

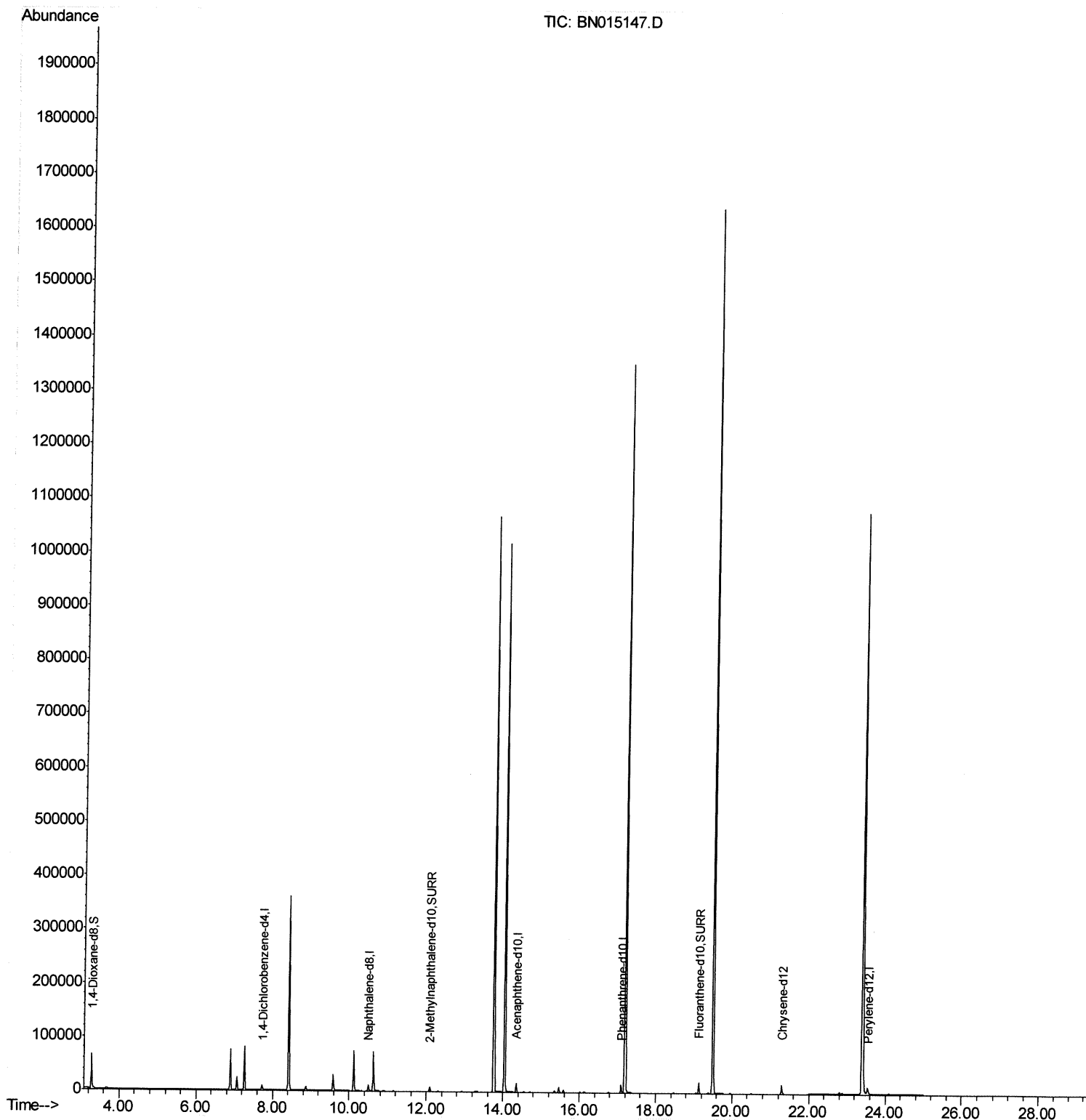
Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN062721\
Data File : BN015147.D
Acq On : 27 Jun 2021 10:55
Operator : CG/JU
Sample : PB137262BL
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
SBLK262

Quant Time: Jun 28 06:38:39 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SFAM-EPA-SIM-BN062121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 22 11:53:52 2021
Response via : Initial Calibration

Manual Integrations
APPROVED

mohammad
6/28/2021 1:28:29 PM



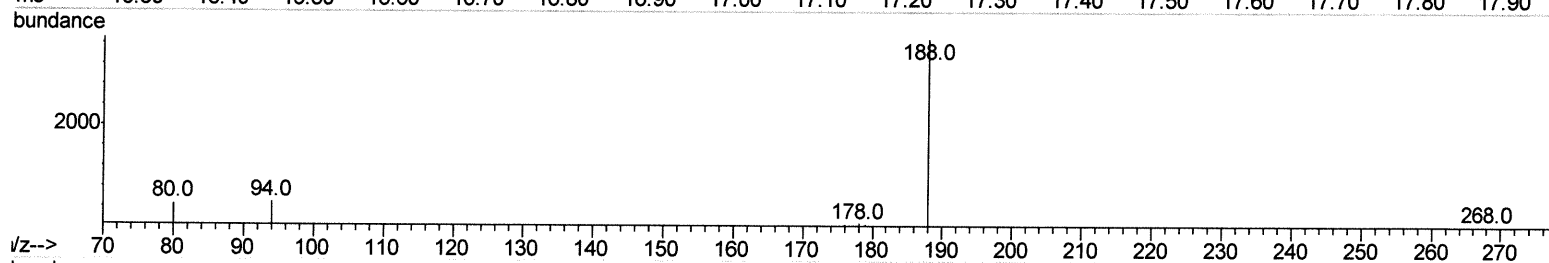
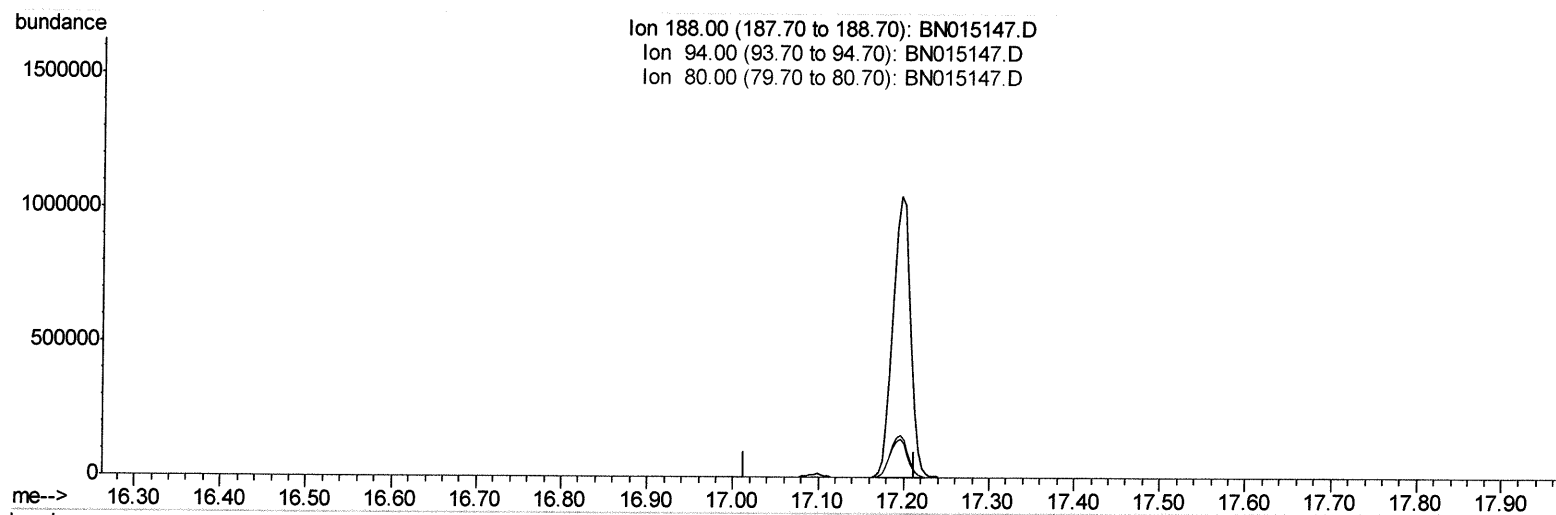
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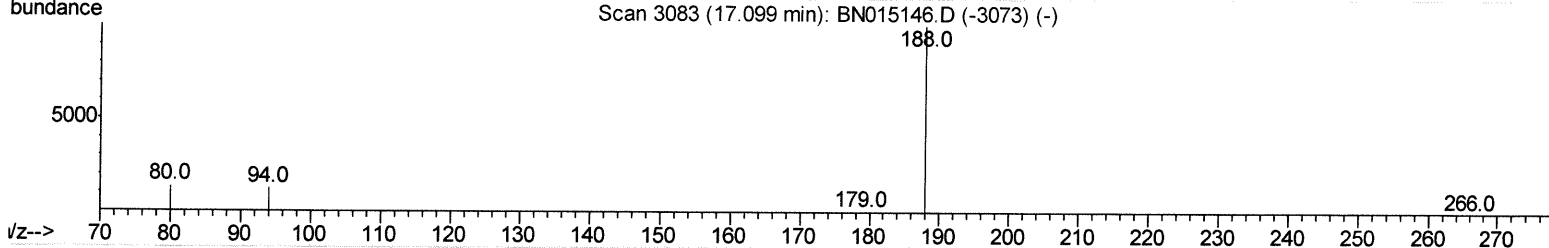
Quant Time: Jun 28 06:36:15 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SFAM-EPA-SIM-BN062121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
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Scan 3083 (17.099 min): BN015146.D (-3073) (-)



TIC: BN015147.D

(13) Phenanthrene-d10 (I)

17.113min (-17.113) 0.00ng/ul

response 0

Ion	Exp%	Act%
188.00	100	0.00
94.00	14.10	0.00#
80.00	14.40	0.00#
0.00	0.00	0.00

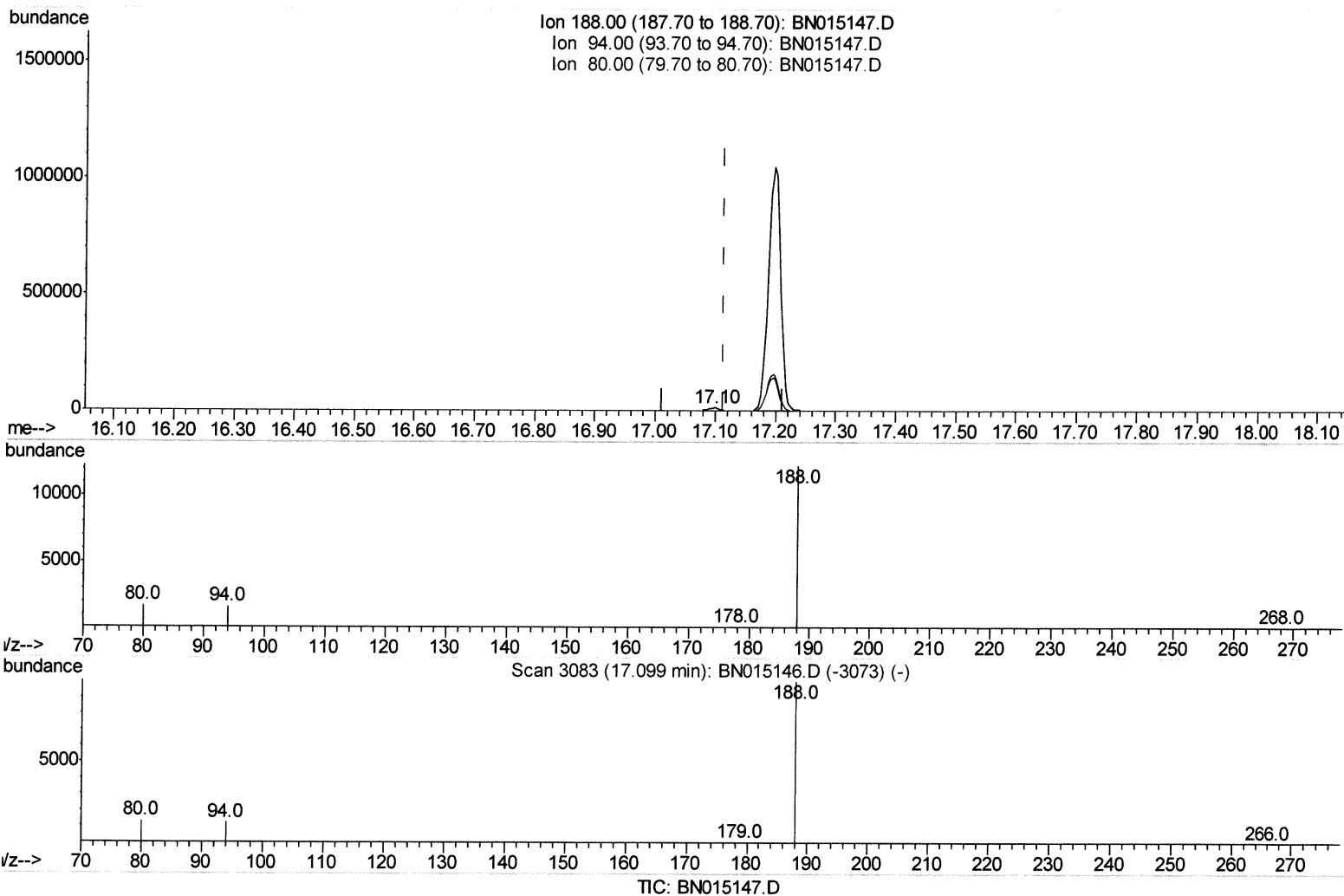
Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN062721\
 Data File : BN015147.D
 Acq On : 27 Jun 2021 10:55
 Operator : CG/JU
 Sample : PB137262BL
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SBLK262

Manual Integrations
 APPROVED

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 6/28/2021 1:28:29 PM

Quant Time: Jun 28 06:36:15 2021
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 22 11:53:52 2021
 Response via : Initial Calibration



(13) Phenanthrene-d10 (I)

17.099min (-0.014) 0.40ng/ul m

ju 6/28/21

response 17368

Ion	Exp%	Act%
188.00	100	100
94.00	14.10	12.76
80.00	14.40	13.38
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN062721\
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Acq On : 27 Jun 2021 10:55
Operator : CG/JU
Sample : PB137262BL
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
SBLK262

Manual Integrations
APPROVED

mohammad
6/28/2021 1:28:29 PM

Quant Time: Jun 28 06:38:39 2021
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	4268	0.40	ng/ul	-0.01
4) Naphthalene-d8	10.50	136	15002	0.40	ng/ul	-0.01
9) Acenaphthene-d10	14.35	164	8504	0.40	ng/ul	-0.02
13) Phenanthrene-d10	17.10	188	17368m>	0.40	ng/ul	>-0.01
17) Chrysene-d12	21.29	240	14478	0.40	ng/ul	-0.02
23) Perylene-d12	23.54	264	14984	0.40	ng/ul	-0.03

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System Monitoring Compounds

3) 1,4-Dioxane-d8	3.29	96	38553	8.01	ng/ul	-0.01
6) 2-Methylnaphthalene-d10	12.09	152	9879	0.42	ng/ul	-0.02
18) Fluoranthene-d10	19.13	212	20934	0.44	ng/ul	-0.02

Target Compounds

Ovalue

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