

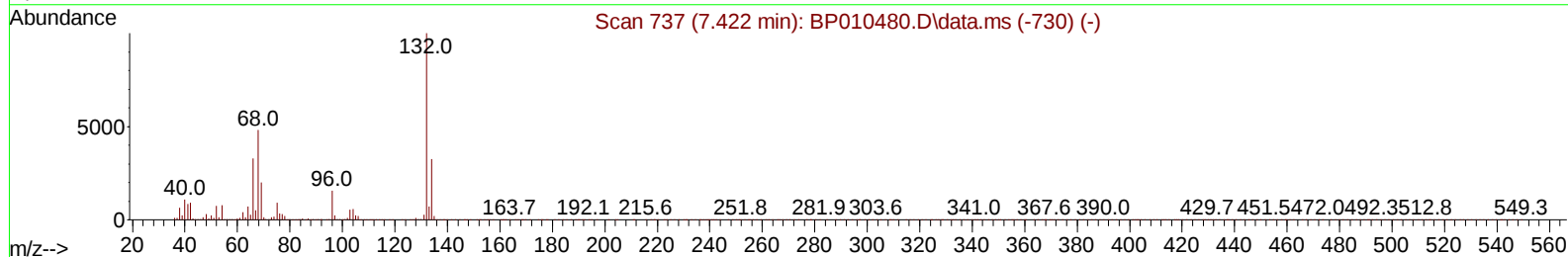
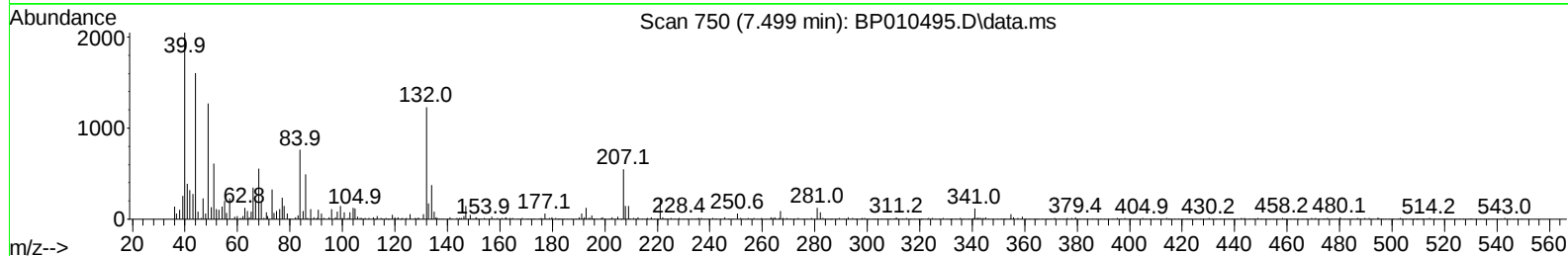
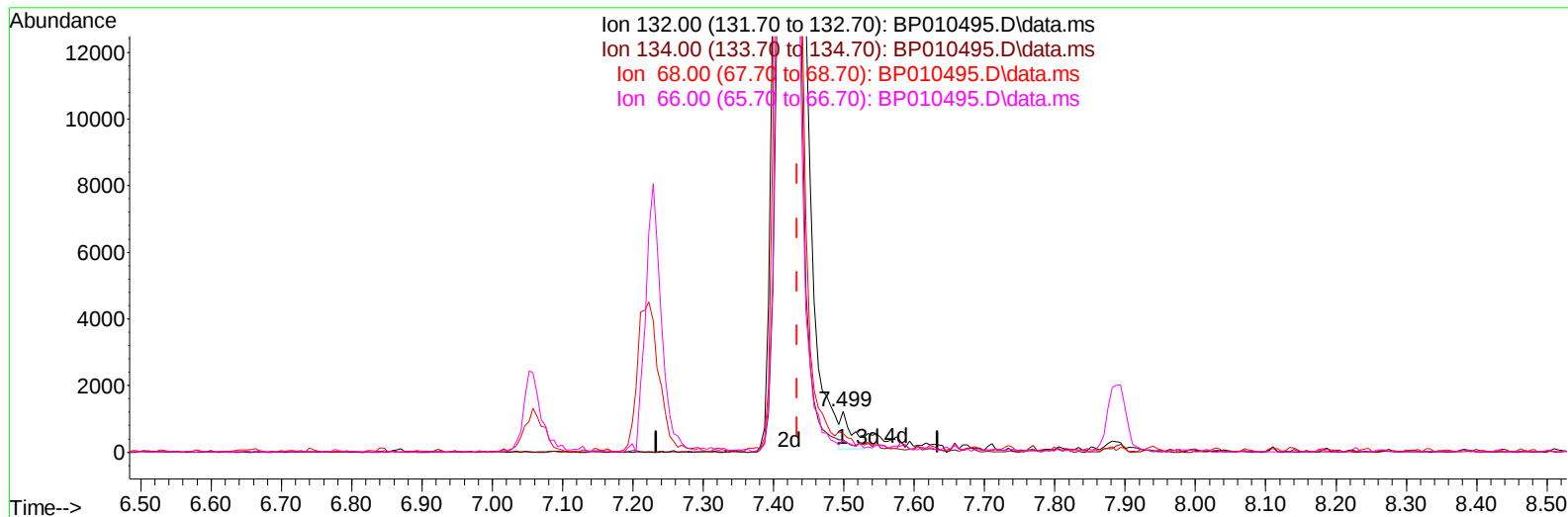
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052322\
 Data File : BP010495.D
 Acq On : 24 May 2022 01:40
 Operator : CG/JU
 Sample : N2950-01
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBTE6

Manual IntegrationsAPPROVED

Quant Time: May 24 02:39:46 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri May 13 15:04:31 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/24/2022
 Supervised By :Yogesh Patel 05/26/2022



TIC: BP010495.D\data.ms

(11) 2-Chlorophenol-d4 (S)

7.499min (+ 0.065) 0.11 ng/u1

response 1204

Ion	Exp%	Act%
132.00	100.00	100.00
134.00	30.90	30.65
68.00	51.40	45.37
66.00	26.30	28.05

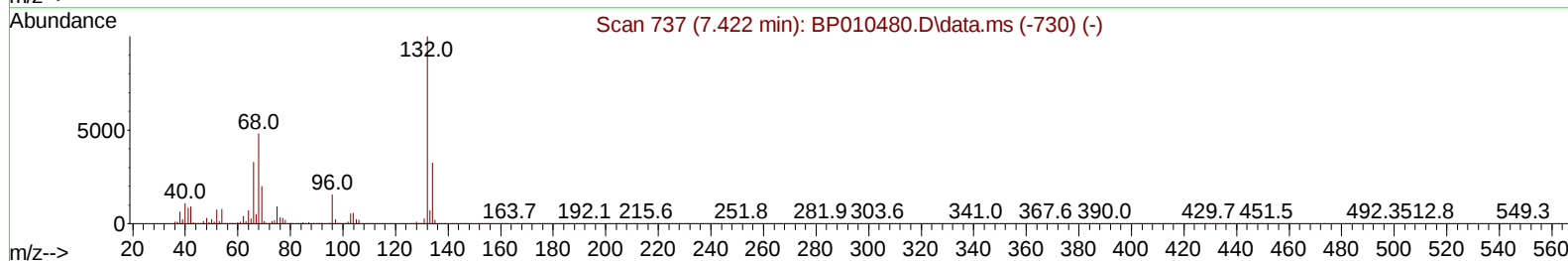
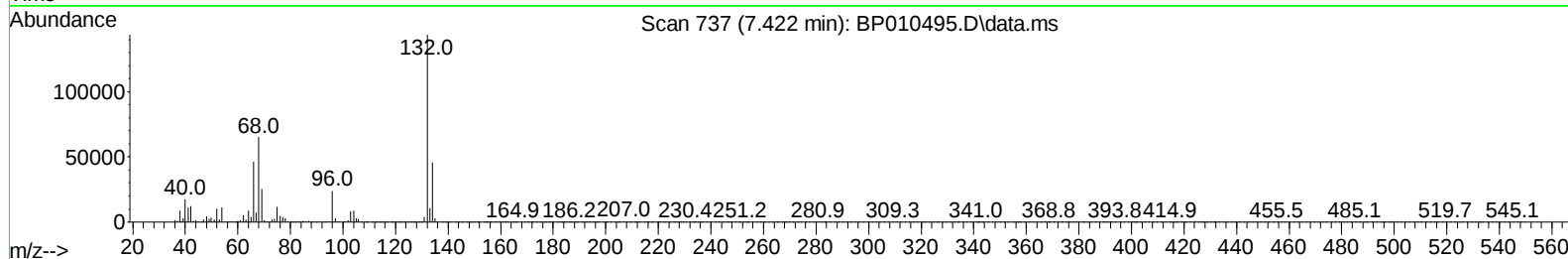
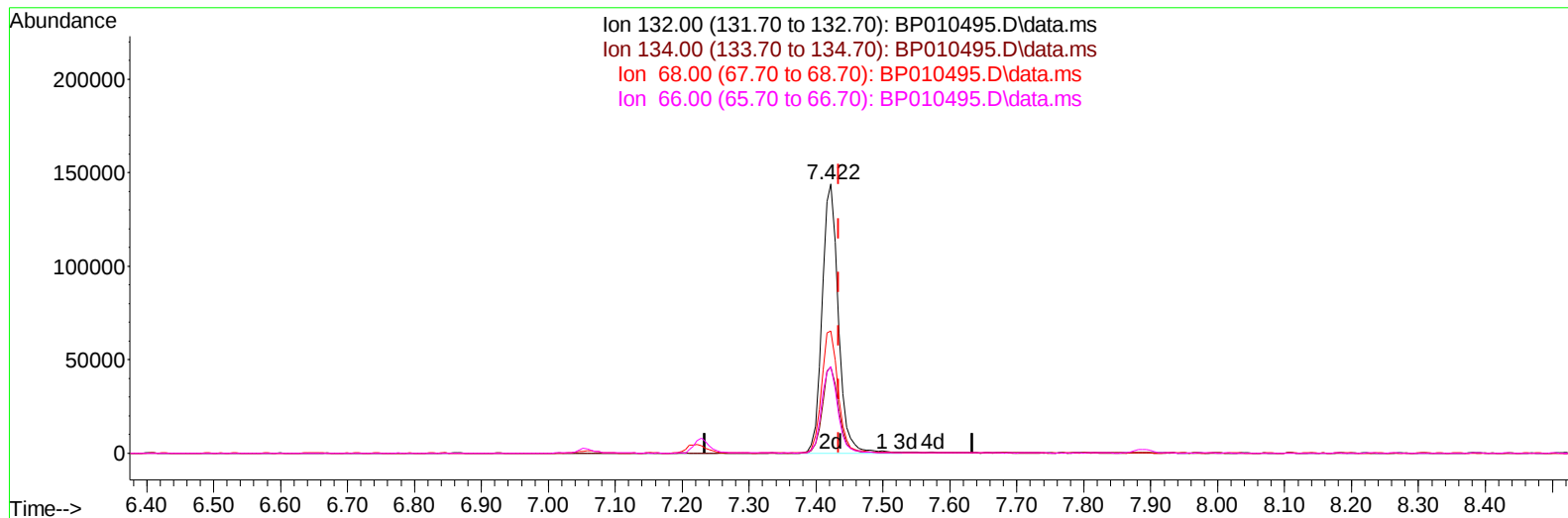
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TIC: BP010495.D\data.ms

(11) 2-Chlorophenol-d4 (S)

7.422min (-0.011) 22.85 ng/ul m

response 242147

Ion	Exp%	Act%
132.00	100.00	100.00
134.00	30.90	32.01
68.00	51.40	45.31
66.00	26.30	32.18#

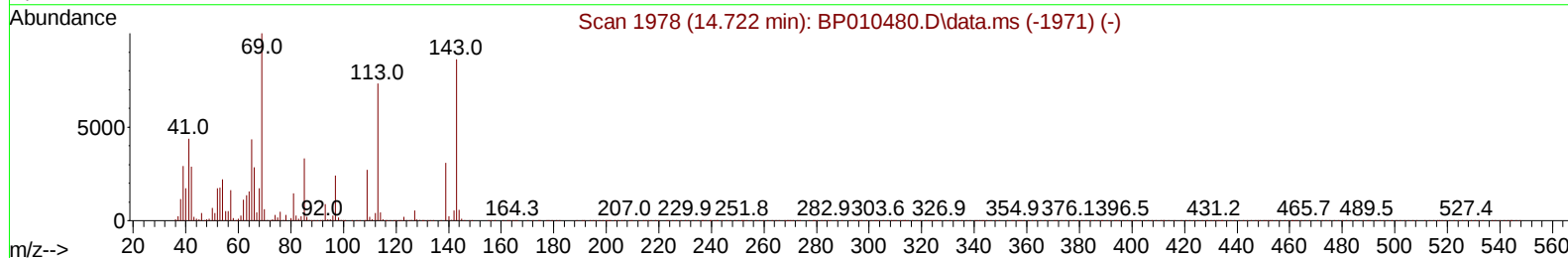
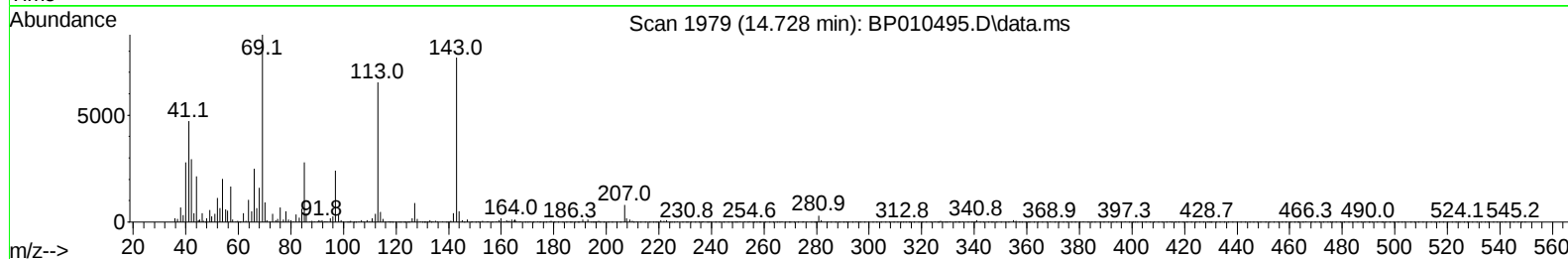
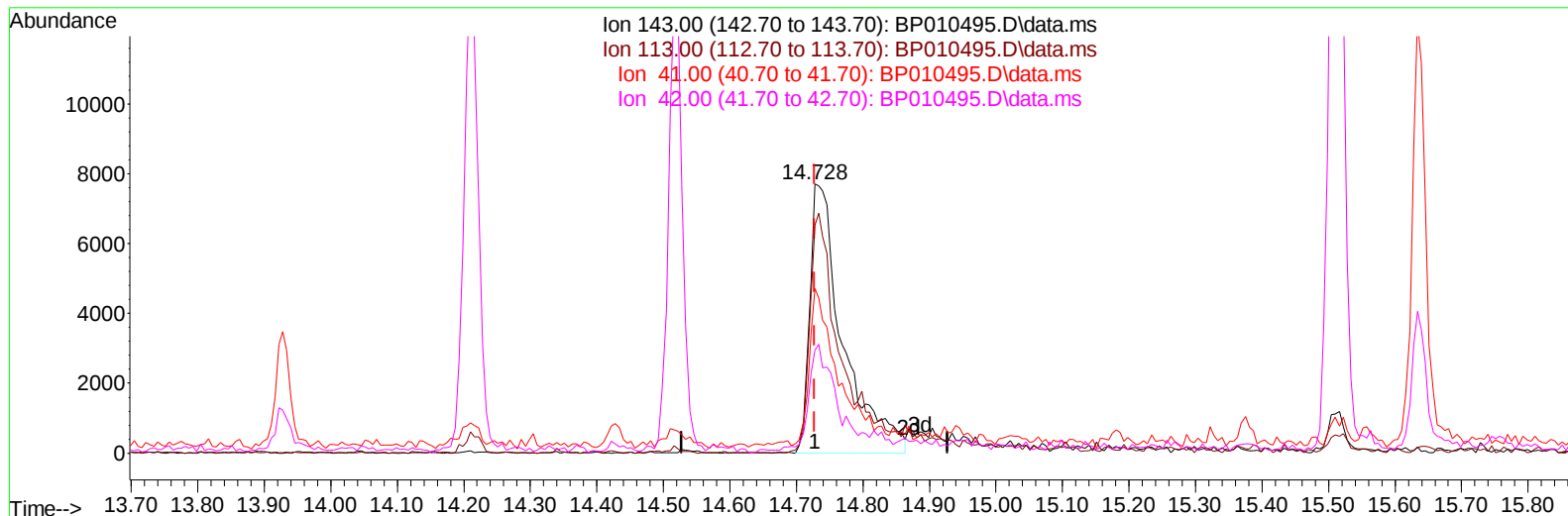
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TIC: BP010495.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.728min (+ 0.000) 3.92 ng/ul m

response 26651

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	84.72#
41.00	23.20	61.13#
42.00	18.00	38.36#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.887	152	164999	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.687	136	753761	20.000	ng/ul	#-0.02
38) Acenaphthene-d10	14.522	164	517858	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.269	188	1097326	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.351	240	946609	20.000	ng/ul	#-0.01
88) Perylene-d12	23.774	264	732680	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	15966	4.125	ng/uL	0.00
4) Pyridine-d5	3.770	84	49768	4.528	ng/ul	0.00
7) Phenol-d5	7.058	99	81042	5.852	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.217	67	270529	29.075	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.422	132	242147m	22.851	ng/ul	-0.01
15) 4-Methylphenol-d8	8.593	113	165162	14.033	ng/ul	-0.02
21) Nitrobenzene-d5	9.046	128	173073	29.668	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.769	143	178051	28.800	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.310	165	293632	25.119	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.828	131	372652	21.440	ng/ul	-0.01
46) Dimethylphthalate-d6	13.928	166	1194225	31.300	ng/ul	-0.01
49) Acenaphthylene-d8	14.210	160	1322755	30.056	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.728	143	26651m	3.924	ng/ul	0.00
60) Fluorene-d10	15.516	176	1048309	31.582	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.640	200	168851	26.311	ng/ul	0.00
73) Anthracene-d10	17.369	188	1815071	36.529	ng/ul	-0.01
81) Pyrene-d10	19.604	212	2114871	42.069	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.622	264	1373199	37.325	ng/ul	-0.01

Target Compounds Qvalue

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

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