

Quantitation Report (Qedit)

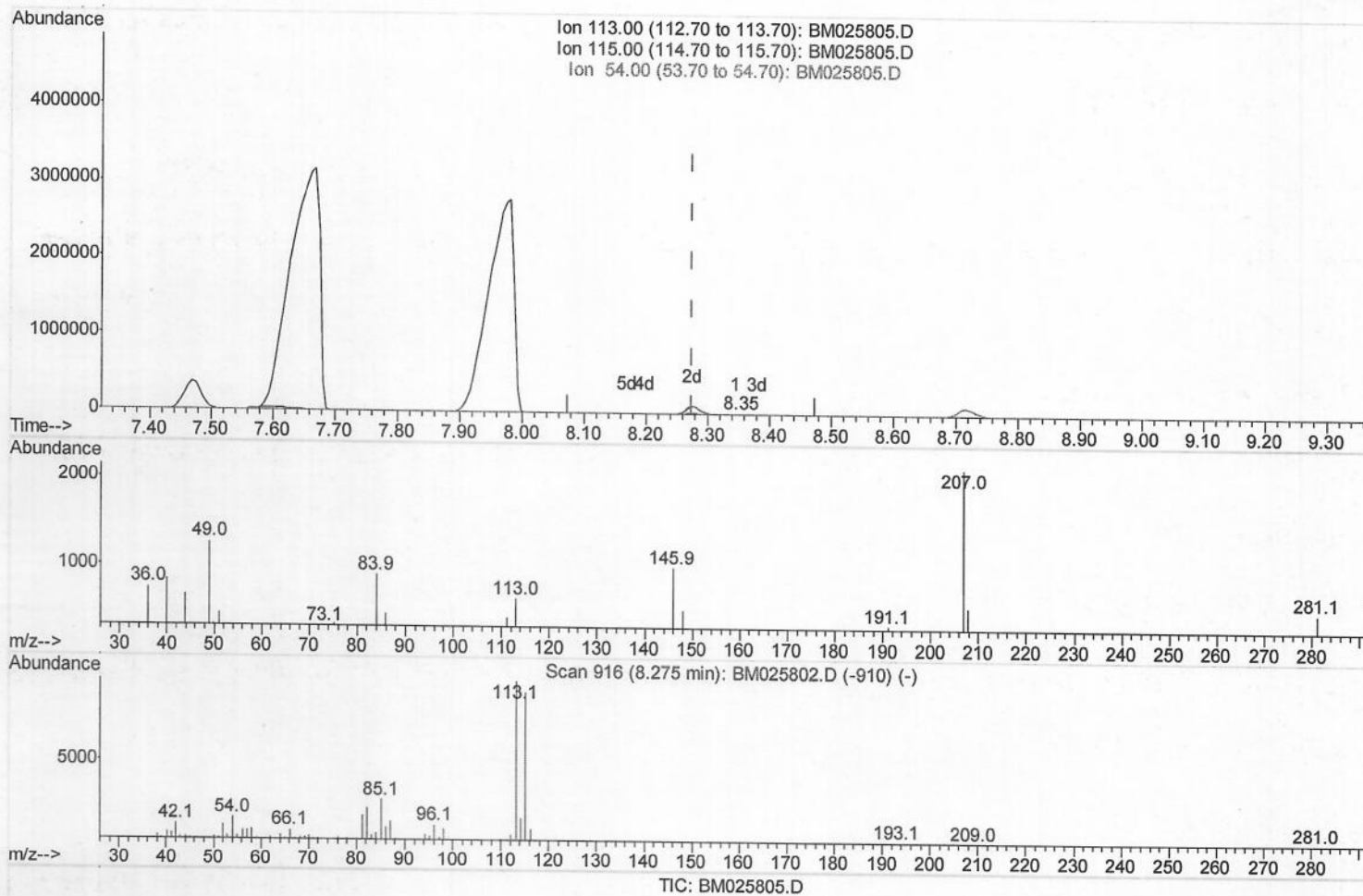
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM032620\
 Data File : BM025805.D
 Acq On : 26 Mar 2020 18:16
 Operator : CG/JU
 Sample : L1968-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0DC7

Manual Integrations
 APPROVED

mohammad
 3/30/2020 8:30:40 AM

Quant Time: Mar 26 18:51:35 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM031420MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Mar 26 18:34:50 2020
 Response via : Initial Calibration



(13) 4-Methylphenol-d8 (S)

8.346min (+0.071) 0.04ng/ul

response 509

Ion	Exp%	Act%
113.00	100	100
115.00	95.30	0.00#
54.00	13.30	0.00#
0.00	0.00	0.00

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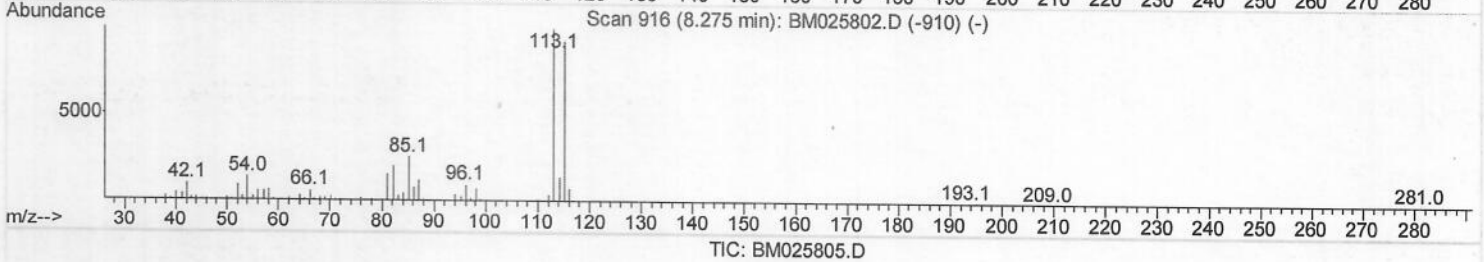
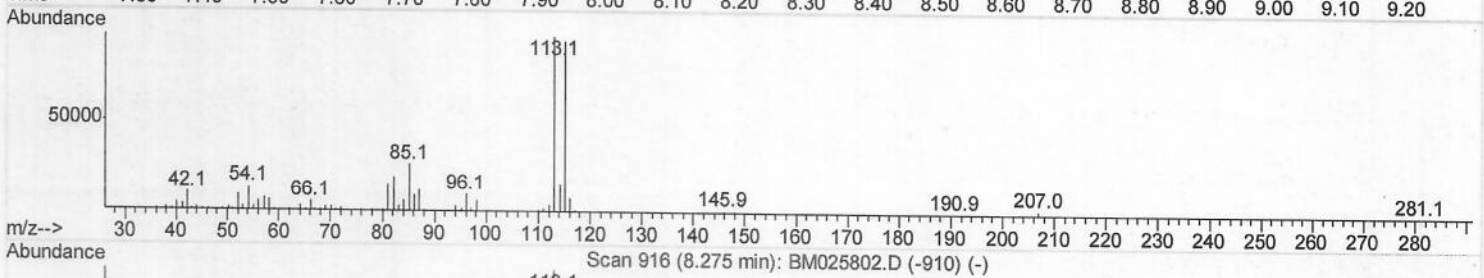
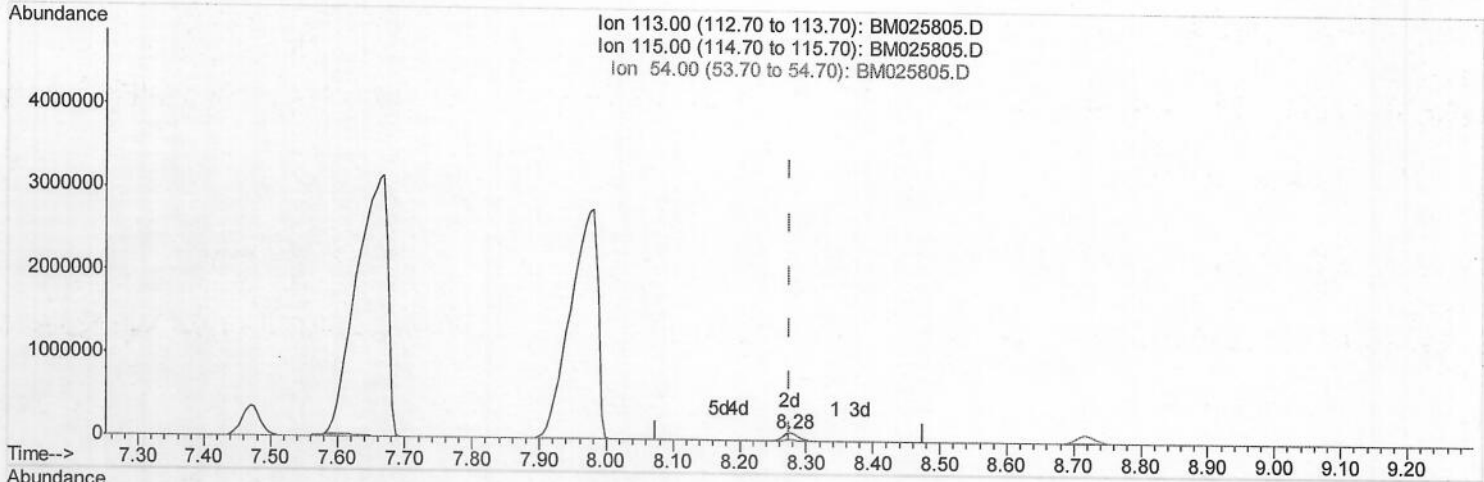
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(13) 4-Methylphenol-d8 (S)

8.275min (+0.000) 11.69ng/ul m >

response 144146

Ion	Exp%	Act%
113.00	100	100
115.00	95.30	97.31
54.00	13.30	13.63
0.00	0.00	0.00

JU
 03/31/2020

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	190127	20.00	ng/ul	0.04
18) Naphthalene-d8	10.34	136	664915	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.20	164	393976	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.95	188	842736	20.00	ng/ul	0.00
78) Chrysene-d12	21.14	240	926750	20.00	ng/ul	0.00
86) Perylene-d12	23.30	264	1034658	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.14	96	4967	1.10	ng/uL	0.00
5) Phenol-d5	6.75	99	72841	4.85	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.92	67	228696	26.70	ng/ul	0.01
9) 2-Chlorophenol-d4	7.11	132	260832	21.42	ng/ul	0.00
13) 4-Methylphenol-d8	8.28	113	144146m	11.69	ng/ul	0.00
19) Nitrobenzene-d5	8.72	128	169564	34.31	ng/ul	0.00
22) 2-Nitrophenol-d4	9.43	143	178022	33.88	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.96	165	308641	28.73	ng/ul	0.00
29) 4-Chloroaniline-d4	10.47	131	18916	1.52	ng/ul	0.00
44) Dimethylphthalate-d6	13.62	166	1015960	33.93	ng/ul	0.00
47) Acenaphthylene-d8	13.89	160	1246198	34.74	ng/ul	0.00
52) 4-Nitrophenol-d4	14.41	143	24132	4.23	ng/ul	0.00
58) Fluorene-d10	15.20	176	905336	34.45	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.33	200	134055	26.07	ng/ul	0.00
71) Anthracene-d10	17.05	188	1408576	35.78	ng/ul	0.00
79) Pyrene-d10	19.34	212	1539600	34.59	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.16	264	1910841	36.07	ng/ul	0.00
Target Compounds						
20) Nitrobenzene	8.78	77	7199801	597.993	ng/ul	98
37) 1,2,4,5-Tetrachlorobenzene	12.38	216	402078	32.268	ng/ul	98
45) Dimethylphthalate	13.67	163	102853	3.300	ng/ul	98
73) 1,2,3,4-Tetrachlorobenzene	12.99	216	543940	42.122	ng/uL	96
74) Pentachlorobenzene	14.53	250	23501	1.635	ng/uL	96

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

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