

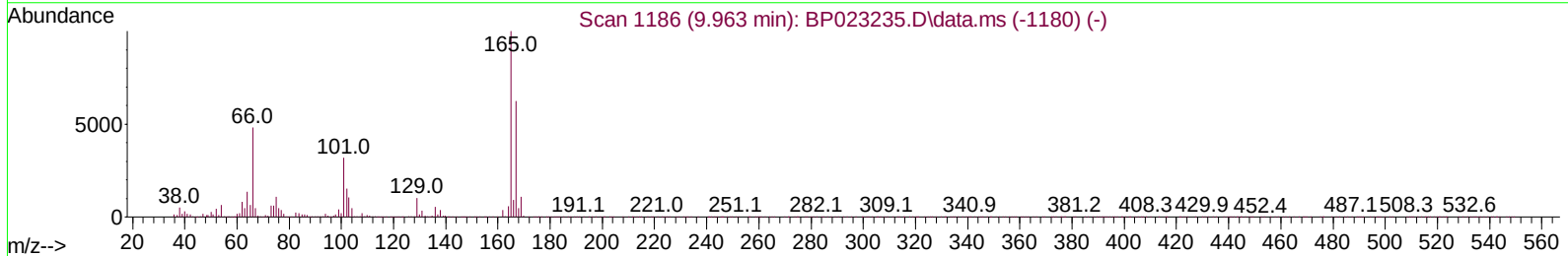
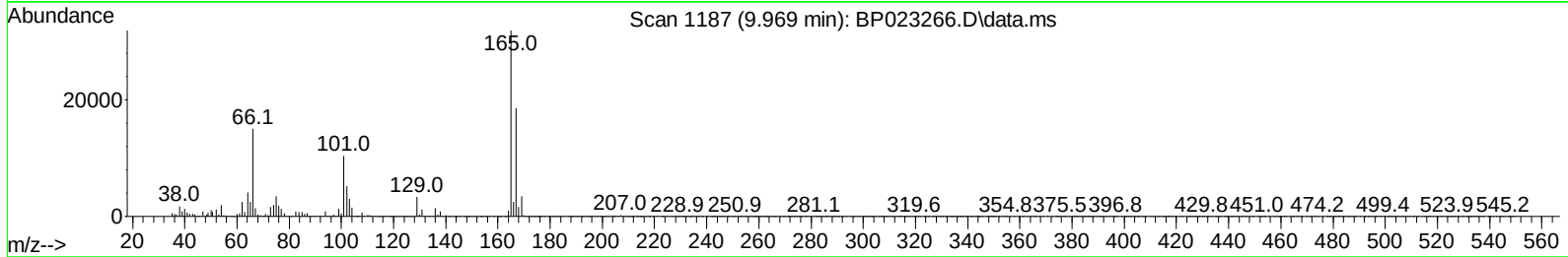
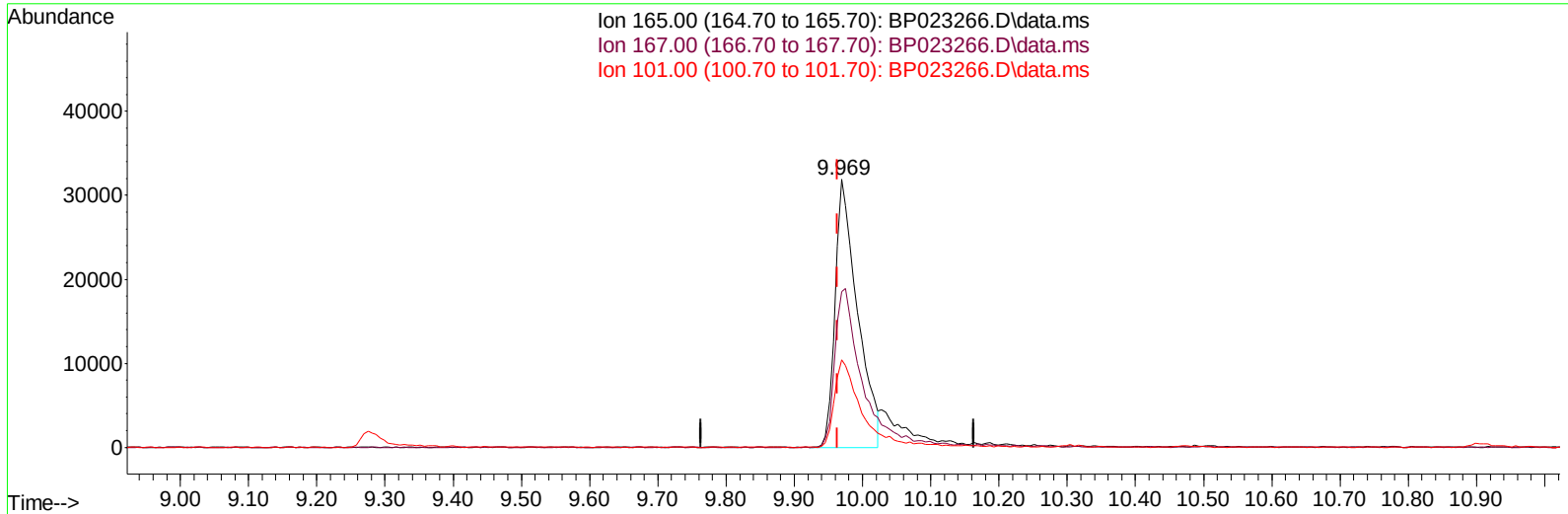
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111924\
 Data File : BP023266.D
 Acq On : 21 Nov 2024 01:50
 Operator : RC/JU
 Sample : P4872-09
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E29X7

Manual IntegrationsAPPROVED

Quant Time: Nov 21 03:15:32 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 20 04:34:35 2024
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/22/2024
 Supervised By :mohammad ahmed 11/22/2024



TIC: BP023266.D\data.ms

(28) 2,4-Dichlorophenol-d3 (S)

9.969min (+ 0.006) 13.86 ng/ul

response 72166

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	65.80	58.15
101.00	32.60	32.58
0.00	0.00	0.00

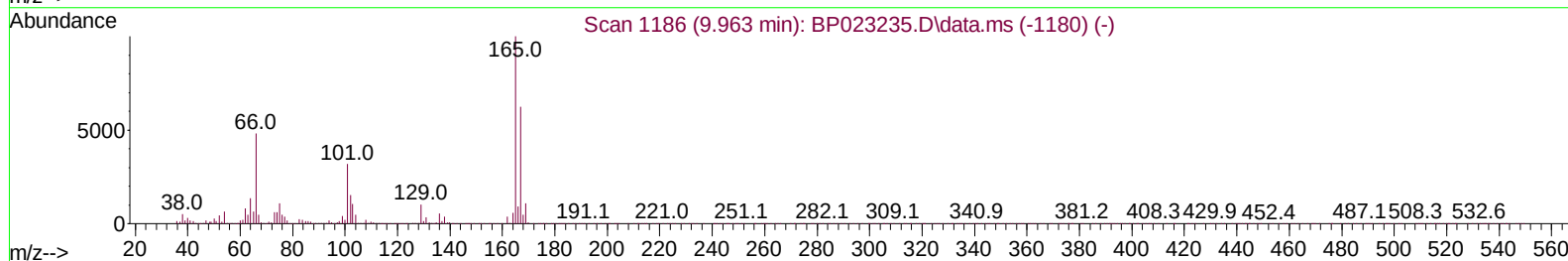
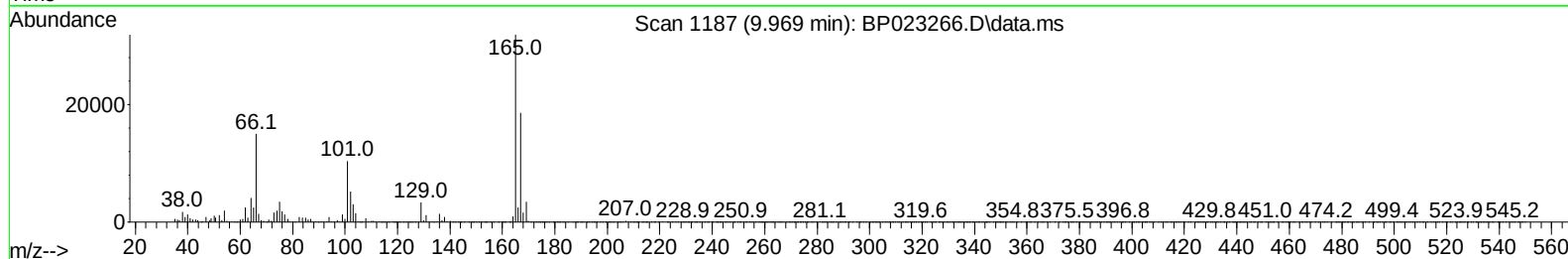
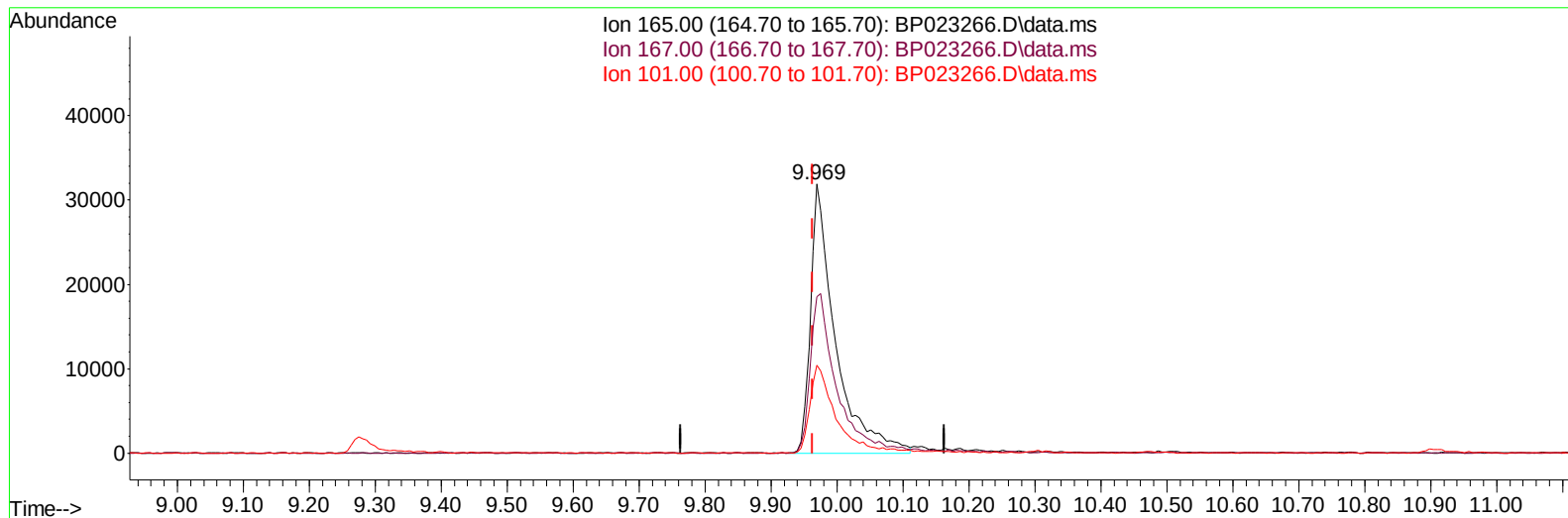
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TIC: BP023266.D\data.ms

(28) 2,4-Dichlorophenol-d3 (S)

9.969min (+ 0.006) 16.02 ng/ul m

response 83387

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	65.80	58.15
101.00	32.60	32.58
0.00	0.00	0.00

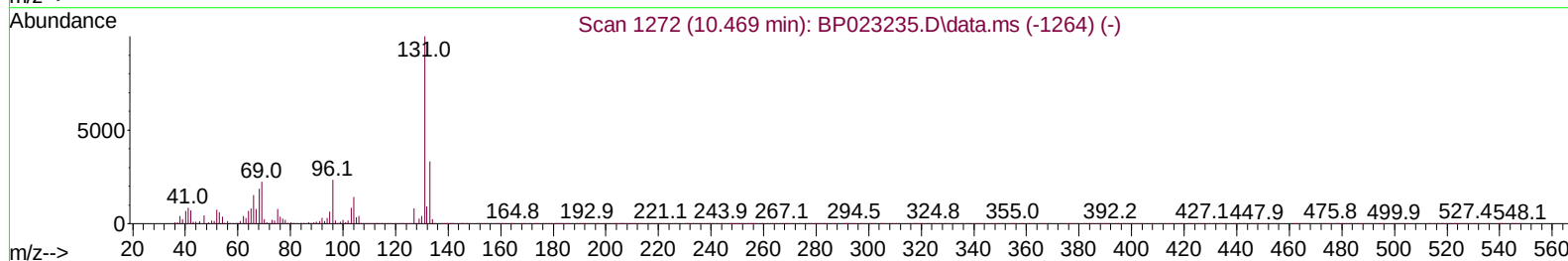
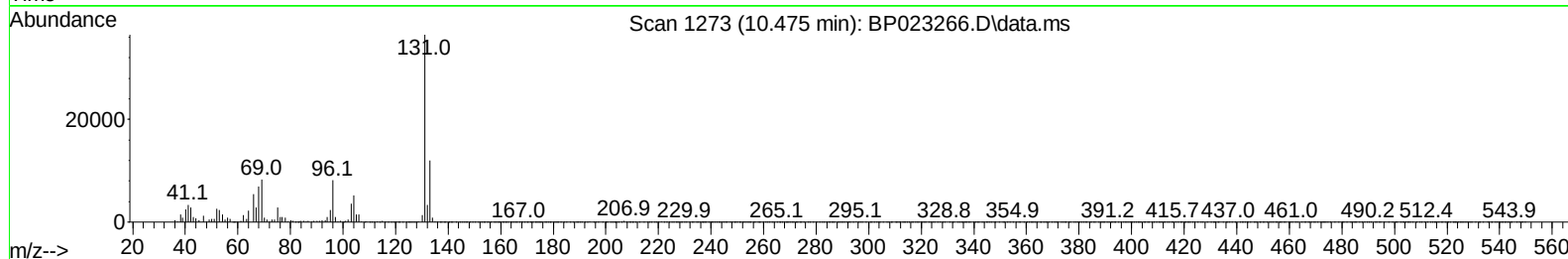
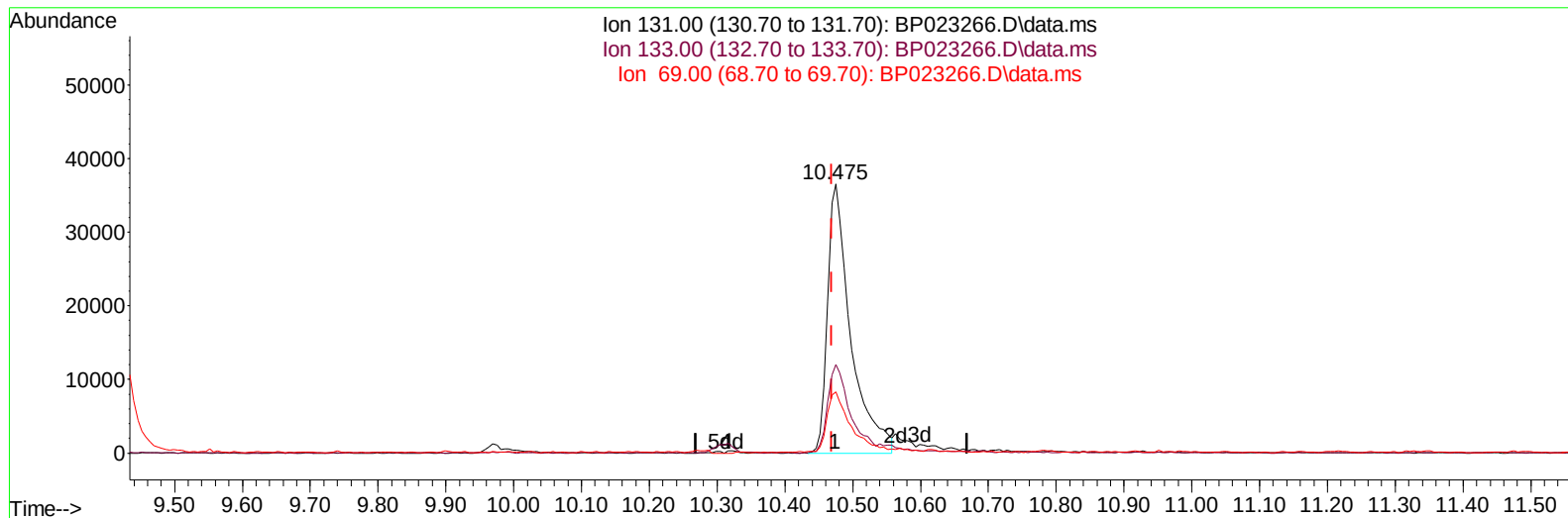
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TIC: BP023266.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.475min (+ 0.006) 13.89 ng/ul

response 86270

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	33.90	32.92
69.00	23.10	22.88
0.00	0.00	0.00

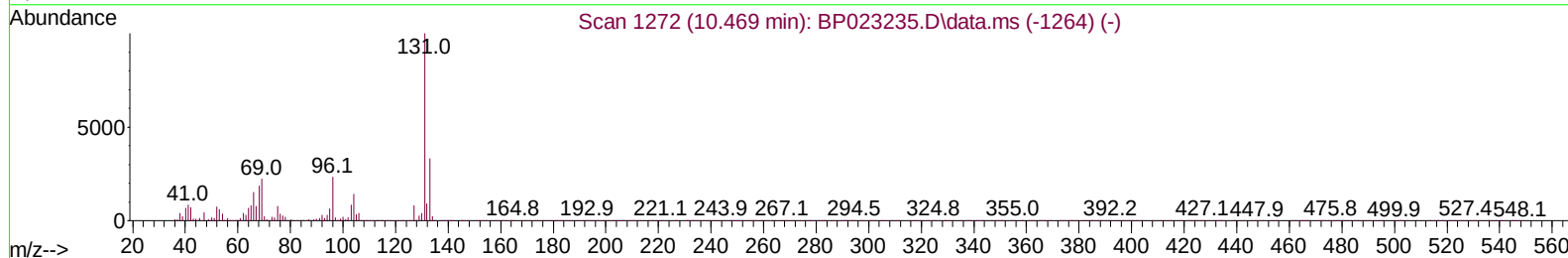
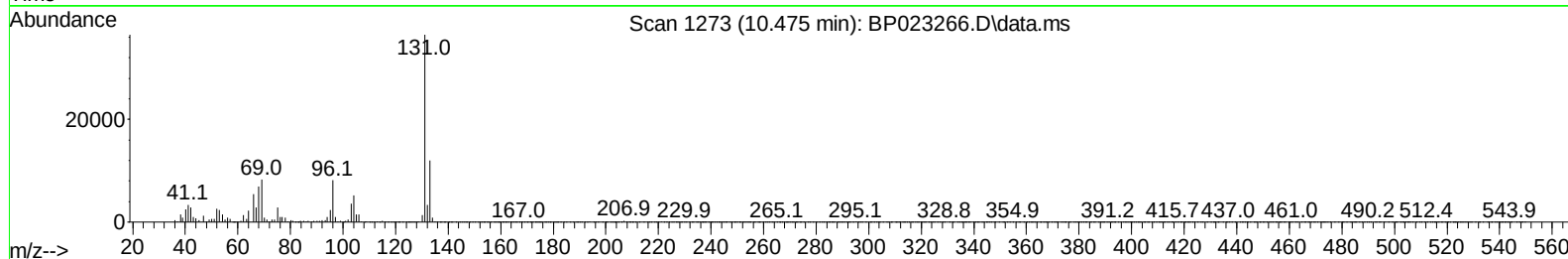
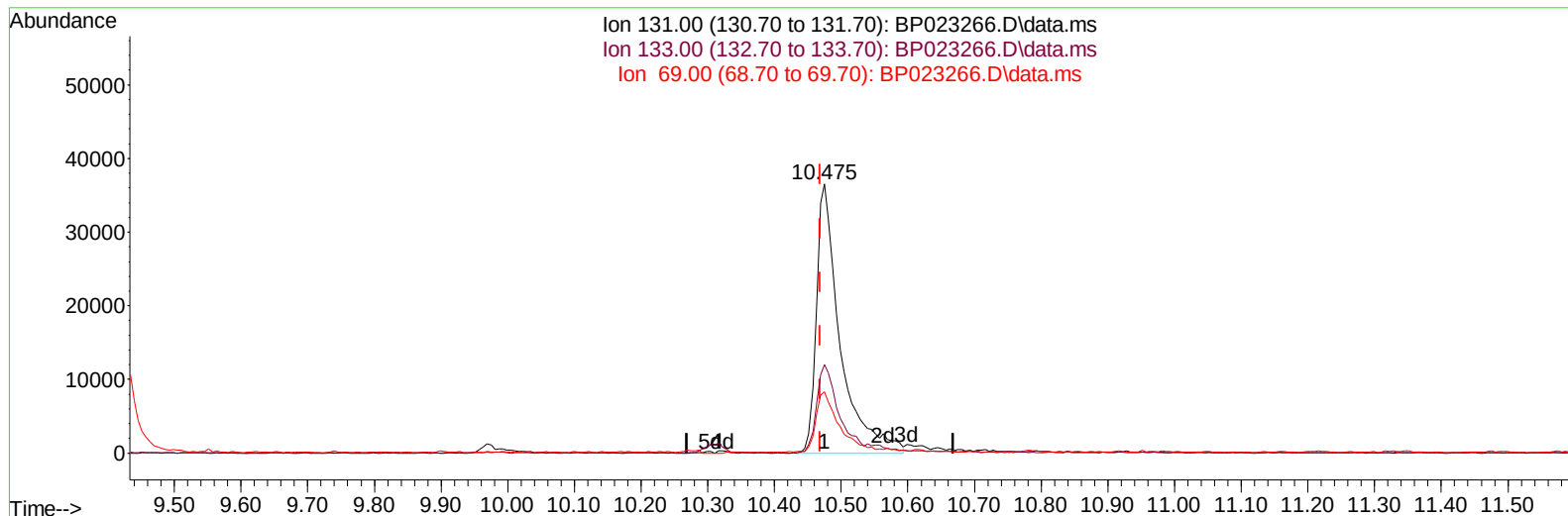
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TIC: BP023266.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.475min (+ 0.006) 14.47 ng/ul m

response 89888

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	33.90	32.92
69.00	23.10	22.88
0.00	0.00	0.00

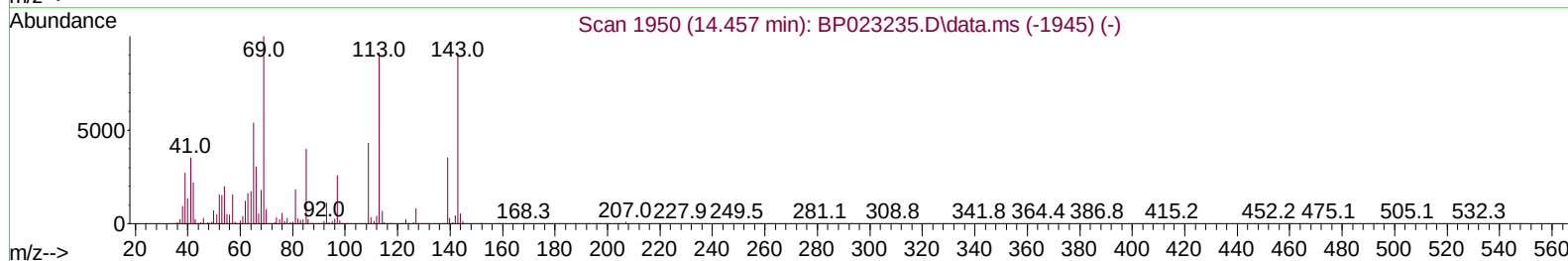
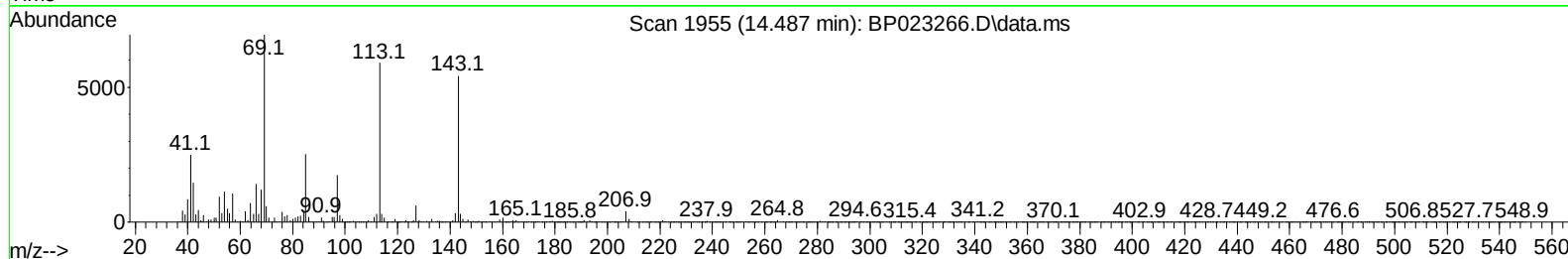
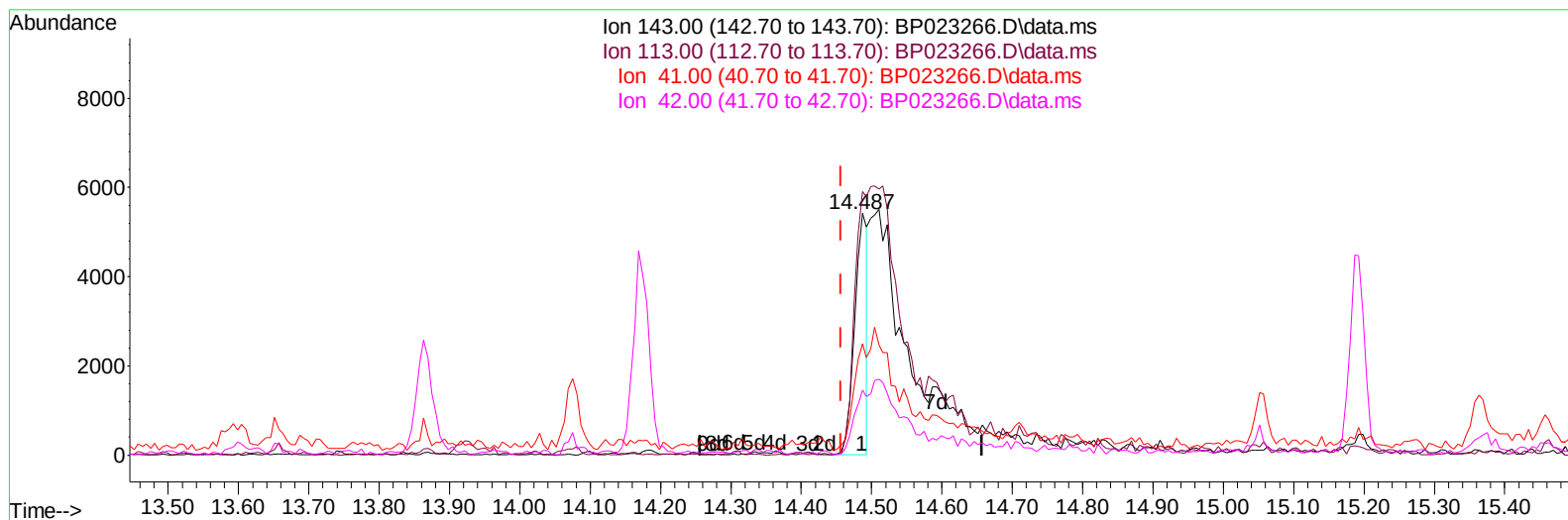
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TIC: BP023266.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.487min (+ 0.030) 2.80 ng/ul

response 6839

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	109.00
41.00	39.20	46.14
42.00	25.90	26.88

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 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/22/2024
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Quant Time: Nov 21 03:17:36 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.564	152	84332	20.000	ng/ul	0.00
20) Naphthalene-d8	10.310	136	299504	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.175	164	225614	20.000	ng/ul	-0.01
64) Phenanthrene-d10	16.987	188	551006	20.000	ng/ul	-0.01
79) Chrysene-d12	21.439	240	690771	20.000	ng/ul	0.02
88) Perylene-d12	24.657	264	822214	20.000	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.134	96	5396	3.076	ng/uL	0.00
4) Pyridine-d5	3.552	84	61030	11.708	ng/ul	0.00
7) Phenol-d5	6.775	99	91022	14.599	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.916	67	55575	15.171	ng/ul	0.00
11) 2-Chlorophenol-d4	7.116	132	67894	14.869	ng/ul	0.00
15) 4-Methylphenol-d8	8.281	113	75648	14.744	ng/ul	0.00
21) Nitrobenzene-d5	8.716	128	33167	15.907	ng/ul	0.00
24) 2-Nitrophenol-d4	9.428	143	38683	15.562	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.969	165	83387m	16.019	ng/ul	0.00
31) 4-Chloroaniline-d4	10.475	131	89888m	14.474	ng/ul	0.00
46) Dimethylphthalate-d6	13.598	166	293485	18.049	ng/ul	0.01
49) Acenaphthylene-d8	13.863	160	299164	17.632	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.510	143	30078m	12.309	ng/ul	0.05
60) Fluorene-d10	15.193	176	260213	18.252	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.363	200	32634	9.887	ng/ul	0.01
73) Anthracene-d10	17.092	188	437410	18.749	ng/ul	0.00
81) Pyrene-d10	19.481	212	596187	18.738	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.433	264	762001	19.456	ng/ul	0.01

Target Compounds Qvalue

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

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