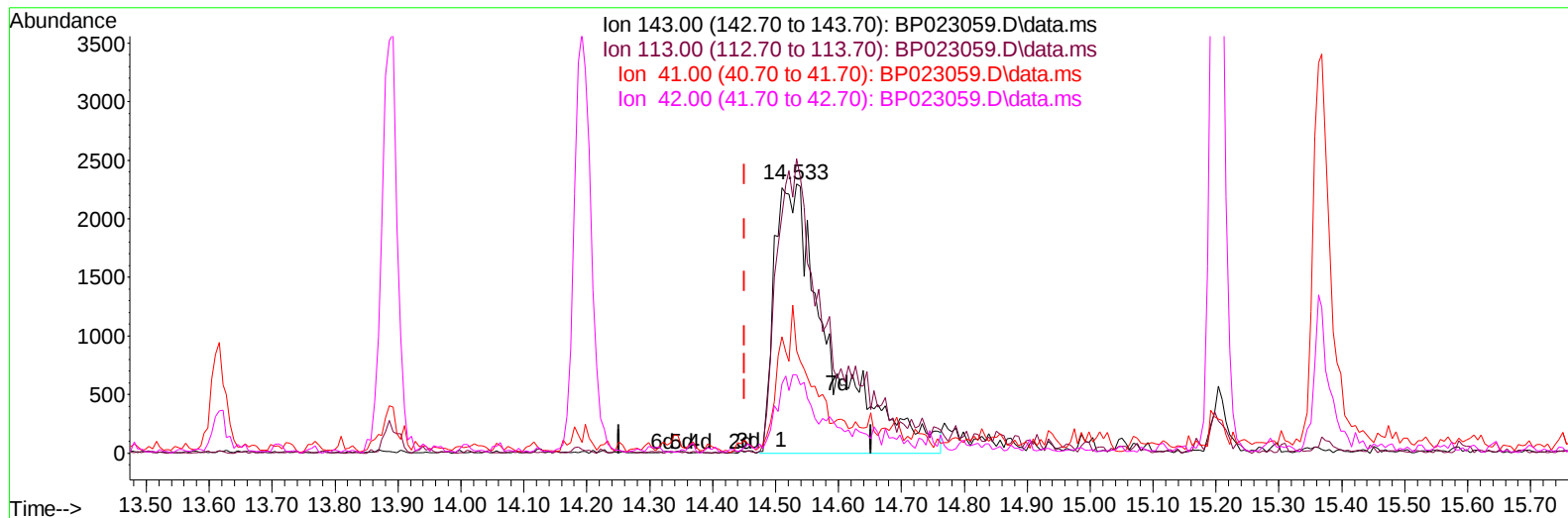
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(54) 4-Nitrophenol-d4 (S)

14.533min (+ 0.082) 6.11 ng/ul m

response 14053

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	109.41
41.00	39.20	38.02
42.00	25.90	29.09

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Instrument :
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Manual IntegrationsAPPROVED

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

Quant Time: Nov 13 03:18:14 2024
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 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 12 23:59:19 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.592	152	66523	20.000	ng/ul	0.00
20) Naphthalene-d8	10.334	136	262126	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.192	164	212470	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.992	188	540081	20.000	ng/ul	-0.02
79) Chrysene-d12	21.439	240	627835	20.000	ng/ul	-0.01
88) Perylene-d12	24.650	264	769283	20.000	ng/ul	-0.04
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.140	96	6090	4.400	ng/uL	0.00
4) Pyridine-d5	3.575	84	13753	3.345	ng/ul	0.02
7) Phenol-d5	6.798	99	42794	8.701	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	6.940	67	85869	29.717	ng/ul	0.00
11) 2-Chlorophenol-d4	7.140	132	93880	26.063	ng/ul	0.01
15) 4-Methylphenol-d8	8.298	113	82726	20.440	ng/ul	0.00
21) Nitrobenzene-d5	8.734	128	55143	30.218	ng/ul	0.00
24) 2-Nitrophenol-d4	9.445	143	64433	29.617	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.986	165	126070	27.671	ng/ul	0.00
31) 4-Chloroaniline-d4	10.498	131	83724	15.404	ng/ul	0.01
46) Dimethylphthalate-d6	13.616	166	508198	33.187	ng/ul	0.00
49) Acenaphthylene-d8	13.886	160	457587	28.638	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.533	143	14053m	6.107	ng/ul	0.08
60) Fluorene-d10	15.204	176	431502	32.139	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.369	200	99082	30.625	ng/ul	0.00
73) Anthracene-d10	17.104	188	839092	36.694	ng/ul	-0.02
81) Pyrene-d10	19.486	212	1125635	38.924	ng/ul	-0.01
92) Benzo(a)pyrene-d12	24.427	264	1430422	39.036	ng/ul	-0.04
Target Compounds						
					Qvalue	
8) Phenol	6.828	94	9313	1.813	ng/ul	93

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

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