

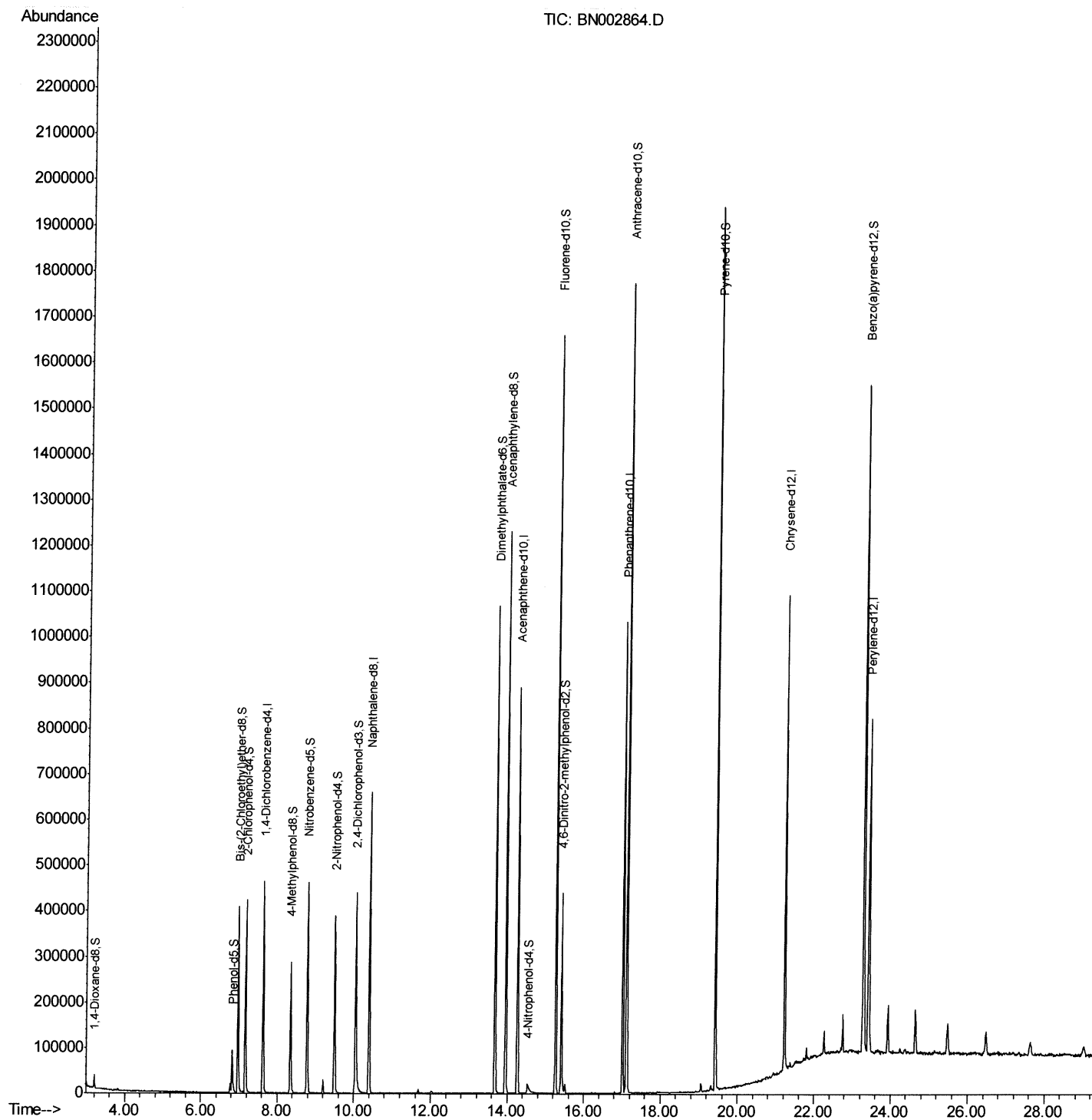
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Data File : BN002864.D
Acq On : 28 Sep 2018 23:17
Operator : MJ/SJ
Sample : J5117-07
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
JLC83

Manual Integrations
APPROVED

Sohil
10/1/2018 4:21:29 PM

Quant Time: Sep 29 06:58:07 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN091418MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Sep 29 06:35:10 2018
Response via : Initial Calibration



Quantitation Report (Qedit)

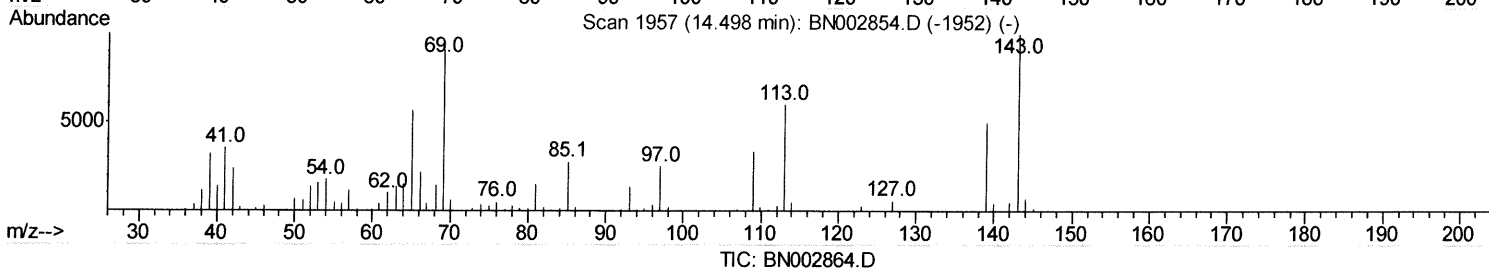
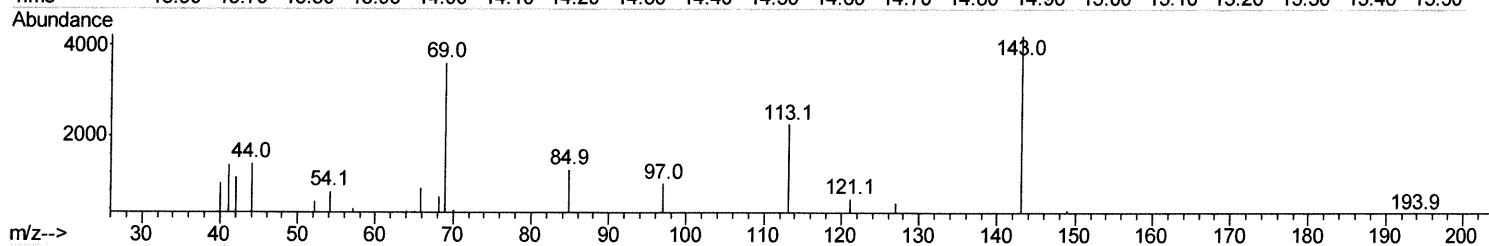
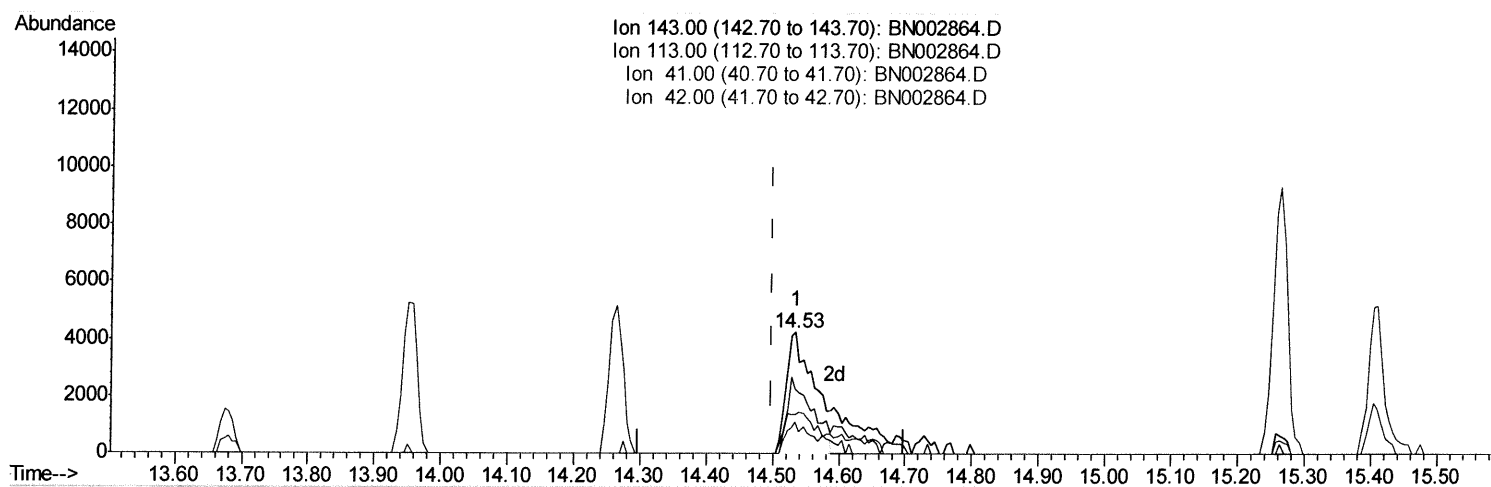
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10/1/2018 4:21:29 PM

Quant Time: Sep 29 06:36:28 2018
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Response via : Initial Calibration



(51) 4-Nitrophenol-d4 (S)

14.533min (+0.035) 3.14ng/ul

response 12285

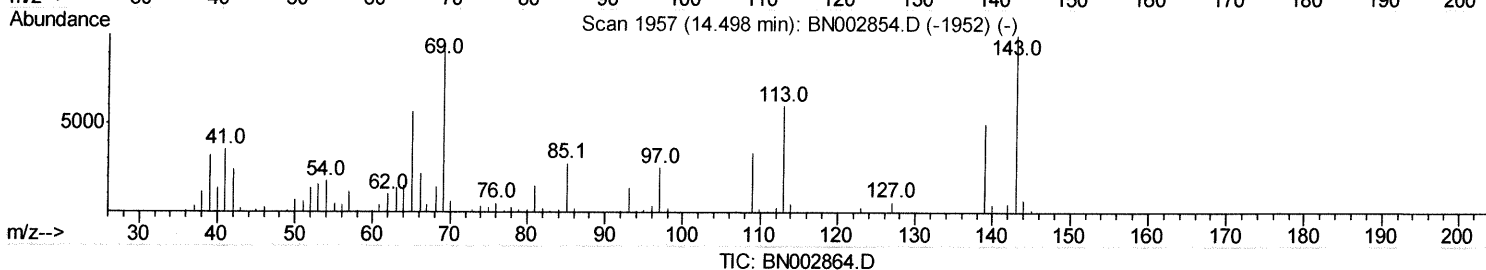
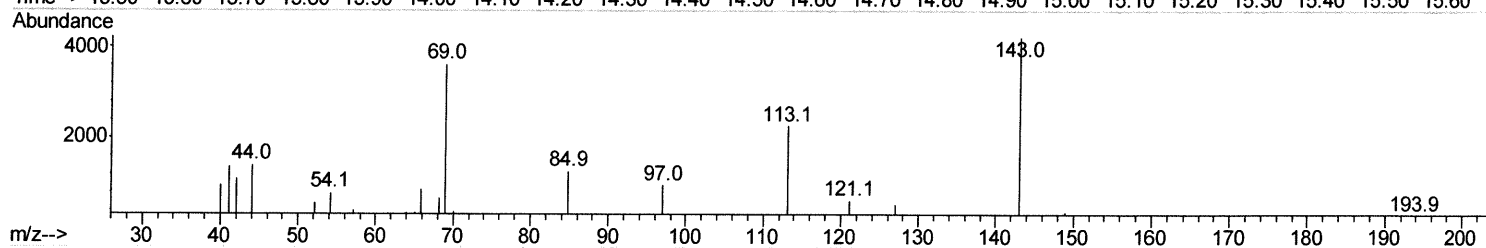
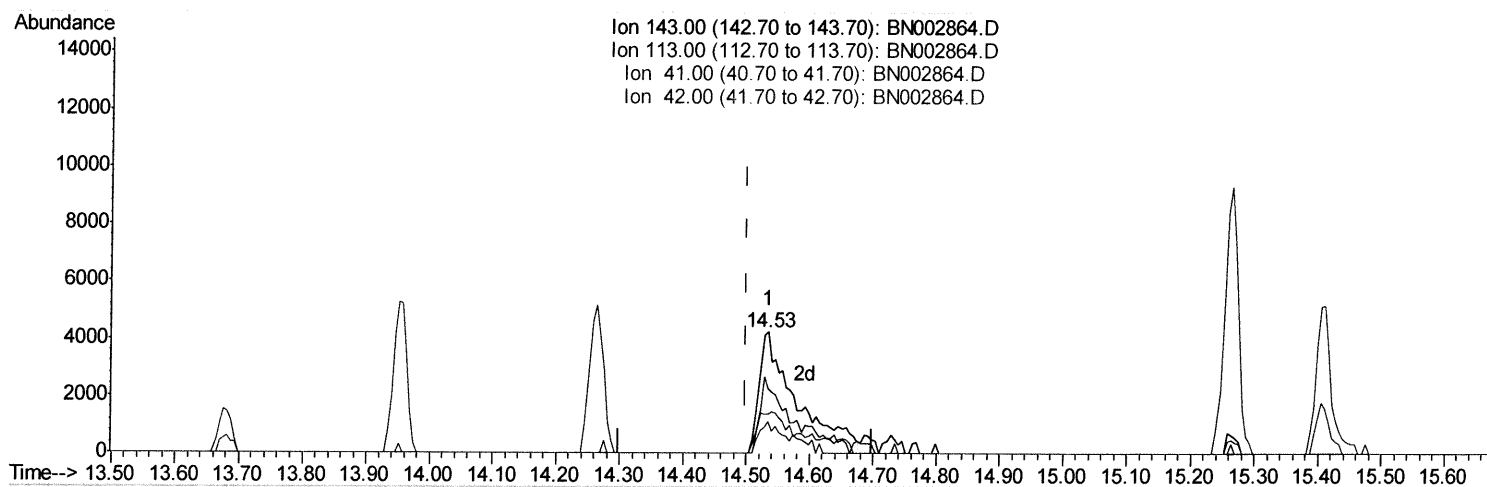
Ion	Exp%	Act%
143.00	100	100
113.00	66.10	53.58
41.00	39.90	32.27
42.00	25.80	25.78

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Operator   : MJ/SJ  
Sample     : J5117-07  
Misc      :  
ALS Vial   : 12      Sample Multiplier: 1
```

Manual Integrations
APPROVED

Sohil
10/1/2018 4:21:29 PM

Quant Time: Sep 29 06:36:28 2018
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Quant Title : SVOA CALIBRATION
QLast Update : Sat Sep 29 06:35:10 2018
Response via : Initial Calibration



(51) 4-Nitrophenol-d4 (S)

14.533min (+0.035) 4.44ng/ul m

response 17369

Ion	Exp%	Act%
143.00	100	100
113.00	66.10	53.58
41.00	39.90	32.27
42.00	25.80	25.78

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN092818\
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 BNA_N
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Manual Integrations
 APPROVED

Sohil
 10/1/2018 4:21:29 PM

Quant Time: Sep 29 06:58:07 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN091418MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Sep 29 06:35:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	132435	20.00	ng/ul	0.00
18) Naphthalene-d8	10.39	136	558184	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.26	164	302196	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.02	188	607177	20.00	ng/ul	0.00
77) Chrysene-d12	21.22	240	525579	20.00	ng/ul	0.00
85) Perylene-d12	23.42	264	540452	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.20	96	14364	5.00	ng/uL	0.00
5) Phenol-d5	6.82	99	67373	5.99	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.97	67	192900	27.62	ng/ul	0.00
9) 2-Chlorophenol-d4	7.17	132	216550	23.39	ng/ul	0.00
13) 4-Methylphenol-d8	8.34	113	124041	14.24	ng/ul	0.00
19) Nitrobenzene-d5	8.78	128	124822	28.87	ng/ul	0.00
22) 2-Nitrophenol-d4	9.49	143	143577	29.34	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.03	165	224088	25.37	ng/ul	0.00
29) 4-Chloroaniline-d4	0.00	131	0d	0.00	ng/ul	0.00
43) Dimethylphthalate-d6	13.68	166	755743	31.95	ng/ul	0.00
46) Acenaphthylene-d8	13.96	160	924718	29.94	ng/ul	0.00
51) 4-Nitrophenol-d4	14.53	143	17369m	4.44	ng/ul	0.04
57) Fluorene-d10	15.26	176	669090	32.29	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.41	200	114264	30.09	ng/ul	0.00
70) Anthracene-d10	17.12	188	1014712	35.00	ng/ul	0.00
78) Pyrene-d10	19.42	212	1003728	36.19	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.29	264	1006172	35.37	ng/ul	0.00

5/10/1/18

Target Compounds Ovalue

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