

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM122319\
Quantitation Report (QT Reviewed)

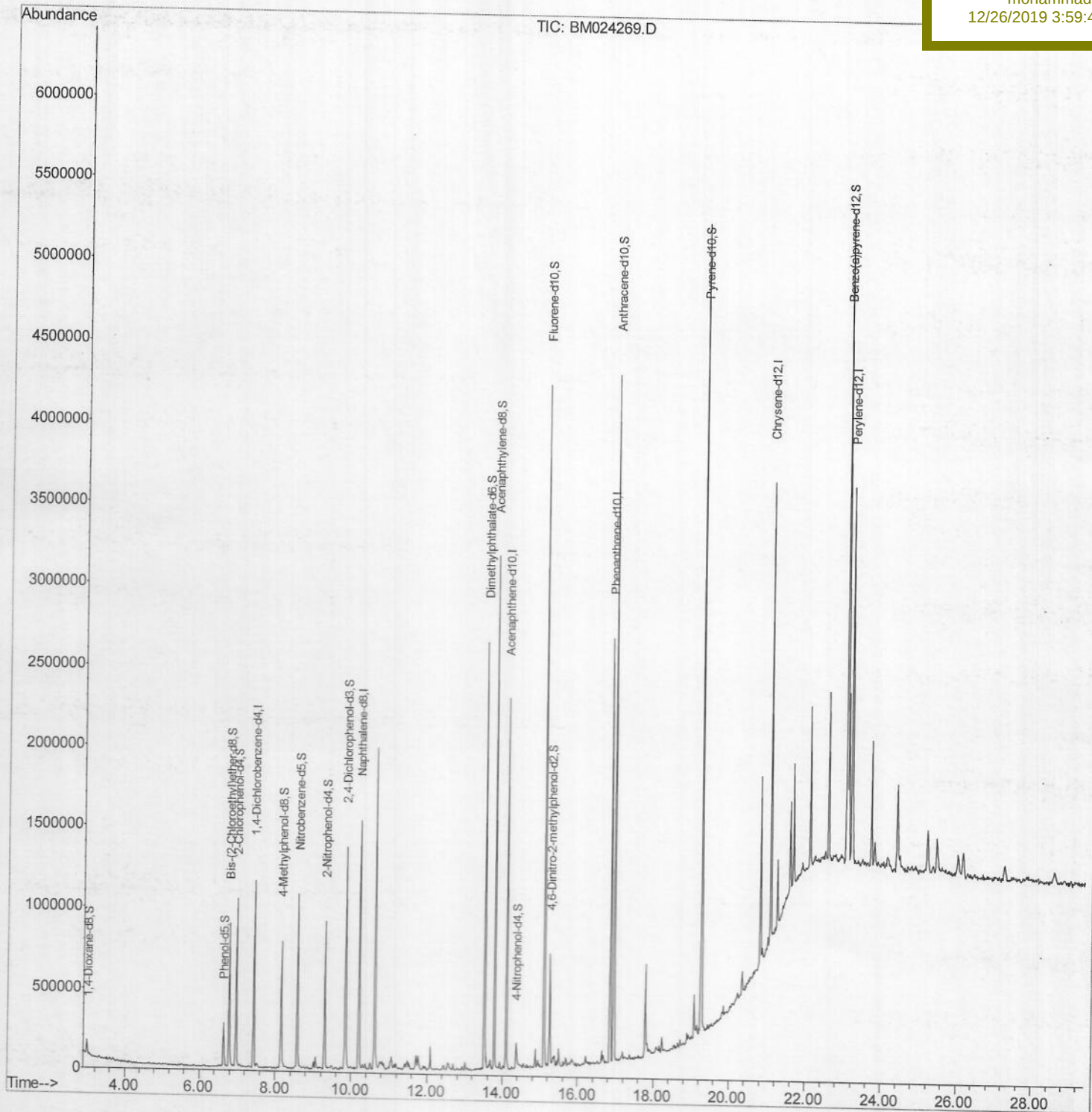
Data File : BM024269.D
Acq On : 25 Dec 2019 00:25
Operator : JU
Sample : K6240-12
Misc :
ALS Vial : 18 Sample Multiplier: 1

Quant Time: Dec 25 03:55:31 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM121719MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Dec 25 03:33:48 2019
Response via : Initial Calibration

Instrument :
BNA_M
Client Sample ID :
BFQW5

Manual Integrations
APPROVED

mohammad
12/26/2019 3:59:46 PM



SOM-EPA-BM121719MA.M Wed Dec 25 03:55:04 2019

Page: 2

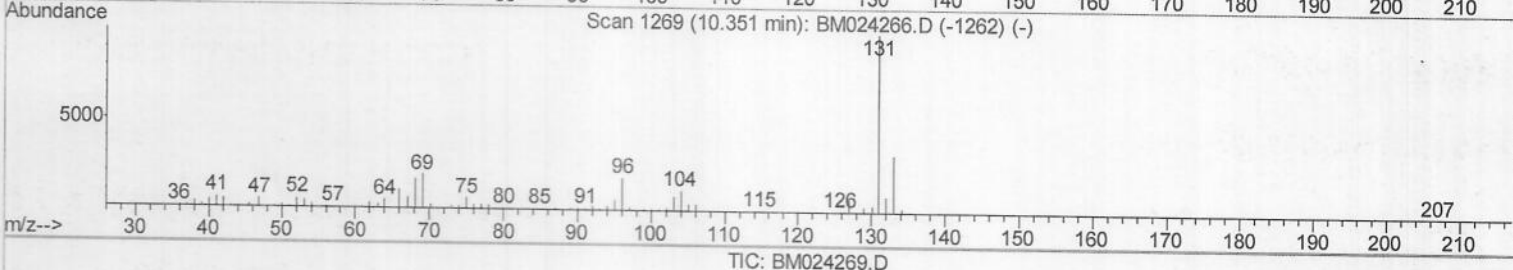
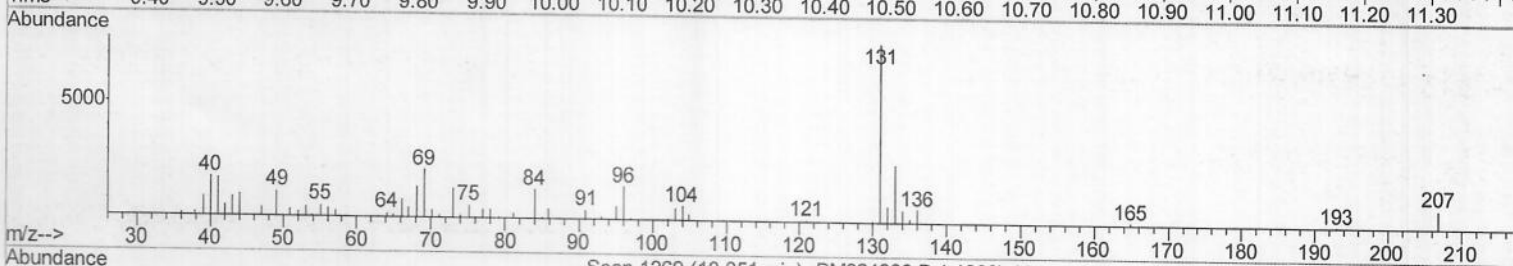
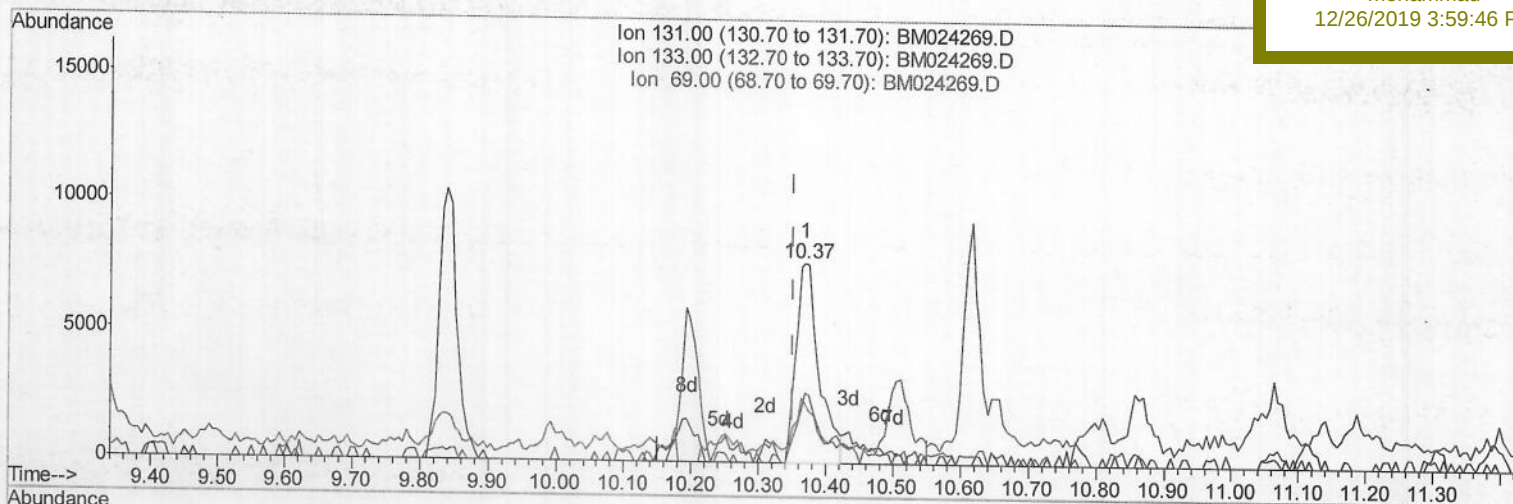
Data File : BM024269.D
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Quant Time: Dec 25 03:36:33 2019
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(29) 4-Chloroaniline-d4 (S)
10.369min (+0.018) 0.79ng/ul
response 18351

Ion	Exp%	Act%
131.00	100	100
133.00	32.50	35.16
69.00	19.00	29.40#
0.00	0.00	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM122319\
Quantitation Report (Qedit)

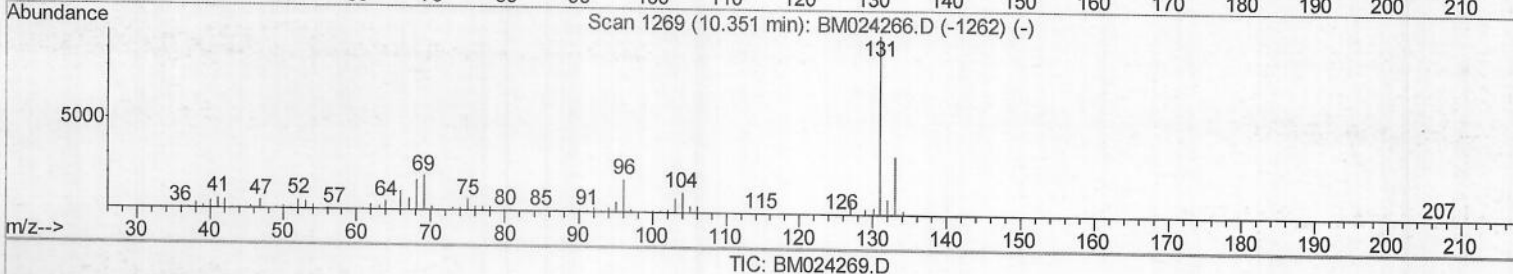
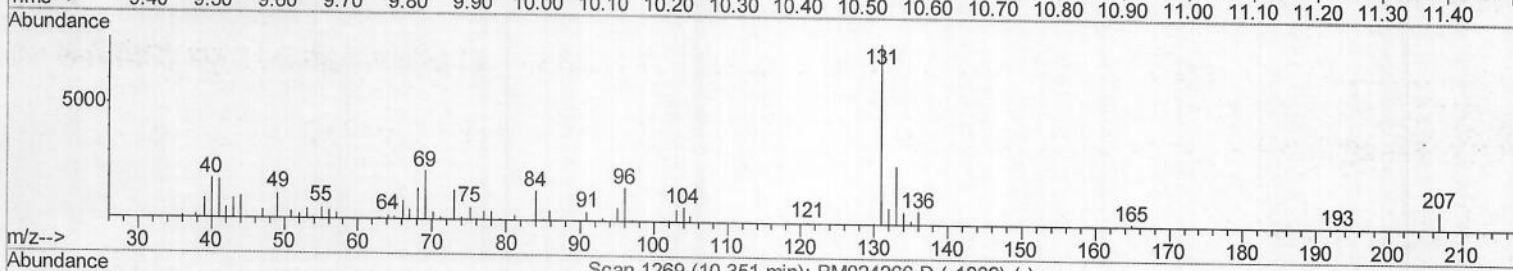
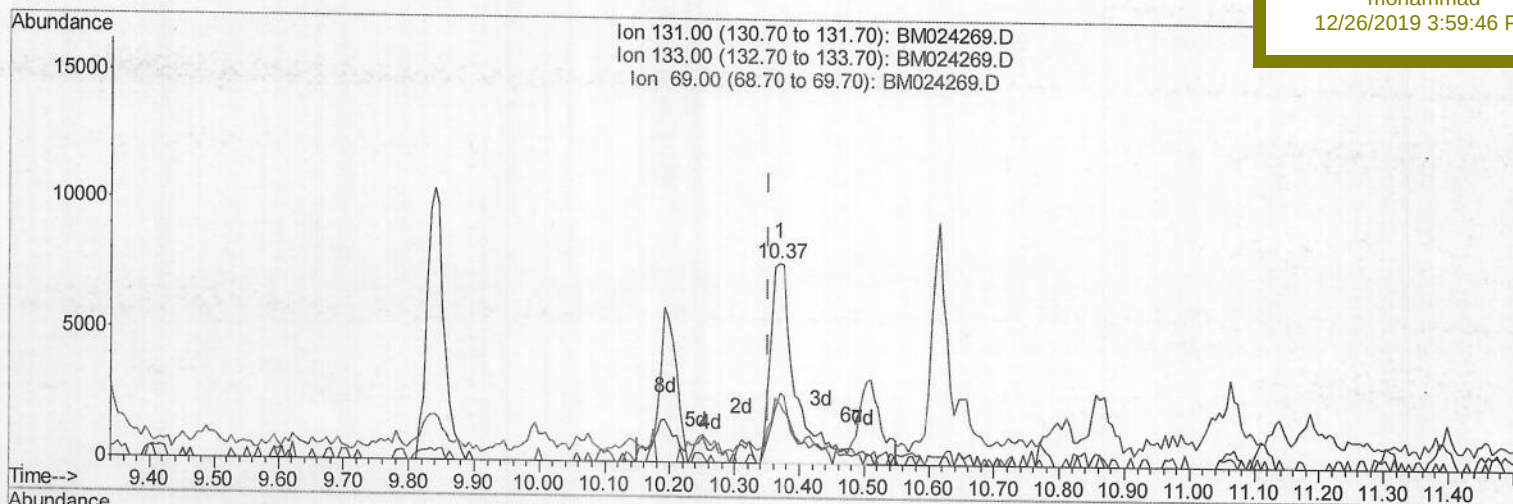
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Manual Integrations
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(29) 4-Chloroaniline-d4 (S)

10.369min (+0.018) 0.93ng/ul m

response 21636

Ion	Exp%	Act%
131.00	100	100
133.00	32.50	35.16
69.00	19.00	29.40#
0.00	0.00	0.00

> JU 12/27/19

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Quantitation Report (QT Reviewed)

Data File : BM024269.D

Acq On : 25 Dec 2019 00:25

Operator : JU

Sample : K6240-12

Misc :

ALS Vial : 18 Sample Multiplier: 1

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM121719MA.M

Quant Title : SVOA CALIBRATION

QLast Update : Wed Dec 25 03:33:48 2019

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Instrument :

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ClientSampleId :

BFQW5

Manual Integrations

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.43	152	303964	20.00	ng/ul	0.00
18) Naphthalene-d8	10.20	136	1244266	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.09	164	712059	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.85	188	1417259	20.00	ng/ul	0.00
78) Chrysene-d12	21.07	240	1256515	20.00	ng/ul	0.00
86) Perylene-d12	23.20	264	1560529	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	2.99	96	45498	6.61	ng/uL	0.00
5) Phenol-d5	6.63	99	170306	5.92	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.78	67	435257	25.22	ng/ul	0.00
9) 2-Chlorophenol-d4	6.97	132	456350	22.23	ng/ul	0.00
13) 4-Methylphenol-d8	8.16	113	316622	13.46	ng/ul	0.00
19) Nitrobenzene-d5	8.58	128	266211	29.77	ng/ul	0.00
22) 2-Nitrophenol-d4	9.30	143	297256	30.24	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.84	165	536698	28.08	ng/ul	0.00
29) 4-Chloroaniline-d4	10.37	131	21636m	0.93	ng/ul	0.02
44) Dimethylphthalate-d6	13.52	166	1744973	33.14	ng/ul	0.00
47) Acenaphthylene-d8	13.78	160	2083322	33.02	ng/ul	0.00
52) 4-Nitrophenol-d4	14.36	143	55417	5.86	ng/ul	0.01
58) Fluorene-d10	15.10	176	1505998	32.63	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.26	200	139848	16.18	ng/ul	0.00
71) Anthracene-d10	16.95	188	2240721	34.75	ng/ul	0.00
79) Pyrene-d10	19.26	212	2336230	39.03	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.06	264	2695142	34.77	ng/ul	0.00

Target Compounds

Qvalue

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

7 JU 12/27/19