

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN122819\

Data File : BN009219.D

Acq On : 28 Dec 2019 09:44

Operator : JU

Sample : PB125532BL

Misc :

ALS Vial : 3 Sample Multiplier: 1

Quant Time: Dec 29 06:24:05 2019

Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN121619.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Sun Dec 29 06:16:40 2019

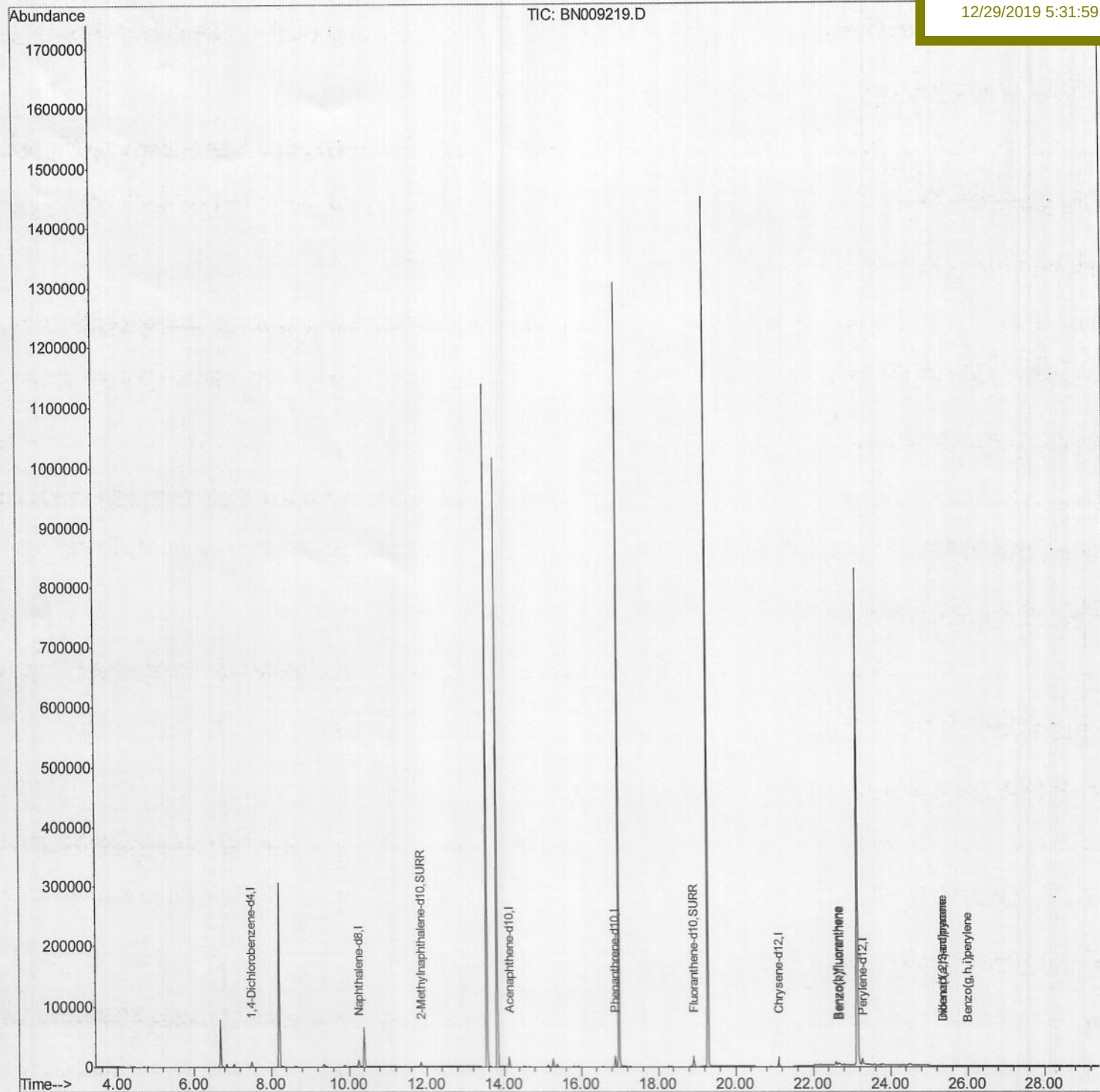
Response via : Initial Calibration

Instrument :

BNA_N

ClientSampleId :

SBLK32

Manual Integrations
APPROVEDmohammad
12/29/2019 5:31:59 PM

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Quant Time: Dec 29 06:19:12 2019

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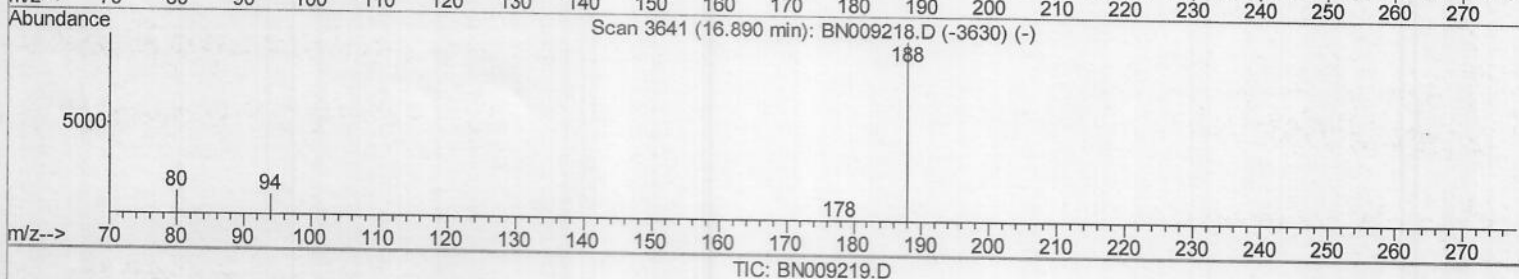
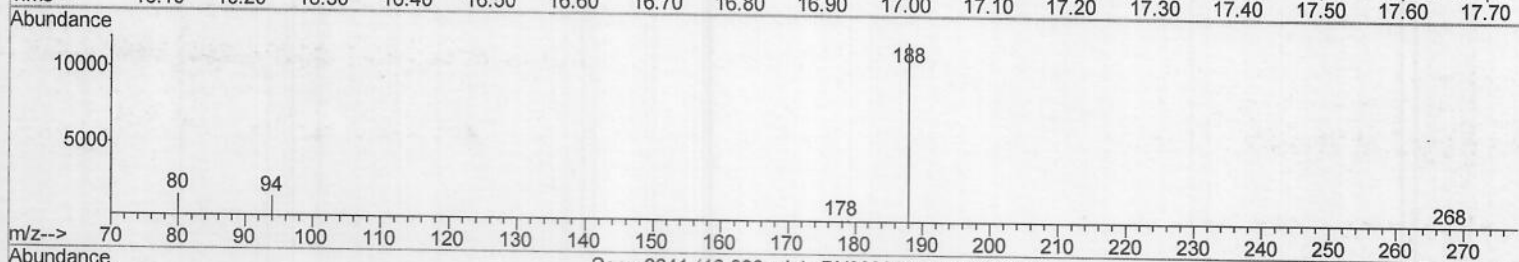
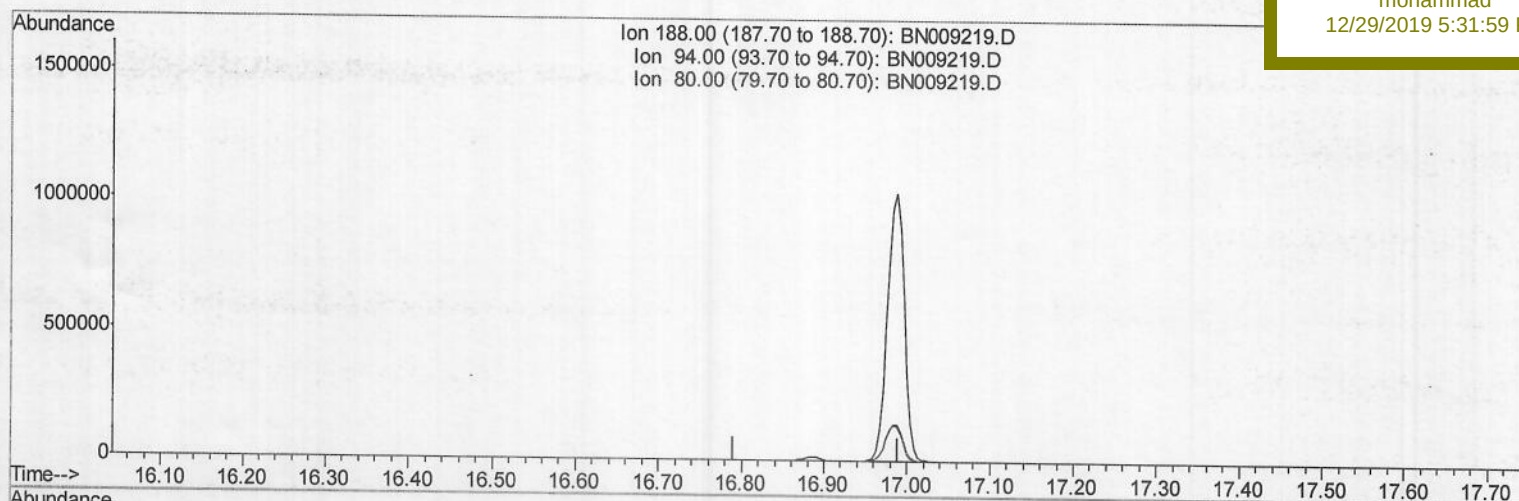
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APPROVEDmohammad
12/29/2019 5:31:59 PM

(10) Phenanthrene-d10 (I)

16.890min (-16.890) 0.00ng/ul

response 0

Ion	Exp%	Act%
188.00	100	0.00
94.00	11.60	0.00#
80.00	13.70	0.00#
0.00	0.00	0.00

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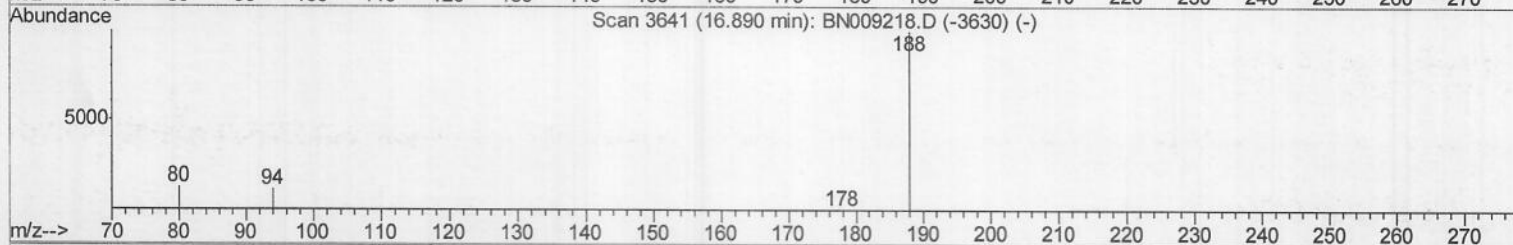
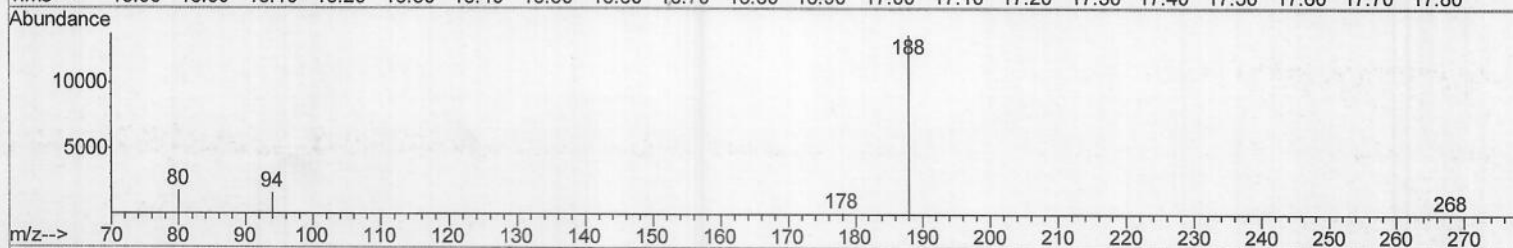
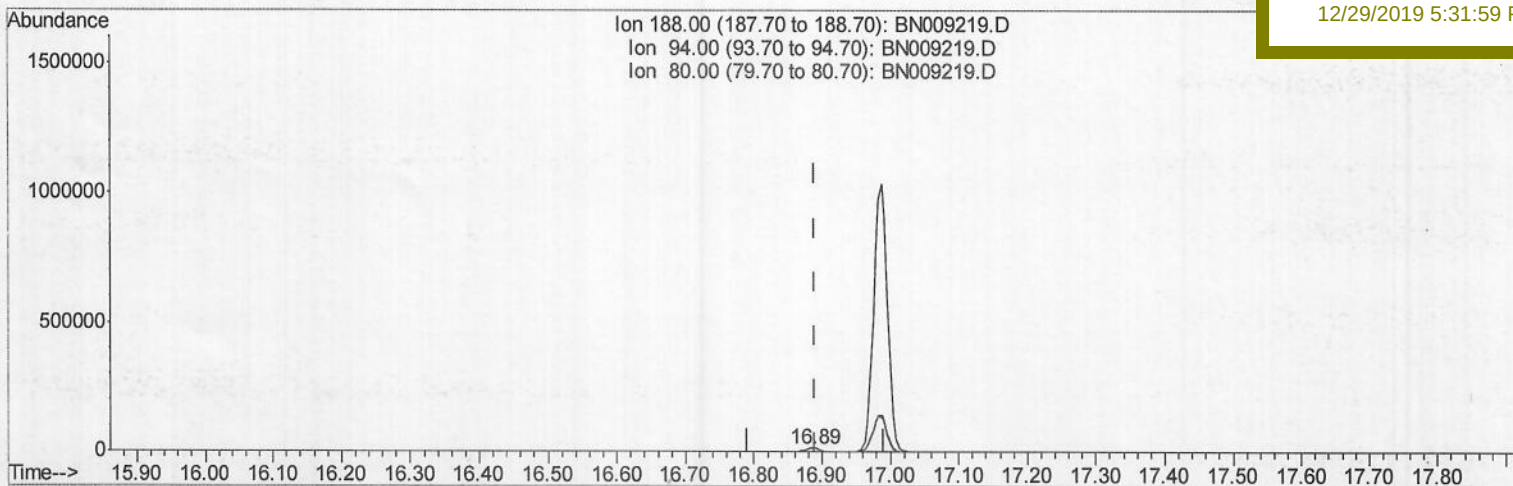
Response via : Initial Calibration

Instrument :

BNA_N

ClientSampleId :

SBLK32

Manual Integrations
APPROVEDmohammad
12/29/2019 5:31:59 PM

TIC: BN009219.D

(10) Phenanthrene-d10 (I)

16.886min (-0.004) 0.40ng/ul m) JU 12/30/19

response 18617

Ion	Exp%	Act%
188.00	100	100
94.00	11.60	12.11
80.00	13.70	13.43
0.00	0.00	0.00

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 Misc :
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Instrument :
 BNA_N
 ClientSampleId :
 SBLK32

Manual Integrations
 APPROVED

mohammad
 12/29/2019 5:31:59 PM

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) 1,4-Dichlorobenzene-d4	7.50	152	3944	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.26	136	14977	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.13	164	8928	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.89	188	18617m	0.40	ng/ul	0.00
16) Chrysene-d12	21.12	240	13303	0.40	ng/ul	0.03
20) Perylene-d12	23.27	264	10722	0.40	ng/ul	0.06

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.85	152	8789	0.33	ng/ul	-0.01
14) Fluoranthene-d10	18.92	212	18837	0.34	ng/ul	0.00

Target Compounds				Ovalue		
21) Benzo(b)fluoranthene	22.65	252	3070	0.061	ng/ul#	55
22) Benzo(k)fluoranthene	22.69	252	2052	0.041	ng/ul#	27
24) Indeno(1,2,3-cd)pyrene	25.36	276	1390	0.029	ng/ul#	89
25) Dibenzo(a,h)anthracene	25.37	278	865	0.023	ng/ul#	21
26) Benzo(a,h,i)perylene	25.99	276	1172	0.029	ng/ul#	62

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