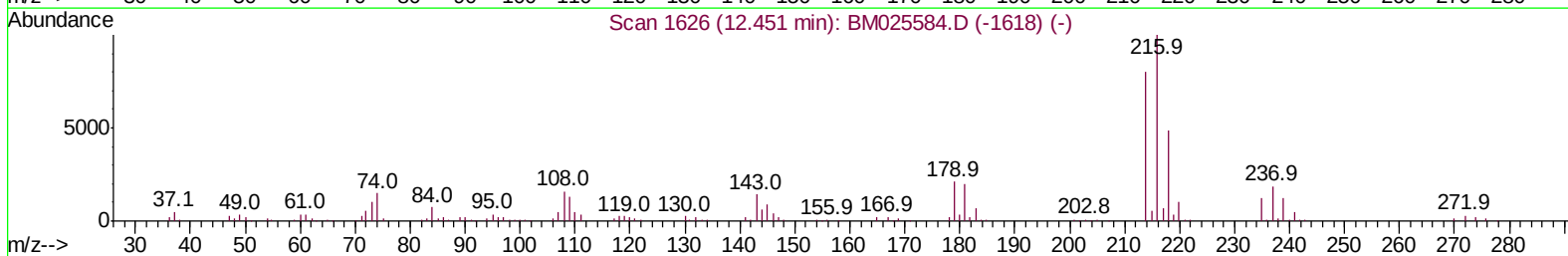
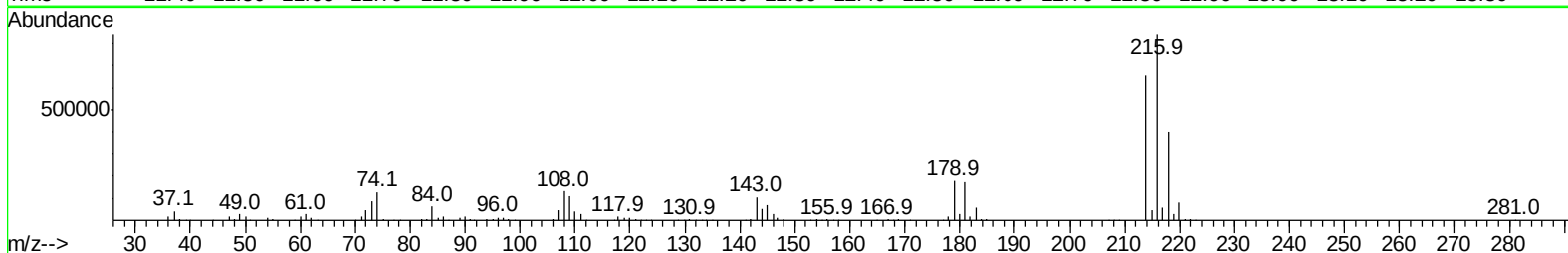
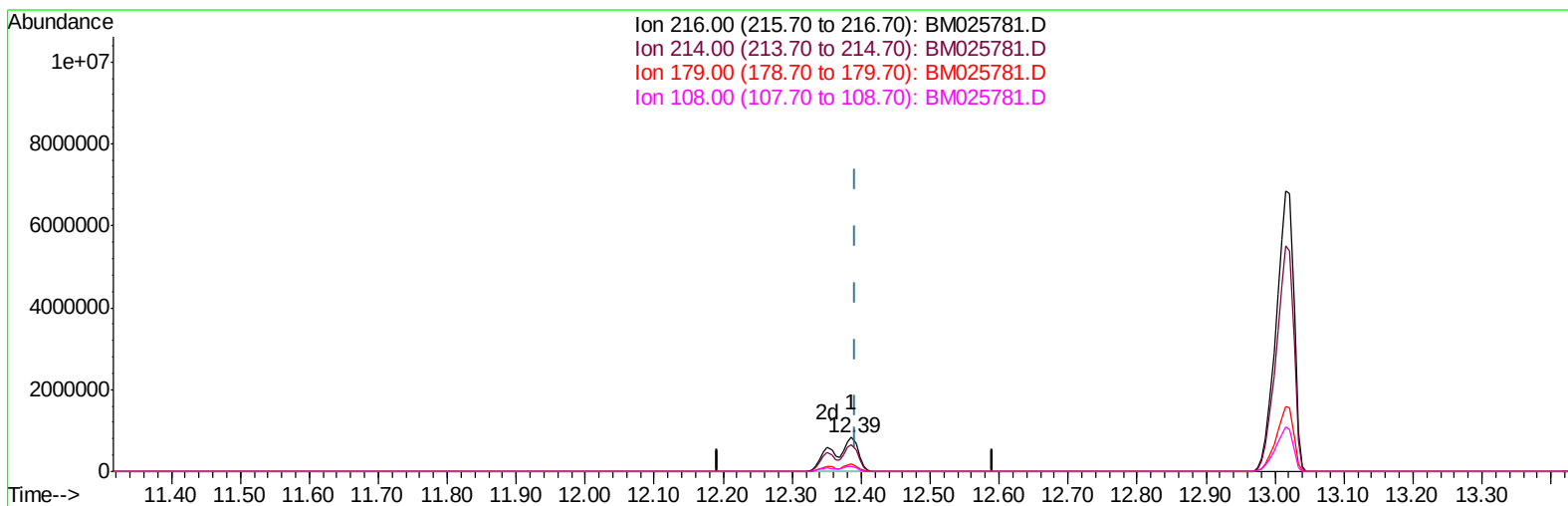


Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032320\
Data File : BM025781.D
Acq On : 25 Mar 2020 23:06
Operator : CG/JU
Sample : L1941-06
Misc :
ALS Vial : 11 Sample Multiplier: 1

Quant Time: Mar 26 01:17:19 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM031420MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Mar 24 00:45:06 2020
Response via : Initial Calibration



TIC: BM025781.D

(37) 1,2,4,5-Tetrachlorobenzene

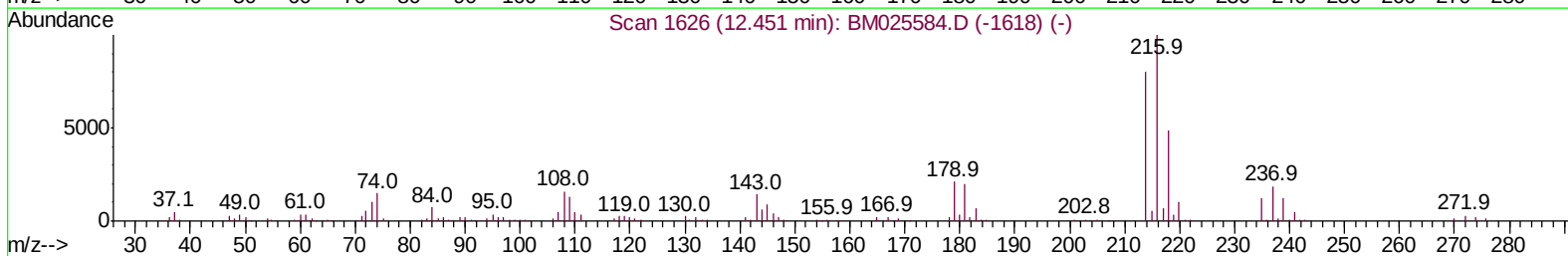
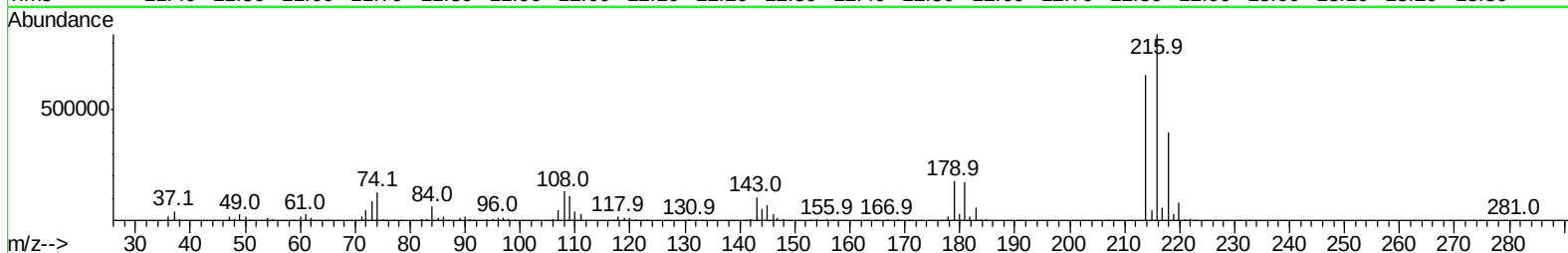
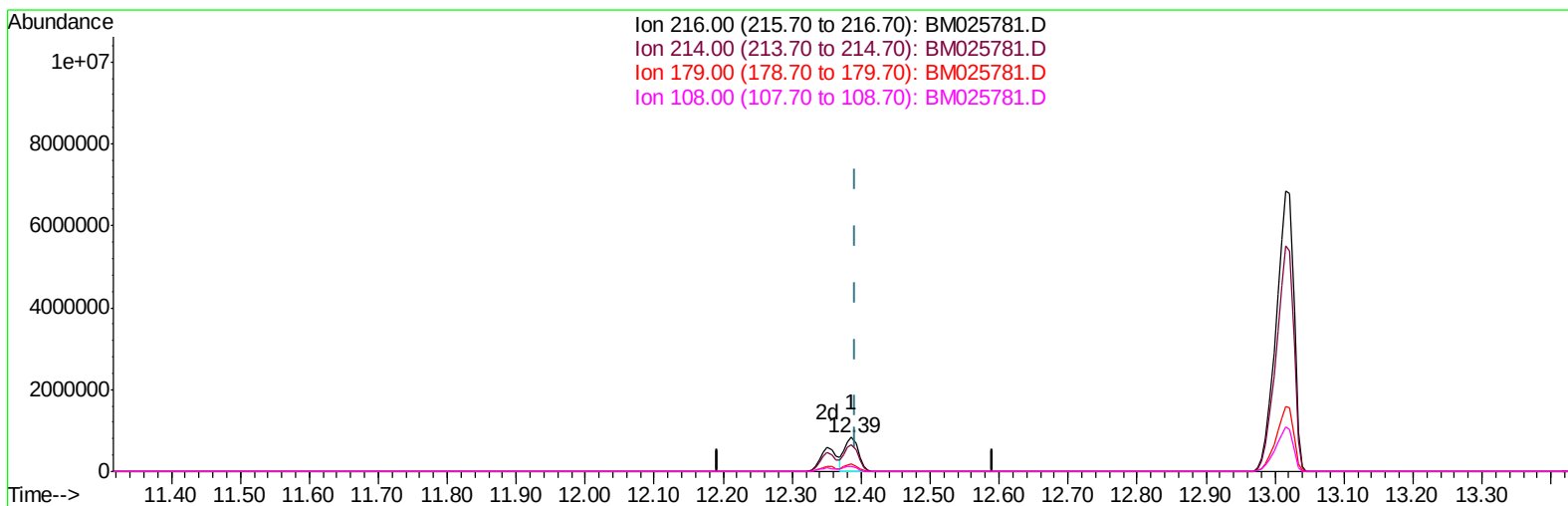
12.386min (-0.006) 135.31ng/ul m

response 2124264

Ion	Exp%	Act%
216.00	100	100
214.00	77.40	78.32
179.00	20.70	20.98
108.00	16.30	15.56

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032320\
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Acq On : 25 Mar 2020 23:06
Operator : CG/JU
Sample : L1941-06
Misc :
ALS Vial : 11 Sample Multiplier: 1

Quant Time: Mar 26 01:17:19 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM031420MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Mar 24 00:45:06 2020
Response via : Initial Calibration



TIC: BM025781.D

(37) 1,2,4,5-Tetrachlorobenzene

12.386min (-0.006) 73.75ng/ul m

response 1157777

Ion	Exp%	Act%
216.00	100	100
214.00	77.40	78.32
179.00	20.70	20.98
108.00	16.30	15.56

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032320\
 Data File : BM025781.D
 Acq On : 25 Mar 2020 23:06
 Operator : CG/JU
 Sample : L1941-06
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Quant Time: Mar 26 10:10:04 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM031420MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Mar 24 00:45:06 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.58	152	217397	20.00	ng/ul	0.00
18) Naphthalene-d8	10.36	136	807906	20.00	ng/ul	0.01
36) Acenaphthene-d10	14.20	164	496364	20.00	ng/ul	-0.01
62) Phenanthrene-d10	16.95	188	1044470	20.00	ng/ul	-0.01
78) Chrysene-d12	21.14	240	1157098	20.00	ng/ul	-0.01
86) Perylene-d12	23.29	264	1302367	20.00	ng/ul	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.15	96	7005	1.35	ng/uL	0.00
5) Phenol-d5	6.75	99	107174	6.25	ng/ul	-0.01
7) Bis-(2-Chloroethyl)ether-d	6.92	67	289845	29.60	ng/ul	0.00
9) 2-Chlorophenol-d4	7.11	132	348718	25.04	ng/ul	0.00
13) 4-Methylphenol-d8	8.27	113	207316	14.71	ng/ul	-0.01
19) Nitrobenzene-d5	8.71	128	218051	36.31	ng/ul	0.00
22) 2-Nitrophenol-d4	9.43	143	240420	37.66	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.97	165	411010	31.48	ng/ul	0.00
29) 4-Chloroaniline-d4	10.48	131	19415	1.28	ng/ul	0.00
44) Dimethylphthalate-d6	13.63	166	1384166	36.69	ng/ul	0.00
47) Acenaphthylene-d8	13.89	160	1724077	38.15	ng/ul	-0.01
52) 4-Nitrophenol-d4	14.41	143	40180	5.59	ng/ul	-0.01
58) Fluorene-d10	15.20	176	1253439	37.85	ng/ul	-0.01
63) 4,6-Dinitro-2-methylphenol	15.33	200	195406	30.66	ng/ul	-0.01
71) Anthracene-d10	17.05	188	1939286	39.75	ng/ul	0.00
79) Pyrene-d10	19.35	212	2112764	38.01	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.16	264	2618703	39.27	ng/ul	-0.01

Target Compounds				Ovalue		
28) Naphthalene	10.40	128	82088	1.866	ng/ul	99
37) 1,2,4,5-Tetrachlorobenzene	12.39	216	1157777m	73.748	ng/ul	
45) Dimethylphthalate	13.67	163	84421	2.150	ng/ul	100
73) 1,2,3,4-Tetrachlorobenzene	13.02	216	12145322	758.857	ng/uL	96
74) Pentachlorobenzene	14.53	250	461586	25.909	ng/uL	99

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

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