

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030320\

Data File : BM025259.D

Acq On : 04 Mar 2020 20:34

Operator : CG/JU

Sample : L1516-08

Misc :

ALS Vial : 55 Sample Multiplier: 1

Quant Time: Mar 05 01:16:20 2020

Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM030320.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Wed Mar 04 08:02:03 2020

Response via : Initial Calibration

Instrument :

BNA_M

ClientSampleId :

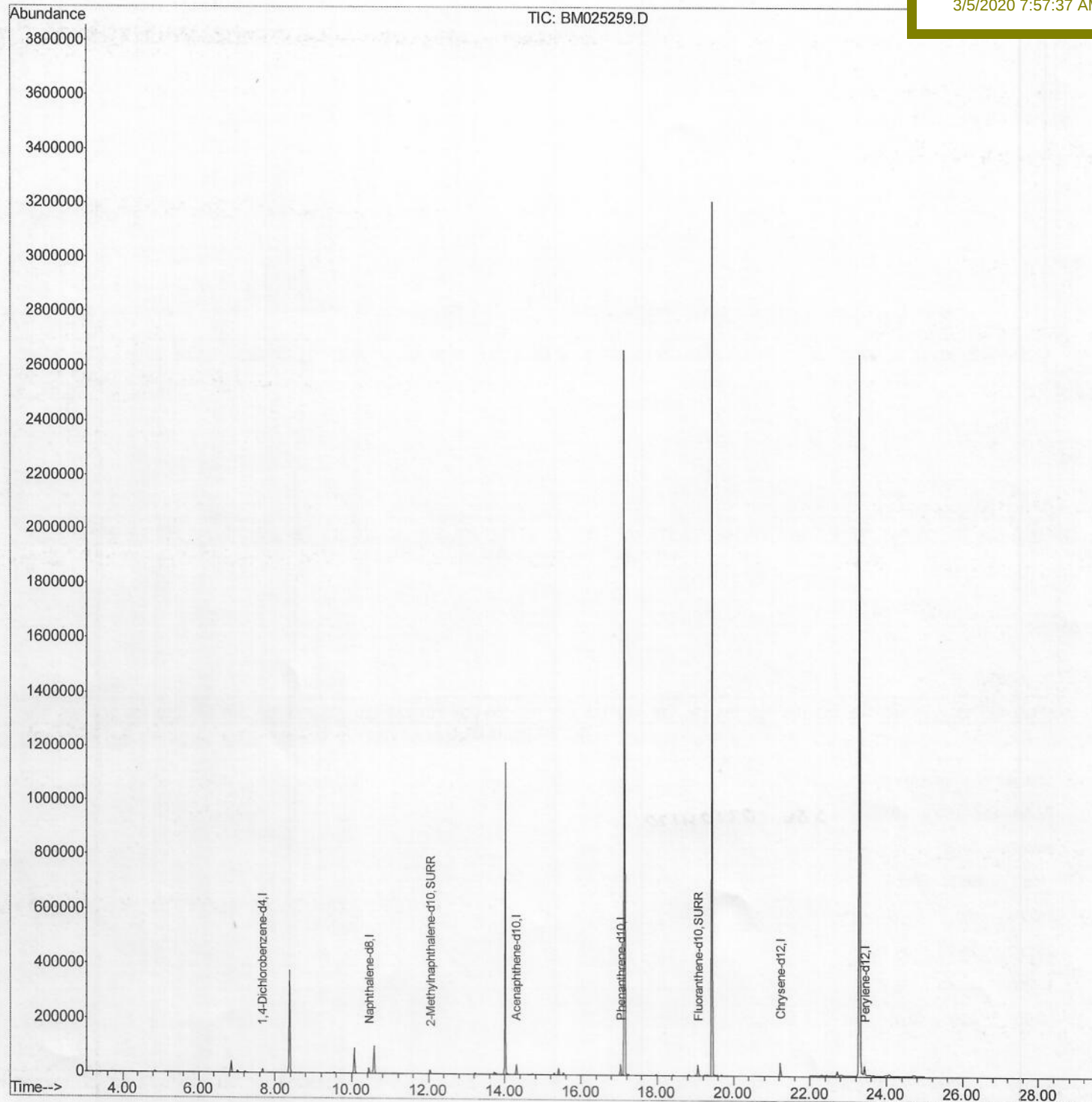
DBD26

Manual Integrations

APPROVED

mohammad

3/5/2020 7:57:37 AM



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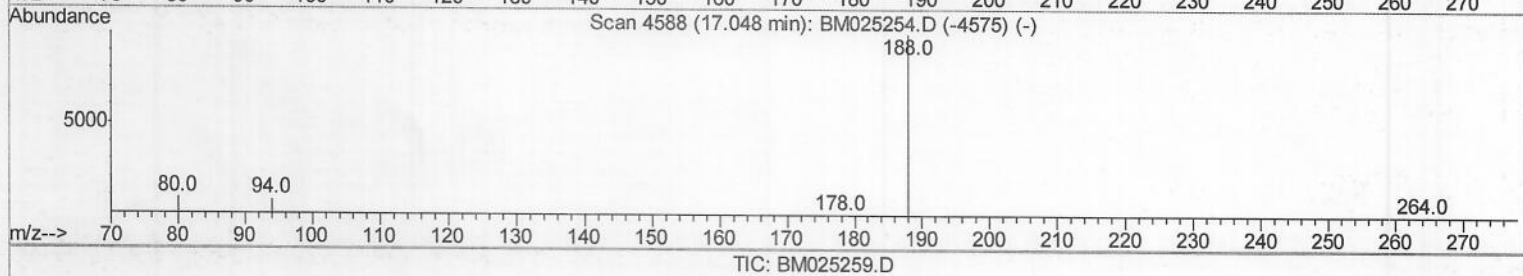
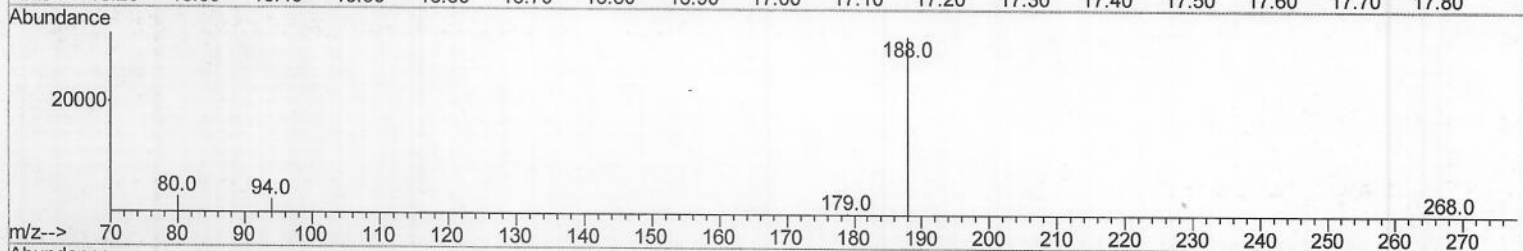
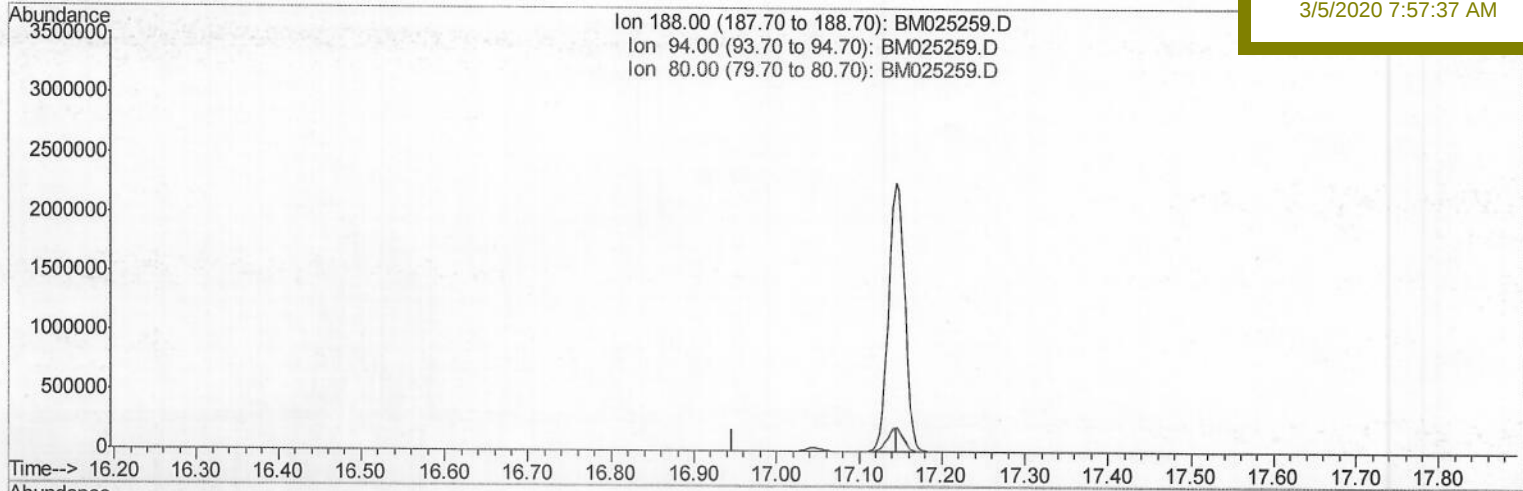
DBD26

Manual Integrations

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mohammad

3/5/2020 7:57:37 AM



(10) Phenanthrene-d10 (I)

17.046min (-17.046) 0.00ng/ul

response 0

Ion	Exp%	Act%
188.00	100	0.00
94.00	26.60	0.00#
80.00	22.60	0.00#
0.00	0.00	0.00

Quantitation Report (Qedit)

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ClientSampleId :

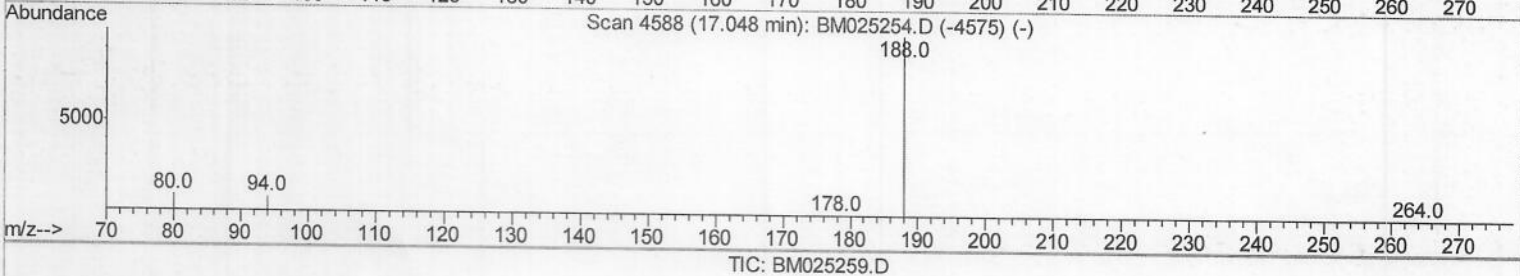
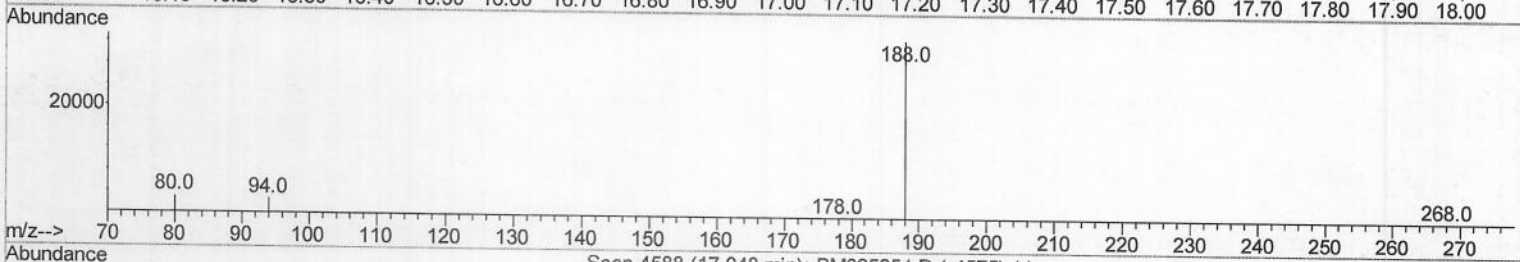
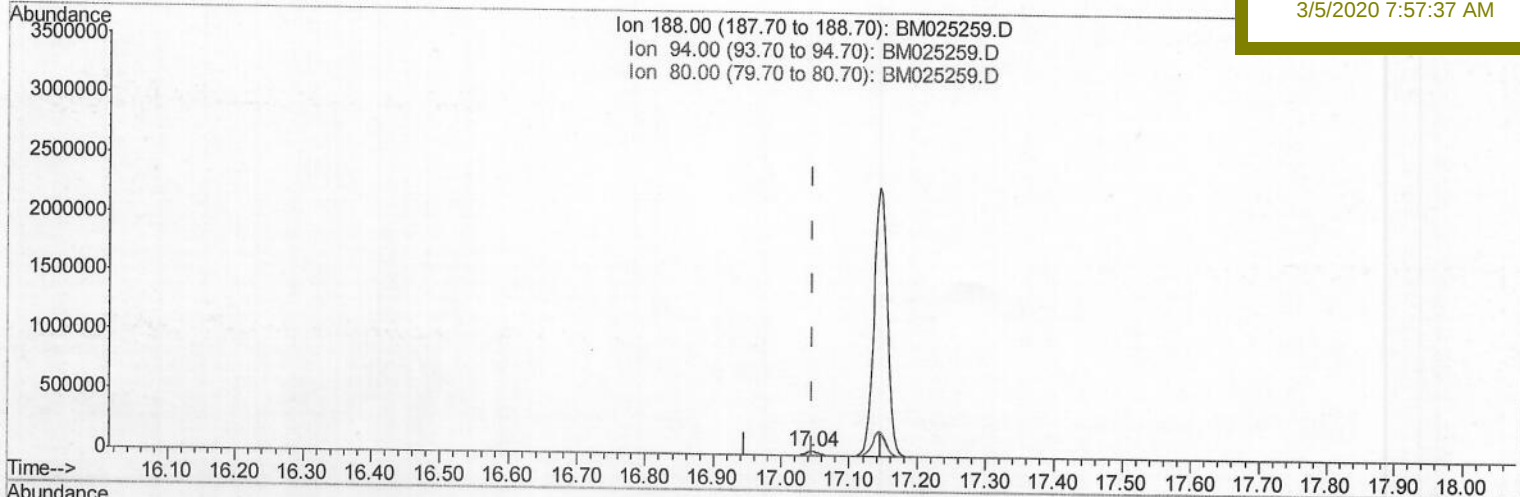
DBD26

Manual Integrations

APPROVED

mohammad

3/5/2020 7:57:37 AM



(10) Phenanthrene-d10 (I)

17.044min (-0.002) 0.40ng/ul m >3V 03105120

response 45057

Ion	Exp%	Act%
188.00	100	100
94.00	26.60	8.12#
80.00	22.60	9.24#
0.00	0.00	0.00

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 Operator : CG/JU
 Sample : L1516-08
 Misc :
 ALS Vial : 55 Sample Multiplier: 1

Instrument :
 BNA_M
 Client Sampled :
 DBD26

Quant Time: Mar 05 01:16:20 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM030320.M
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 QLast Update : Wed Mar 04 08:02:03 2020
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Manual Integrations
 APPROVED

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 3/5/2020 7:57:37 AM

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) 1,4-Dichlorobenzene-d4	7.68	152	5939	0.40	ng/ul	0.00
2) Naphthalene-d8	10.44	136	26049	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	18927	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	45057m)	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	44648	0.40	ng/ul	0.00
20) Perylene-d12	23.43	264	43277	0.40	ng/ul	0.00

SV 03/05/20

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev (Min)
4) 2-Methylnaphthalene-d10	12.04	152	14695	0.30	ng/ul	0.00
14) Fluoranthene-d10	19.07	212	42942	0.29	ng/ul	0.00

Target Compounds	Ovalue

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