

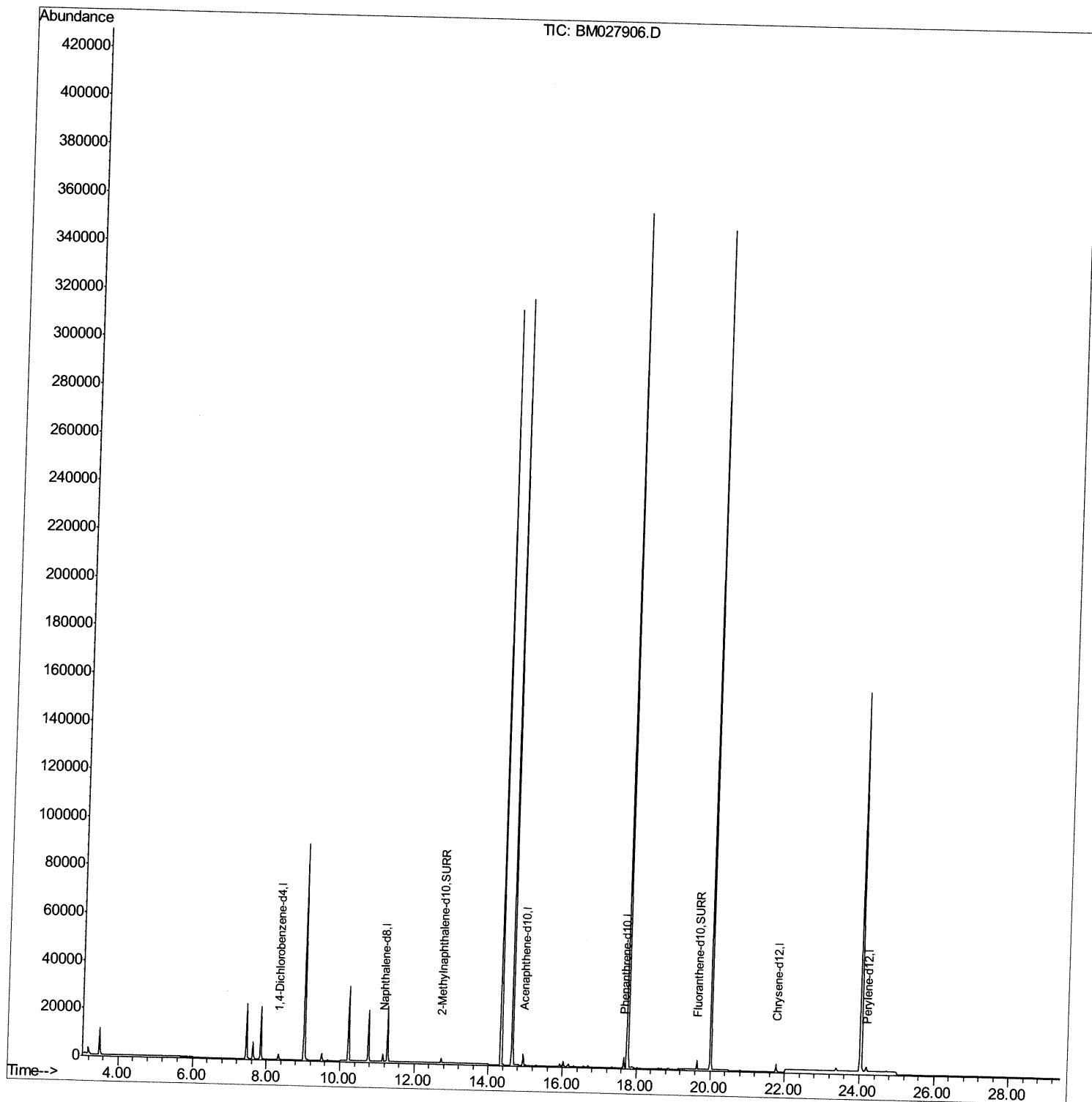
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM112320\
Data File : BM027906.D
Acq On : 23 Nov 2020 10:43
Operator : JU/CG
Sample : PB133151BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SBLK51

Manual Integrations
APPROVED

mohammad
11/25/2020 12:44:07 PM

Quant Time: Nov 23 11:38:46 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM112020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Nov 23 11:35:49 2020
Response via : Initial Calibration



Quantitation Report (Qedit)

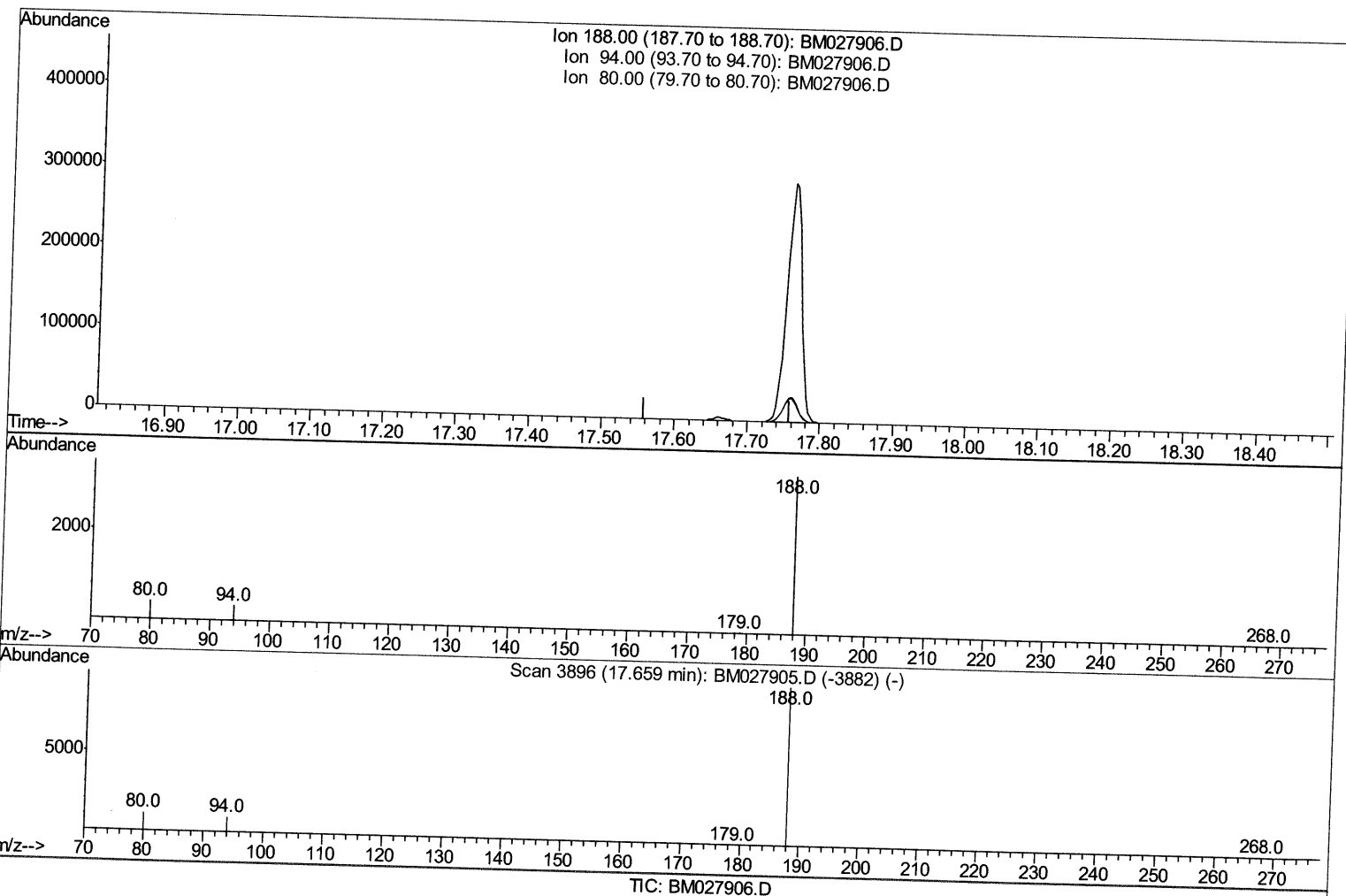
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(10) Phenanthrene-d10 (I)
 17.659min (-17.659) 0.00ng/ul
 response 0

Ion	Exp%	Act%
188.00	100	0.00
94.00	9.30	0.00#
80.00	10.20	0.00#
0.00	0.00	0.00

Quantitation Report (Qedit)

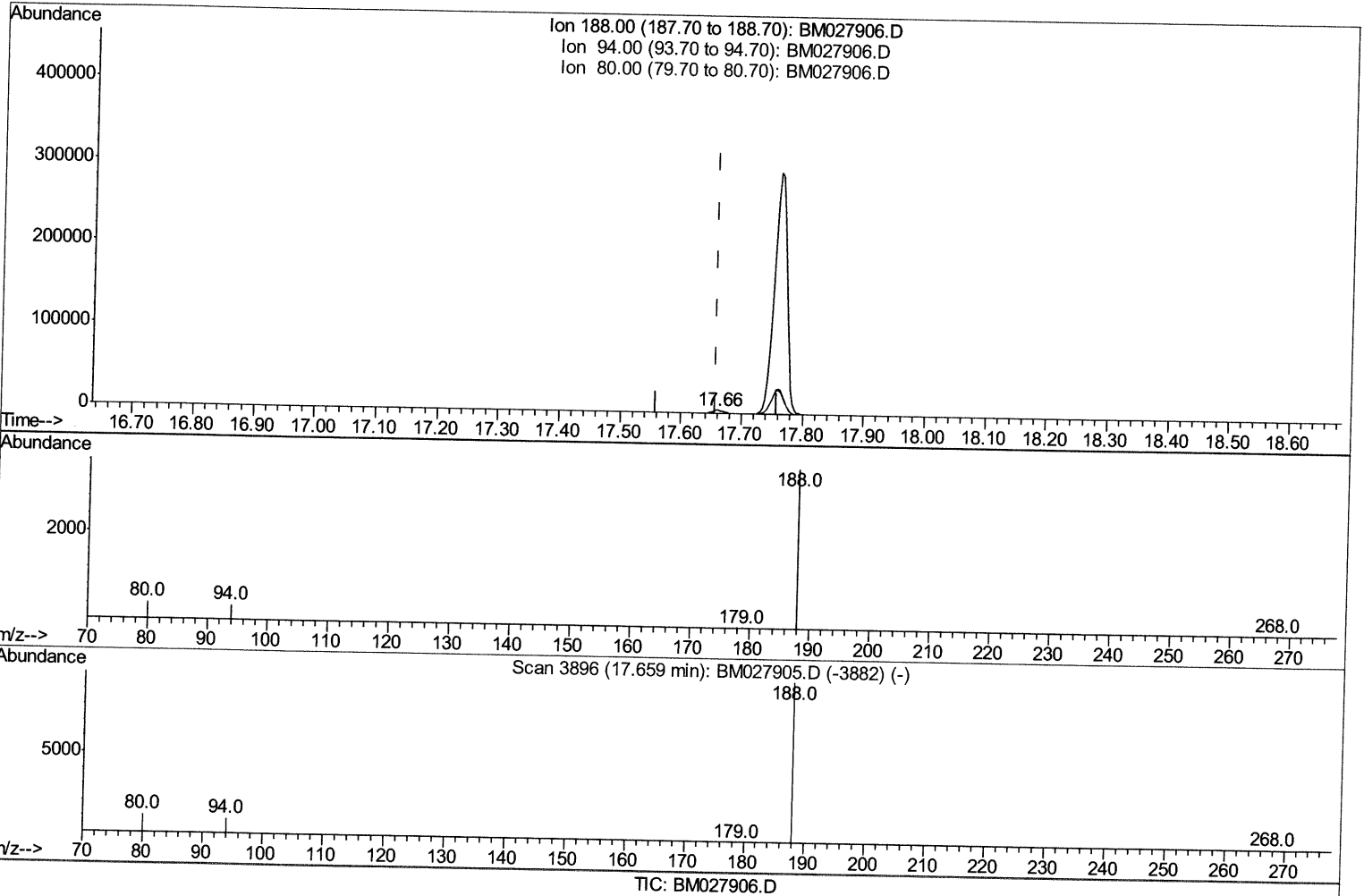
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(10) Phenanthrene-d10 (I)

17.662min (+0.003) 0.40ng/ul m 11/25/20 JU

response 5088

Ion	Exp%	Act%
188.00	100	100
94.00	9.30	10.16
80.00	10.20	11.71
0.00	0.00	0.00

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.33	152	1037	0.40	ng/ul	0.00
2) Naphthalene-d8	11.17	136	4189	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.94	164	2687	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.66	188	5088m >	0.40	ng/ul >	0.00
16) Chrysene-d12	21.77	240	2851	0.40	ng/ul	0.00
20) Perylene-d12	24.19	264	2370	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.73	152	2403	0.34	ng/ul	0.00
14) Fluoranthene-d10	19.65	212	4065	0.30	ng/ul	0.00

11/25/20 JU

Target Compounds

Ovalue

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