

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122319\

Data File : BM024285.D

Acq On : 25 Dec 2019 12:42

Operator : JU

Sample : K6254-07

Misc :

ALS Vial : 34 Sample Multiplier: 1

Quant Time: Dec 25 14:26:46 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

ClientSampleId :

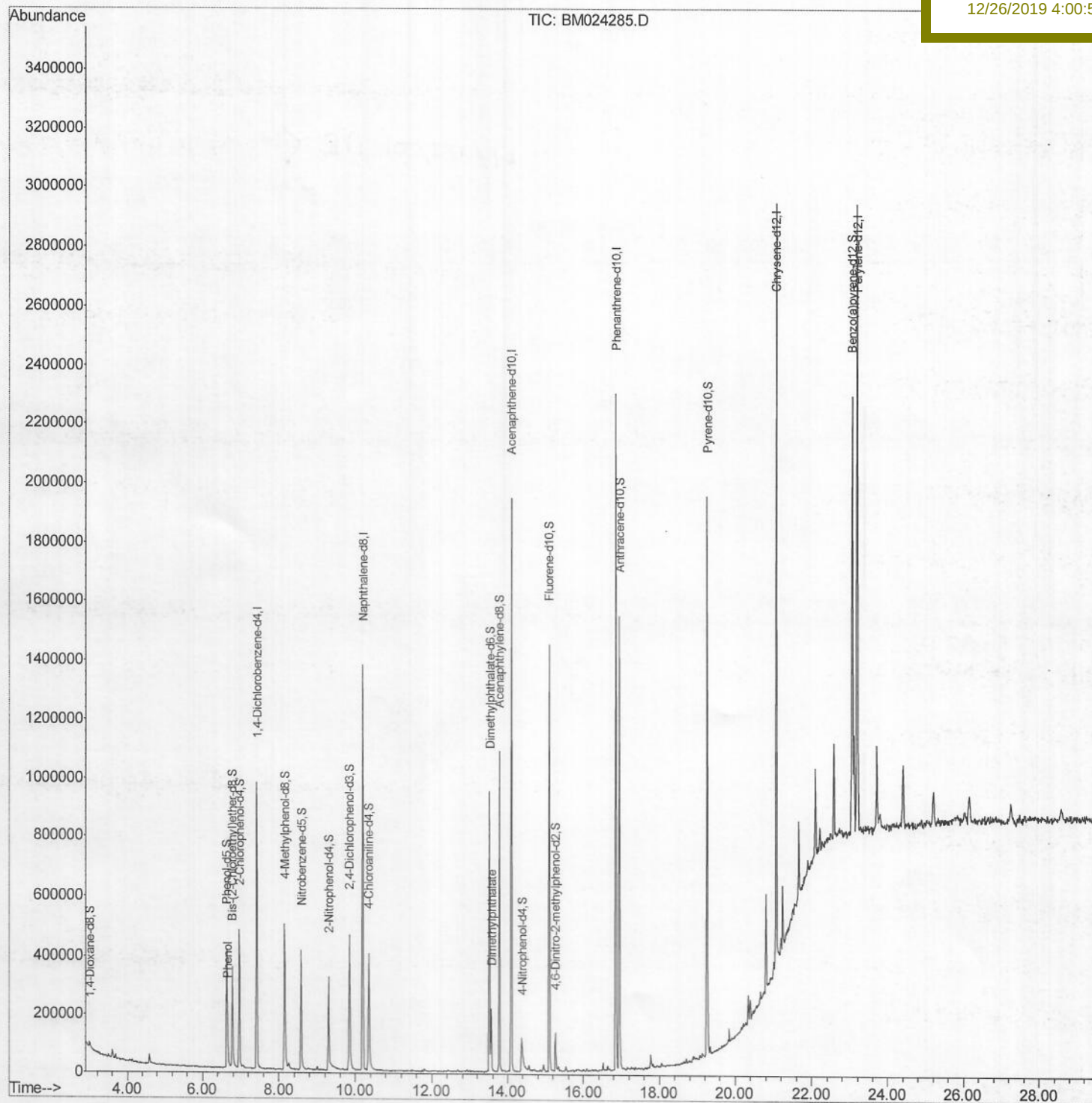
C0C21

Manual Integrations

APPROVED

mohammad

12/26/2019 4:00:54 PM



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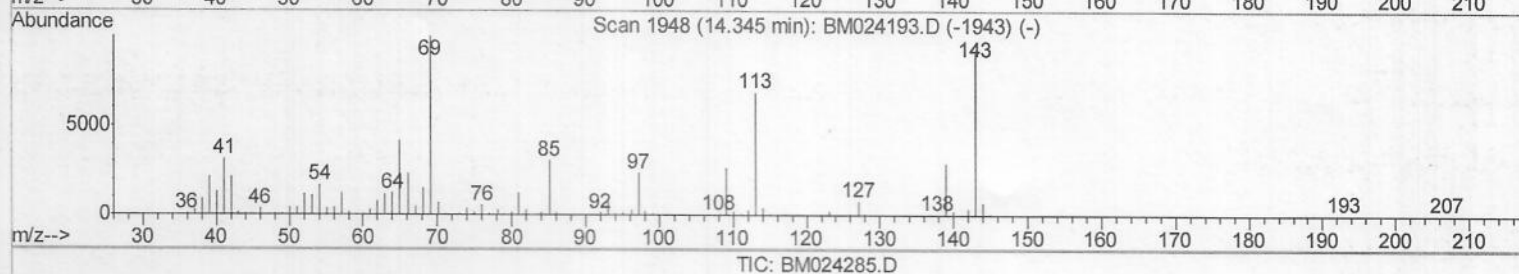
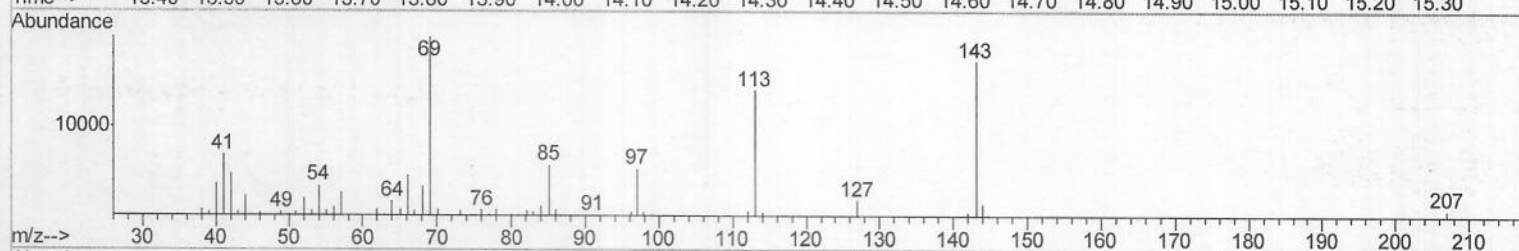
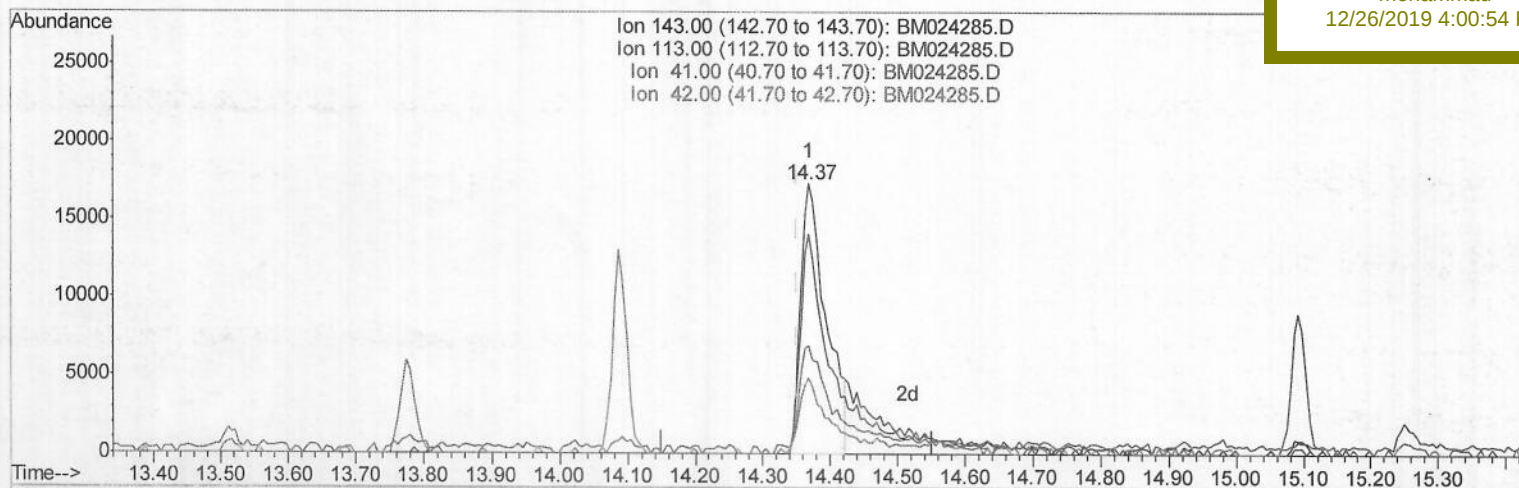
C0C21

Manual Integrations

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(52) 4-Nitrophenol-d4 (S)

14.369min (+0.018) 5.45ng/ul

response 45486

Ion	Exp%	Act%
143.00	100	100
113.00	73.60	81.26
41.00	33.90	40.20
42.00	22.80	28.21#

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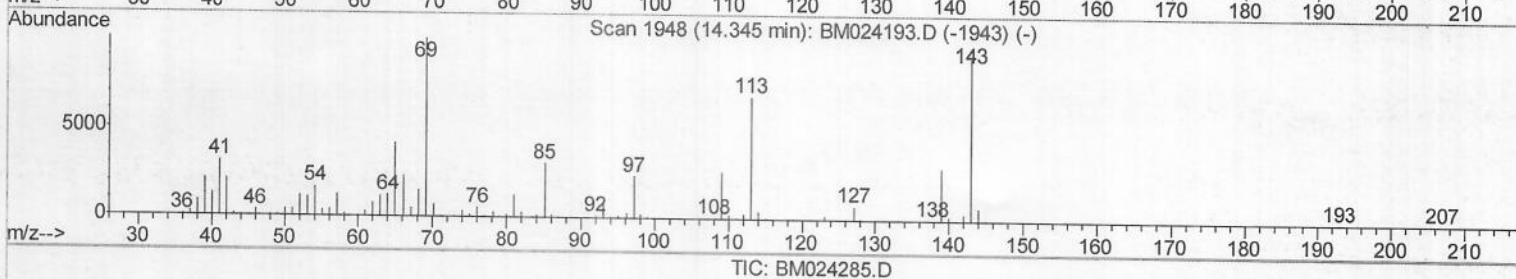
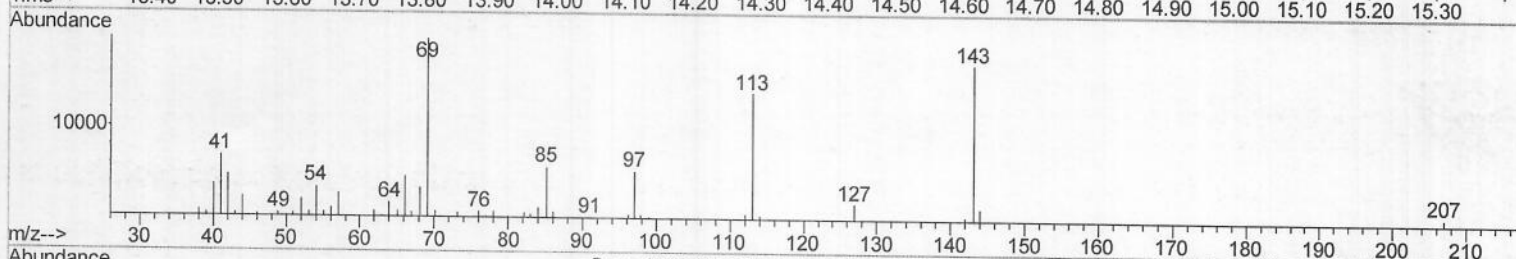
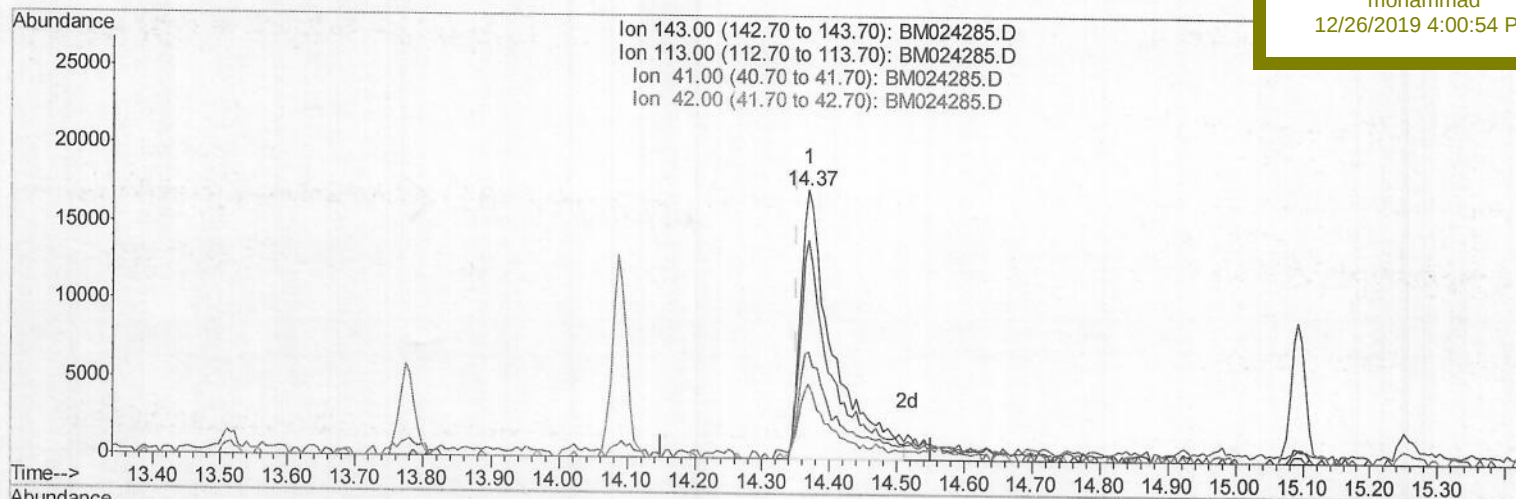
C0C21

Manual Integrations

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mohammad

12/26/2019 4:00:54 PM



(52) 4-Nitrophenol-d4 (S)

14.369min (+0.018) 7.02ng/ul m

response 58568

Ion	Exp%	Act%
143.00	100	100
113.00	73.60	81.26
41.00	33.90	40.20
42.00	22.80	28.21#

> JU 12/27/19

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BNA_M

Client Sampled :

C0C21

Manual Integrations

APPROVED

mohammad

12/26/2019 4:00:54 PM

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) 1,4-Dichlorobenzene-d4	7.42	152	258284	20.00	ng/ul	0.00
18) Naphthalene-d8	10.19	136	1080860	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.09	164	628189	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.85	188	1250661	20.00	ng/ul	0.00
78) Chrysene-d12	21.06	240	1140743	20.00	ng/ul	0.00
86) Perylene-d12	23.19	264	1336280	20.00	ng/ul	-0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.00	96	13400	2.29	ng/uL	0.00
5) Phenol-d5	6.63	99	265186	10.84	ng/ul	0.00
7) Bis-(2-Chloroethyl) ether-d	6.78	67	179358	12.23	ng/ul	0.00
9) 2-Chlorophenol-d4	6.96	132	215494	12.36	ng/ul	0.00
13) 4-Methylphenol-d8	8.15	113	210583	10.53	ng/ul	0.00
19) Nitrobenzene-d5	8.58	128	98368	12.66	ng/ul	0.00
22) 2-Nitrophenol-d4	9.30	143	105051	12.30	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.84	165	192726	11.61	ng/ul	0.00
29) 4-Chloroaniline-d4	10.35	131	253211	12.56	ng/ul	0.00
44) Dimethylphthalate-d6	13.52	166	631305	13.59	ng/ul	0.00
47) Acenaphthylene-d8	13.77	160	730807	13.13	ng/ul	0.00
52) 4-Nitrophenol-d4	14.37	143	58568m	7.02	ng/ul	0.02
58) Fluorene-d10	15.09	176	533380	13.10	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.26	200	29811	3.91	ng/ul	0.00
71) Anthracene-d10	16.94	188	811154	14.25	ng/ul	0.00
79) Pyrene-d10	19.26	212	899612	16.56	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.06	264	1016188	15.31	ng/ul	-0.01

Target Compounds

					Ovalue	
6) Phenol	6.66	94	47051	1.863	ng/ul	93
45) Dimethylphthalate	13.56	163	135943	2.816	ng/ul	99

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

> 5012/27/19