

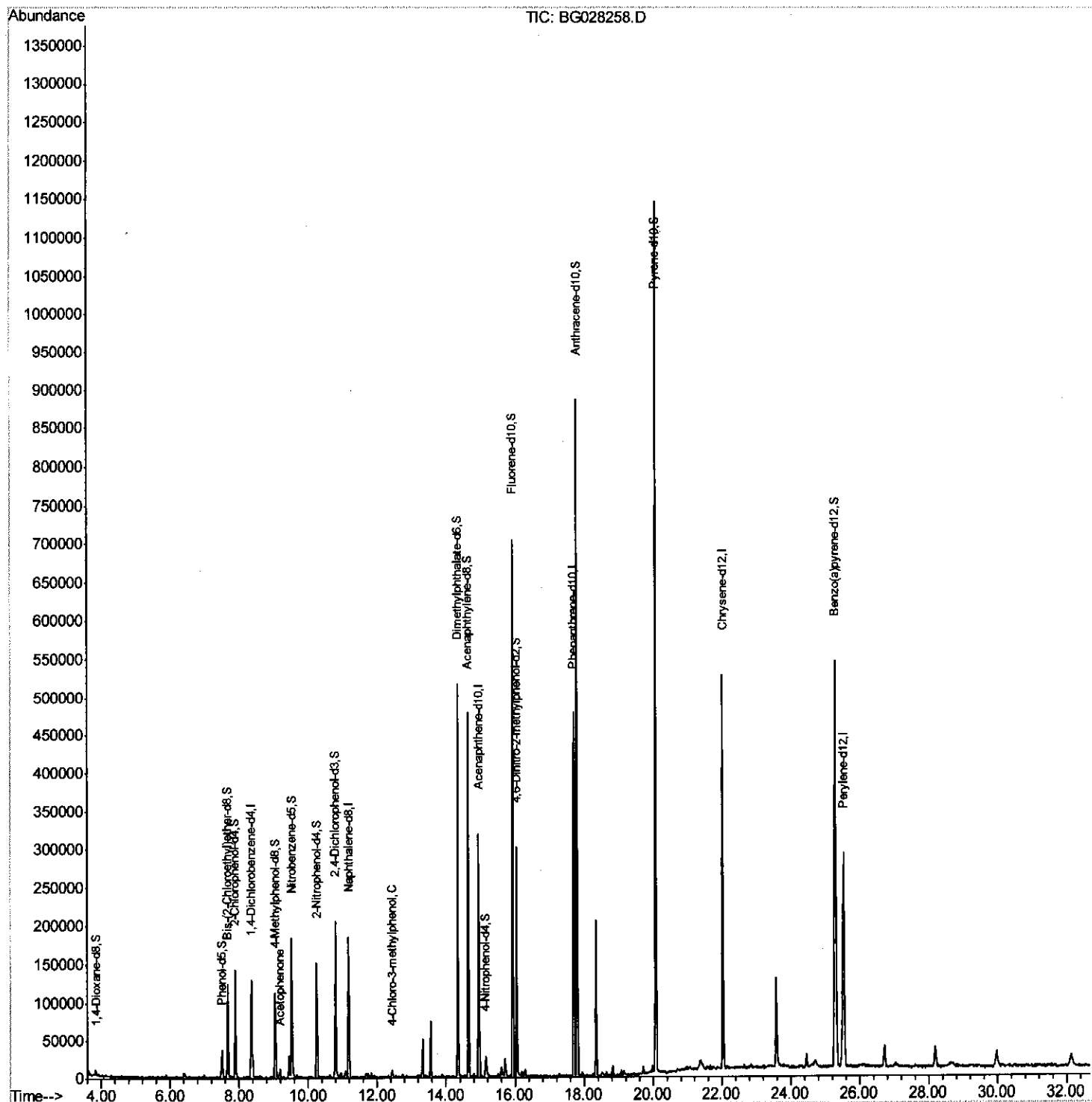
Data Path : Z:\HPCHEM1\BNA G\DATA\BG081017\
Data File : BG028258.D
Acq On : 10 Aug 2017 14:47
Operator : SJ/JU
Sample : I4630-03
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampled :
C0AA9

Manual Integrations
APPROVED

Sohil
8/14/2017 4:02:07 PM

Quant Time: Aug 11 01:26:37 2017
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM-EPA-BG080217.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Aug 11 01:09:14 2017
Response via : Initial Calibration



Quantitation Report (Qedit)

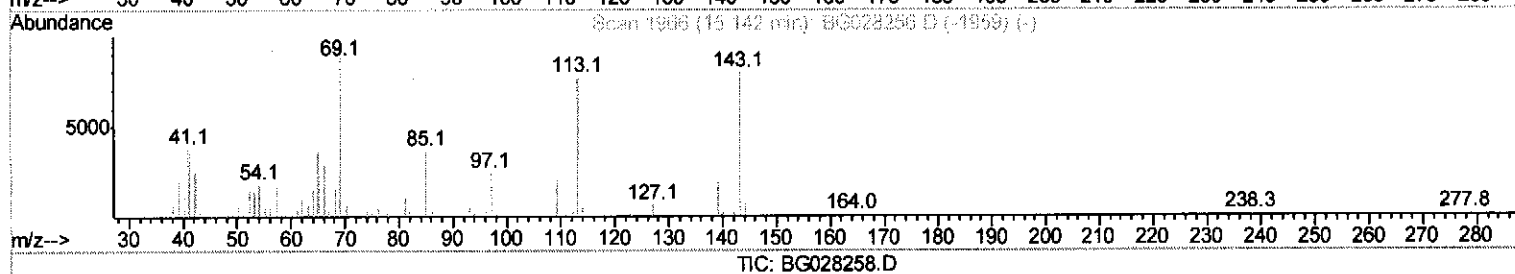
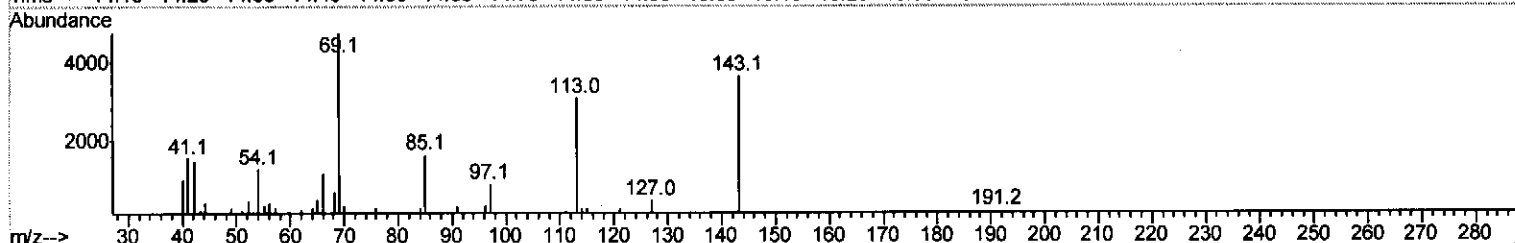
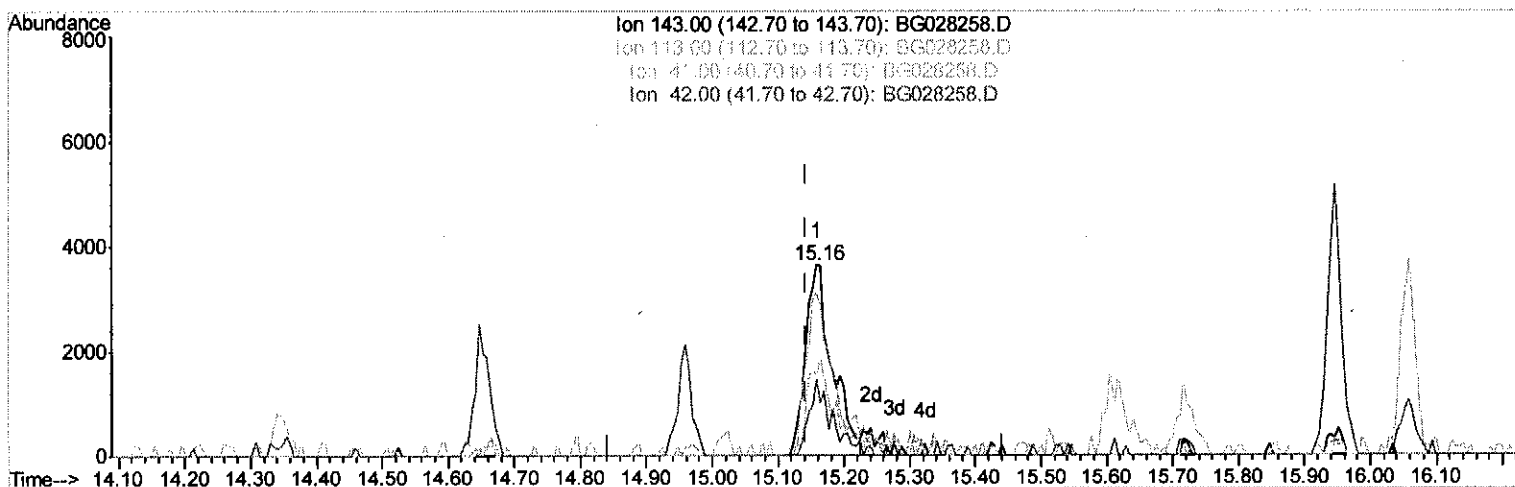
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Quant Time: Aug 11 01:09:59 2017
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Quant Title : SVOA CALIBRATION
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Response via : Initial Calibration



(51) 4-Nitrophenol-d4 (S)

15.159min (+0.017) 6.72ng/ul

response 10302

Ion	Exp%	Act%
143.00	100	100
113.00	117.30	84.42#
41.00	58.00	43.12#
42.00	42.20	40.34

Quantitation Report (Qedit)

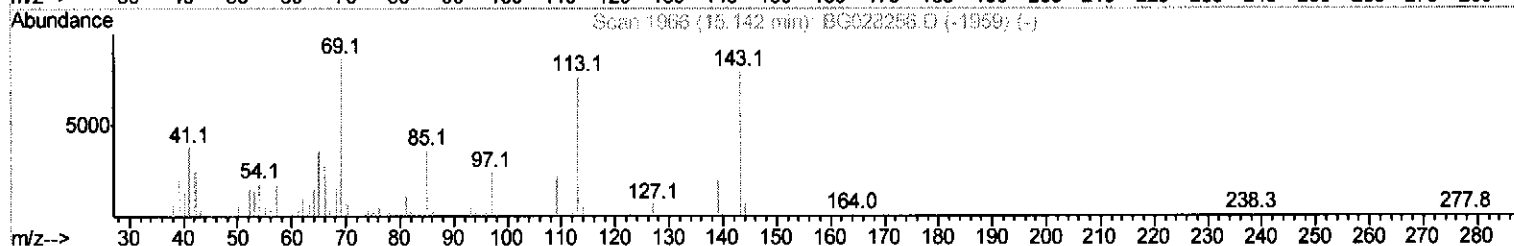
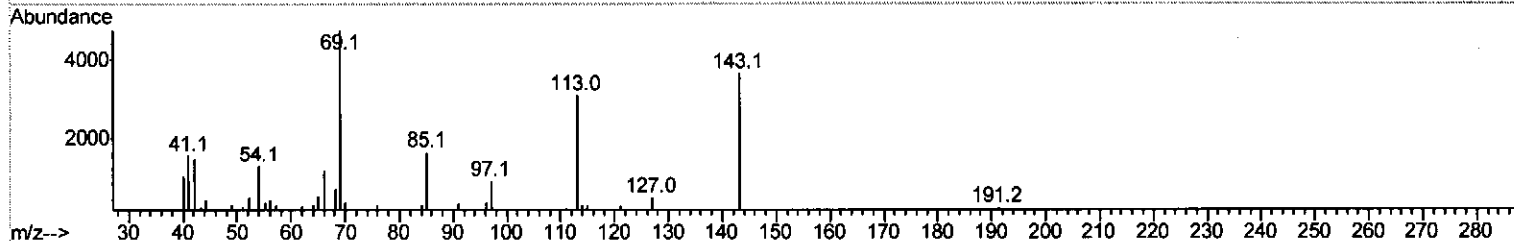
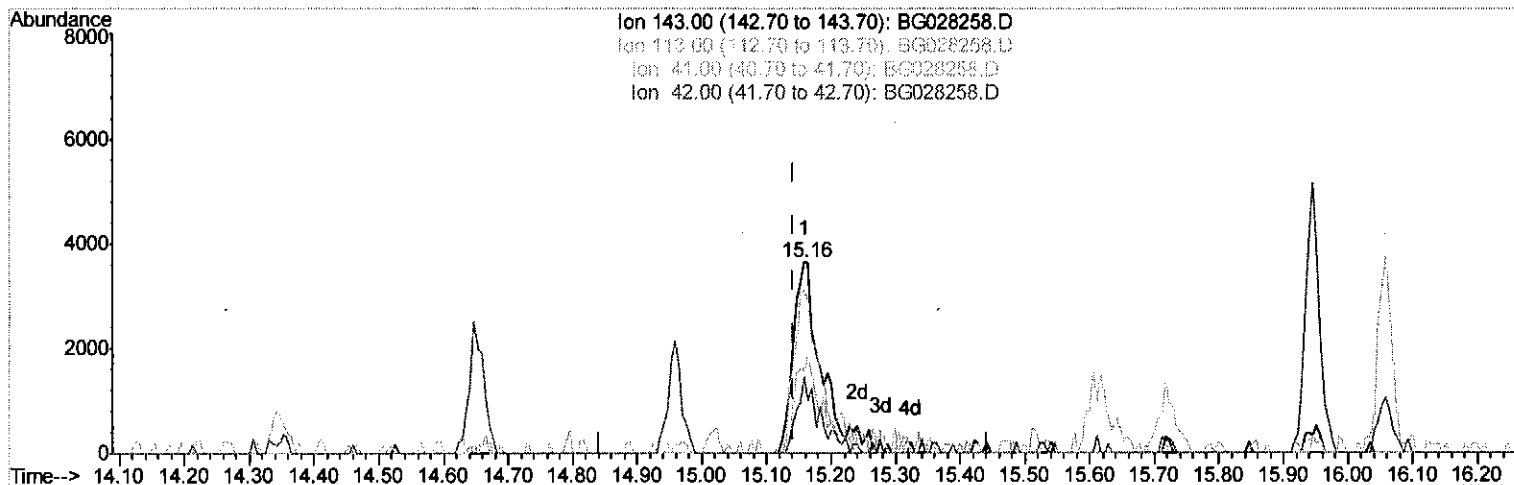
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BNA_G
ClientSampleId :
C0AA9

Manual Integrations
APPROVED

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8/14/2017 4:02:07 PM

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Quant Title : SVOA CALIBRATION
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Response via : Initial Calibration



TIC: BG028258.D

(51) 4-Nitrophenol-d4 (S)

15.159min (+0.017) 7.24ng/ul m

response 11109

Ion	Exp%	Act%
143.00	100	100
113.00	117.30	84.42#
41.00	58.00	43.12#
42.00	42.20	40.34

Data Path : Z:\HPCHEM1\BNA G\DATA\BG081017\
 Data File : BG028258.D
 Acq On : 10 Aug 2017 14:47
 Operator : SJ/JU
 Sample : I4630-03
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0AA9

Manual Integrations
 APPROVED

Sohil
 8/14/2017 4:02:07 PM

Quant Time: Aug 11 01:26:37 2017
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 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Aug 11 01:09:14 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.35	152	36145	20.00	ng/ul	0.00
18) Naphthalene-d8	11.17	136	150411	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.96	164	108447	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.71	188	302878	20.00	ng/ul	0.00
75) Chrysene-d12	22.04	240	353450	20.00	ng/ul	0.00
83) Perylene-d12	25.53	264	344243	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.84	96	3599	4.64	ng/uL	0.00
5) Phenol-d5	7.50	99	20868	5.26	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.67	67	67309	26.77	ng/ul	0.00
9) 2-Chlorophenol-d4	7.88	132	59347	23.70	ng/ul	0.00
13) 4-Methylphenol-d8	9.05	113	41734	12.58	ng/ul	0.00
19) Nitrobenzene-d5	9.52	128	38600	32.82	ng/ul	0.00
22) 2-Nitrophenol-d4	10.24	143	46396	35.98	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.79	165	77456	29.03	ng/ul	0.00
29) 4-Chloroaniline-d4	11.29	131	79	0.03	ng/ul	0.00
43) Dimethylphthalate-d6	14.35	166	337243	34.96	ng/ul	0.00
46) Acenaphthylene-d8	14.65	160	348934	32.09	ng/ul	0.00
51) 4-Nitrophenol-d4	15.16	143	11109m)	7.24	ng/ul	0.02
57) Fluorene-d10	15.95	176	288817	33.38	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	16.06	200	64291	32.68	ng/ul	0.00
70) Anthracene-d10	17.81	188	543328	37.41	ng/ul	0.00
76) Pyrene-d10	20.08	212	649307	35.82	ng/ul	0.00
87) Benzo(a)pyrene-d12	25.30	264	574532	35.65	ng/ul	0.00
Target Compounds						
14) Acetophenone	9.18	105	6646	1.405	ng/ul#	82
33) 4-Chloro-3-methylphenol	12.44	107	5133	1.656	ng/ul#	86

0.02 > 8/14/17

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