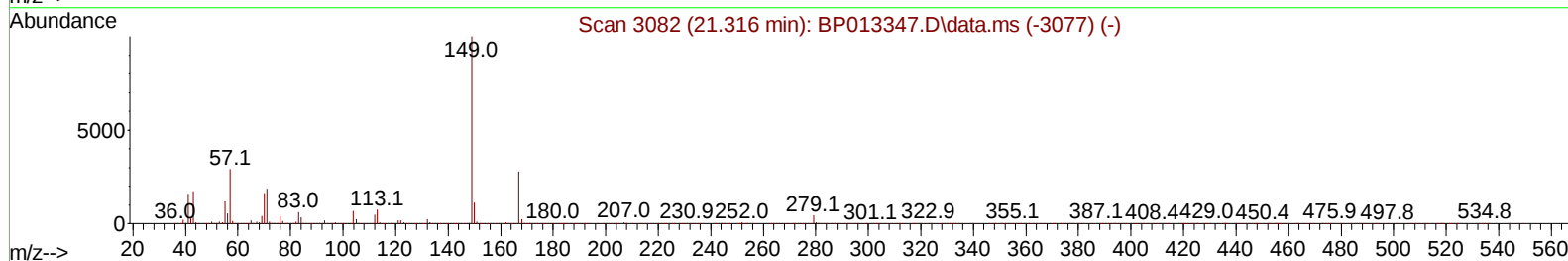
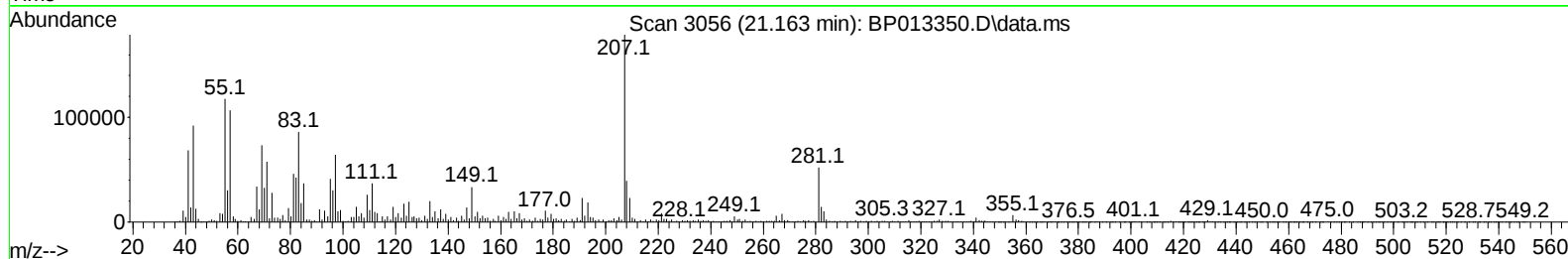
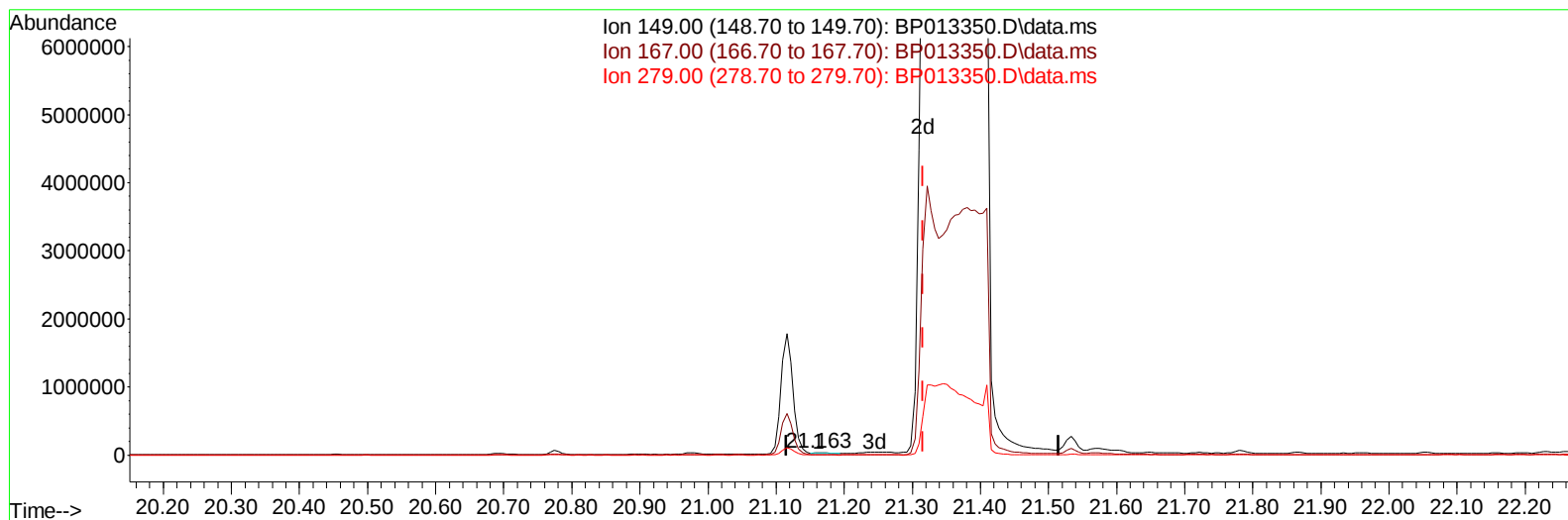


Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122822\
Data File : BP013350.D
Acq On : 28 Dec 2022 13:30
Operator : CG/JU
Sample : N6128-04
Misc :
ALS Vial : 5 Sample Multiplier: 1

Quant Time: Dec 29 00:17:47 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Dec 28 00:16:35 2022
Response via : Initial Calibration



TIC: BP013350.D\data.ms

(86) Bis(2-ethylhexyl)phthalate

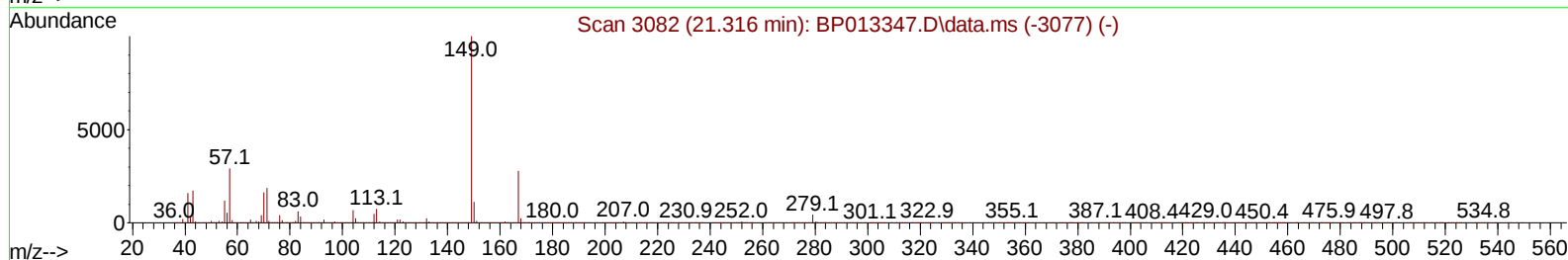
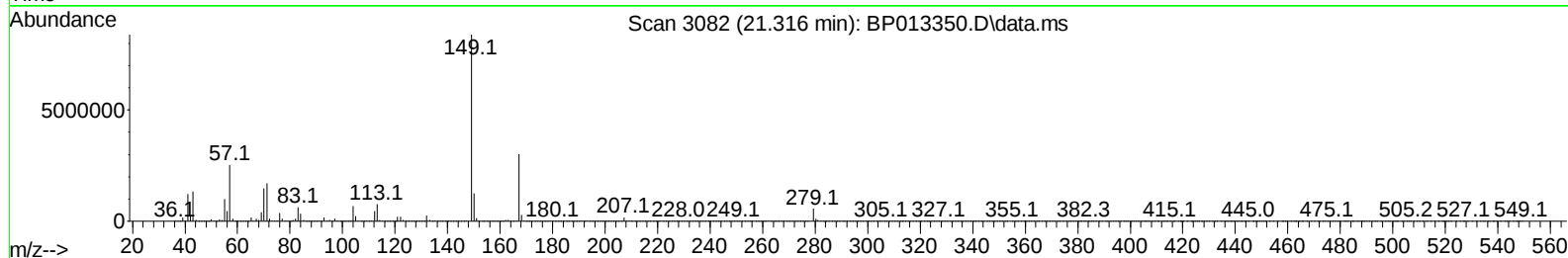
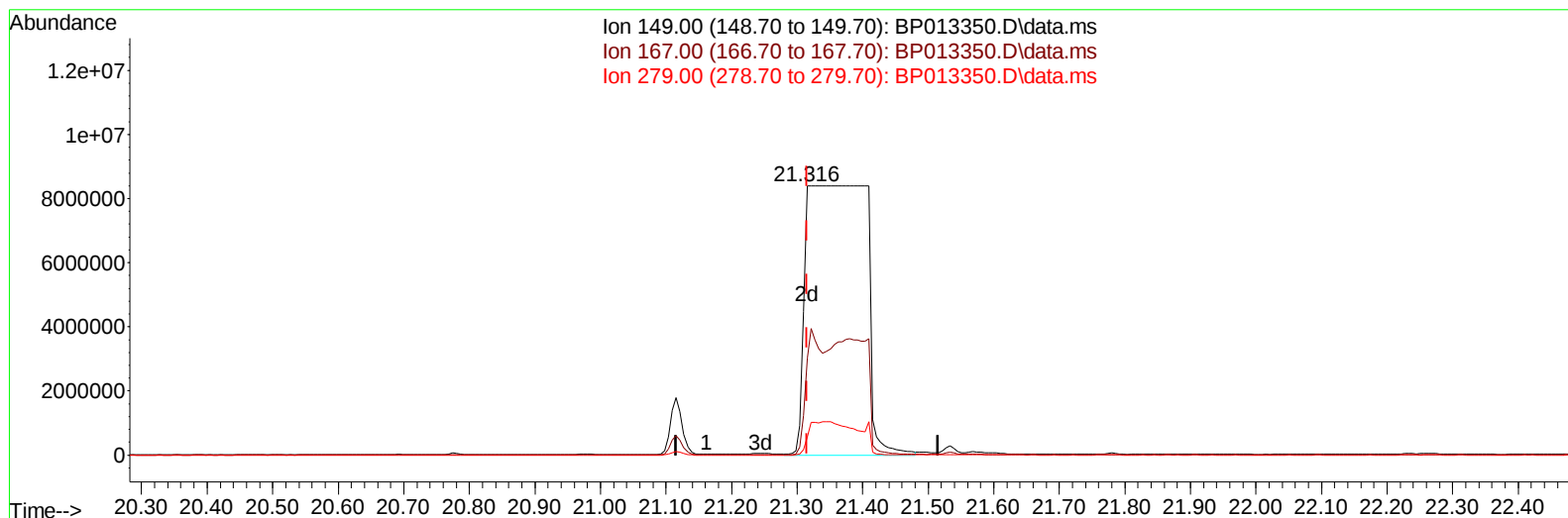
21.163min (-0.153) 0.13 ng/u1

response 11251

Ion	Exp%	Act%
149.00	100.00	100.00
167.00	28.10	25.79
279.00	4.80	3.97
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122822\
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Sample : N6128-04
Misc :
ALS Vial : 5 Sample Multiplier: 1

Quant Time: Dec 29 00:17:47 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION
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TIC: BP013350.D\data.ms

(86) Bis(2-ethylhexyl)phthalate

21.316min (-0.000) 627.57 ng/ul m

response 53545990

Ion	Exp%	Act%
149.00	100.00	100.00
167.00	28.10	36.05#
279.00	4.80	6.77#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122822\
 Data File : BP013350.D
 Acq On : 28 Dec 2022 13:30
 Operator : CG/JU
 Sample : N6128-04
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Quant Time: Dec 29 00:17:47 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
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 QLast Update : Wed Dec 28 00:16:35 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	568264	20.000	ng/ul	0.00
20) Naphthalene-d8	10.751	136	2467678	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.581	164	1654622	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.339	188	3743927	20.000	ng/ul	0.00
79) Chrysene-d12	21.427	240	2329783	20.000	ng/ul	0.00
88) Perylene-d12	23.904	264	2757213	20.000	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.322	96	41651	3.134	ng/uL	0.00
4) Pyridine-d5	3.752	84	606335	16.060	ng/ul	0.00
7) Phenol-d5	7.093	99	1045670	22.591	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.263	67	655875	23.490	ng/ul	0.00
11) 2-Chlorophenol-d4	7.463	132	829142	22.755	ng/ul	0.00
15) 4-Methylphenol-d8	8.646	113	876888	23.121	ng/ul	0.00
21) Nitrobenzene-d5	9.104	128	414426	22.858	ng/ul	0.00
24) 2-Nitrophenol-d4	9.828	143	469798	23.016	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.369	165	913513	23.041	ng/ul	0.00
31) 4-Chloroaniline-d4	10.887	131	1204328	21.517	ng/ul	0.00
46) Dimethylphthalate-d6	13.987	166	2840677	22.338	ng/ul	0.00
49) Acenaphthylene-d8	14.275	160	3248607	23.249	ng/ul	0.00
54) 4-Nitrophenol-d4	14.787	143	405515	17.728	ng/ul	0.00
60) Fluorene-d10	15.575	176	2521325	23.436	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.710	200	372426	15.458	ng/ul	0.02
73) Anthracene-d10	17.445	188	3748815	22.195	ng/ul	0.01
81) Pyrene-d10	19.674	212	3923109	29.336	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.745	264	3256935	23.699	ng/ul	0.01
Target Compounds						
					Qvalue	
86) Bis(2-ethylhexyl)phtha...	21.316	149	53545990m	627.572	ng/ul	

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

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