

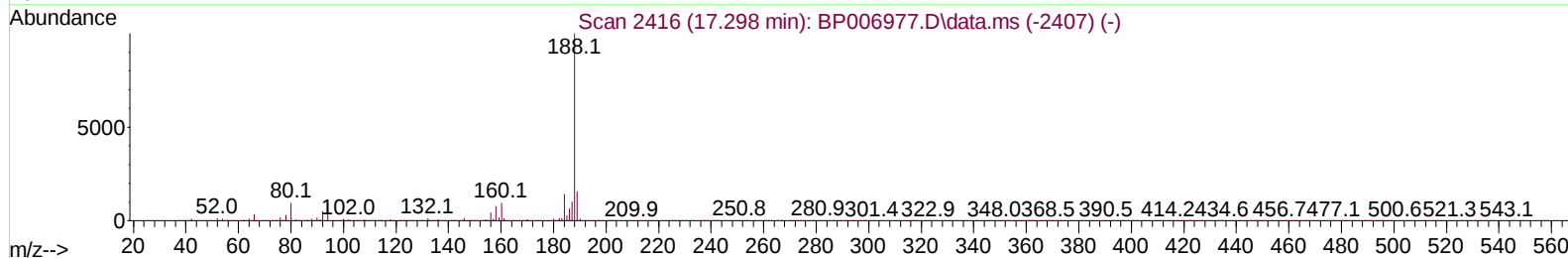
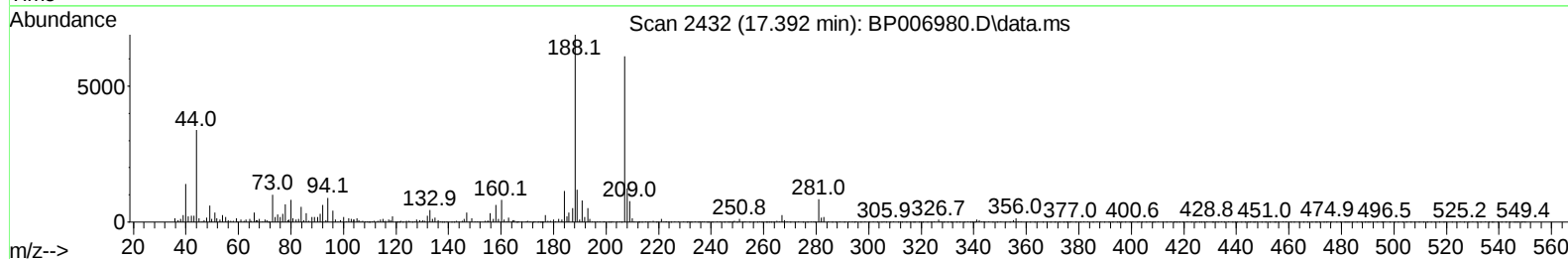
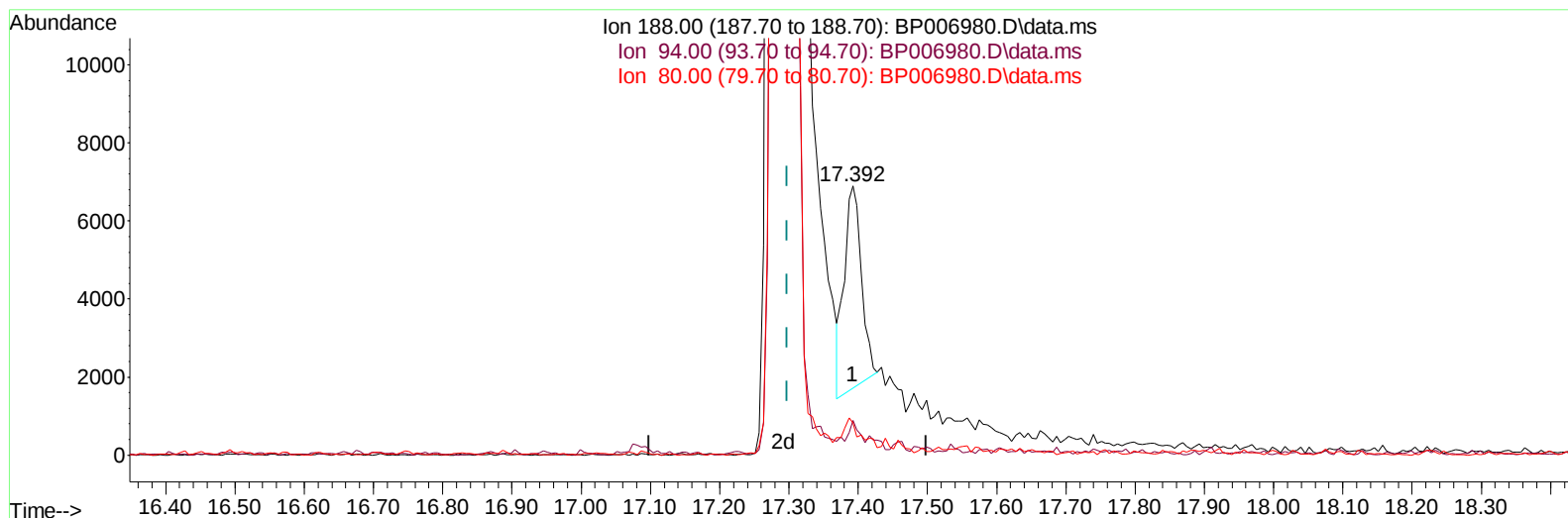
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091521\
Data File : BP006980.D
Acq On : 15 Sep 2021 17:56
Operator : CG/JU
Sample : M3708-10
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SVOC-GPC2-BLANK

Manual Integrations
APPROVED

mohammad
9/16/2021 11:44:01 AM

Quant Time: Sep 15 18:26:18 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP090821.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Sep 15 17:01:19 2021
Response via : Initial Calibration



TIC: BP006980.D\data.ms

(64) Phenanthrene-d10 (I)

17.392min (+ 0.094) 20.00 ng/u1

response 9079

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	11.50	12.90
80.00	13.50	13.97
0.00	0.00	0.00

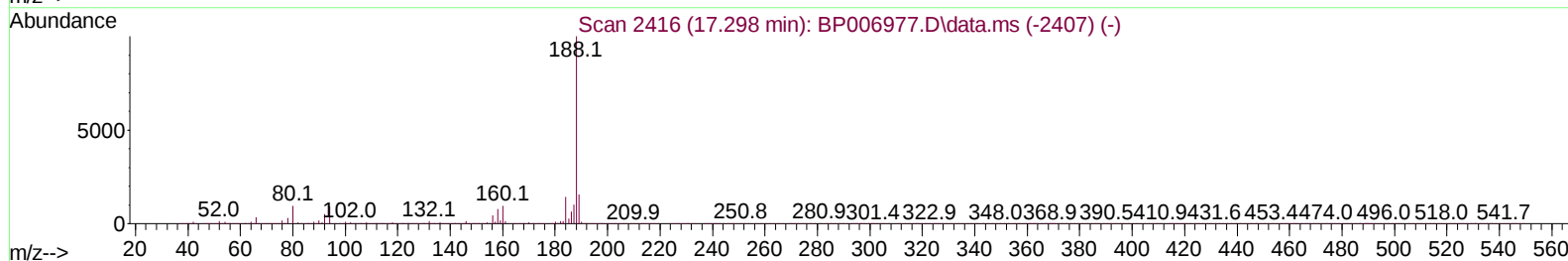
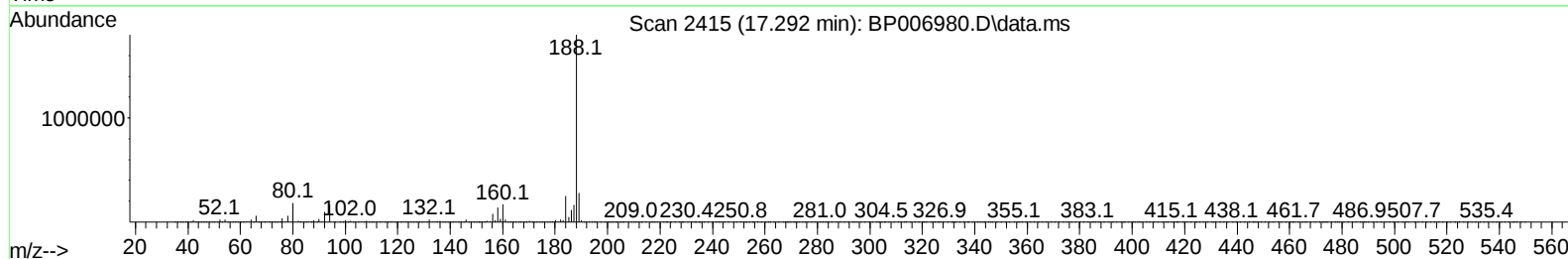
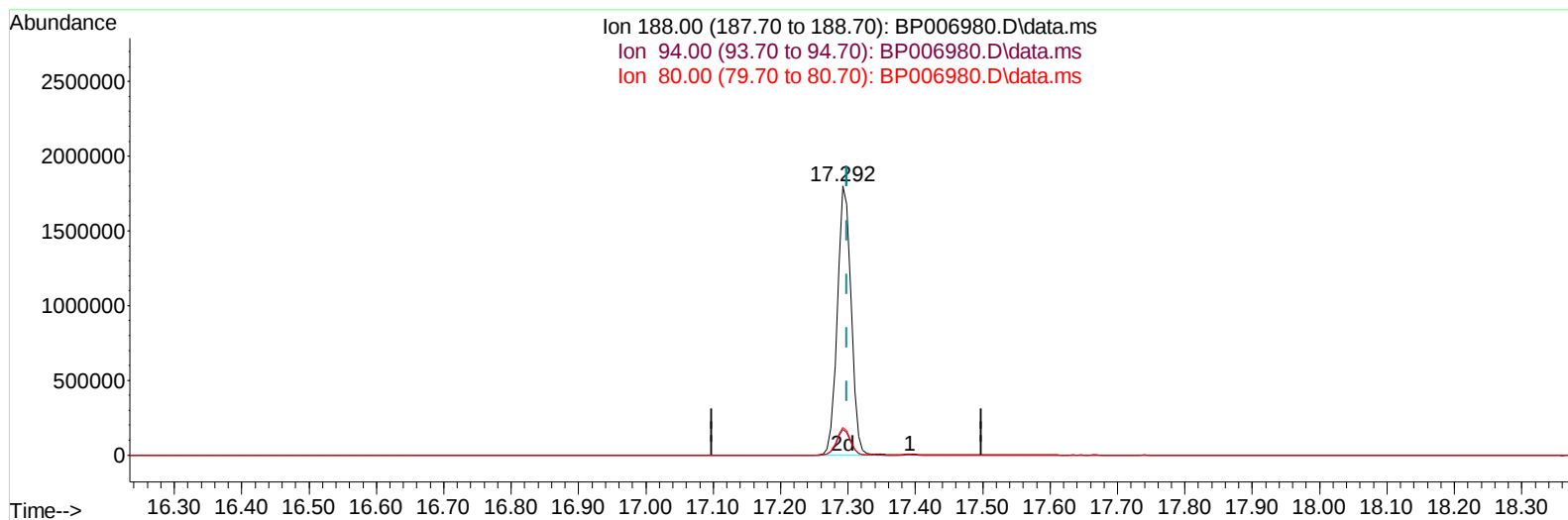
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TIC: BP006980.D\data.ms

(64) Phenanthrene-d10 (I)

17.292min (-0.006) 20.00 ng/ul m

response 2558346

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	11.50	9.59
80.00	13.50	10.28#
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	7.922	152	418834	20.000	ng/ul	0.00
20) Naphthalene-d8	10.722	136	1732212	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.546	164	1190055	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.292	188	2558346m	20.000	ng/ul	0.00
79) Chrysene-d12	21.375	240	2213750	20.000	ng/ul	0.00
88) Perylene-d12	23.798	264	2042290	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	0.000	96	0d	0.000	ng/uL	
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	0.000	99	0d	0.000	ng/ul	
9) Bis-(2-Chloroethyl)eth...	0.000	67	0d	0.000	ng/ul	
11) 2-Chlorophenol-d4	0.000	132	0d	0.000	ng/ul	
15) 4-Methylphenol-d8	0.000	113	0d	0.000	ng/ul	
21) Nitrobenzene-d5	0.000	128	0d	0.000	ng/ul	
24) 2-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
28) 2,4-Dichlorophenol-d3	0.000	165	0d	0.000	ng/ul	
31) 4-Chloroaniline-d4	0.000	131	0d	0.000	ng/ul	
46) Dimethylphthalate-d6	0.000	166	0d	0.000	ng/ul	
49) Acenaphthylene-d8	0.000	160	0d	0.000	ng/ul	
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
60) Fluorene-d10	0.000	176	0d	0.000	ng/ul	
65) 4,6-Dinitro-2-methylph...	0.000	200	0d	0.000	ng/ul	
73) Anthracene-d10	0.000	188	0d	0.000	ng/ul	
81) Pyrene-d10	0.000	212	0d	0.000	ng/ul	
92) Benzo(a)pyrene-d12	0.000	264	0d	0.000	ng/ul	

Target Compounds Qvalue

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

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