

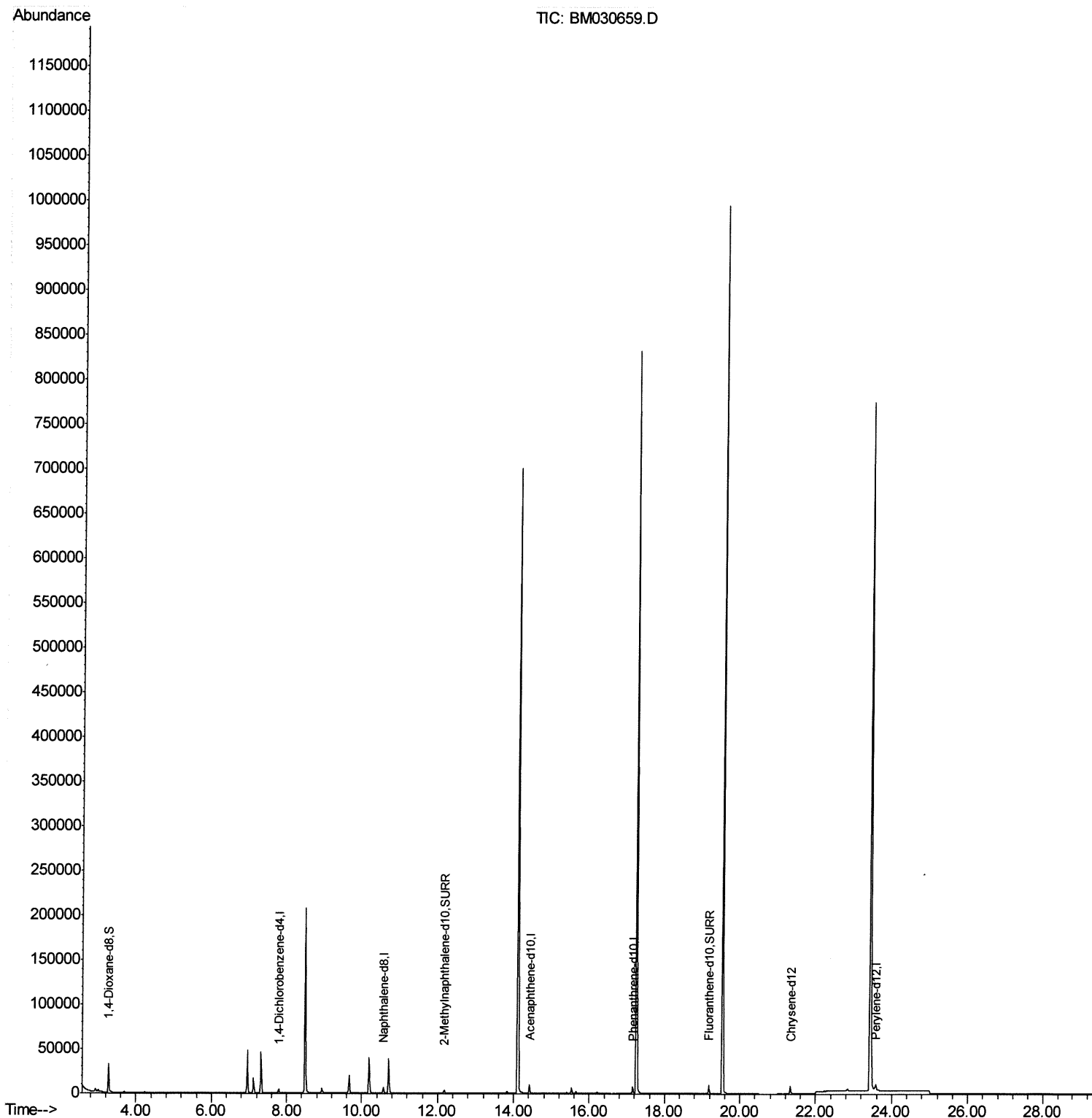
Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062821\
Data File : BM030659.D
Acq On : 28 Jun 2021 18:37
Operator : CG/JU
Sample : PB137370BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SBLK370

Manual Integrations
APPROVED

mohammad
6/29/2021 2:31:26 PM

Quant Time: Jun 28 19:08:35 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SFAM-EPA-SIM-BM062821.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 28 14:33:12 2021
Response via : Initial Calibration



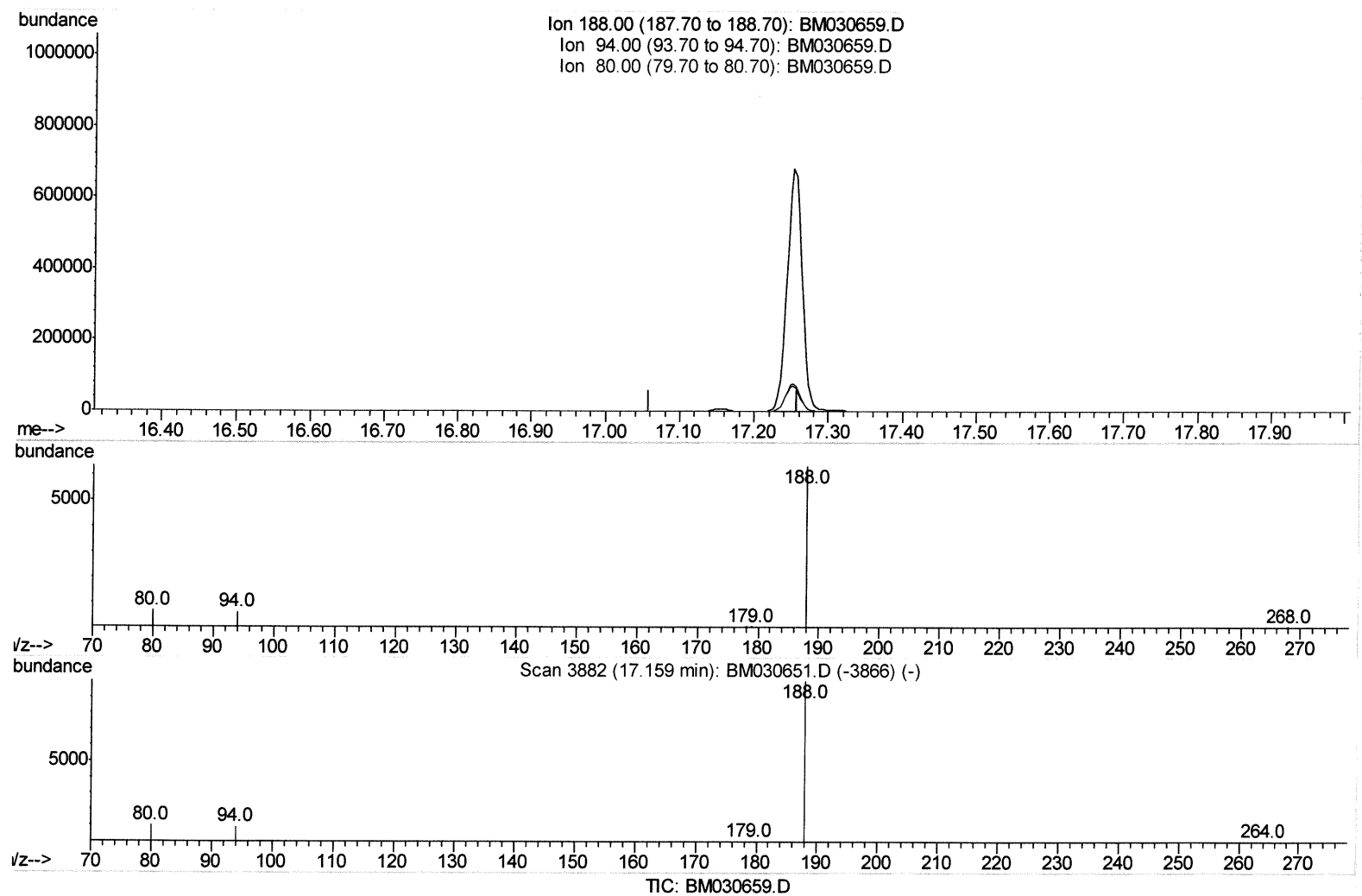
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(13) Phenanthrene-d10 (I)

17.159min (-17.159) 0.00ng/ul

response 0

Ion	Exp%	Act%
188.00	100	0.00
94.00	10.80	0.00#
80.00	11.40	0.00#
0.00	0.00	0.00

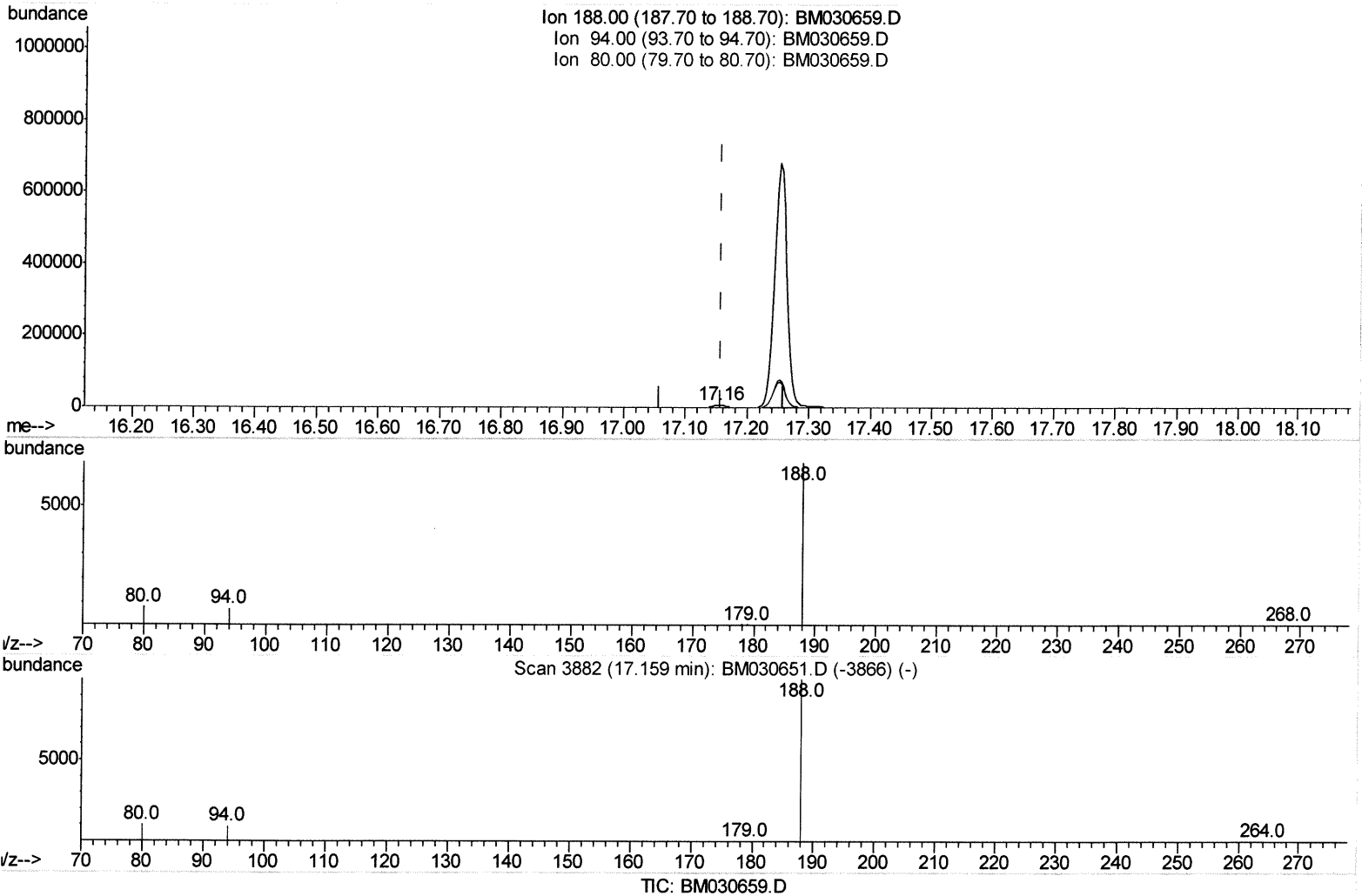
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(13) Phenanthrene-d10 (I)

17.155min (-0.004) 0.40ng/ul m *JY 6/30/21*

response 9927

Ion	Exp%	Act%
188.00	100	100
94.00	10.80	10.38
80.00	11.40	11.31
0.00	0.00	0.00

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Manual Integrations
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.79	152	2331	0.40	ng/ul	0.00
4) Naphthalene-d8	10.57	136	8766	0.40	ng/ul	0.00
9) Acenaphthene-d10	14.42	164	5397	0.40	ng/ul	0.00
13) Phenanthrene-d10	17.16	188	9927m	0.40	ng/ul	0.00
17) Chrysene-d12	21.33	240	8353	0.40	ng/ul	0.00
23) Perylene-d12	23.58	264	10249	0.40	ng/ul	0.00

> J4 b) 20/21

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.28	96	20148	7.93	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.17	152	5982	0.43	ng/ul	0.00
18) Fluoranthene-d10	19.18	212	11357	0.47	ng/ul	0.00

Target Compounds	Ovalue

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