

Quantitation Report (Qedit)

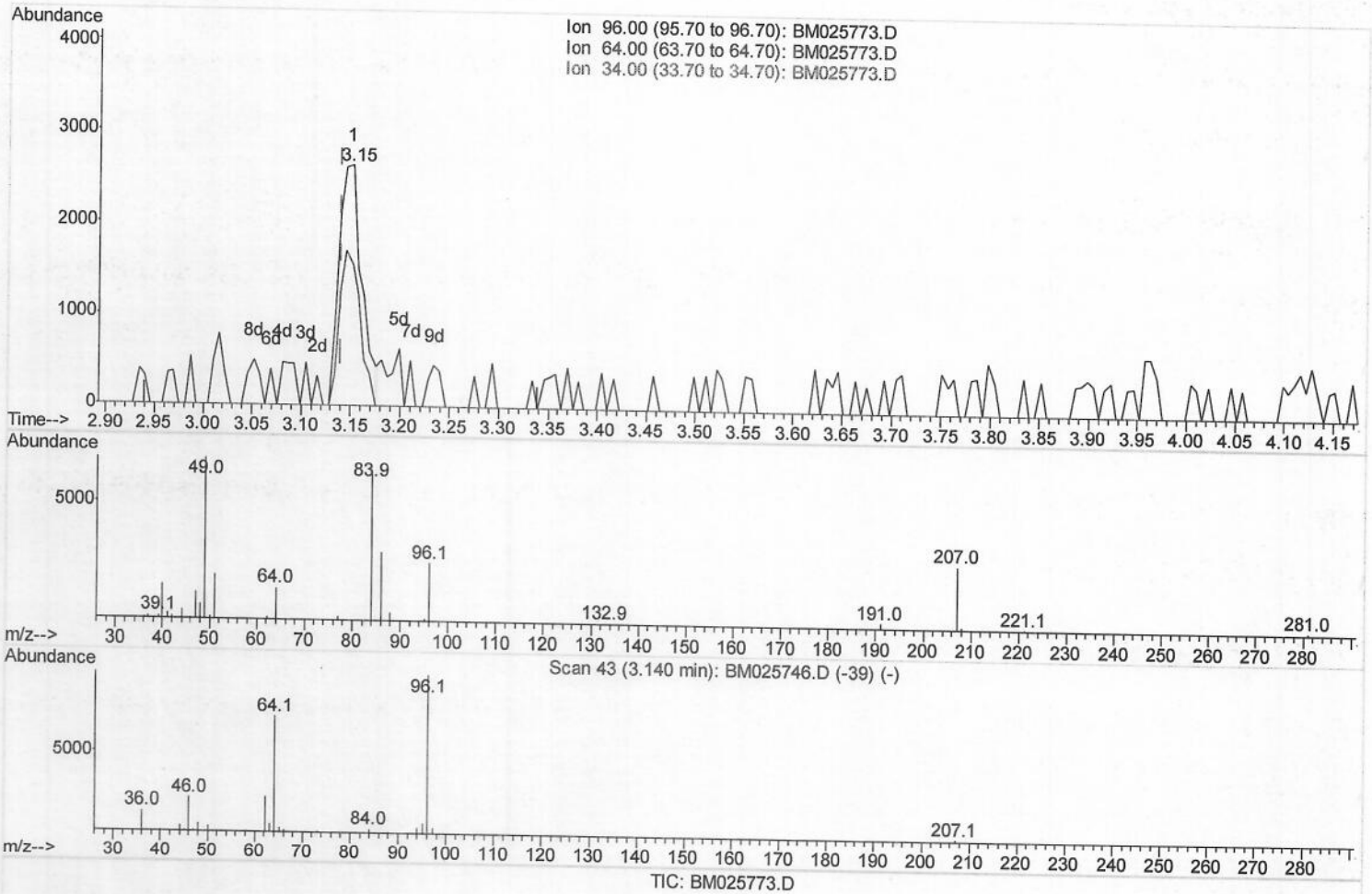
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM032320\
 Data File : BM025773.D
 Acq On : 25 Mar 2020 18:19
 Operator : CG/JU
 Sample : PB127634BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK34

Manual Integrations
 APPROVED

mohammad
 3/27/2020 8:15:11 AM

Quant Time: Mar 26 00:38:10 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM031420MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Mar 24 00:45:06 2020
 Response via : Initial Calibration



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

3.152min (+0.012) 0.95ng/uL

response 4239

Ion	Exp%	Act%
96.00	100	100
64.00	75.60	58.62#
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

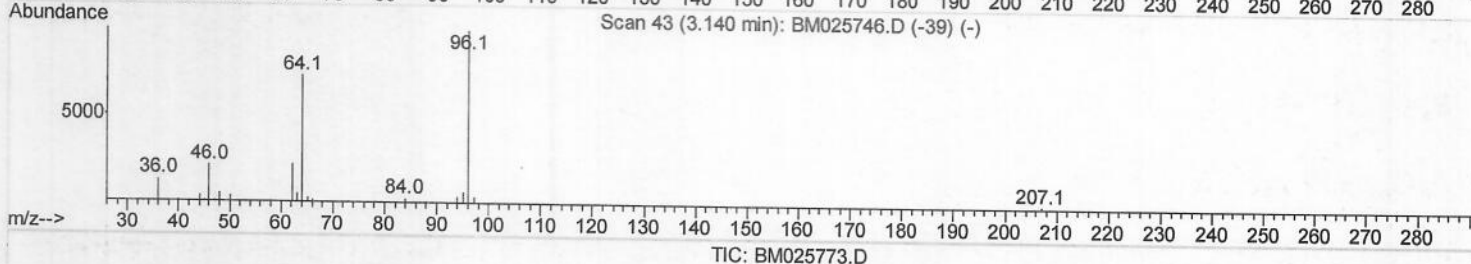
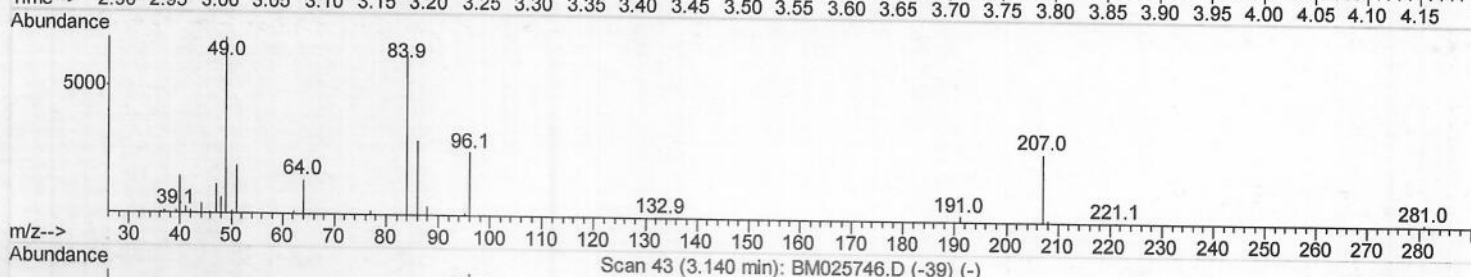
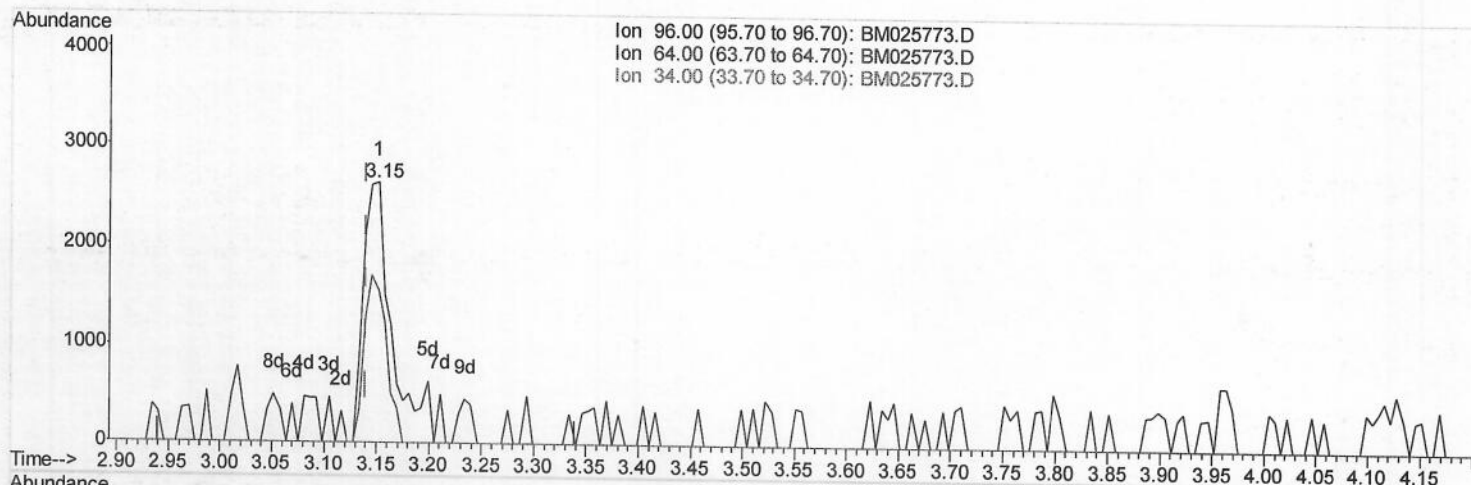
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(3) 1,4-Dioxane-d8 (S)

3.152min (+0.012) 1.10ng/uL m >

response 4883

Ion	Exp%	Act%
96.00	100	100
64.00	75.60	58.62#
34.00	0.00	0.00
0.00	0.00	0.00

JU
 03/28/2020

Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM032320\
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 Acq On : 25 Mar 2020 18:19
 Operator : CG/JU
 Sample : PB127634BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
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 Client Sampled :
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Manual Integrations
 APPROVED

mohammad
 3/27/2020 8:15:11 AM

Quant Time: Mar 26 00:39:00 2020

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM031420MA.M

Quant Title : SVOA CALIBRATION

QLast Update : Tue Mar 24 00:45:06 2020

Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	187066	20.00	ng/ul	0.00
18) Naphthalene-d8	10.33	136	754110	20.00	ng/ul	-0.01
36) Acenaphthene-d10	14.20	164	472508	20.00	ng/ul	-0.01
62) Phenanthrene-d10	16.95	188	1025064	20.00	ng/ul	-0.01
78) Chrysene-d12	21.14	240	1095839	20.00	ng/ul	-0.01
86) Perylene-d12	23.30	264	1220914	20.00	ng/ul	-0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.15	96	4883m	1.10	ng/uL	0.01
5) Phenol-d5	6.75	99	60181	4.08	ng/ul	-0.01
7) Bis-(2-Chloroethyl)ether-d	6.91	67	171884	20.40	ng/ul	-0.01
9) 2-Chlorophenol-d4	7.10	132	208642	17.41	ng/ul	-0.01
13) 4-Methylphenol-d8	8.27	113	116868	9.63	ng/ul	-0.01
19) Nitrobenzene-d5	8.71	128	131980	23.55	ng/ul	0.00
22) 2-Nitrophenol-d4	9.43	143	137025	22.99	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.96	165	241492	19.82	ng/ul	0.00
29) 4-Chloroaniline-d4	10.47	131	15773	1.11	ng/ul	0.00
44) Dimethylphthalate-d6	13.62	166	844071	23.50	ng/ul	-0.01
47) Acenaphthylene-d8	13.89	160	1030174	23.94	ng/ul	-0.01
52) 4-Nitrophenol-d4	14.42	143	21610	3.16	ng/ul	0.00
58) Fluorene-d10	15.20	176	766189	24.31	ng/ul	-0.01
63) 4,6-Dinitro-2-methylphenol	15.33	200	111788	17.87	ng/ul	-0.01
71) Anthracene-d10	17.05	188	1245978	26.02	ng/ul	0.00
79) Pyrene-d10	19.35	212	1358842	25.82	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.16	264	1638689	26.22	ng/ul	-0.01

Target Compounds

45) Dimethylphthalate	13.67	163	79724	2.133	ng/ul	Ovalue 99
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(#) = qualifier out of range (m) = manual integration (+) = signals summed

JU
 03/28/2020

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