

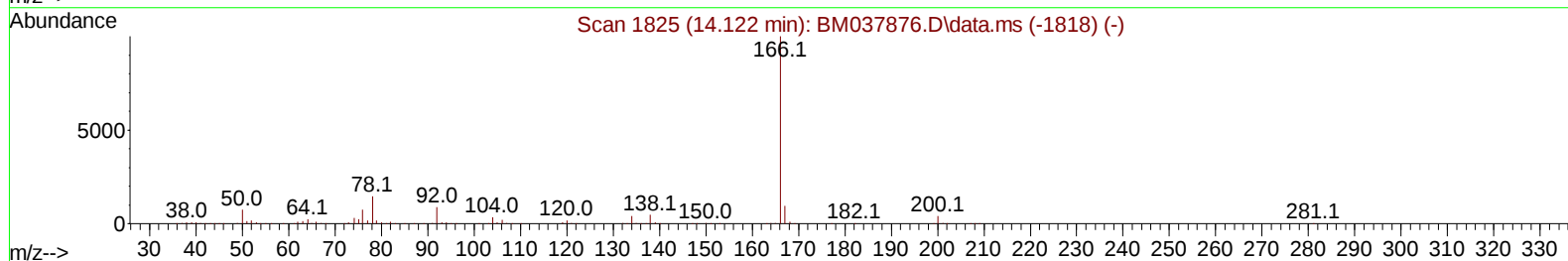
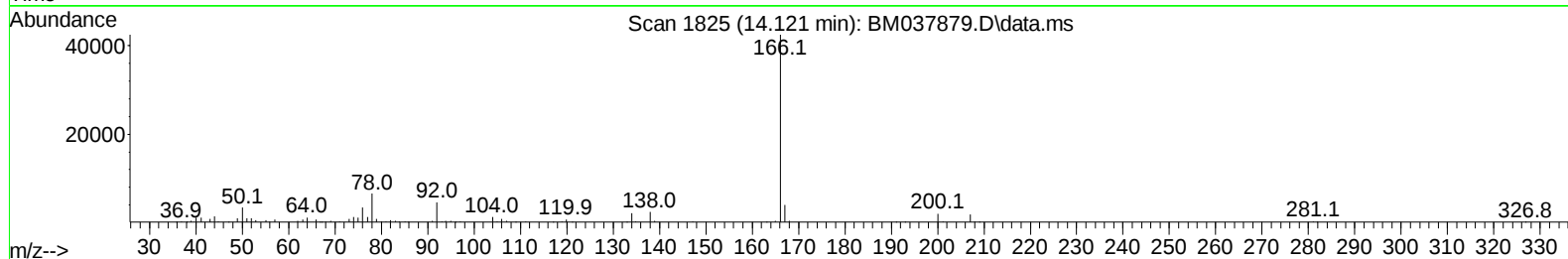
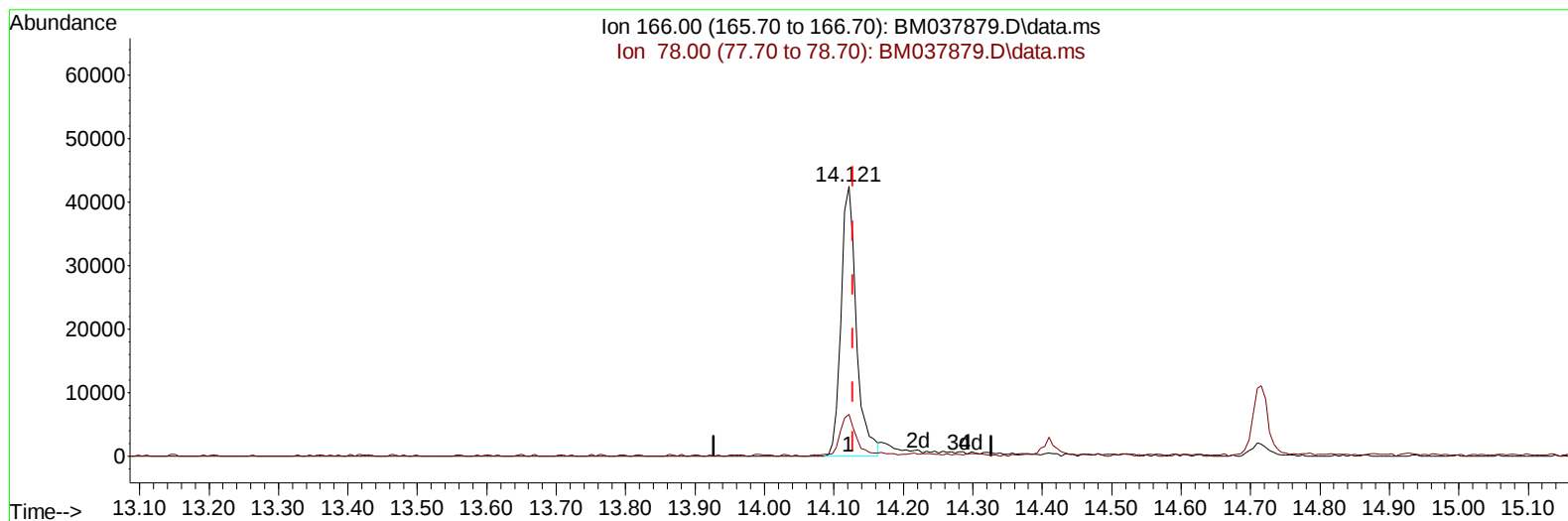
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120822\
 Data File : BM037879.D
 Acq On : 08 Dec 2022 10:23
 Operator : CG/JU
 Sample : N5738-17MEDL 10X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 GBNJ6MEDL

Manual IntegrationsAPPROVED

Quant Time: Dec 08 11:23:20 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Dec 08 01:49:06 2022
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 12/08/2022
 Supervised By :mohammad ahmed 12/08/2022



TIC: BM037879.D\data.ms

(46) Dimethylphthalate-d6 (S)

14.121min (-0.006) 2.61 ng/u1

response 64445

Ion	Exp%	Act%
166.00	100.00	100.00
78.00	14.10	15.41
0.00	0.00	0.00
0.00	0.00	0.00

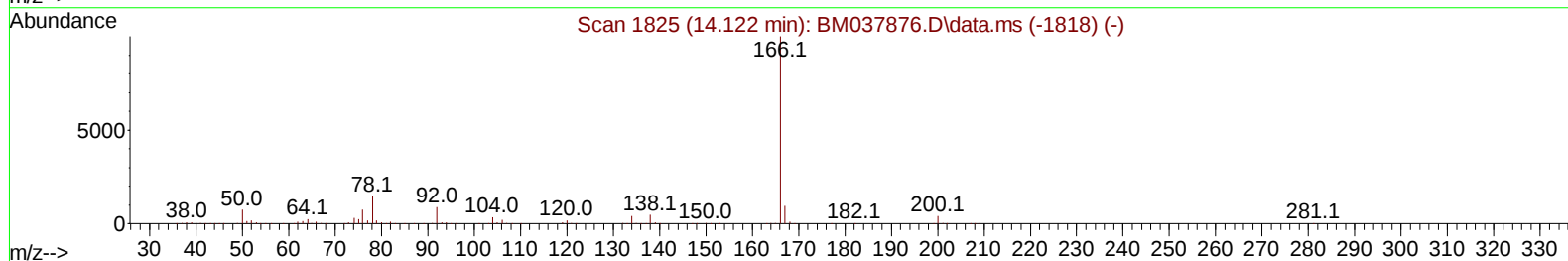
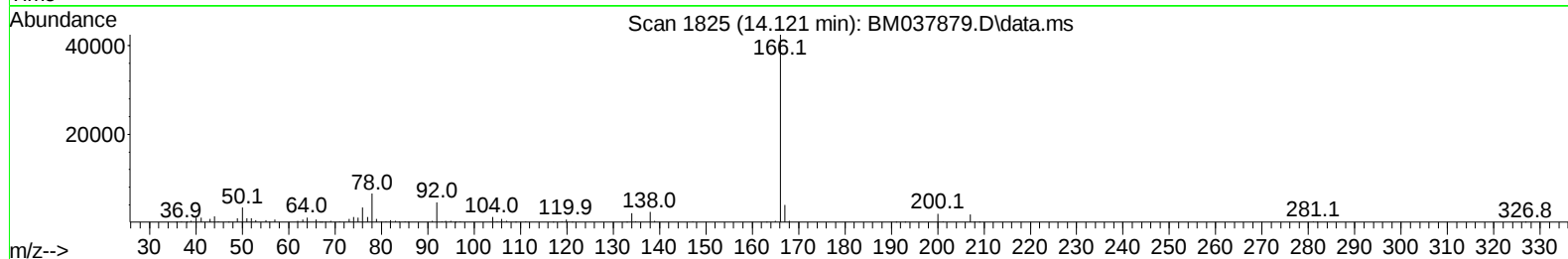
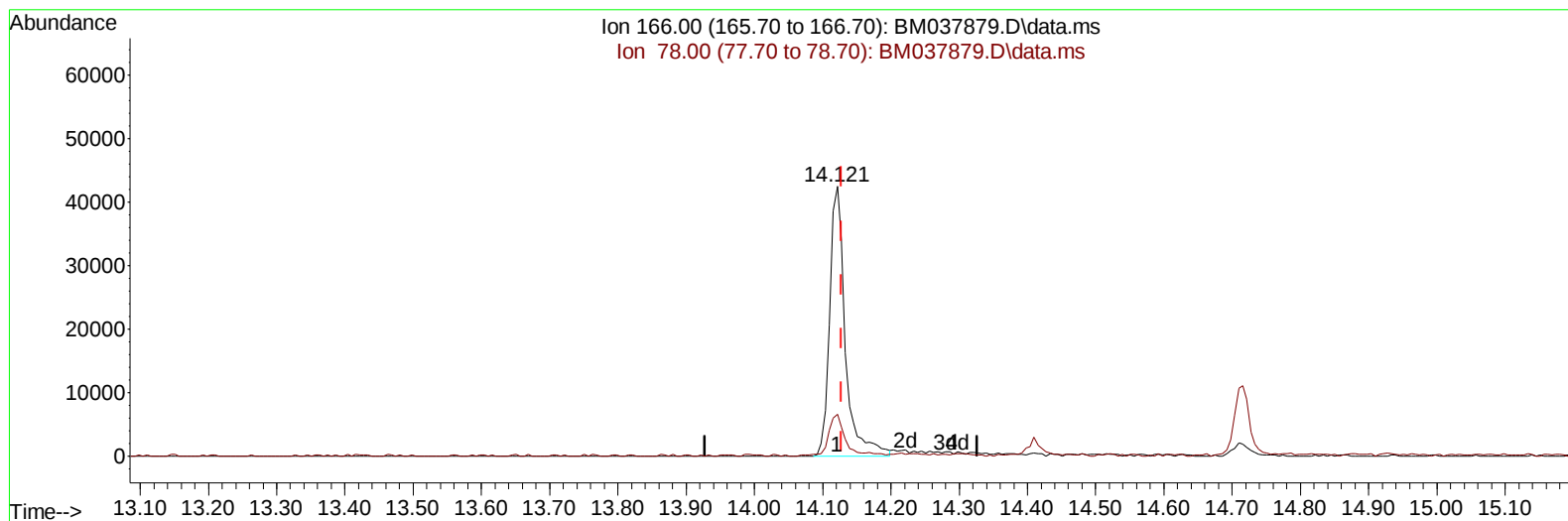
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TIC: BM037879.D\data.ms

(46) Dimethylphthalate-d6 (S)

14.121min (-0.006) 2.74 ng/ul m

response 67597

Ion	Exp%	Act%
166.00	100.00	100.00
78.00	14.10	15.41
0.00	0.00	0.00
0.00	0.00	0.00

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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	8.087	152		175458	20.000	ng/ul	0.00
20) Naphthalene-d8	10.910	136		672301	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.716	164		345737	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.451	188		776019	20.000	ng/ul	0.00
79) Chrysene-d12	21.621	240		801432	20.000	ng/ul	0.00
88) Perylene-d12	24.103	264		879104	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	0.000	96		0d	0.000	ng/uL	
4) Pyridine-d5	0.000	84		0d	0.000	ng/ul	
7) Phenol-d5	7.240	99		37633	2.639	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.410	67		26670	3.072	ng/ul	0.00
11) 2-Chlorophenol-d4	7.616	132		32099	2.744	ng/ul	0.00
15) 4-Methylphenol-d8	8.792	113		23884	2.146	ng/ul	0.00
21) Nitrobenzene-d5	9.257	128		14944	2.807	ng/ul	0.00
24) 2-Nitrophenol-d4	9.986	143		13479	2.416	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.522	165		23661	2.233	ng/ul	0.00
31) 4-Chloroaniline-d4	11.051	131		10643	0.728	ng/ul	0.00
46) Dimethylphthalate-d6	14.121	166		67597m	2.742	ng/ul	0.00
49) Acenaphthylene-d8	14.410	160		82129	2.726	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143		0d	0.000	ng/ul	
60) Fluorene-d10	15.704	176		68211	3.106	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.815	200		6080	1.465	ng/ul	0.00
73) Anthracene-d10	17.551	188		115955	3.202	ng/ul	0.00
81) Pyrene-d10	19.833	212		153631	3.191	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.938	264		144412	3.289	ng/ul	0.00
Target Compounds							
						Qvalue	
71) Pentachlorophenol	17.104	266		453012	80.242	ng/ul	99

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

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