

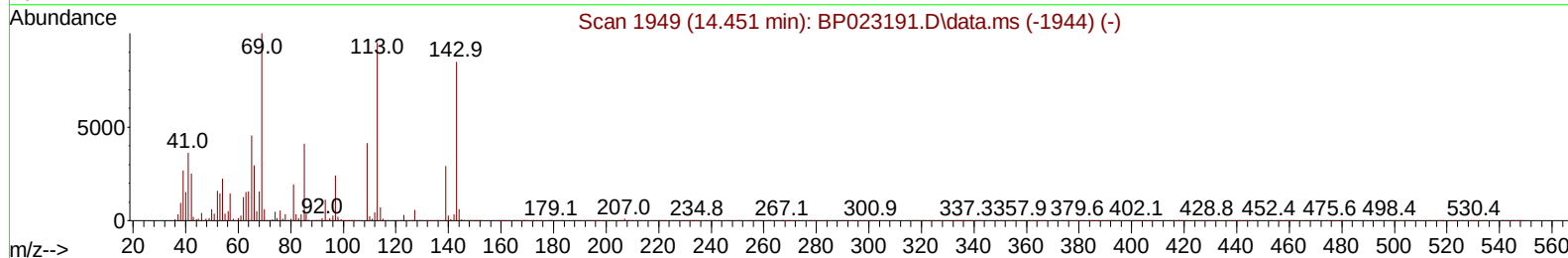
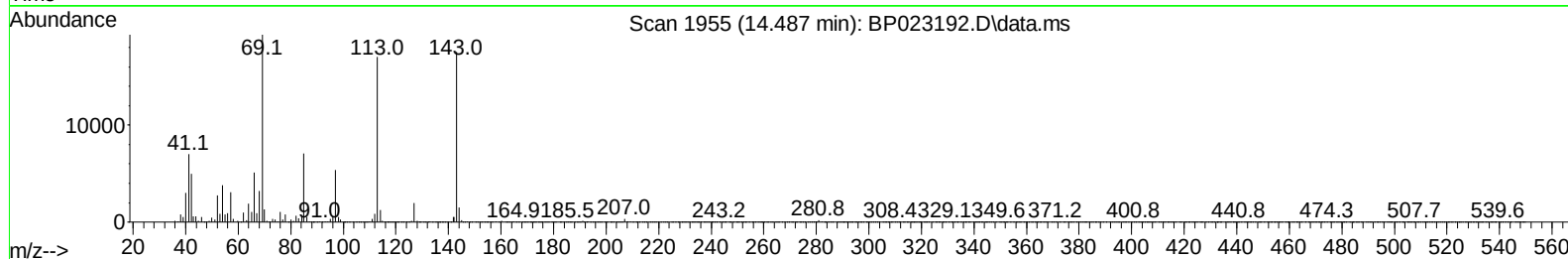
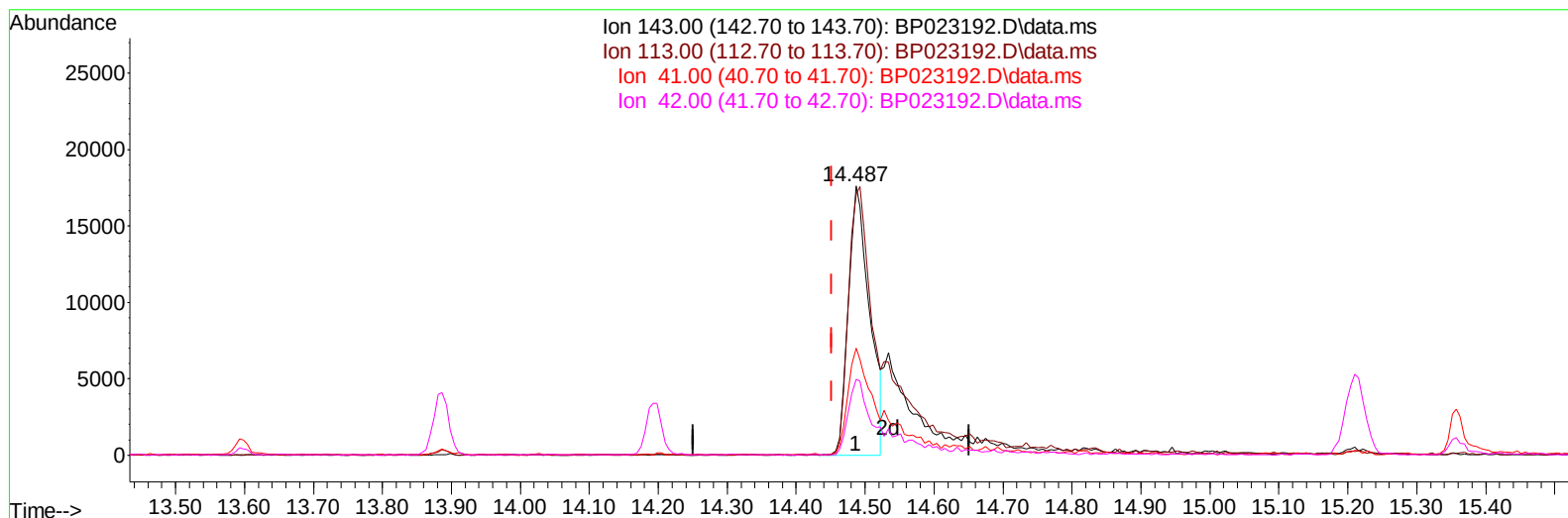
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111824\
 Data File : BP023192.D
 Acq On : 18 Nov 2024 11:09
 Operator : RC/JU
 Sample : PB164978BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK978

Manual IntegrationsAPPROVED

Quant Time: Nov 18 11:34:38 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 11/19/2024
 Supervised By :mohammad ahmed 11/19/2024



TIC: BP023192.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.487min (+ 0.035) 19.27 ng/ul

response 37516

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	96.91
41.00	39.20	39.77
42.00	25.90	28.27

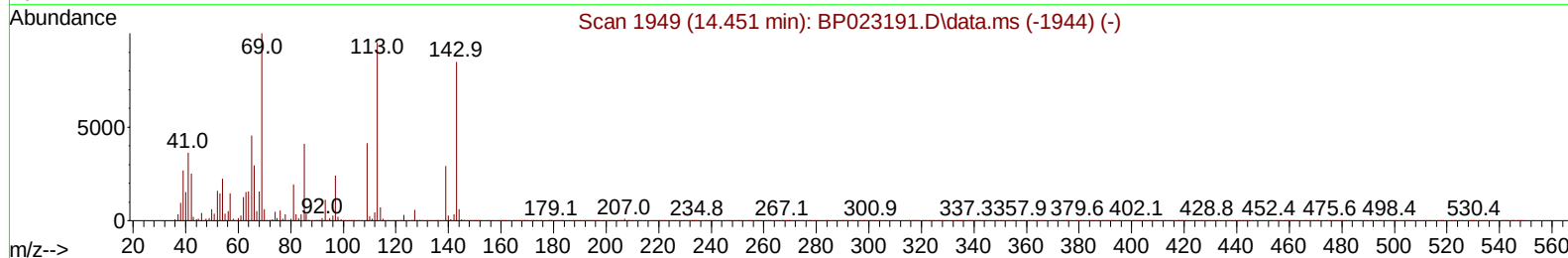
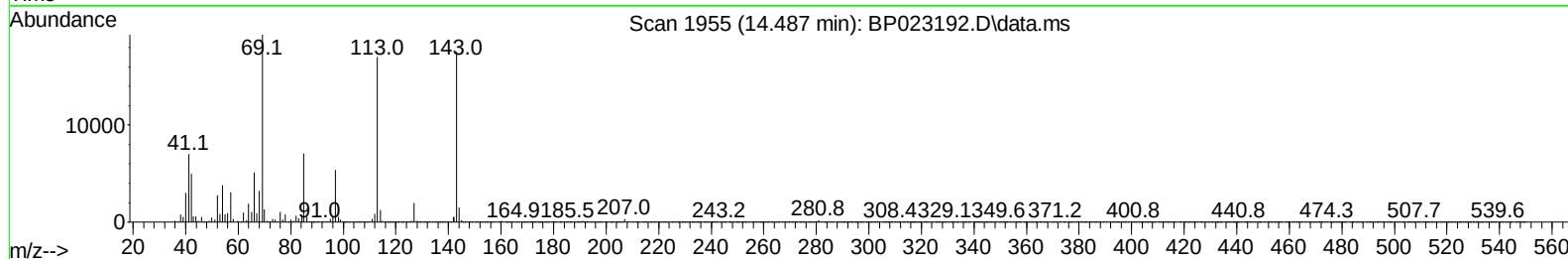
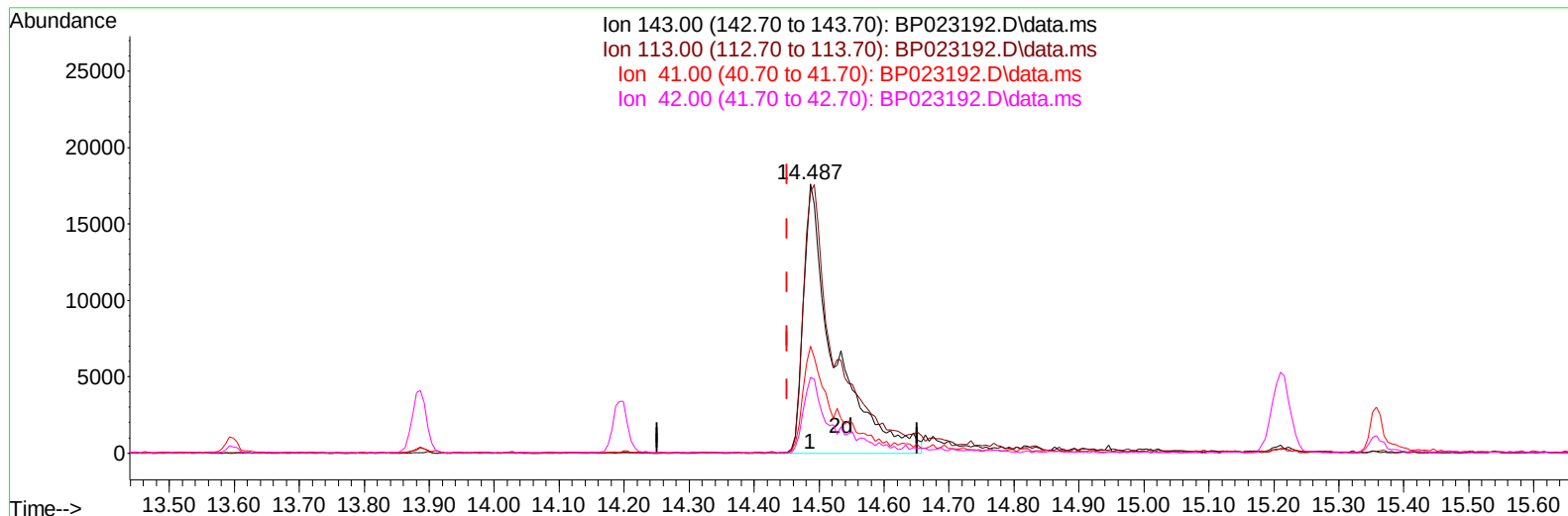
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TIC: BP023192.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.487min (+ 0.035) 29.75 ng/ul m

response 57919

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	96.91
41.00	39.20	39.77
42.00	25.90	28.27

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/19/2024
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Quant Time: Nov 18 11:44:17 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.575	152	56440	20.000	ng/ul	0.00
20) Naphthalene-d8	10.322	136	222306	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.192	164	179750	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.004	188	464908	20.000	ng/ul	-0.01
79) Chrysene-d12	21.445	240	489255	20.000	ng/ul	0.00
88) Perylene-d12	24.651	264	539150	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.134	96	9542	8.126	ng/uL	0.00
4) Pyridine-d5	3.546	84	118300	33.911	ng/ul	0.00
7) Phenol-d5	6.781	99	141897	34.006	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.922	67	90298	36.832	ng/ul	0.00
11) 2-Chlorophenol-d4	7.116	132	112269	36.737	ng/ul	-0.01
15) 4-Methylphenol-d8	8.287	113	114230	33.266	ng/ul	0.00
21) Nitrobenzene-d5	8.710	128	56928	36.784	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.428	143	67364	36.511	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.975	165	126218	32.666	ng/ul	0.00
31) 4-Chloroaniline-d4	10.475	131	152231	33.025	ng/ul	0.00
46) Dimethylphthalate-d6	13.598	166	500040	38.598	ng/ul	-0.01
49) Acenaphthylene-d8	13.887	160	514624	38.070	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.487	143	57919m	29.749	ng/ul	0.04
60) Fluorene-d10	15.210	176	442230	38.934	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.357	200	74405	26.716	ng/ul	0.00
73) Anthracene-d10	17.104	188	790747	40.171	ng/ul	-0.01
81) Pyrene-d10	19.486	212	991301	43.989	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.433	264	1060020	41.275	ng/ul	-0.01

Target Compounds Qvalue

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

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