

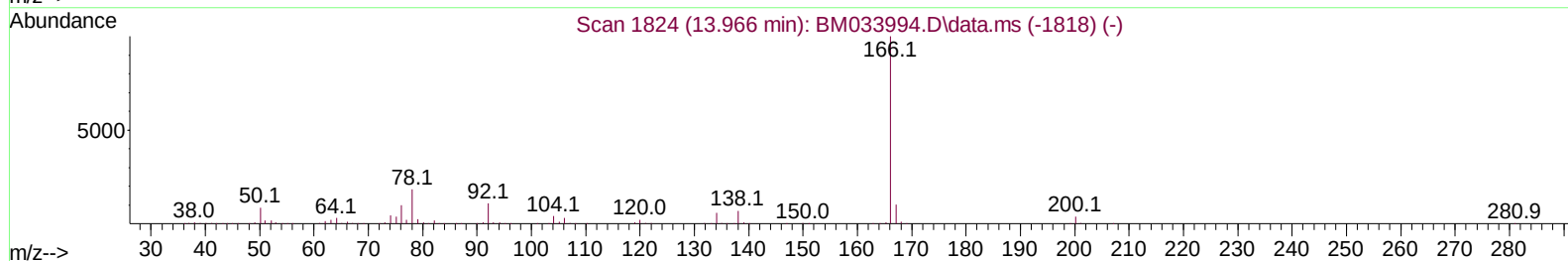
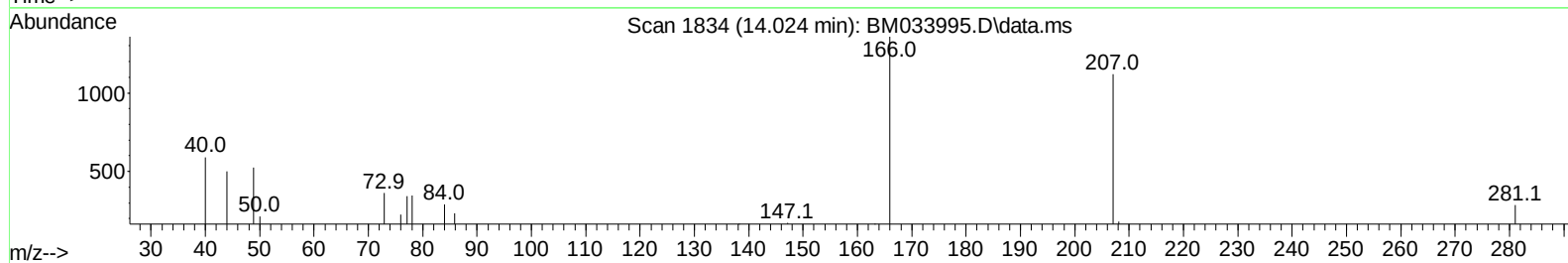
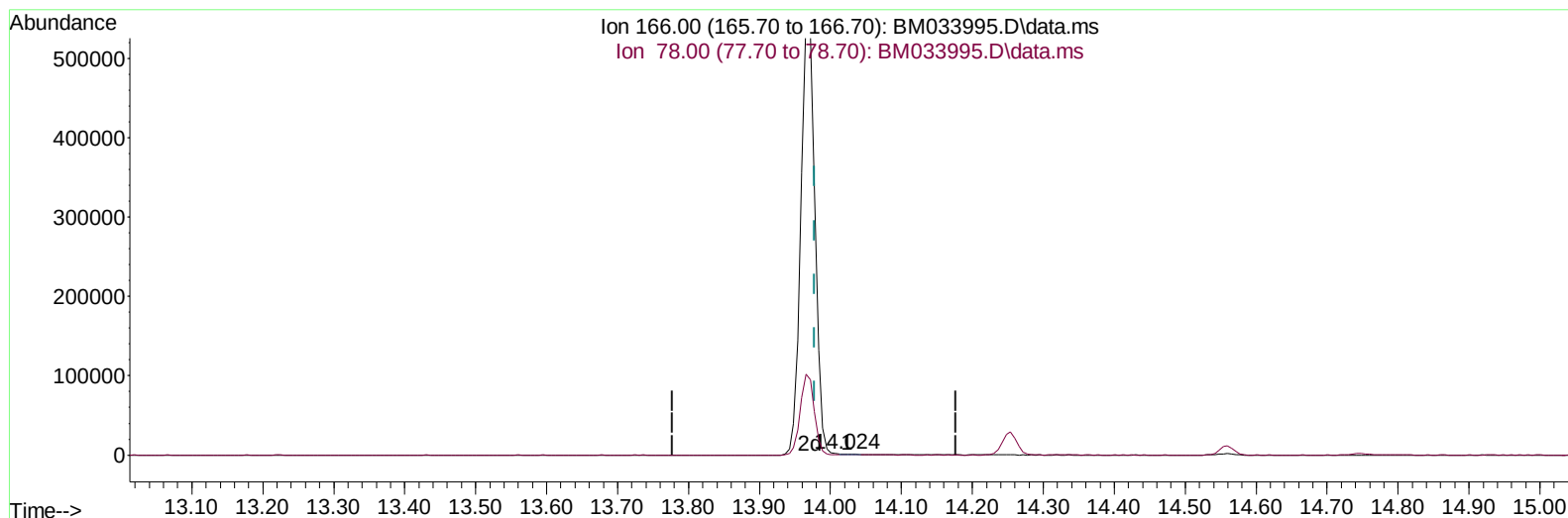
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020322\
Data File : BM033995.D
Acq On : 03 Feb 2022 11:01
Operator : CG/JU
Sample : PB142413BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SBLK413

Manual IntegrationsAPPROVED

Quant Time: Feb 04 02:59:37 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Feb 01 02:16:28 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 02/04/2022
Supervised By :mohammad ahmed 02/04/2022



TIC: BM033995.D\data.ms

(46) Dimethylphthalate-d6 (S)

14.024min (+ 0.047) 0.04 ng/u1

response 910

Ion	Exp%	Act%
166.00	100.00	100.00
78.00	21.40	25.55
0.00	0.00	0.00
0.00	0.00	0.00

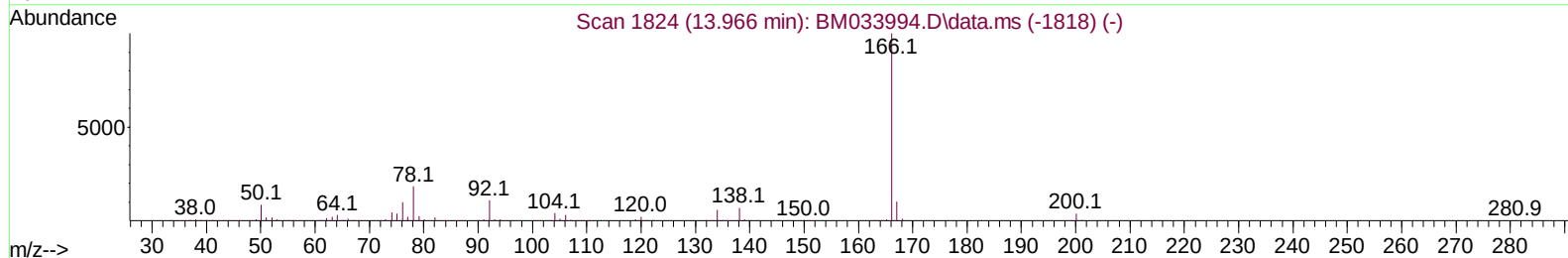
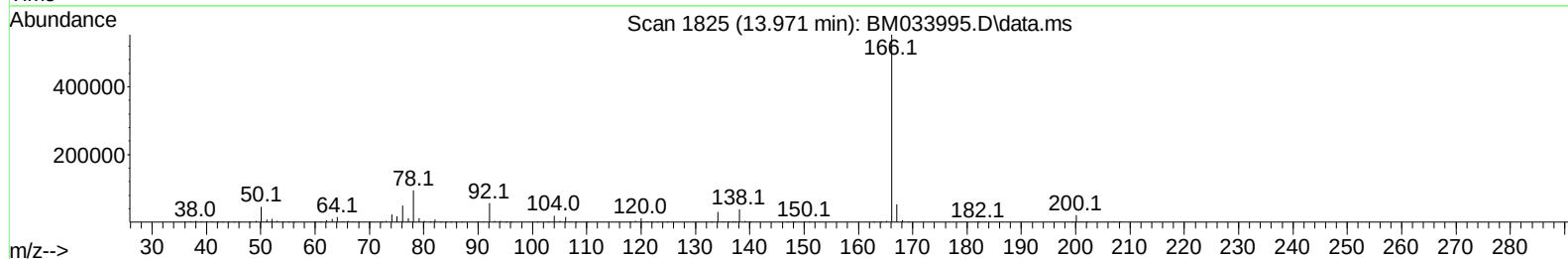
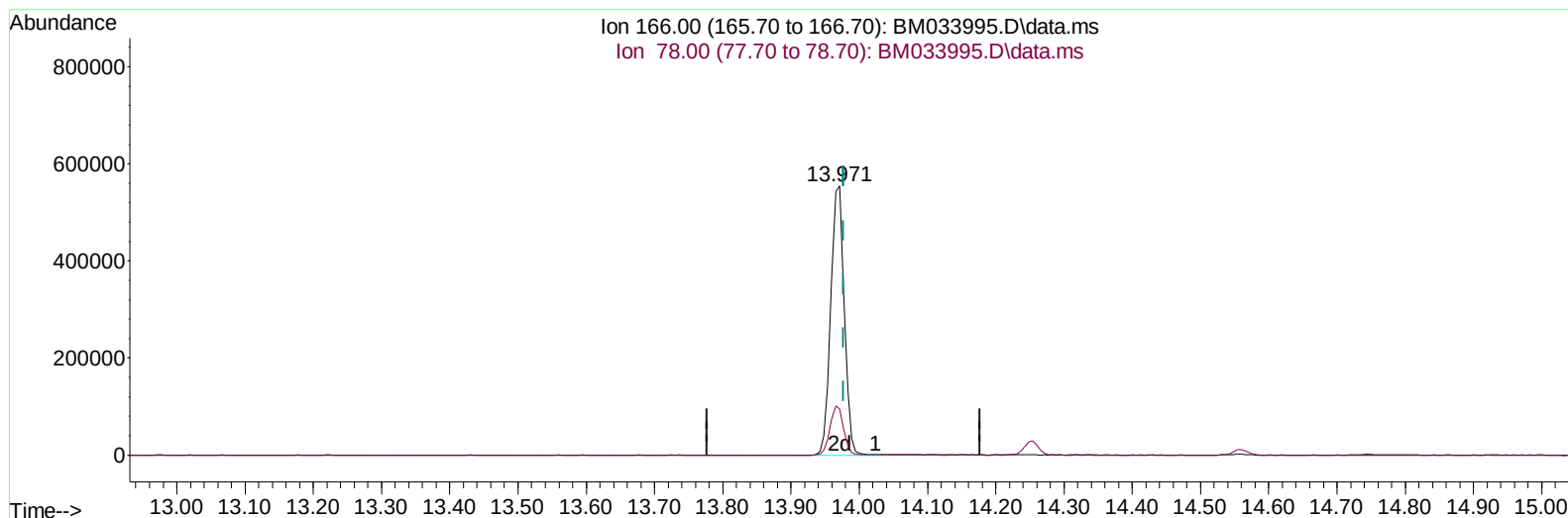
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(46) Dimethylphthalate-d6 (S)

13.971min (-0.006) 34.41 ng/ul m

response 765767

Ion	Exp%	Act%
166.00	100.00	100.00
78.00	21.40	17.07#
0.00	0.00	0.00
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.925	152	103500	20.000	ng/ul	0.00
20) Naphthalene-d8	10.730	136	453373	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.560	164	297817	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.295	188	604511	20.000	ng/ul	-0.01
79) Chrysene-d12	21.459	240	481290	20.000	ng/ul	0.00
88) Perylene-d12	23.847	264	431616	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.284	96	15687	5.978	ng/uL	0.00
4) Pyridine-d5	3.713	84	203573	27.792	ng/ul	0.00
7) Phenol-d5	7.078	99	262630	30.289	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.242	67	171443	32.432	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.448	132	217943	31.373	ng/ul	-0.01
15) 4-Methylphenol-d8	8.631	113	223319	32.502	ng/ul	-0.01
21) Nitrobenzene-d5	9.083	128	110642	30.414	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.807	143	119454	30.407	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.348	165	206895	29.103	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.866	131	304275	30.995	ng/ul	0.00
46) Dimethylphthalate-d6	13.971	166	765767m	34.409	ng/ul	0.00
49) Acenaphthylene-d8	14.254	160	849627	31.165	ng/ul	0.00
54) 4-Nitrophenol-d4	14.748	143	98953	26.025	ng/ul	0.00
60) Fluorene-d10	15.548	176	602081	32.400	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.660	200	96362	25.391	ng/ul	-0.01
73) Anthracene-d10	17.395	188	984767	34.016	ng/ul	0.00
81) Pyrene-d10	19.677	212	1039088	36.569	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.694	264	776872	34.268	ng/ul	-0.01

Target Compounds Qvalue

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

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