

Quantitation Report (Qedit)

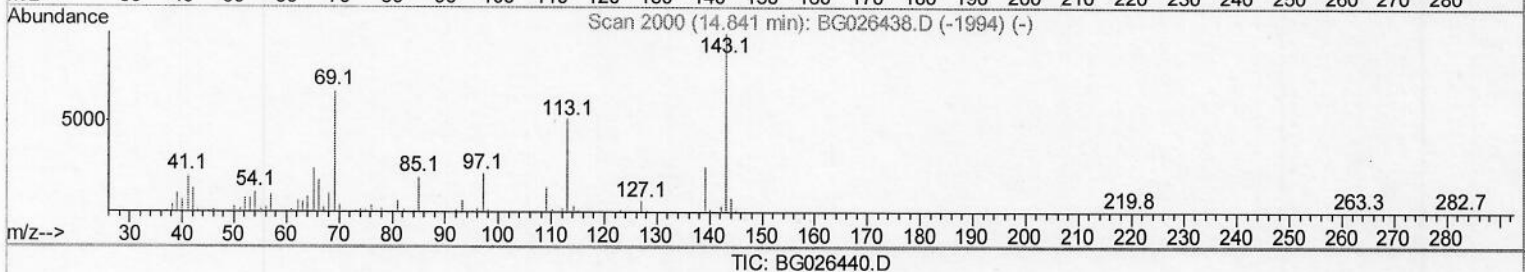
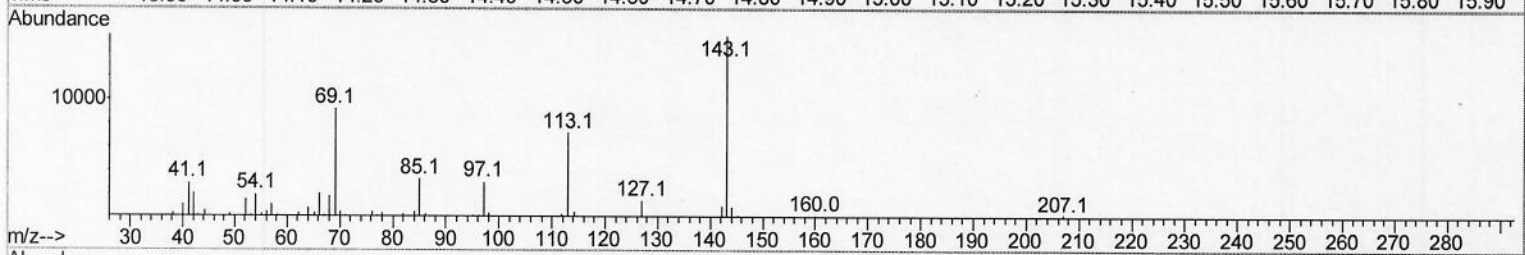
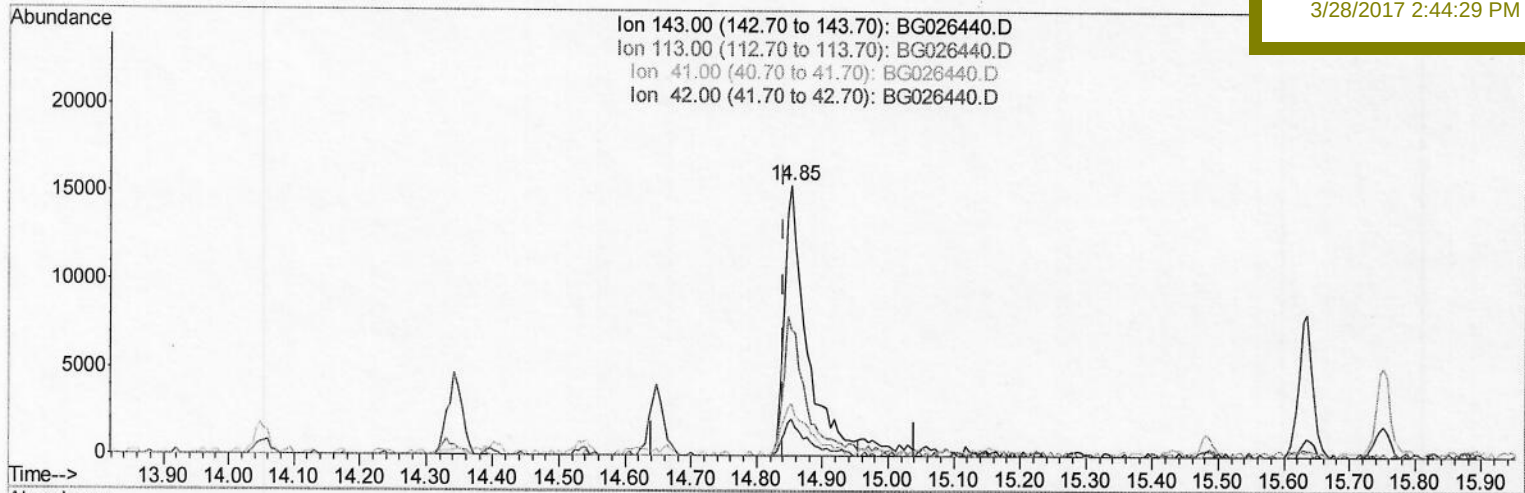
Data Path : Z:\HPCHEM1\BNA G\Data\BG032717\
 Data File : BG026440.D
 Acq On : 27 Mar 2017 19:52
 Operator : SJ/MA
 Sample : I2358-12
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Quant Time: Mar 28 01:23:58 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM-EPA-BG032717MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Mar 28 01:23:51 2017
 Response via : Initial Calibration

Instrument :
 BNA_G
 ClientSampleId :
 C0BL1

Manual Integrations
 APPROVED

mohammad
 3/28/2017 2:44:29 PM



(51) 4-Nitrophenol-d4 (S)

14.853min (+0.012) 6.06ng/ul m

response 36754

Ion	Exp%	Act%
143.00	100	100
113.00	75.10	46.79#
41.00	28.80	19.41#
42.00	19.10	13.62#

S-J
 03/28/2017

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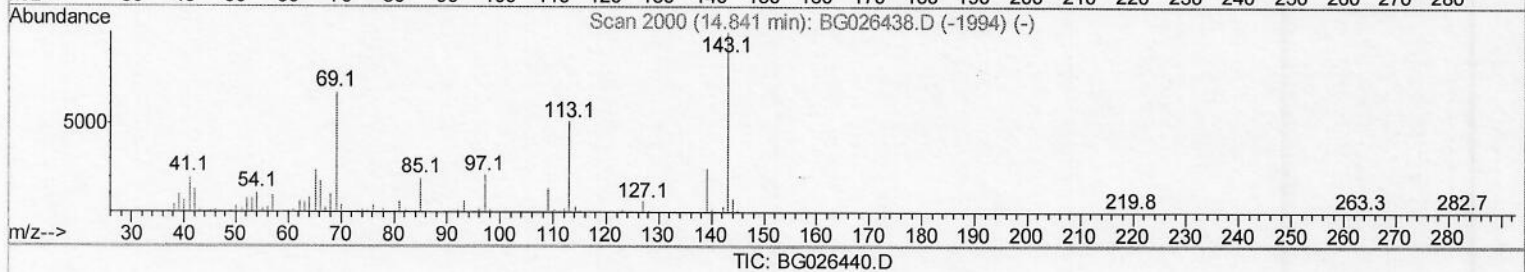
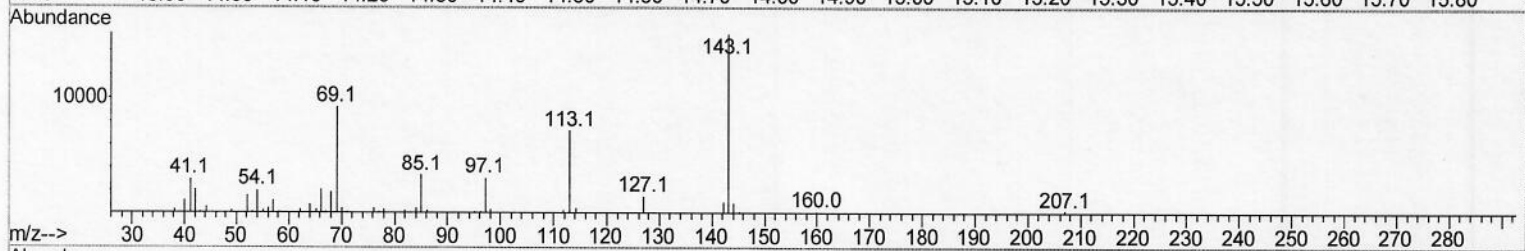
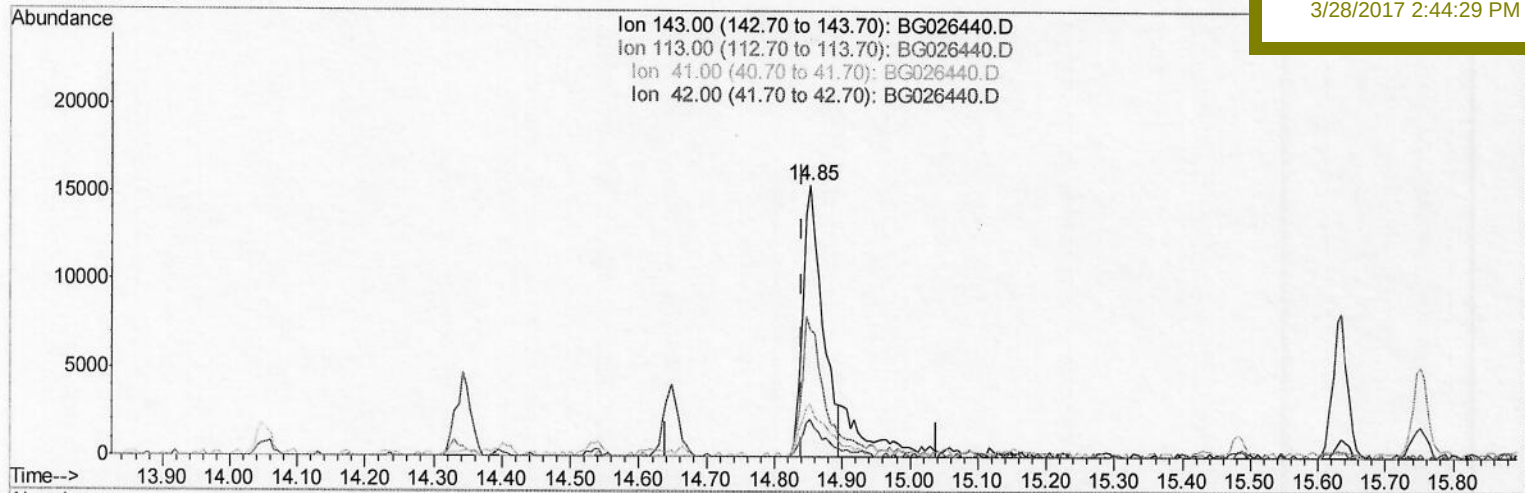
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(51) 4-Nitrophenol-d4 (S)

14.853min (+0.012) 5.17ng/ul

response 31345

Ion	Exp%	Act%
143.00	100	100
113.00	75.10	46.79#
41.00	28.80	19.41#
42.00	19.10	13.62#

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.04	152	167870	20.00	nq/ul	0.00
18) Naphthalene-d8	10.84	136	697096	20.00	nq/ul	0.00
35) Acenaphthene-d10	14.65	164	402811	20.00	nq/ul	0.00
61) Phenanthrene-d10	17.38	188	973040	20.00	nq/ul	0.00
77) Chrysene-d12	21.63	240	1130432	20.00	nq/ul	0.00
85) Perylene-d12	24.80	264	1106246	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.47	96	20168	5.75	nq/uL	0.00
5) Phenol-d5	7.21	99	86647	7.01	nq/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.37	67	216827	31.30	nq/ul	0.00
9) 2-Chlorophenol-d4	7.58	132	277336	25.02	nq/ul	0.00
13) 4-Methylphenol-d8	8.74	113	158734	15.94	nq/ul	0.00
19) Nitrobenzene-d5	9.20	128	161500	32.42	nq/ul	0.00
22) 2-Nitrophenol-d4	9.92	143	193312	33.11	nq/ul	0.00
26) 2,4-Dichlorophenol-d3	10.46	165	312029	29.12	nq/ul	0.00
29) 4-Chloroaniline-d4	10.99	131	10198	0.96	nq/ul	0.01
43) Dimethylphthalate-d6	14.05	166	1077562	34.55	nq/ul	0.00
46) Acenaphthylene-d8	14.34	160	1283642	33.03	nq/ul	0.00
51) 4-Nitrophenol-d4	14.85	143	36754m	6.06	nq/ul	0.01
57) Fluorene-d10	15.63	176	986543	34.82	nq/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.75	200	197391	31.55	nq/ul	0.00
70) Anthracene-d10	17.48	188	1550601	35.09	nq/ul	0.00
78) Pyrene-d10	19.75	212	1765484	35.01	nq/ul	0.00
89) Benzo(a)pyrene-d12	24.59	264	1773636	37.32	ng/ul	0.00

Target Compounds Ovalue

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

S.S
 03/28/2017

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