

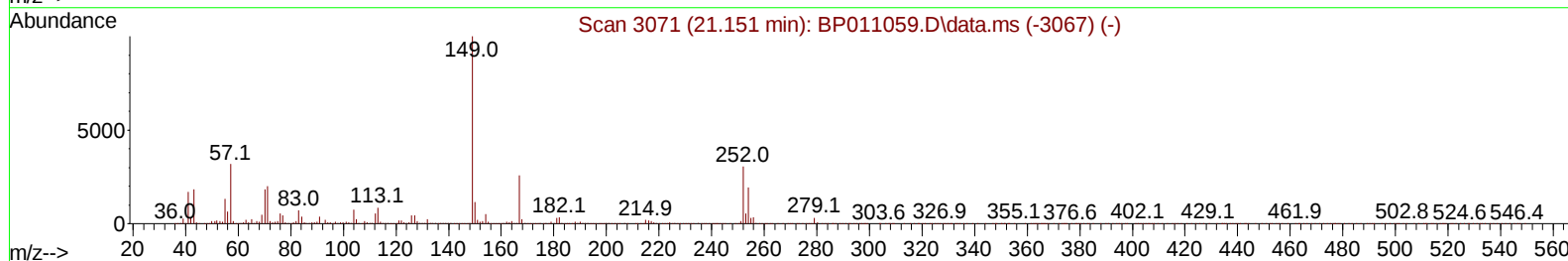
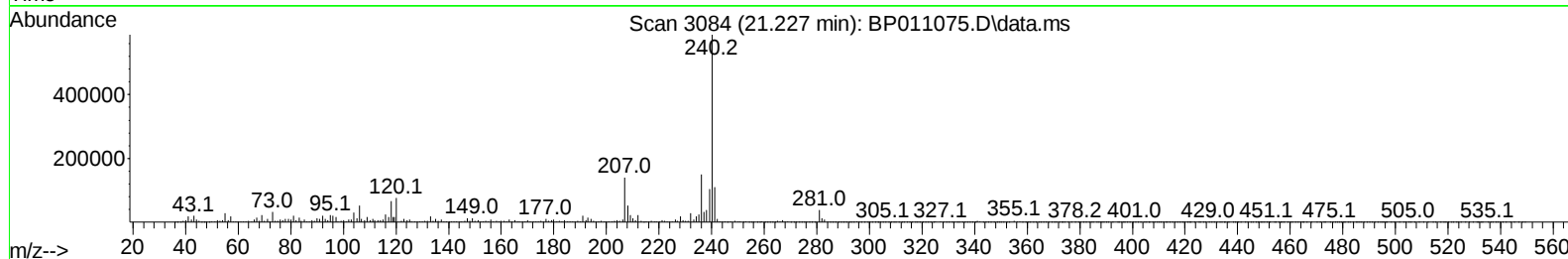
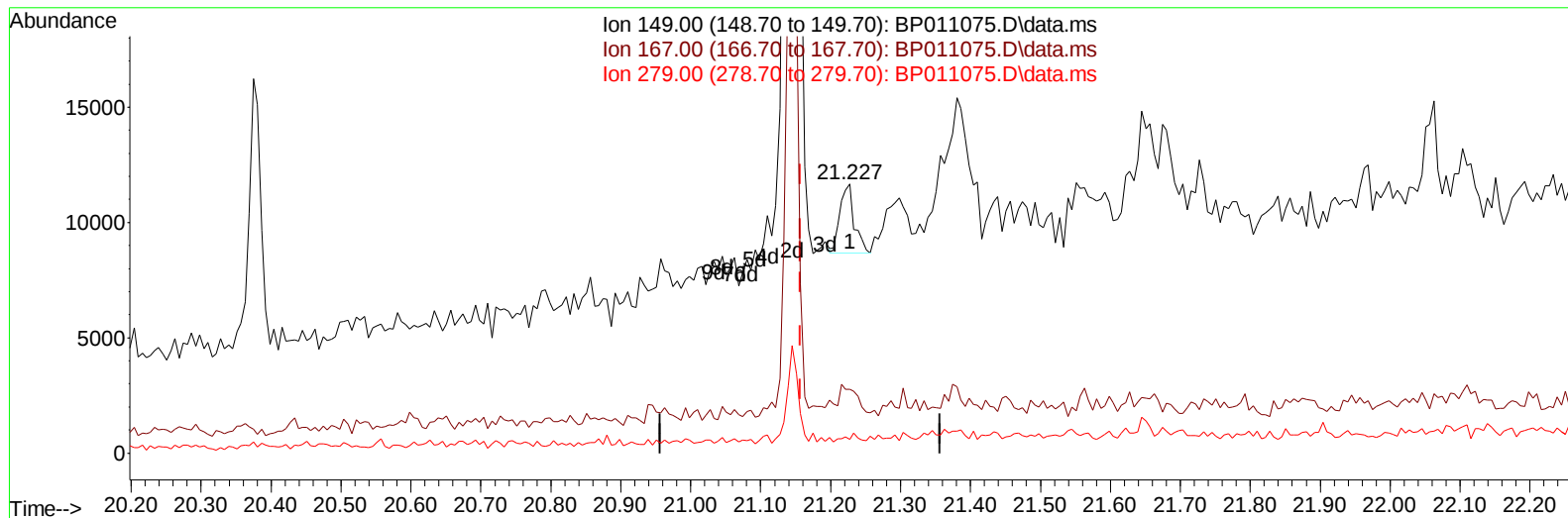
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072122\
 Data File : BP011075.D
 Acq On : 21 Jul 2022 23:06
 Operator : CG/JU
 Sample : N3709-15
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 H0B29

Manual IntegrationsAPPROVED

Quant Time: Jul 22 00:44:50 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jul 19 03:25:54 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 07/22/2022
 Supervised By :mohammad ahmed 07/22/2022



TIC: BP011075.D\data.ms

(86) Bis(2-ethylhexyl)phthalate

21.227min (+ 0.071) 0.08 ng/u1

response 4332

Ion	Exp%	Act%
149.00	100.00	100.00
167.00	28.20	23.81
279.00	5.30	5.15
0.00	0.00	0.00

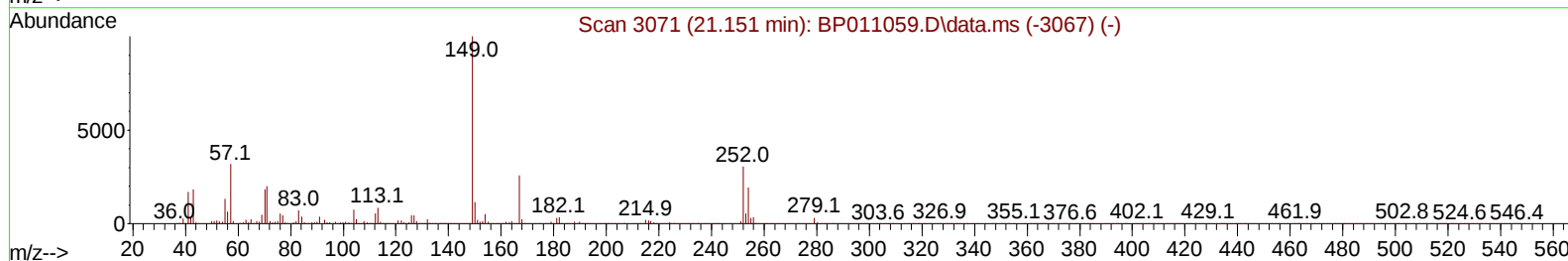
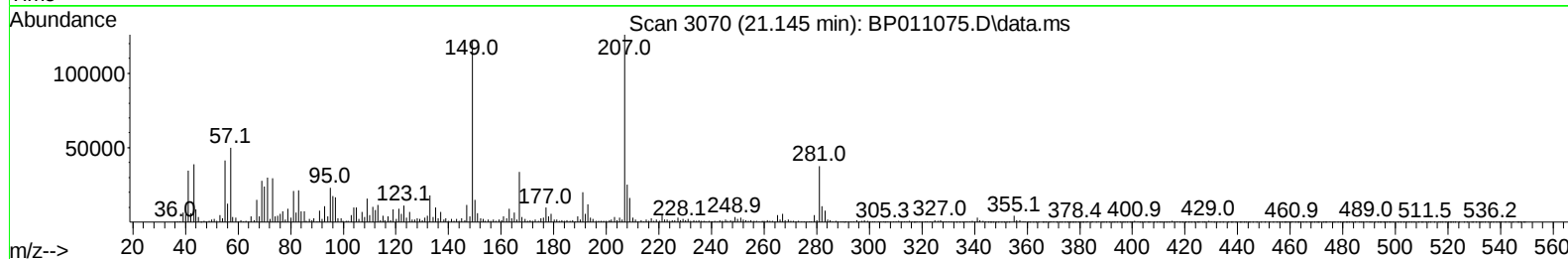
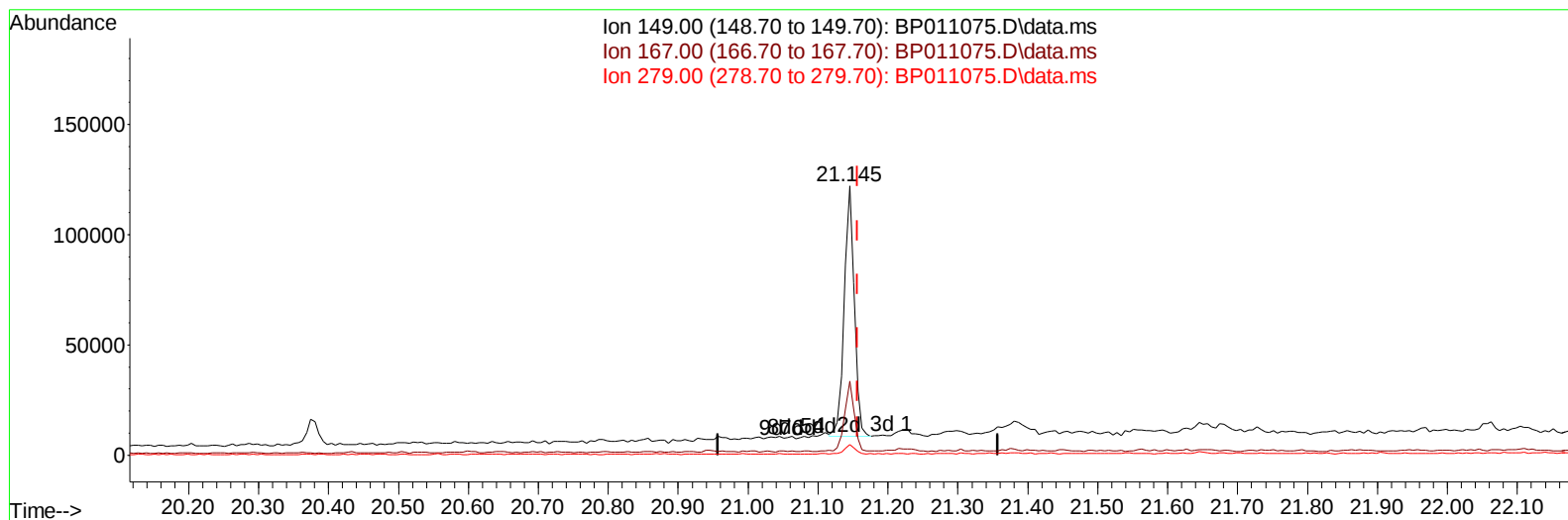
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(86) Bis(2-ethylhexyl)phthalate

21.145min (-0.012) 2.01 ng/ul m

response 114017

Ion	Exp%	Act%
149.00	100.00	100.00
167.00	28.20	27.48
279.00	5.30	3.82#
0.00	0.00	0.00

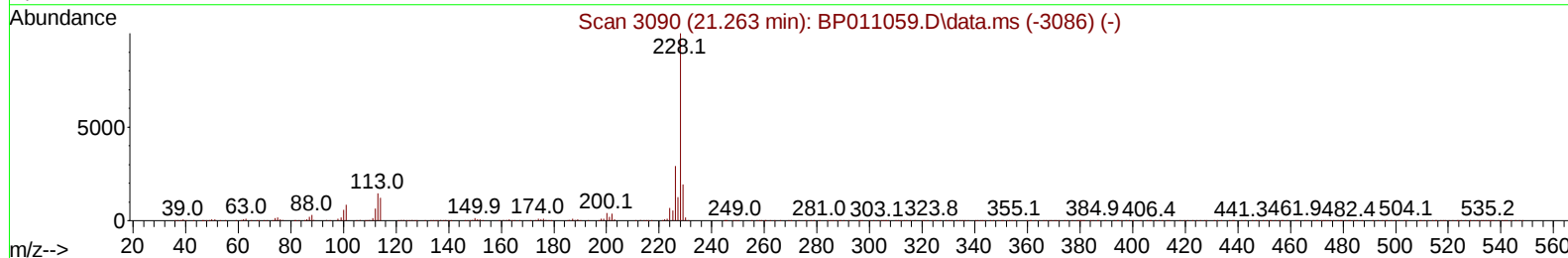
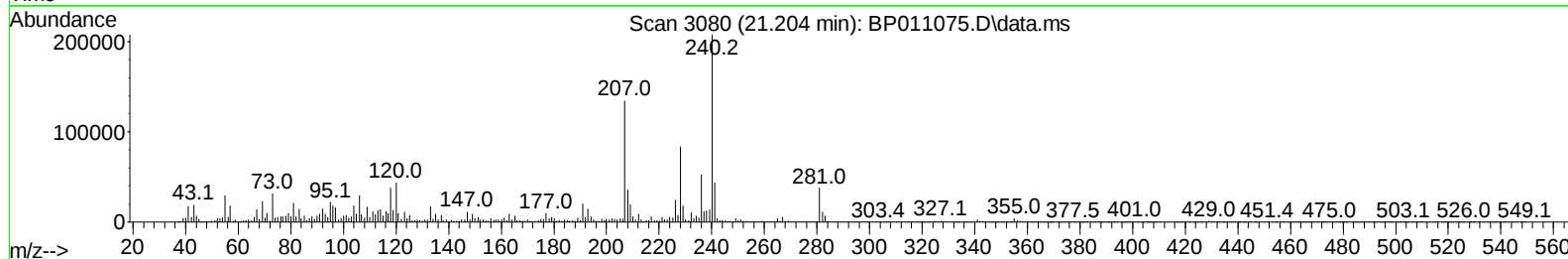
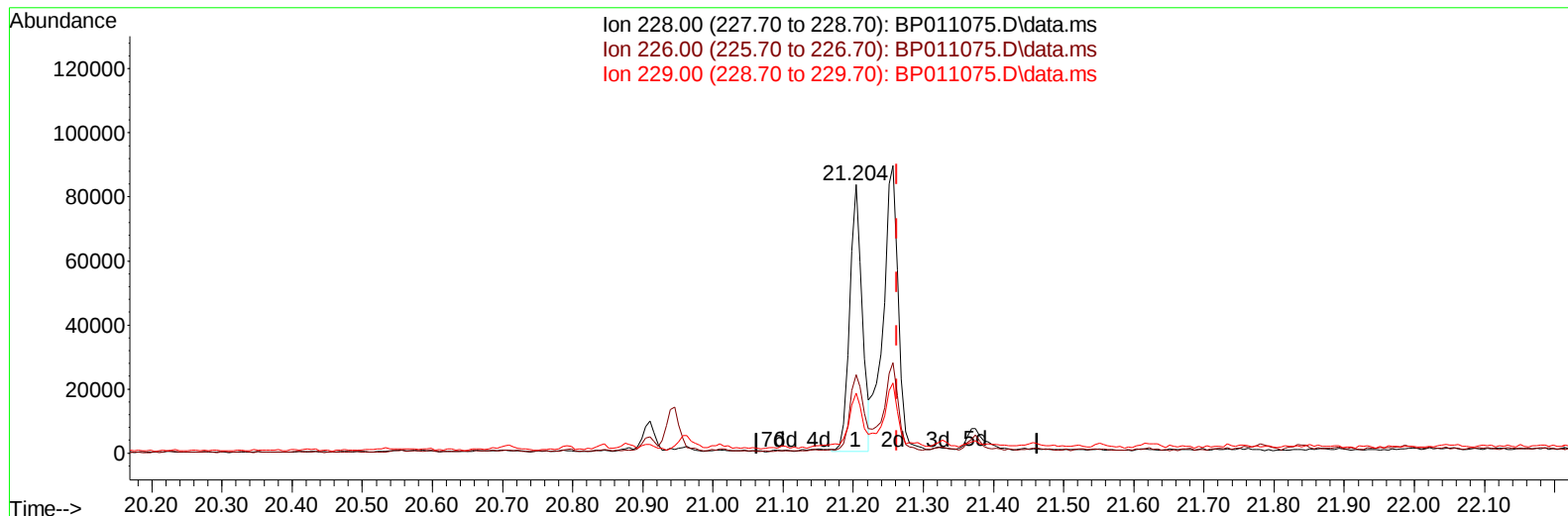
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TIC: BP011075.D\data.ms

(87) Chrysene

21.204min (-0.059) 1.28 ng/u1

response 102723

Ion	Exp%	Act%
228.00	100.00	100.00
226.00	29.50	29.24
229.00	19.50	22.25
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	7.669	152	434775	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.463	136	1778436	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.334	164	938616	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.098	188	1643603	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.222	240	1344946	20.000	ng/ul	# 0.00
88) Perylene-d12	23.569	264	1471270	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.175	96	22957	1.904	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	6.846	99	420689	11.788	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.011	67	303213	13.212	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.205	132	360455	12.499	ng/ul	0.00
15) 4-Methylphenol-d8	8.387	113	281487	9.852	ng/ul	-0.01
21) Nitrobenzene-d5	8.828	128	173748	13.406	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.552	143	152425	11.886	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.087	165	284725	12.140	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.610	131	289545	7.489	ng/ul	0.00
46) Dimethylphthalate-d6	13.751	166	911485	14.208	ng/ul	-0.01
49) Acenaphthylene-d8	14.022	160	1144000	15.114	ng/ul	0.00
54) 4-Nitrophenol-d4	14.551	143	44758	3.447	ng/ul	0.00
60) Fluorene-d10	15.334	176	786942	14.432	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.457	200	31189	3.835	ng/ul	0.00
73) Anthracene-d10	17.198	188	1036373	14.832	ng/ul	0.00
81) Pyrene-d10	19.451	212	1116763	15.810	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.416	264	1119180	15.838	ng/ul	-0.01
Target Compounds						
					Qvalue	
80) Fluoranthene	19.128	202	173682	2.102	ng/ul#	88
82) Pyrene	19.480	202	199869	2.345	ng/ul#	86
85) Benzo(a)anthracene	21.204	228	100954	1.233	ng/ul	93
86) Bis(2-ethylhexyl)phtha...	21.145	149	114017m	2.010	ng/ul	
87) Chrysene	21.257	228	134788m	1.676	ng/ul	
90) Benzo(b)fluoranthene	22.851	252	172711	1.901	ng/ul#	76
93) Benzo(a)pyrene	23.463	252	127773	1.465	ng/ul#	82
96) Benzo(g,h,i)perylene	26.739	276	134076	1.553	ng/ul#	82

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

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