

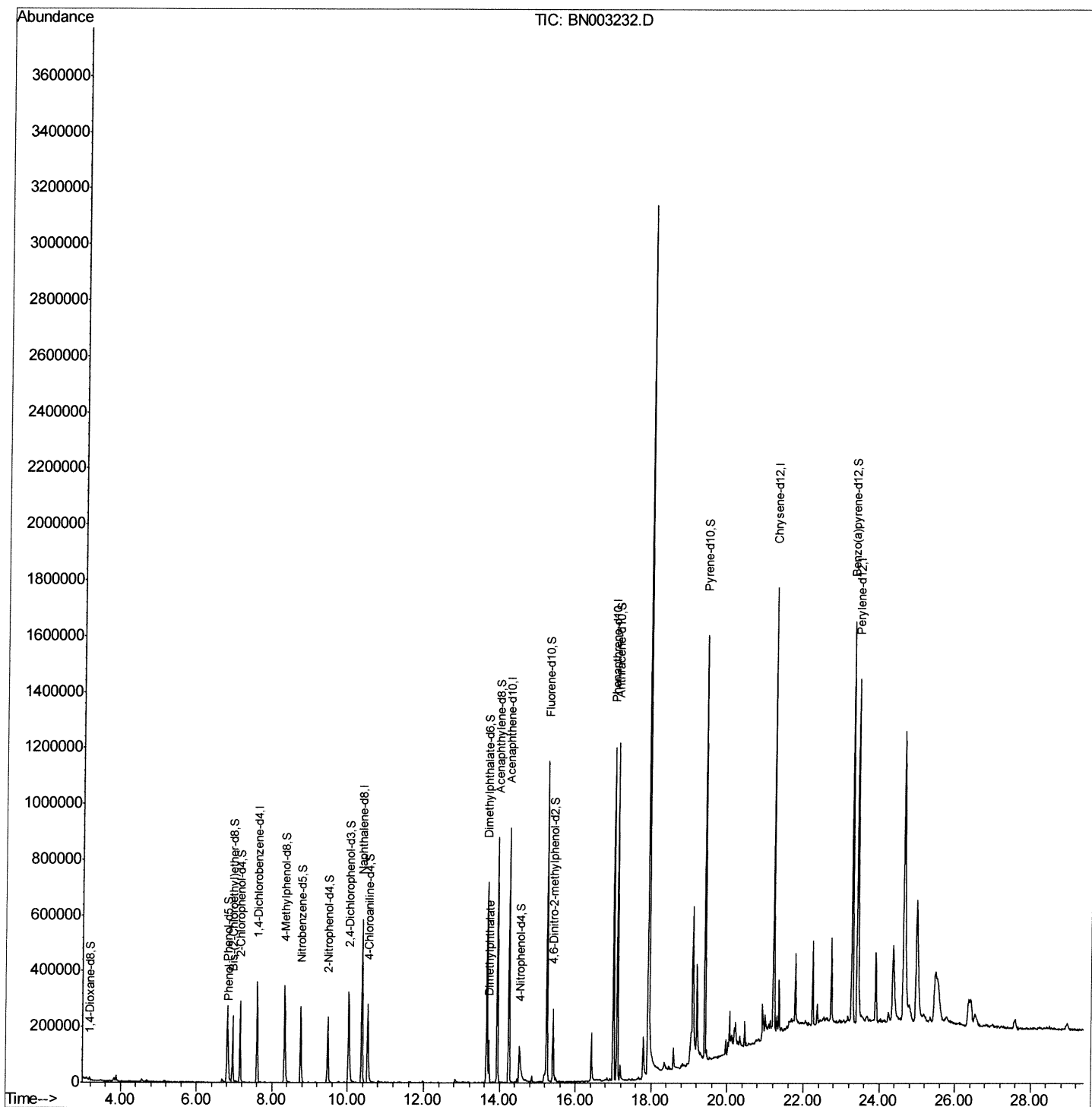
Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN101918\
Data File : BN003232.D
Acq On : 20 Oct 2018 06:12
Operator : MJ/SJ
Sample : J5567-08
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
C0AH6

Manual Integrations
APPROVED

Sohil
10/22/2018 4:53:11 PM

Quant Time: Oct 22 04:17:49 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN101918MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Oct 22 04:05:35 2018
Response via : Initial Calibration



Quantitation Report (Qedit)

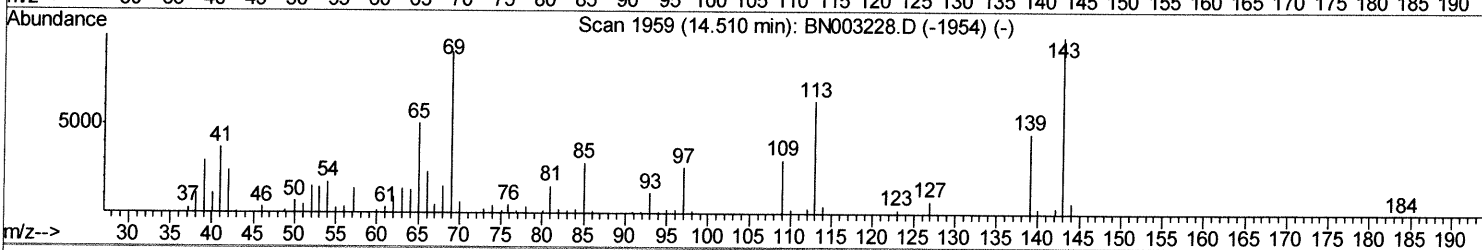
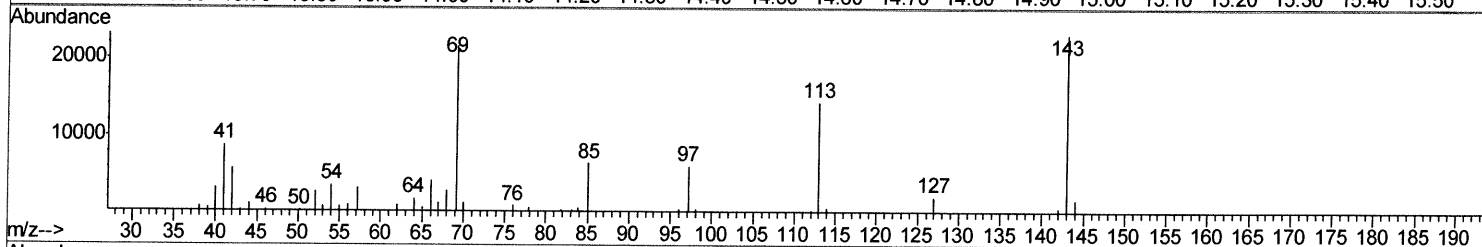
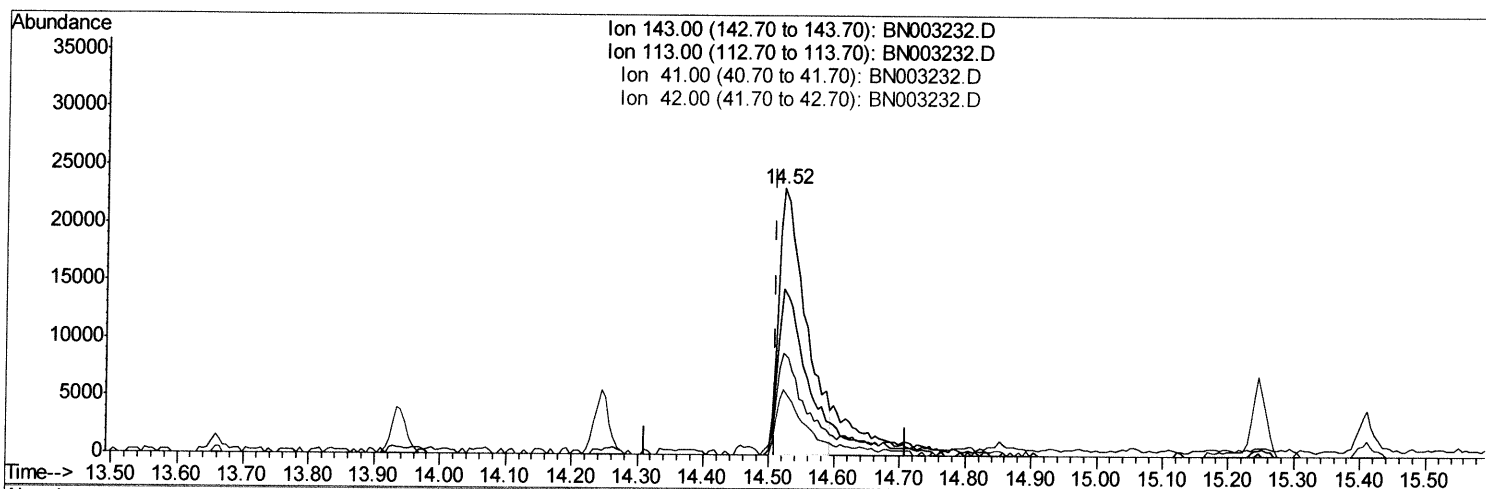
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TIC: BN003232.D

(51) 4-Nitrophenol-d4 (S)

14.522min (+0.012) 18.36ng/ul

response 66911

Ion	Exp%	Act%
143.00	100	100
113.00	60.10	62.25
41.00	34.20	38.15
42.00	23.00	24.71

Quantitation Report (Qedit)

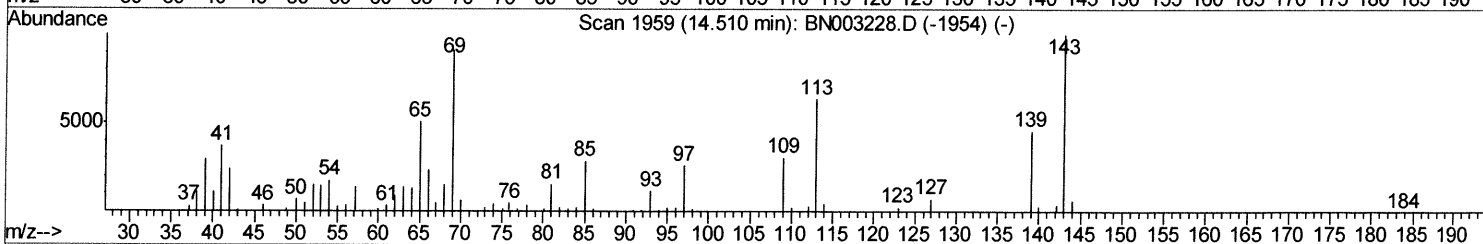
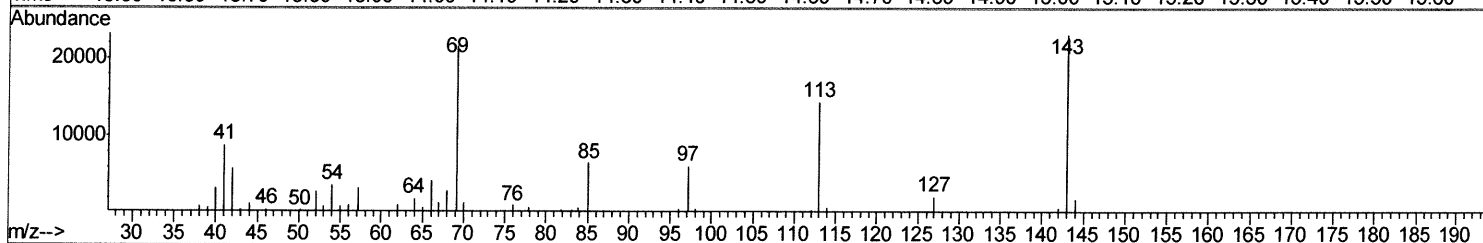
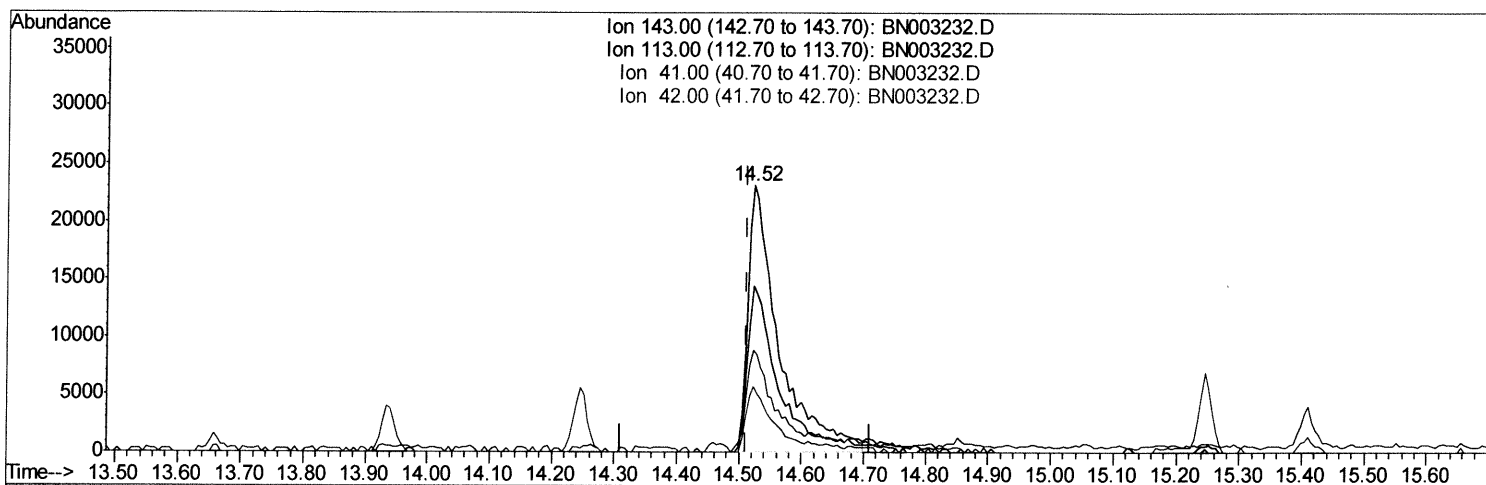
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Manual Integrations
APPROVED

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TIC: BN003232.D

(51) 4-Nitrophenol-d4 (S)

14.522min (+0.012) 22.05ng/ul m

SS 10/24/18

response 80363

Ion	Exp%	Act%
143.00	100	100
113.00	60.10	62.25
41.00	34.20	38.15
42.00	23.00	24.71

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN101918\
 Data File : BN003232.D
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 Operator : MJ/SJ
 Sample : J5567-08
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AH6

Manual Integrations
 APPROVED

Sohil
 10/22/2018 4:53:11 PM

Quant Time: Oct 22 04:17:49 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN101918MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Oct 22 04:05:35 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	102145	20.00	ng/ul	0.00
18) Naphthalene-d8	10.37	136	497155	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.25	164	310419	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.00	188	716814	20.00	ng/ul	0.00
77) Chrysene-d12	21.22	240	782013	20.00	ng/ul	0.00
85) Perylene-d12	23.42	264	740280	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.18	96	6062	3.06	ng/uL	0.00
5) Phenol-d5	6.81	99	182646	18.88	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.95	67	109314	19.31	ng/ul	0.00
9) 2-Chlorophenol-d4	7.15	132	150403	19.93	ng/ul	0.00
13) 4-Methylphenol-d8	8.33	113	155615	19.76	ng/ul	0.00
19) Nitrobenzene-d5	8.76	128	77296	20.04	ng/ul	0.00
22) 2-Nitrophenol-d4	9.48	143	90604	20.24	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.02	165	159286	19.63	ng/ul	0.00
29) 4-Chloroaniline-d4	10.53	131	189002	28.88	ng/ul	0.00
43) Dimethylphthalate-d6	13.66	166	504818	19.08	ng/ul	0.00
46) Acenaphthylene-d8	13.94	160	653707	20.01	ng/ul	0.00
51) 4-Nitrophenol-d4	14.52	143	80363m	22.05	ng/ul	0.01
57) Fluorene-d10	15.25	176	466366	20.13	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.41	200	73965	15.36	ng/ul	0.00
70) Anthracene-d10	17.10	188	710892	20.03	ng/ul	0.00
78) Pyrene-d10	19.41	212	789221	20.82	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.28	264	815508	20.38	ng/ul	0.00

SJ 10/24/08

Target Compounds

					Qvalue
6) Phenol	6.84	94	58445	6.004 ng/ul	98
44) Dimethylphthalate	13.71	163	104046	4.024 ng/ul	97

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