

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN101918\
Quantitation Report (QT Reviewed)

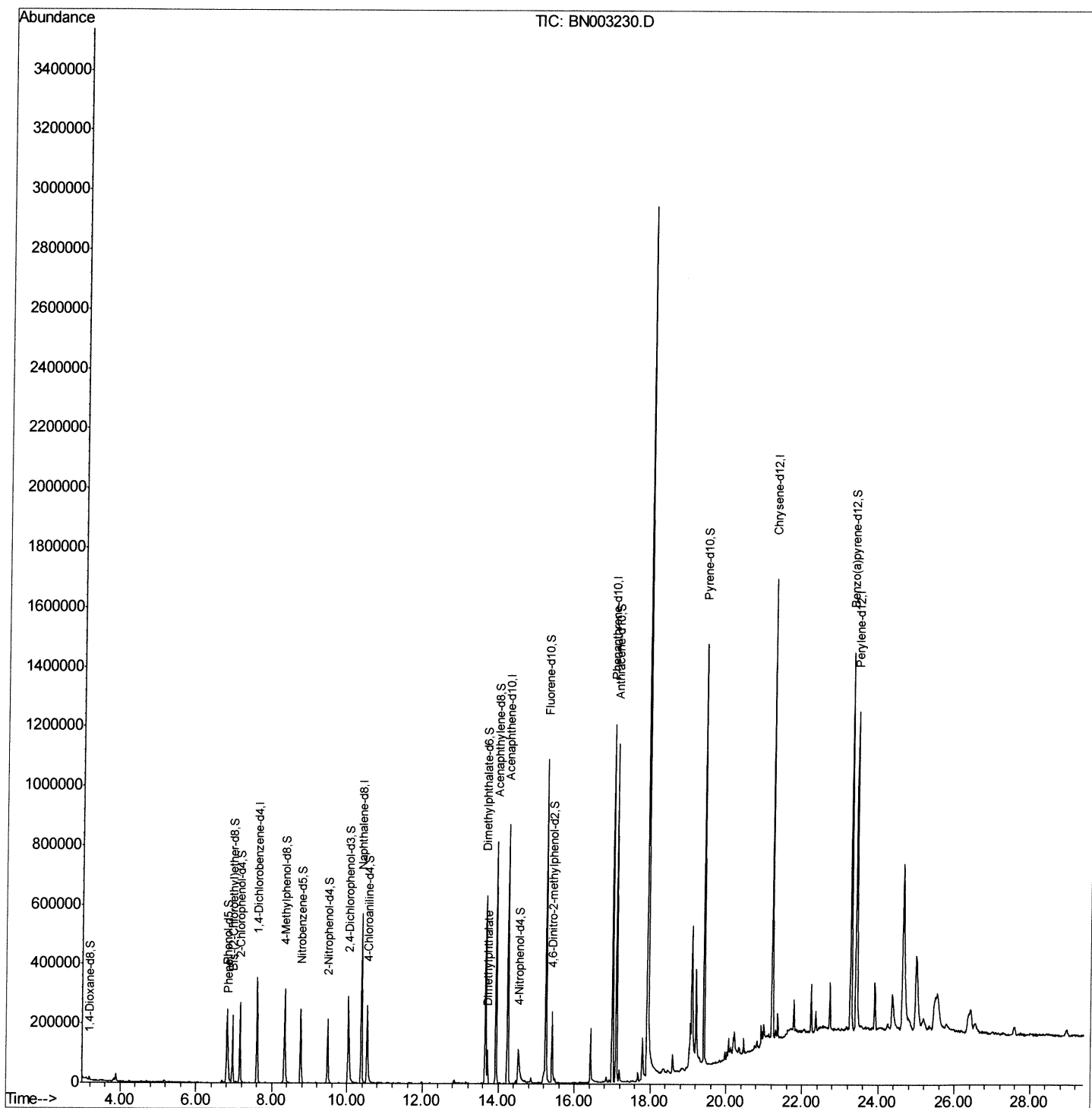
Data File : BN003230.D
Acq On : 20 Oct 2018 05:01
Operator : MJ/SJ
Sample : J5567-20
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
C0AJ6

Manual Integrations
APPROVED

Sohil
10/22/2018 4:53:06 PM

Quant Time: Oct 20 06:49:00 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN101918MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Oct 20 06:06:40 2018
Response via : Initial Calibration



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN101918\
Quantitation Report (Qedit)

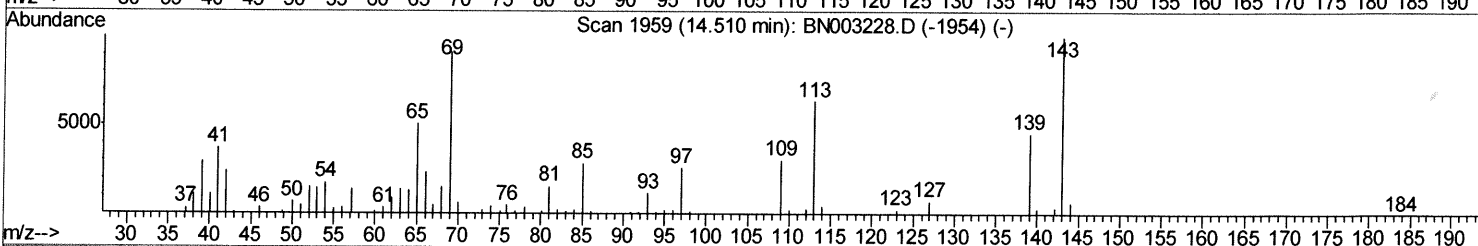
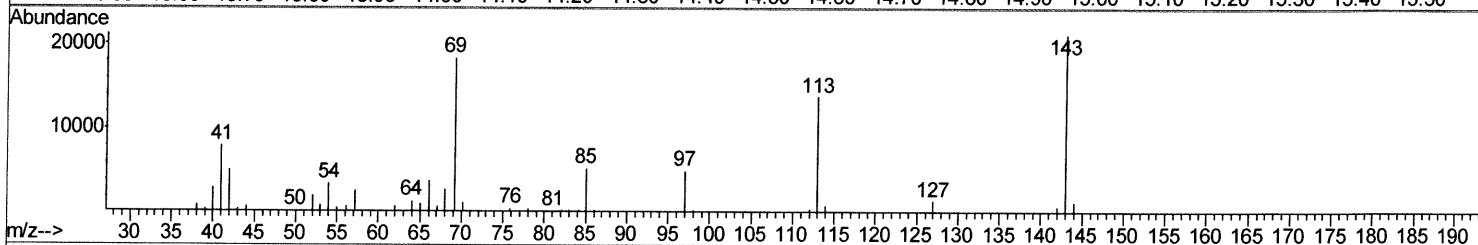
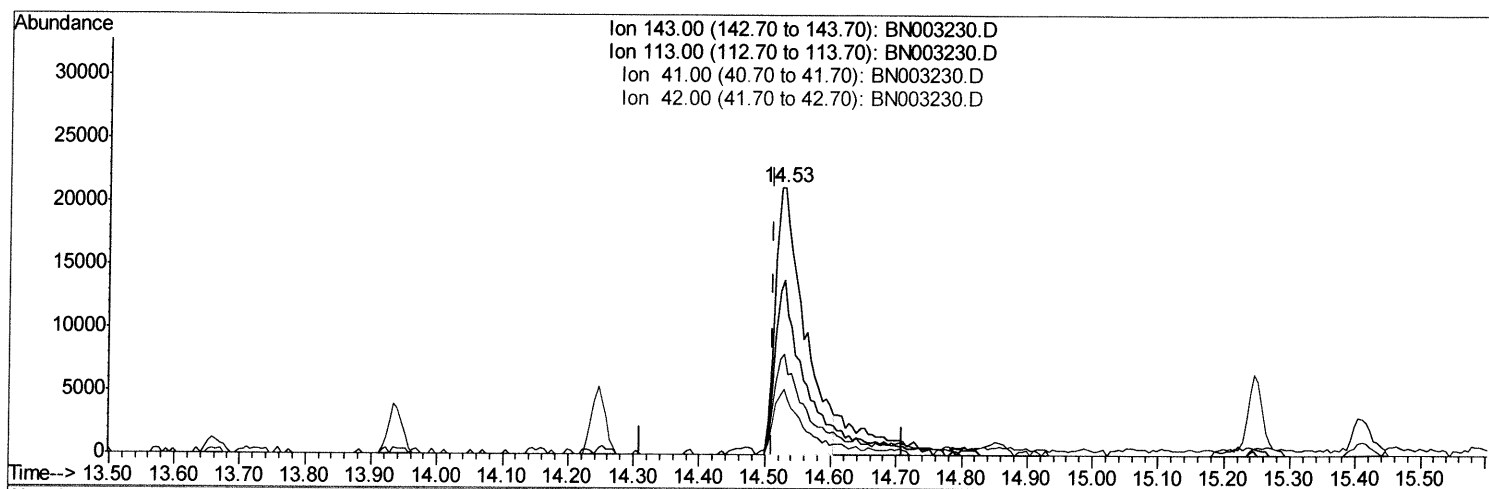
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10/22/2018 4:53:06 PM

Quant Time: Oct 20 06:44:33 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN101918MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Oct 20 06:06:40 2018
Response via : Initial Calibration



TIC: BN003230.D

(51) 4-Nitrophenol-d4 (S)

14.528min (+0.017) 17.93ng/ul

response 64305

Ion	Exp%	Act%
143.00	100	100
113.00	60.10	65.39
41.00	34.20	37.78
42.00	23.00	24.44

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Quantitation Report (Qedit)

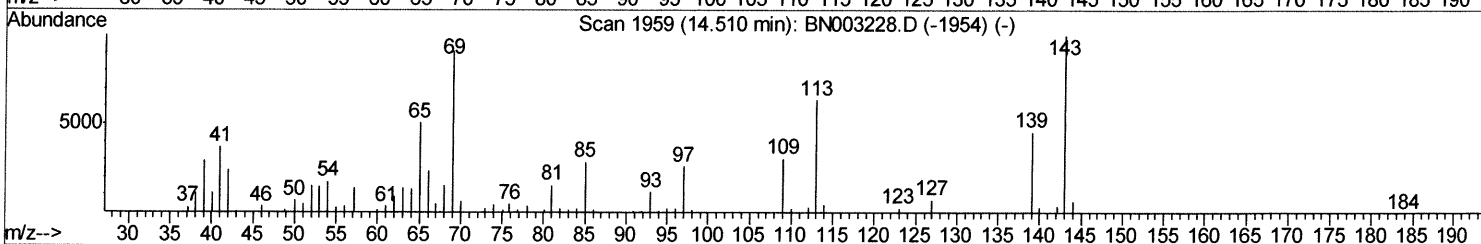
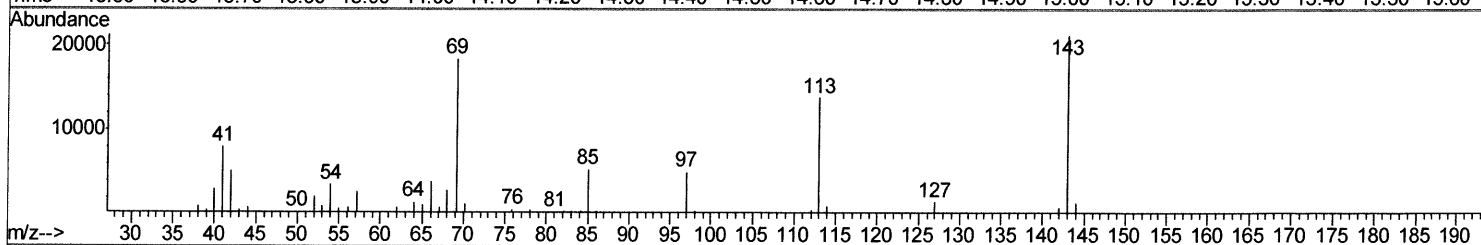
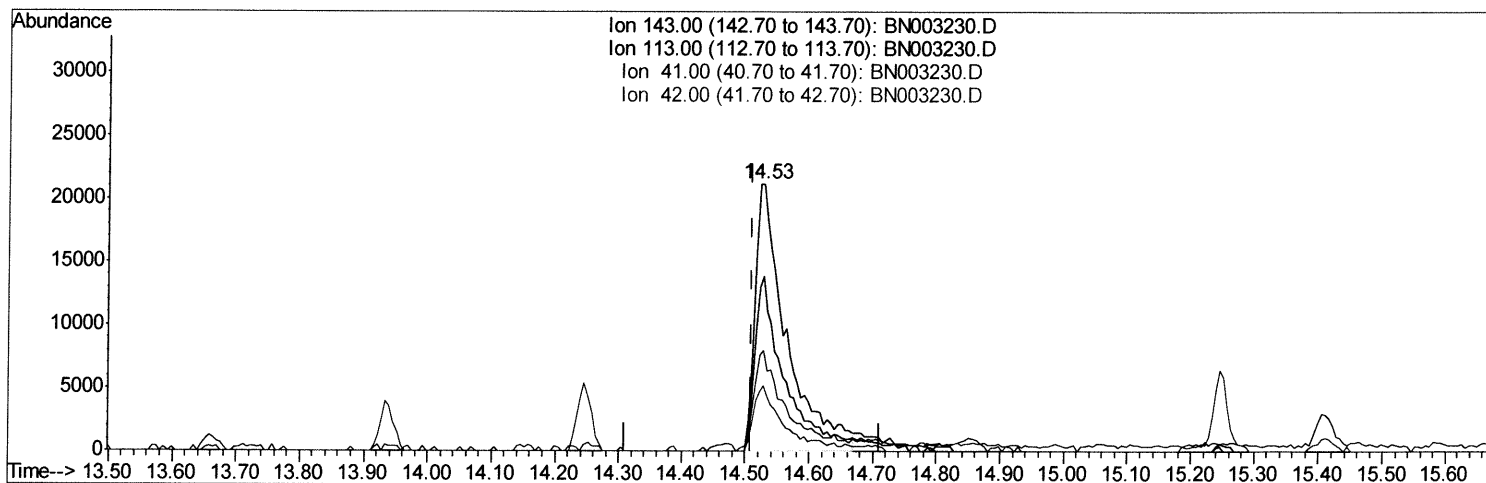
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Manual Integrations
APPROVED

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TIC: BN003230.D

(51) 4-Nitrophenol-d4 (S)

14.528min (+0.017) 20.27ng/ul m

5/10/24/18

response 72685

Ion	Exp%	Act%
143.00	100	100
113.00	60.10	65.39
41.00	34.20	37.78
42.00	23.00	24.44

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Manual Integrations
APPROVED

Sohil
10/22/2018 4:53:06 PM

Quant Time: Oct 20 06:49:00 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN101918MA.M
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Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	100516	20.00	ng/ul	0.00
18) Naphthalene-d8	10.37	136	493256	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.25	164	305450	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.00	188	705193	20.00	ng/ul	0.00
77) Chrysene-d12	21.22	240	782177	20.00	ng/ul	0.00
85) Perylene-d12	23.42	264	750683	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.18	96	6667	3.42	ng/uL	0.00
5) Phenol-d5	6.81	99	172792	18.15	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.95	67	105648	18.96	ng/ul	0.00
9) 2-Chlorophenol-d4	7.15	132	144398	19.44	ng/ul	0.00
13) 4-Methylphenol-d8	8.33	113	148585	19.17	ng/ul	0.00
19) Nitrobenzene-d5	8.76	128	71856	18.77	ng/ul	0.00
22) 2-Nitrophenol-d4	9.48	143	85325	19.22	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.03	165	152323	18.92	ng/ul	0.00
29) 4-Chloroaniline-d4	10.53	131	184386	28.40	ng/ul	0.00
43) Dimethylphthalate-d6	13.66	166	473321	18.18	ng/ul	0.00
46) Acenaphthylene-d8	13.94	160	605729	18.85	ng/ul	0.00
51) 4-Nitrophenol-d4	14.53	143	72685m	20.27	ng/ul	0.02
57) Fluorene-d10	15.25	176	433859	19.03	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.41	200	70237	14.83	ng/ul	0.00
70) Anthracene-d10	17.10	188	666963	19.10	ng/ul	0.00
78) Pyrene-d10	19.41	212	738782	19.49	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.28	264	781528	19.26	ng/ul	0.00

Target Compounds				Qvalue	
6) Phenol	6.83	94	56790	5.929	ng/ul 95
44) Dimethylphthalate	13.71	163	82806	3.255	ng/ul 99

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

5519/24/18