

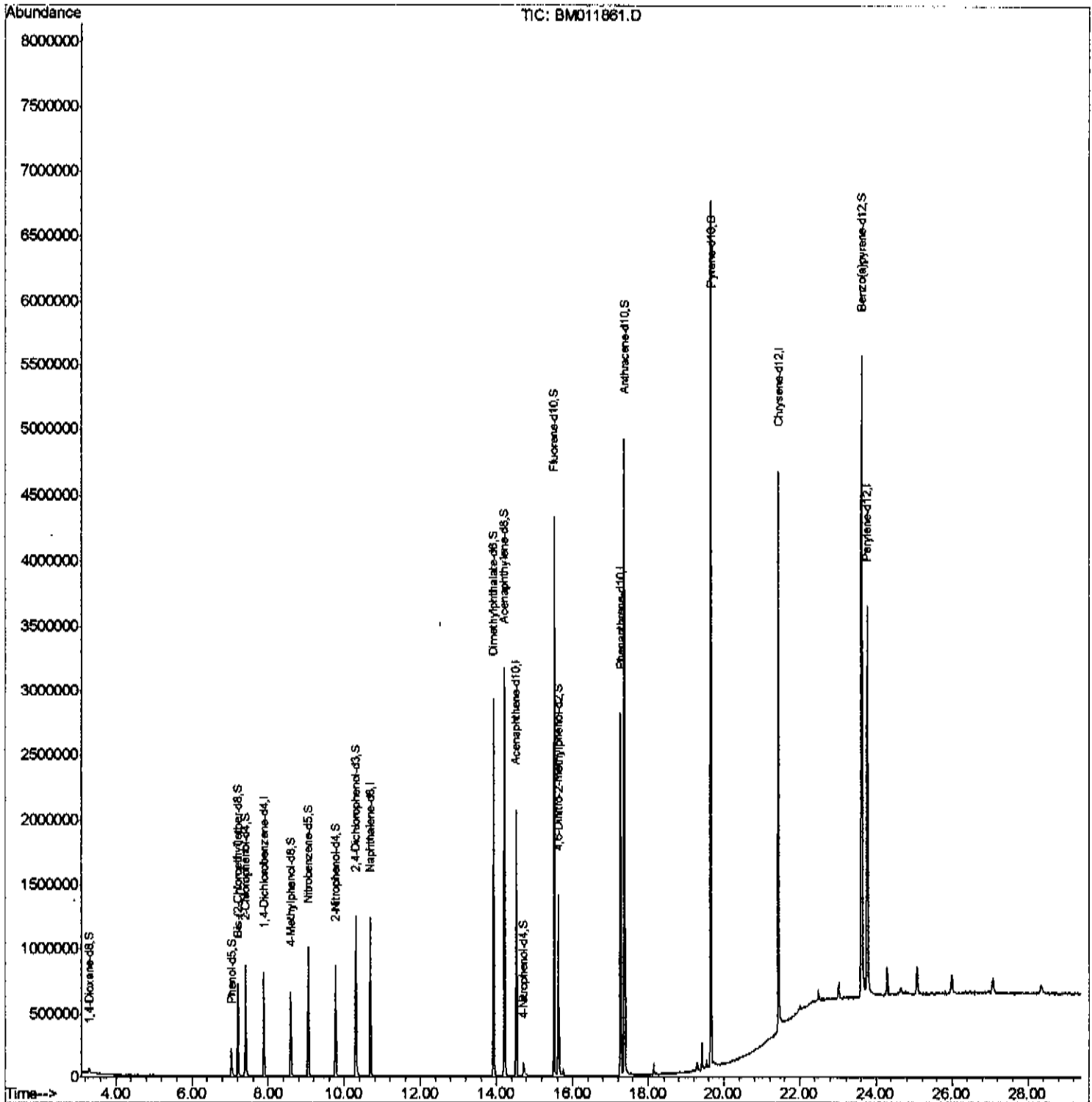
Data Path : Z:\HPCHEM1\BNA M\DATA\BM100317\
Data File : BM011861.D
Acq On : 03 Oct 2017 21:20
Operator : SJ/JU
Sample : I5525-09
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampled :
C0AA9

Manual Integrations
APPROVED

Sohil
10/4/2017 7:28:06 PM

Quant Time: Oct 04 04:23:28 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM100317.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Oct 03 15:47:54 2017
Response via : Initial Calibration



Quantitation Report (Qedit)

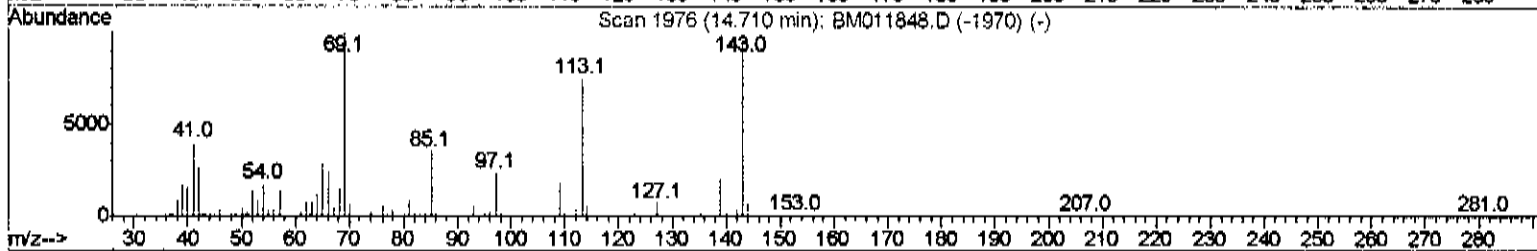
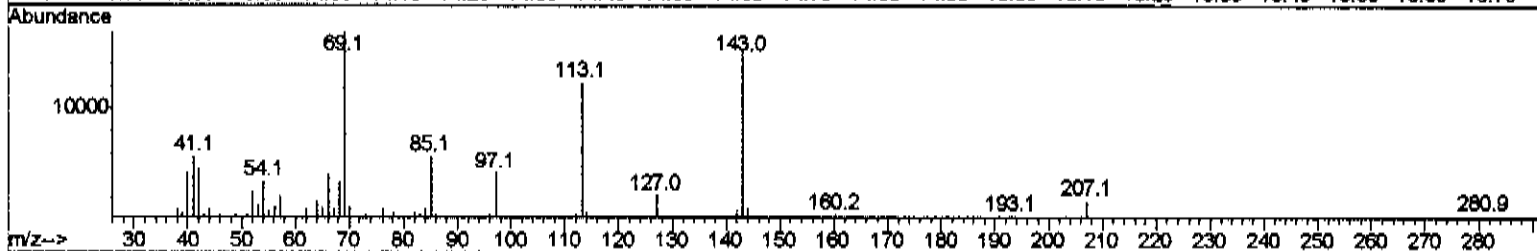
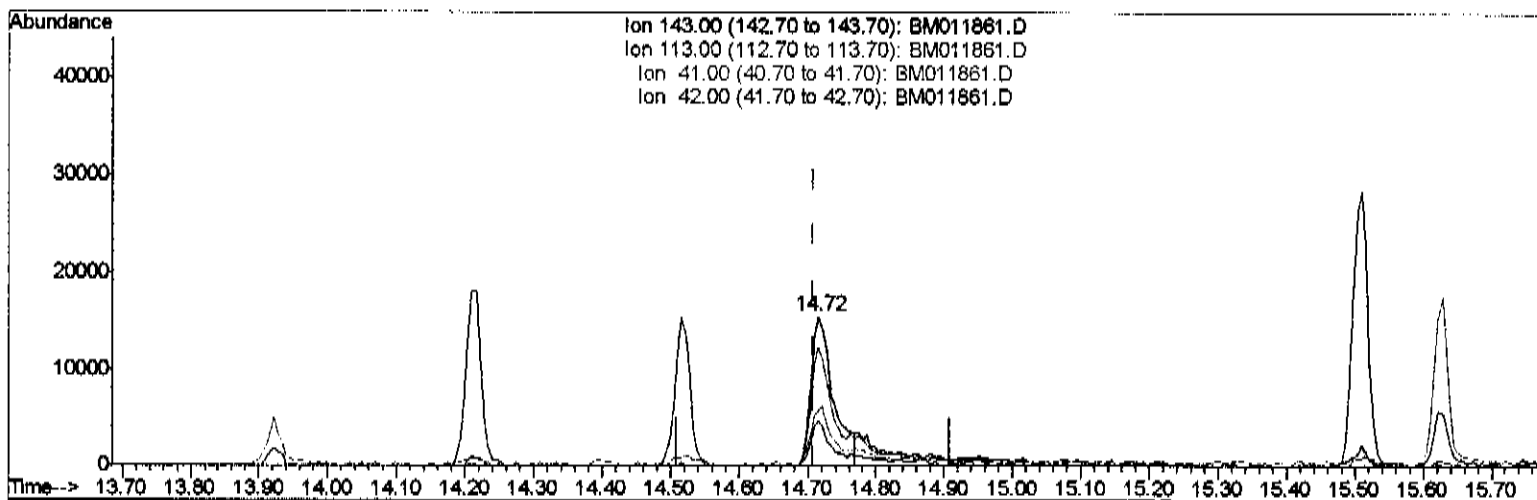
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TIC: BM011861.D

(51) 4-Nitrophenol-d4 (S)

14.715min (+0.006) 3.45ng/ul

response 35516

Ion	Exp%	Act%
143.00	100	100
113.00	114.60	79.59#
41.00	34.80	37.29
42.00	20.50	30.05#

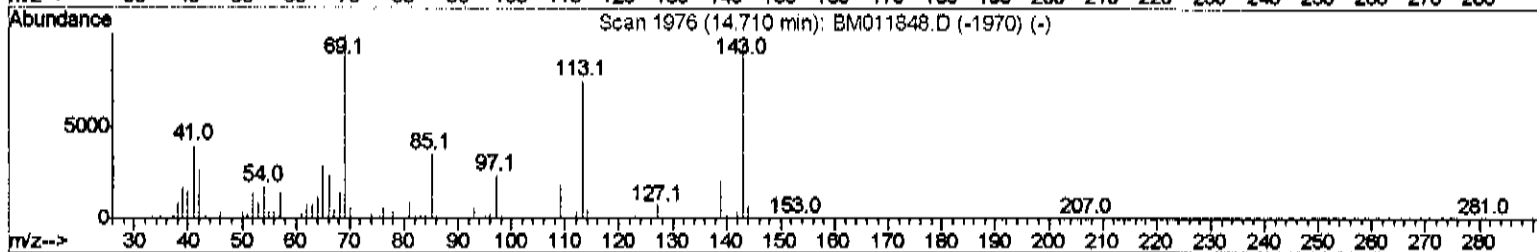
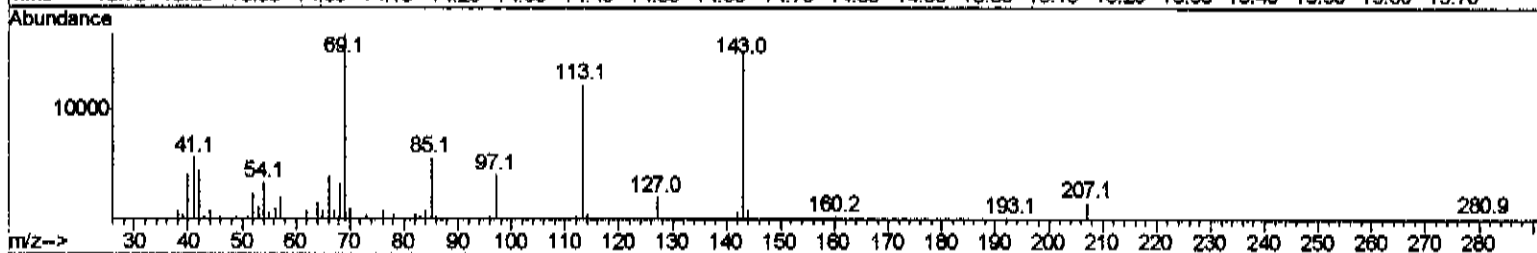
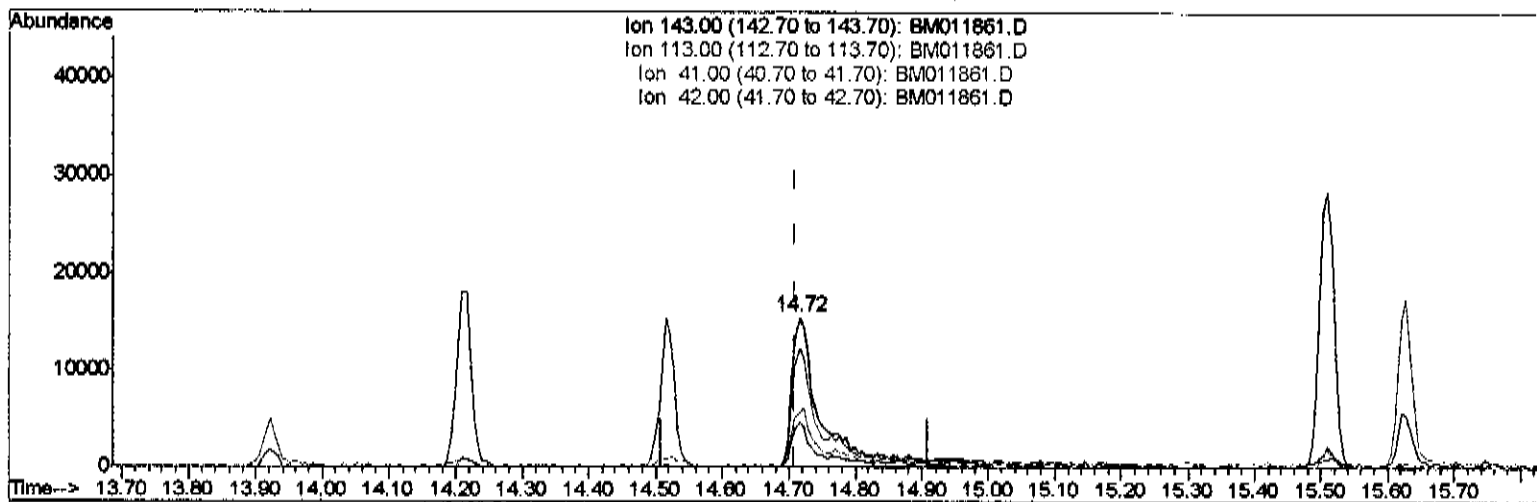
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TIC: BM011861.D

(51) 4-Nitrophenol-d4 (S)

14.715min (+0.006) 4.14ng/ul m

response 42585

Ion	Exp%	Act%
143.00	100	100
113.00	114.60	79.59#
41.00	34.80	37.29
42.00	20.50	30.05#

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Manual Integrations
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.87	152	230776	20.00	ng/ul	0.00
18) Naphthalene-d8	10.68	136	1036444	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.52	164	679655	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.26	188	1705063	20.00	ng/ul	0.00
75) Chrysene-d12	21.43	240	2065148	20.00	ng/ul	0.00
83) Perylene-d12	23.76	264	2228081	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.29	96	24517	5.02	ng/uL	0.00
5) Phenol-d5	7.03	99	137708	6.94	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.20	67	346757	29.51	ng/ul	0.00
9) 2-Chlorophenol-d4	7.40	132	403290	25.32	ng/ul	0.00
13) 4-Methylphenol-d8	8.58	113	264579	15.42	ng/ul	0.00
19) Nitrobenzene-d5	9.04	128	233655	31.66	ng/ul	0.00
22) 2-Nitrophenol-d4	9.76	143	271235	32.63	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.30	165	489509	29.14	ng/ul	0.00
29) 4-Chloroaniline-d4	10.84	131	3853	0.21	ng/ul	0.02
43) Dimethylphthalate-d6	13.92	166	1951083	33.12	ng/ul	0.00
46) Acenaphthylene-d8	14.21	160	2232513	32.25	ng/ul	0.00
51) 4-Nitrophenol-d4	14.72	143	42585ms	4.14	ng/ul	0.00
57) Fluorene-d10	15.51	176	1686813	32.45	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.63	200	307879	26.84	ng/ul	0.00
70) Anthracene-d10	17.36	188	2780571	33.09	ng/ul	0.00
76) Pyrene-d10	19.64	212	3453376	37.07	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.61	264	3498975	32.64	ng/ul	0.00

Target Compounds Ovalue

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