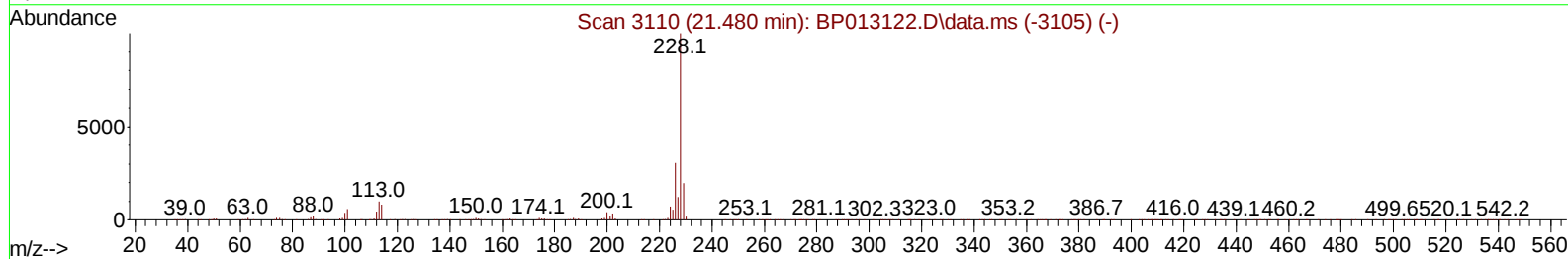
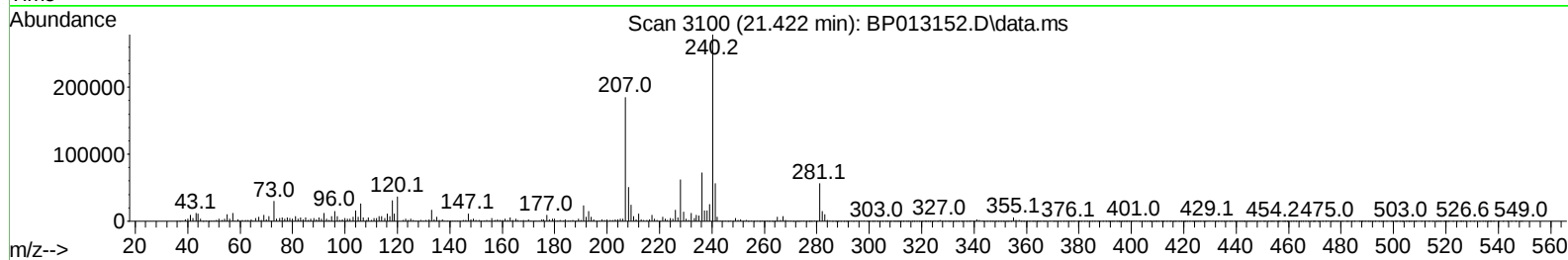
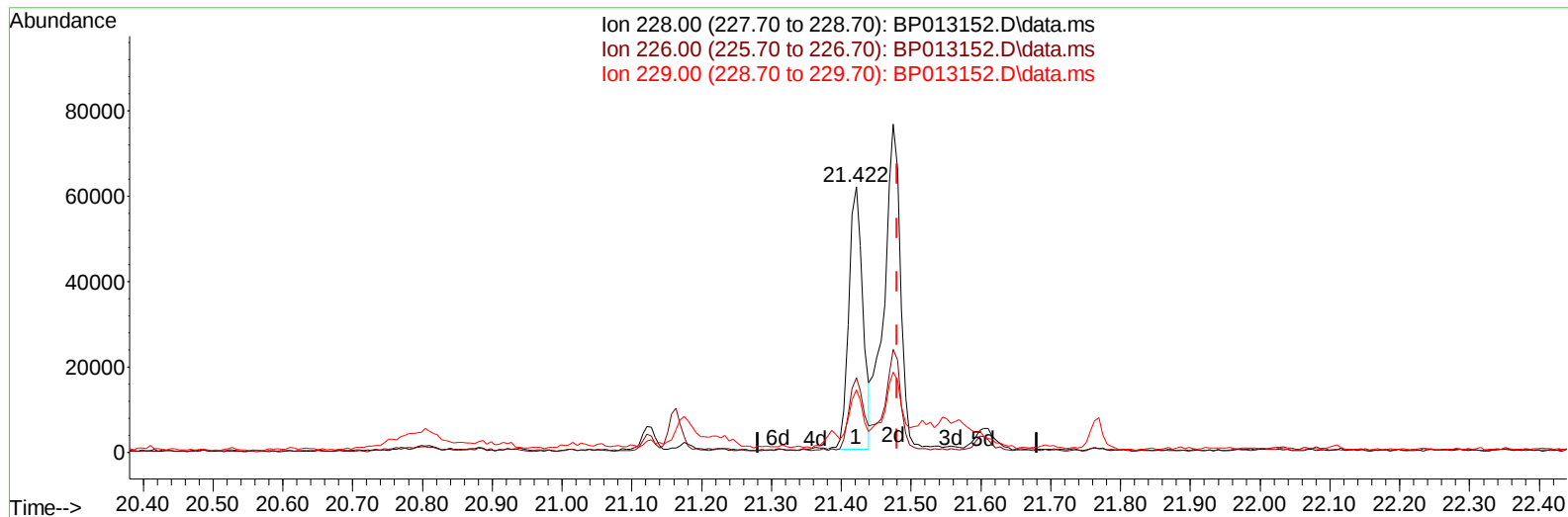


Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013152.D
Acq On : 20 Dec 2022 06:16
Operator : CG/JU
Sample : N6006-09
Misc :
ALS Vial : 32 Sample Multiplier: 1

Quant Time: Dec 20 06:52:02 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Dec 19 21:43:11 2022
Response via : Initial Calibration



TIC: BP013152.D\data.ms

(87) Chrysene

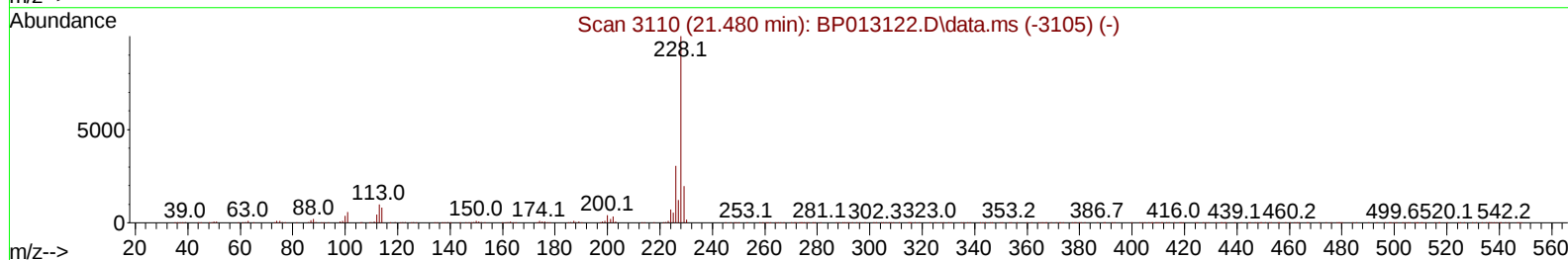
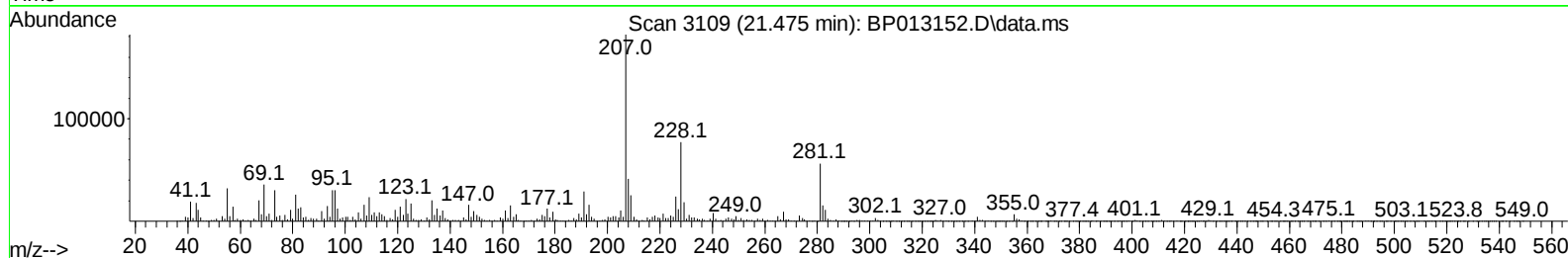
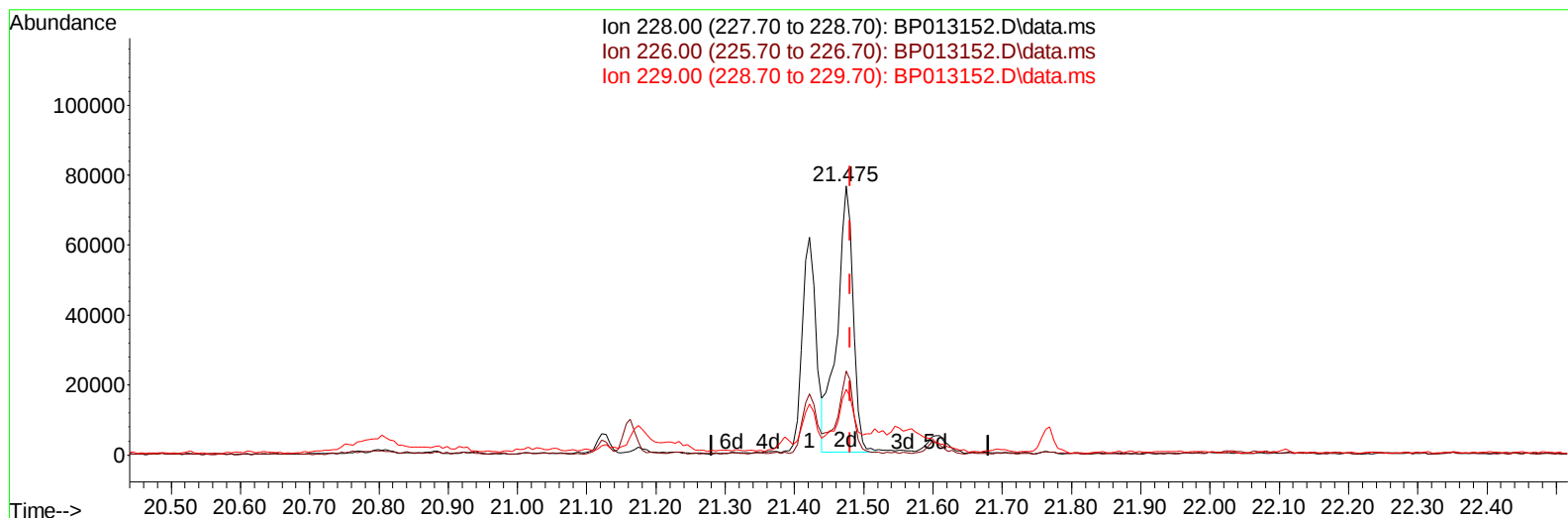
21.422min (-0.059) 0.75 ng/u1

response 86373

Ion	Exp%	Act%
228.00	100.00	100.00
226.00	30.60	28.04
229.00	20.00	23.43
0.00	0.00	0.00

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TIC: BP013152.D\data.ms

(87) Chrysene

21.475min (-0.006) 1.08 ng/ul m

response 123942

Ion	Exp%	Act%
228.00	100.00	100.00
226.00	30.60	31.36
229.00	20.00	24.46#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
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 Operator : CG/JU
 Sample : N6006-09
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Quant Time: Dec 20 06:52:02 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.952	152	475158	20.000	ng/ul	0.00
20) Naphthalene-d8	10.769	136	2020397	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.593	164	1248930	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.351	188	2439159	20.000	ng/ul	0.00
79) Chrysene-d12	21.439	240	1826518	20.000	ng/ul	0.00
88) Perylene-d12	23.916	264	2204627	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.334	96	97826	8.803	ng/uL	0.00
4) Pyridine-d5	3.764	84	926842	29.360	ng/ul	0.00
7) Phenol-d5	7.111	99	2139343	55.276	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.281	67	1327120	56.843	ng/ul	0.00
11) 2-Chlorophenol-d4	7.481	132	1722639	56.541	ng/ul	0.00
15) 4-Methylphenol-d8	8.663	113	1640438	51.729	ng/ul	0.00
21) Nitrobenzene-d5	9.122	128	889543	59.925	ng/ul	0.00
24) 2-Nitrophenol-d4	9.846	143	1030609	61.669	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.387	165	1821727	56.121	ng/ul	0.00
31) 4-Chloroaniline-d4	10.905	131	1677567	36.607	ng/ul	0.00
46) Dimethylphthalate-d6	14.004	166	5440885	56.684	ng/ul	0.00
49) Acenaphthylene-d8	14.293	160	6185507	58.646	ng/ul	0.00
54) 4-Nitrophenol-d4	14.793	143	761887	44.127	ng/ul	0.00
60) Fluorene-d10	15.592	176	4531055	55.798	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.710	200	779720	49.675	ng/ul	0.00
73) Anthracene-d10	17.451	188	6454751	58.657	ng/ul	0.00
81) Pyrene-d10	19.681	212	6829788	65.144	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.757	264	6722842	61.179	ng/ul	0.00
Target Compounds						
					Qvalue	
80) Fluoranthene	19.357	202	146068	1.228	ng/ul	97
82) Pyrene	19.710	202	133065	1.084	ng/ul	98
87) Chrysene	21.475	228	123942m	1.081	ng/ul	
90) Benzo(b)fluoranthene	23.157	252	201078	1.430	ng/ul#	92

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

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