

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM122319\

Quantitation Report (QT Reviewed)

Data File : BM024278.D

Acq On : 25 Dec 2019 05:47

Operator : JU

Sample : K6254-01

Misc :

ALS Vial : 27 Sample Multiplier: 1

Quant Time: Dec 25 06:54:31 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

ClientSampled :

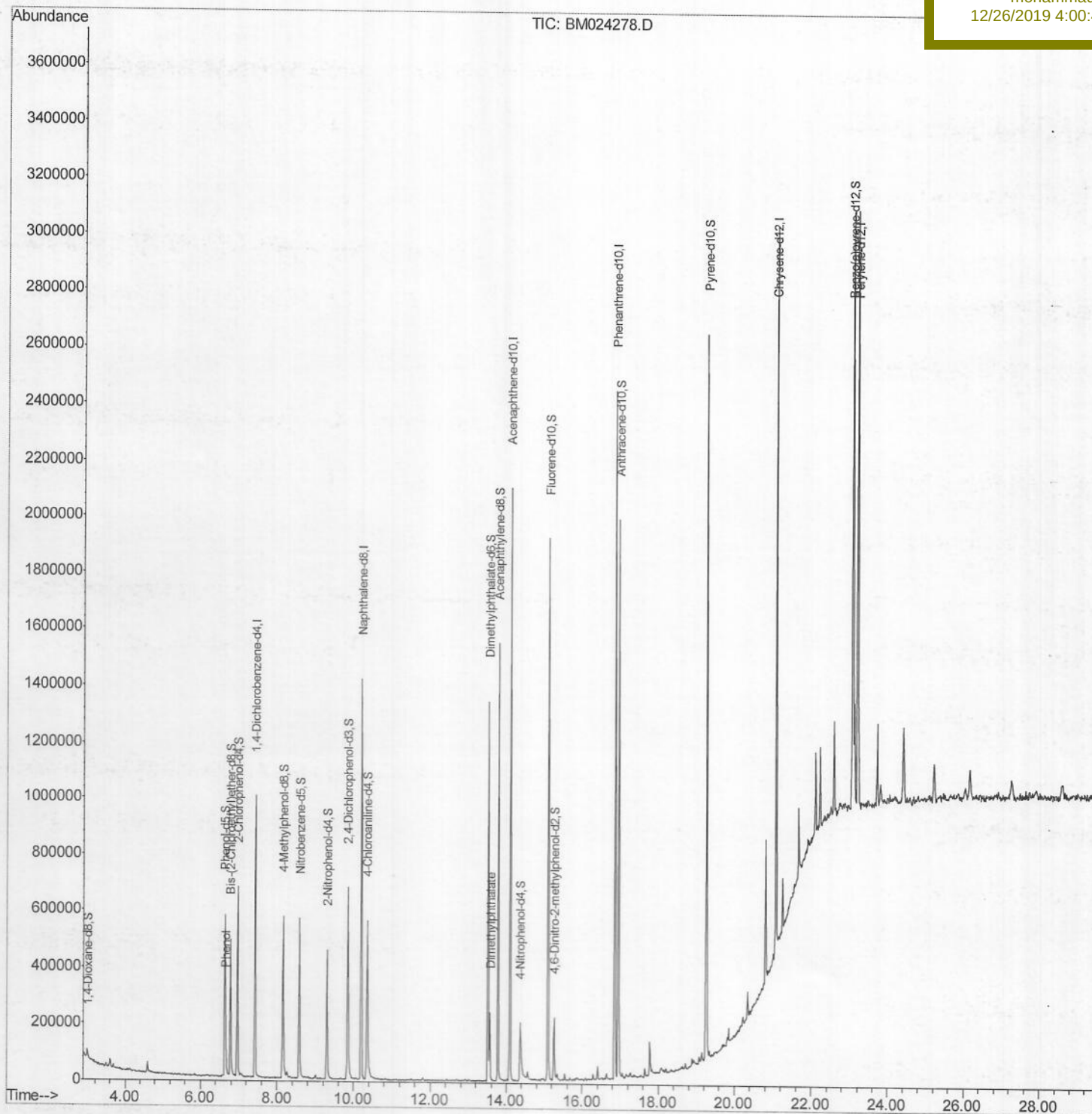
C0C05

Manual Integrations

APPROVED

mohammad

12/26/2019 4:00:40 PM



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

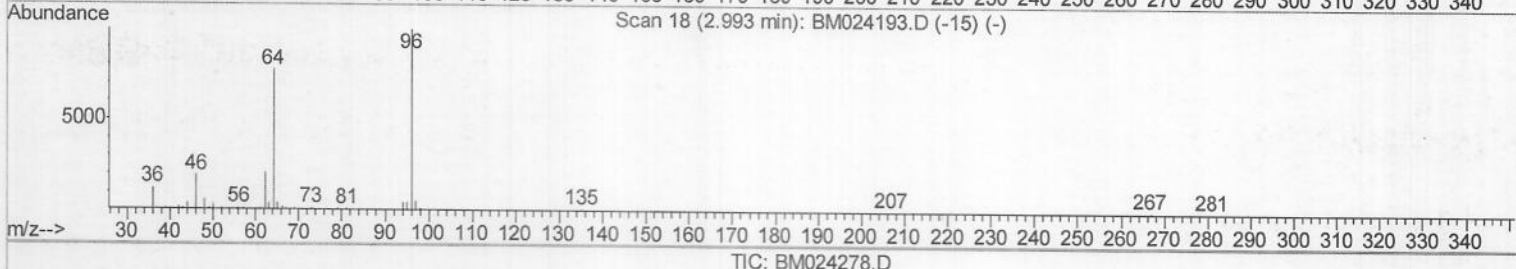
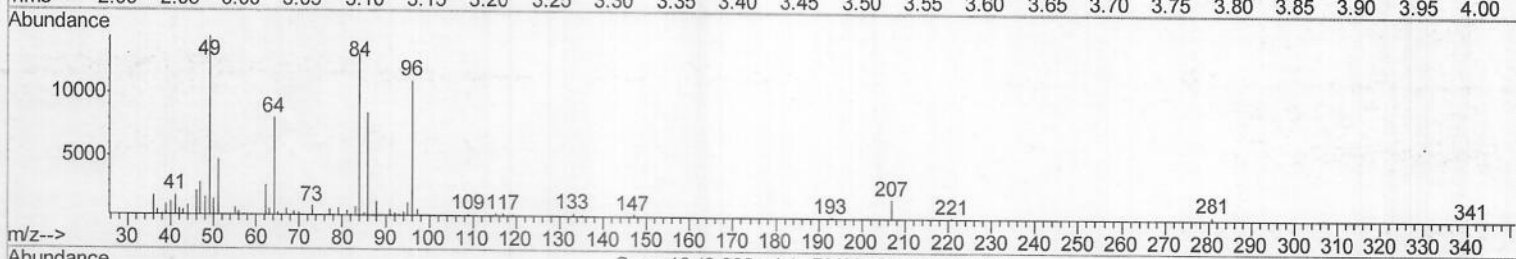
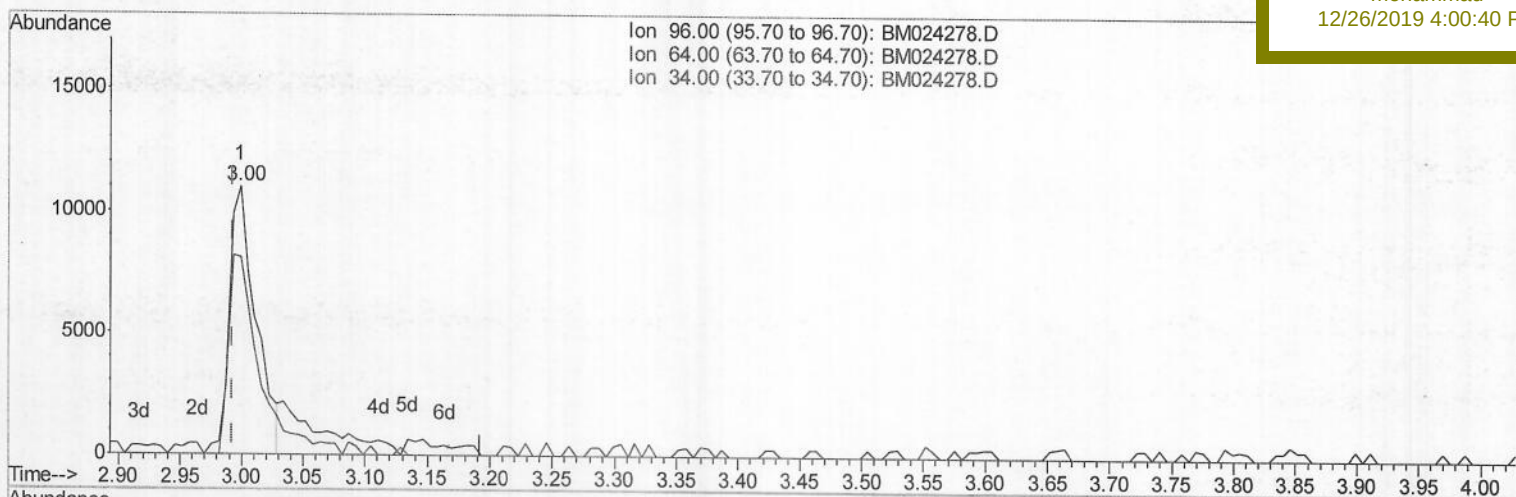
Data File : BM024278.D
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ClientSampleId :
C0C05

Manual Integrations
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(3) 1,4-Dioxane-d8 (S)

2.999min (+0.006) 2.66ng/uL

response 16550

Ion	Exp%	Act%
96.00	100	100
64.00	78.60	73.50
34.00	0.00	0.00
0.00	0.00	0.00

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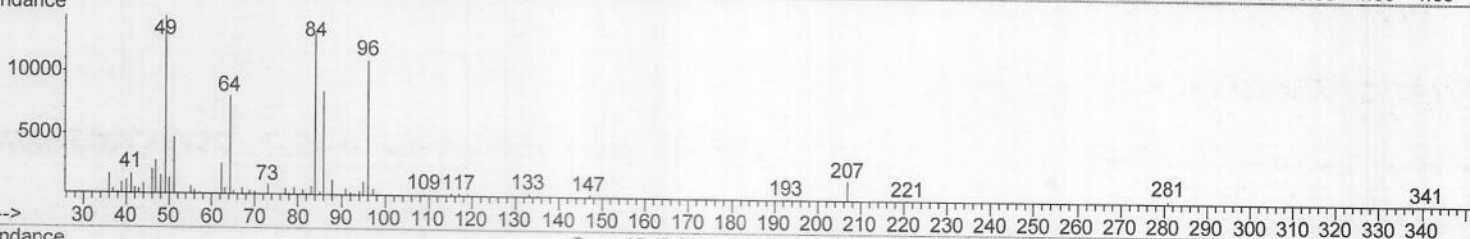
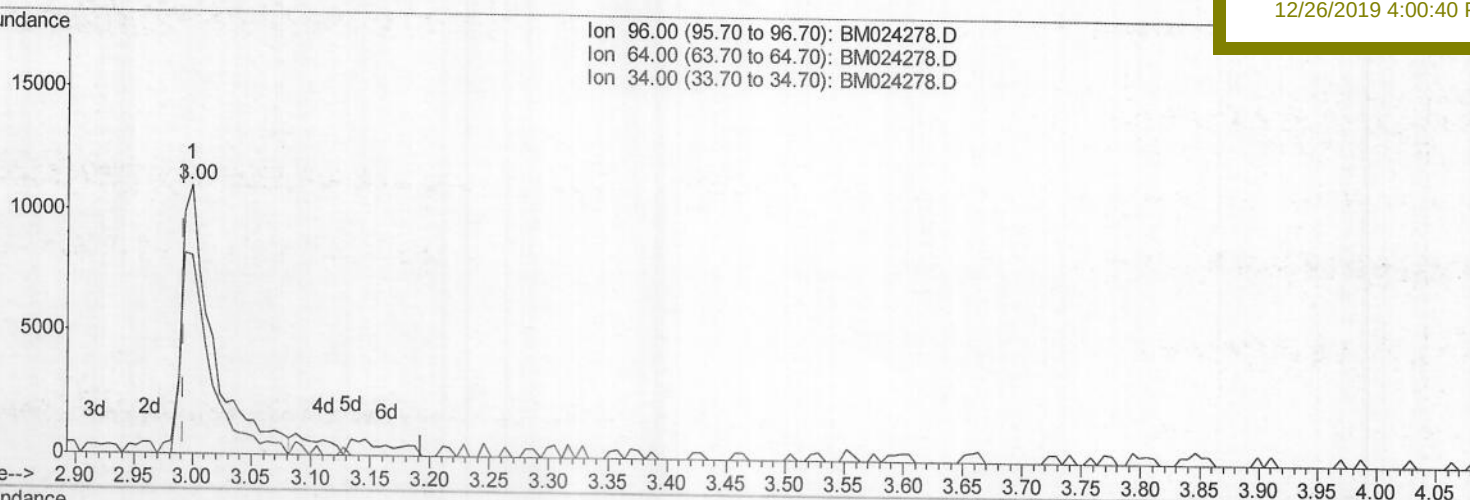
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Instrument :
BNA_M
ClientSampleId :
C0C05

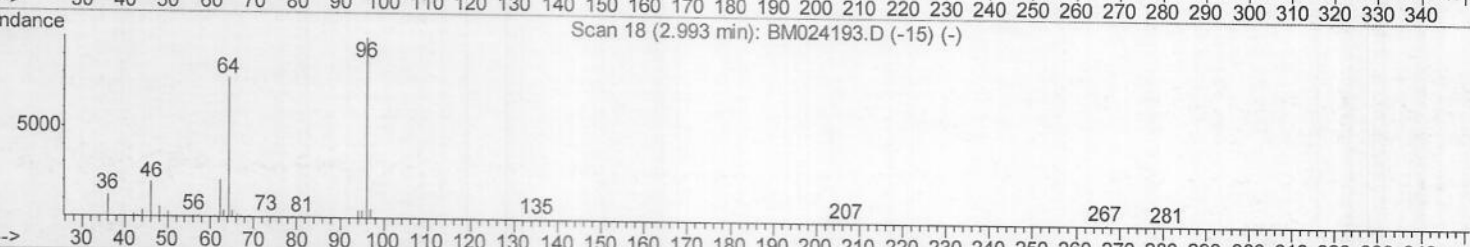
Manual Integrations
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12/26/2019 4:00:40 PM

Ion 96.00 (95.70 to 96.70): BM024278.D
Ion 64.00 (63.70 to 64.70): BM024278.D
Ion 34.00 (33.70 to 34.70): BM024278.D



Scan 18 (2.993 min): BM024193.D (-15) (-)



TIC: BM024278.D

(3) 1,4-Dioxane-d8 (S)

2.999min (+0.006) 3.28ng/uL m

response 20364

Ion	Exp%	Act%
96.00	100	100
64.00	78.60	73.50
34.00	0.00	0.00
0.00	0.00	0.00

>JU 12/27/19

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C0C05

Manual Integrations

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) 1,4-Dichlorobenzene-d4	7.42	152	274621	20.00	ng/ul	0.00
18) Naphthalene-d8	10.19	136	1142745	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.09	164	656600	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.85	188	1304245	20.00	ng/ul	0.00
78) Chrysene-d12	21.07	240	1170758	20.00	ng/ul	0.00
86) Perylene-d12	23.19	264	1415741	20.00	ng/ul	-0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.00	96	20364m	3.28	ng/uL	0.00
5) Phenol-d5	6.63	99	369828	14.22	ng/ul	0.00
7) Bis-(2-Chloroethyl) ether-d	6.78	67	248884	15.96	ng/ul	0.00
9) 2-Chlorophenol-d4	6.97	132	294799	15.90	ng/ul	0.00
13) 4-Methylphenol-d8	8.16	113	242353	11.40	ng/ul	0.00
19) Nitrobenzene-d5	8.58	128	141763	17.26	ng/ul	0.00
22) 2-Nitrophenol-d4	9.30	143	153953	17.05	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.84	165	268572	15.30	ng/ul	0.00
29) 4-Chloroaniline-d4	10.35	131	353359	16.58	ng/ul	0.00
44) Dimethylphthalate-d6	13.52	166	872278	17.97	ng/ul	0.00
47) Acenaphthylene-d8	13.78	160	1034415	17.78	ng/ul	0.00
52) 4-Nitrophenol-d4	14.36	143	88535	10.16	ng/ul	0.01
58) Fluorene-d10	15.10	176	732155	17.20	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.26	200	47975	6.03	ng/ul	0.00
71) Anthracene-d10	16.95	188	1090986	18.38	ng/ul	0.00
79) Pyrene-d10	19.26	212	1228841	22.03	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.06	264	1386630	19.72	ng/ul	0.00

Target Compounds

6) Phenol	6.66	94	62255	2.319	ng/ul	95
45) Dimethylphthalate	13.56	163	154529	3.062	ng/ul	99

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

> JU 12/27/19