

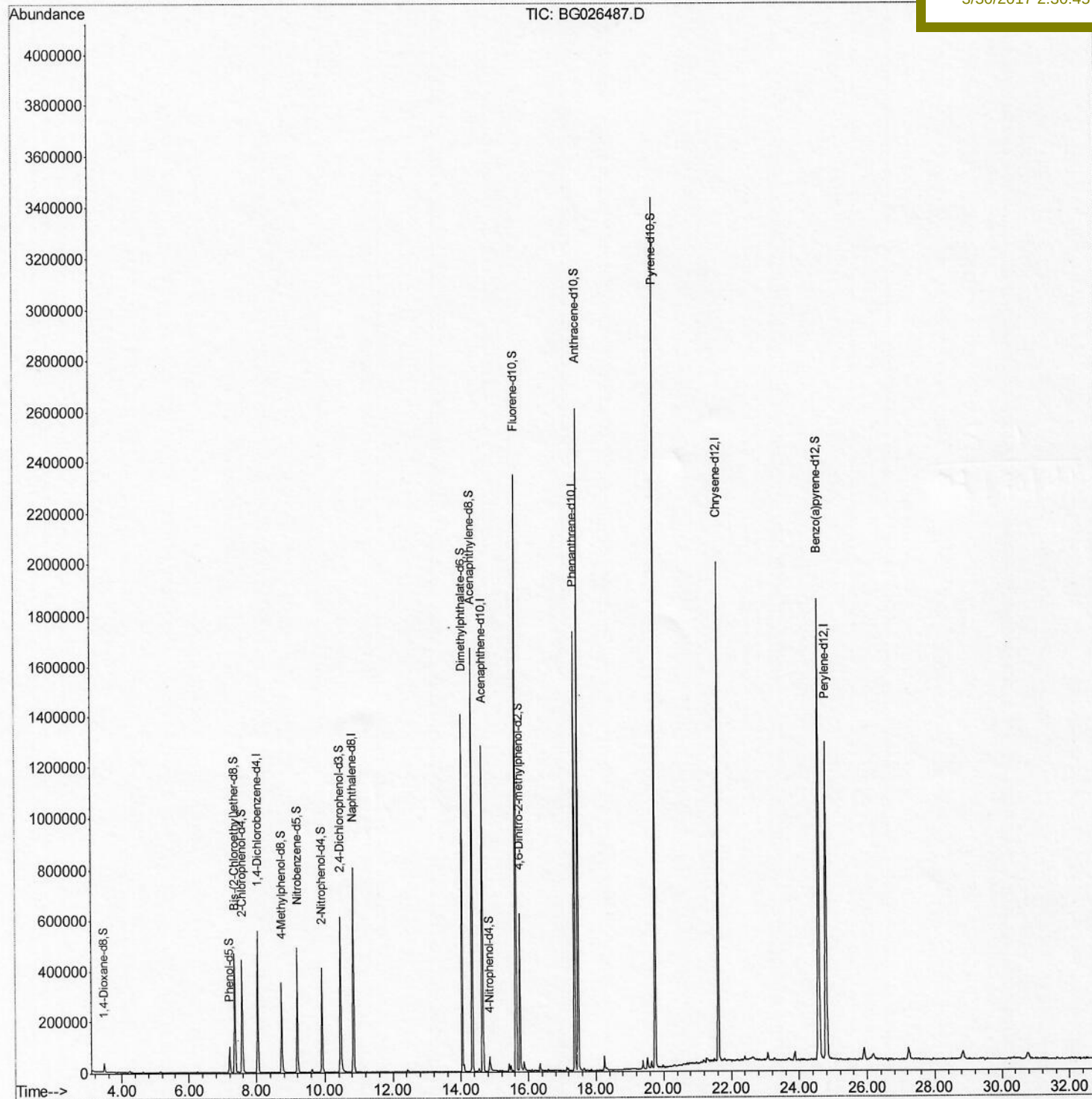
Data Path : Z:\HPCHEM1\BNA G\DATA\BG032917\
Data File : BG026487.D
Acq On : 29 Mar 2017 22:02
Operator : SJ/MA
Sample : I2397-06
Misc :
ALS Vial : 11 Sample Multiplier: 1

Quant Time: Mar 30 03:12:37 2017
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM-EPA-BG032717MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Mar 30 02:19:03 2017
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
G9Y90

Manual Integrations
APPROVED

mohammad
3/30/2017 2:30:45 PM



Quantitation Report (Qedit)

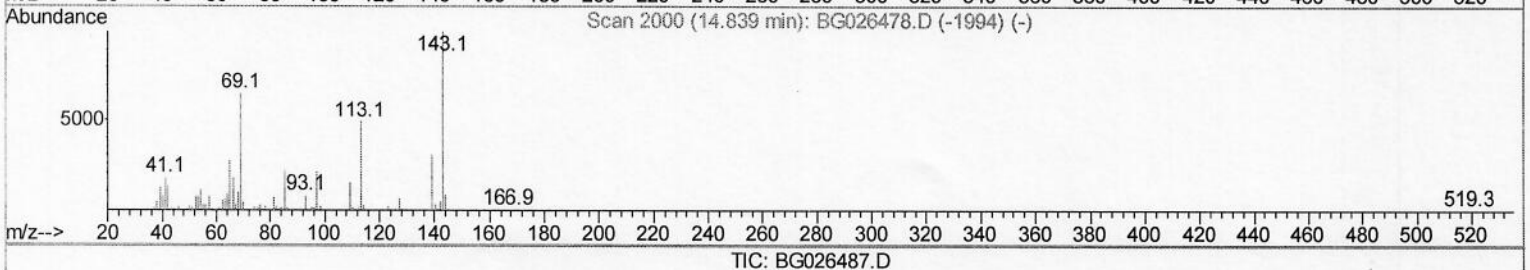
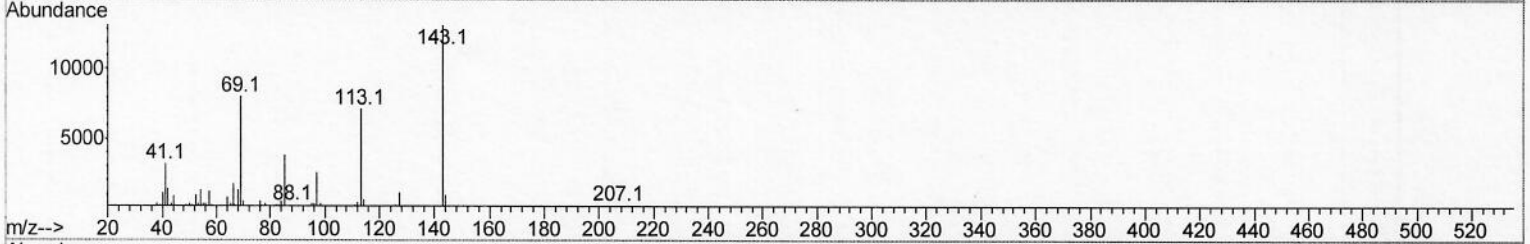
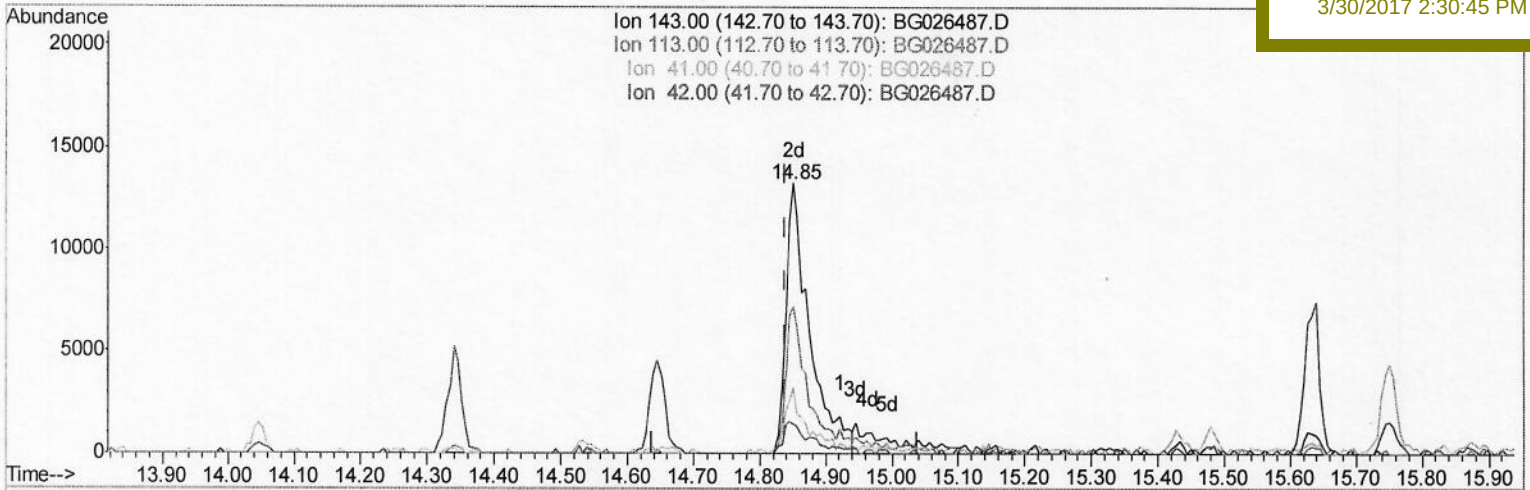
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BNA_G
ClientSampled :
G9Y90

Manual Integrations
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TIC: BG026487.D

(51) 4-Nitrophenol-d4 (S)

14.851min (+0.013) 4.68ng/ul m

response 32948

Ion	Exp%	Act%
143.00	100	100
113.00	75.10	53.91#
41.00	28.80	24.08
42.00	19.10	10.75#

SJ
03/30/2017

Quantitation Report (Qedit)

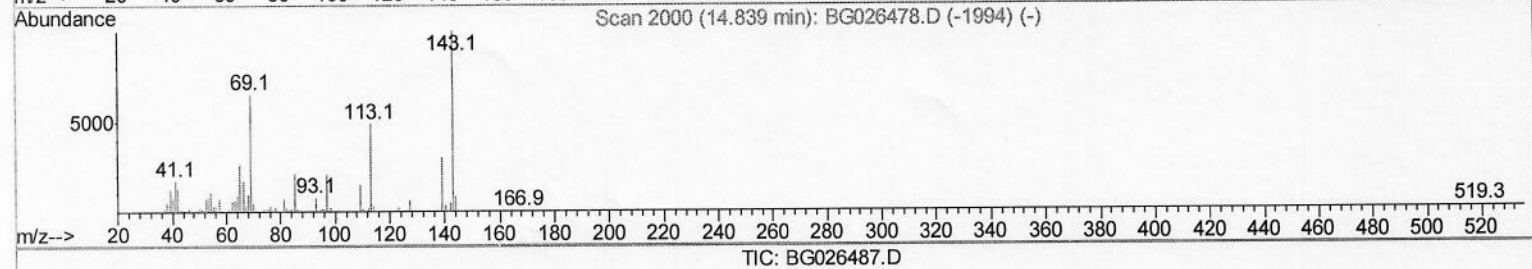
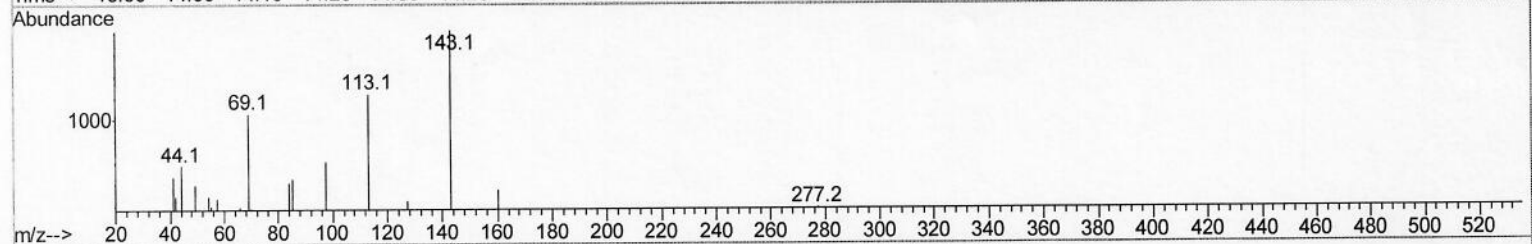
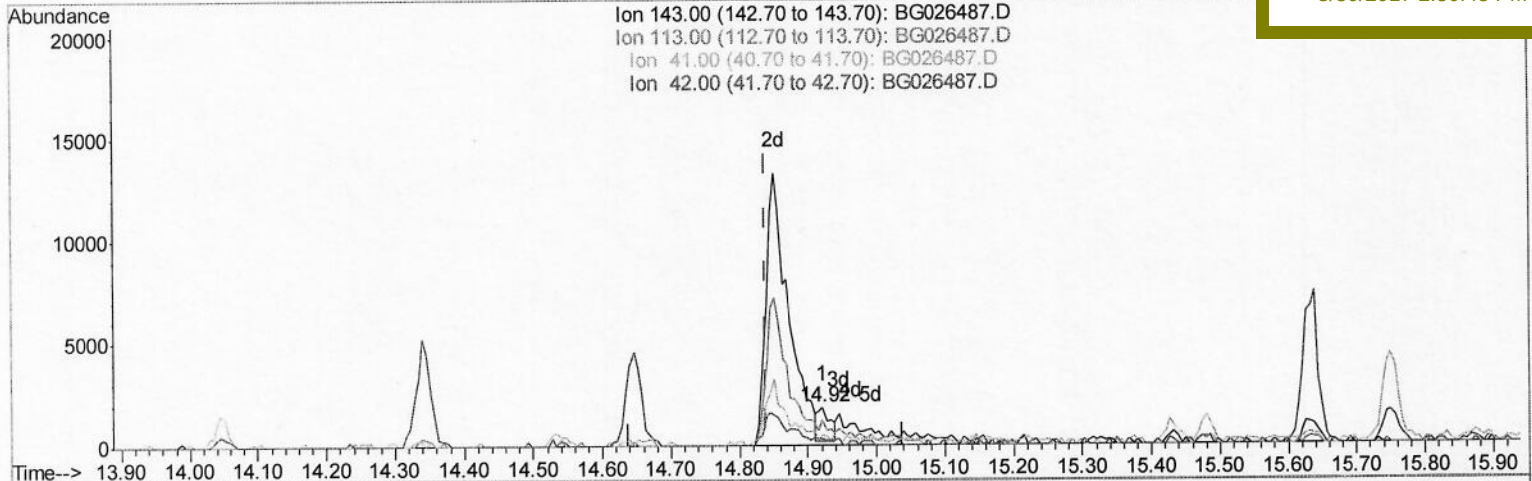
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 Sample : I2397-06
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Instrument :
 BNA_G
 ClientSampleId :
 G9Y90

Quant Time: Mar 30 02:19:45 2017
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Manual Integrations
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(51) 4-Nitrophenol-d4 (S)

14.922min (+0.083) 0.30ng/ul

response 2128

Ion	Exp%	Act%
143.00	100	100
113.00	75.10	67.89
41.00	28.80	25.89
42.00	19.10	15.61

Data Path : Z:\HPCHEM1\BNA G\DATA\BG032917\
 Data File : BG026487.D
 Acq On : 29 Mar 2017 22:02
 Operator : SJ/MA
 Sample : I2397-06
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 G9Y90

Quant Time: Mar 30 03:12:37 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM-EPA-BG032717MA.M
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Manual Integrations
 APPROVED

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) 1,4-Dichlorobenzene-d4	8.04	152	186842	20.00	ng/ul	0.00
18) Naphthalene-d8	10.84	136	785473	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.65	164	467736	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.38	188	1124513	20.00	ng/ul	0.00
77) Chrysene-d12	21.63	240	1253625	20.00	ng/ul	0.00
85) Perylene-d12	24.80	264	1255946	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.47	96	22100	5.66	ng/uL	0.00
5) Phenol-d5	7.21	99	83798	6.09	ng/ul	0.00
7) Bis-(2-Chloroethyl) ether-d	7.36	67	219742	28.50	ng/ul	0.00
9) 2-Chlorophenol-d4	7.58	132	279049	22.62	ng/ul	0.00
13) 4-Methylphenol-d8	8.74	113	161971	14.61	ng/ul	0.00
19) Nitrobenzene-d5	9.20	128	164253	29.26	ng/ul	0.00
22) 2-Nitrophenol-d4	9.92	143	199506	30.32	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.46	165	311218	25.78	ng/ul	0.00
29) 4-Chloroaniline-d4	0.00	131	0d	0.00	ng/ul	0.00
43) Dimethylphthalate-d6	14.05	166	1110143	30.65	ng/ul	0.00
46) Acenaphthylene-d8	14.34	160	1327873	29.43	ng/ul	0.00
51) 4-Nitrophenol-d4	14.85	143	32948m	4.68	ng/ul	0.01
57) Fluorene-d10	15.63	176	1033152	31.40	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.75	200	198525	27.45	ng/ul	0.00
70) Anthracene-d10	17.48	188	1625846	31.84	ng/ul	0.00
78) Pyrene-d10	19.76	212	1873009	33.50	ng/ul	0.00
89) Benzo(a)pyrene-d12	24.59	264	1847488	34.24	ng/ul	0.00

Target Compounds Ovalue

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

SJ
 03/30/2017