

Quantitation Report (Qedit)

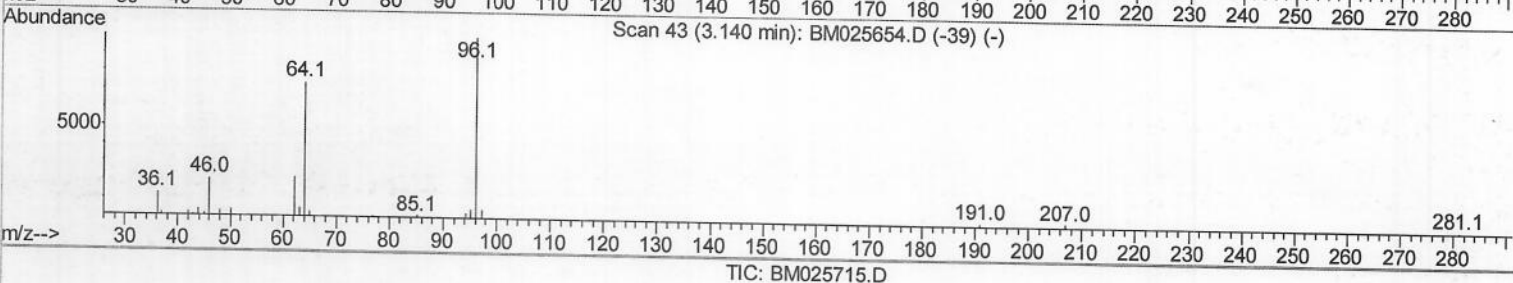
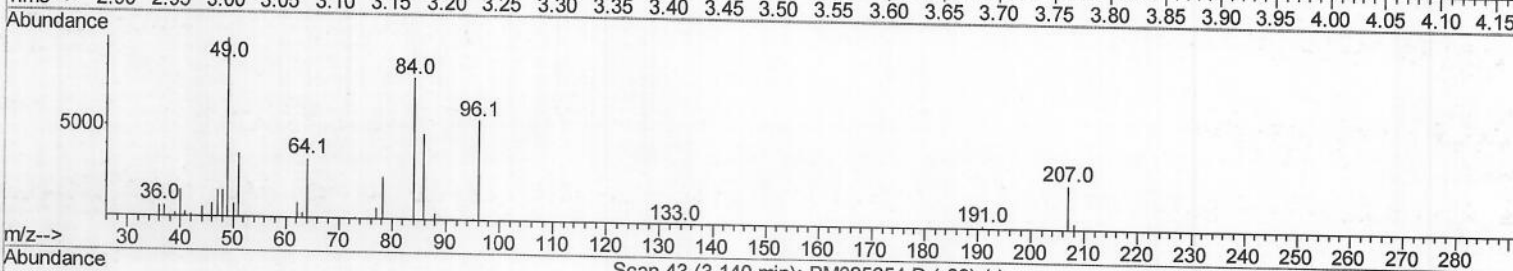
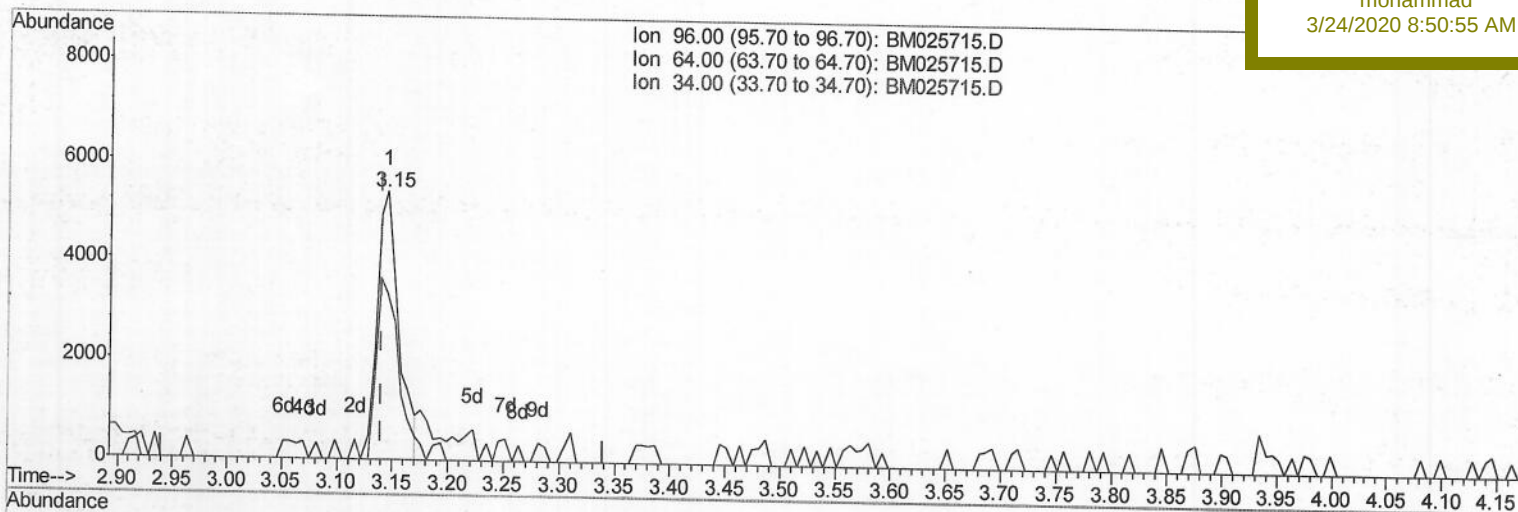
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM032320\
 Data File : BM025715.D
 Acq On : 24 Mar 2020 05:42
 Operator : CG/JU
 Sample : L1890-17
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Quant Time: Mar 24 07:02:55 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM031420MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Mar 24 04:07:11 2020
 Response via : Initial Calibration

Instrument :
 BNA_M
 ClientSampleId :
 C0DH4

Manual Integrations
 APPROVED

mohammad
 3/24/2020 8:50:55 AM



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

3.146min (+0.006) 1.58ng/uL

response 7167

Ion	Exp%	Act%
96.00	100	100
64.00	75.60	61.32
34.00	0.00	0.00
0.00	0.00	0.00

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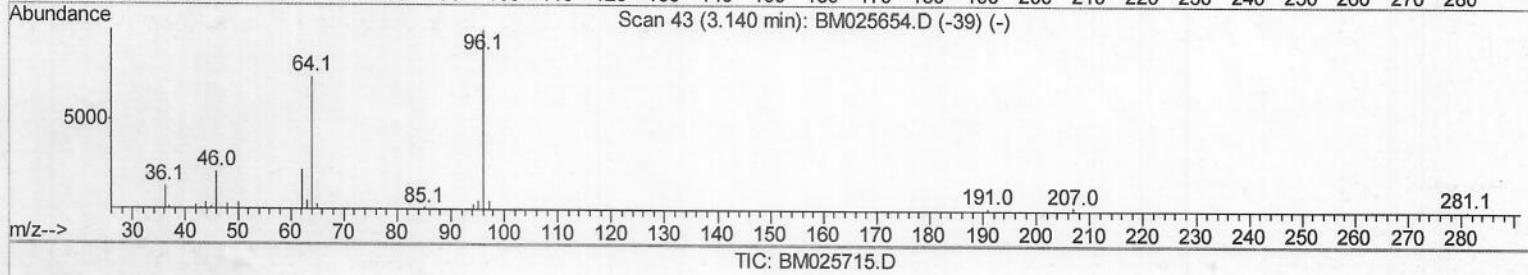
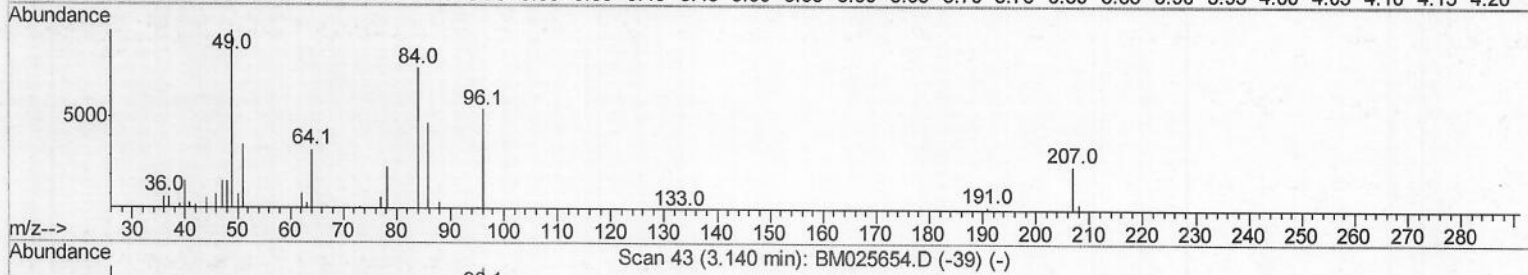
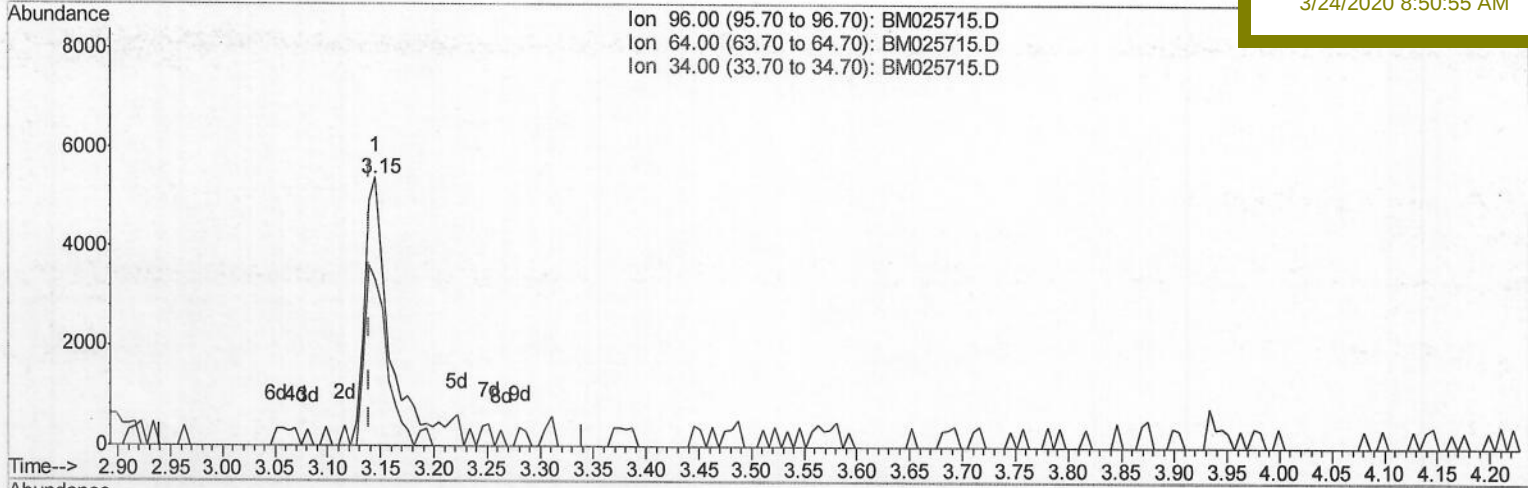
C0DH4

Manual Integrations

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

3.146min (+0.006) 1.97ng/uL m

response 8903

Ion Exp% Act%

96.00 100 100

64.00 75.60 61.32

34.00 0.00 0.00

0.00 0.00 0.00

JU
03/26/2020

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) 1,4-Dichlorobenzene-d4	7.58	152	189841	20.00	ng/ul	0.00
18) Naphthalene-d8	10.34	136	750151	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.21	164	468716	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.96	188	1033671	20.00	ng/ul	0.00
78) Chrysene-d12	21.15	240	1131241	20.00	ng/ul	0.00
86) Perylene-d12	23.30	264	1265763	20.00	ng/ul	-0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.15	96	8903m7	1.97	ng/uL	0.00
5) Phenol-d5	6.76	99	146161	9.76	ng/ul	0.00
7) Bis-(2-Chloroethyl) ether-d	6.92	67	270618	31.64	ng/ul	0.00
9) 2-Chlorophenol-d4	7.11	132	349228	28.72	ng/ul	0.00
13) 4-Methylphenol-d8	8.28	113	259444	21.07	ng/ul	0.00
19) Nitrobenzene-d5	8.72	128	204229	36.63	ng/ul	0.00
22) 2-Nitrophenol-d4	9.43	143	224840	37.93	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.96	165	435322	35.91	ng/ul	0.00
29) 4-Chloroaniline-d4	10.47	131	39224	2.79	ng/ul	0.00
44) Dimethylphthalate-d6	13.63	166	1447953	40.65	ng/ul	0.00
47) Acenaphthylene-d8	13.90	160	1742279	40.82	ng/ul	0.00
52) 4-Nitrophenol-d4	14.42	143	64842	9.56	ng/ul	0.00
58) Fluorene-d10	15.21	176	1276189	40.81	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.33	200	207967	32.98	ng/ul	0.00
71) Anthracene-d10	17.06	188	2053163	42.53	ng/ul	0.00
79) Pyrene-d10	19.36	212	2287710	42.10	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.17	264	2833435	43.72	ng/ul	0.00

Target Compounds

					Ovalue	
6) Phenol	6.78	94	189288	12.098	ng/ul	96
10) 2-Chlorophenol	7.14	128	32952	2.574	ng/ul	93
32) Caprolactam	11.25	113	72543	18.241	ng/ul	87
45) Dimethylphthalate	13.67	163	48764	1.315	ng/ul	99

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

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