

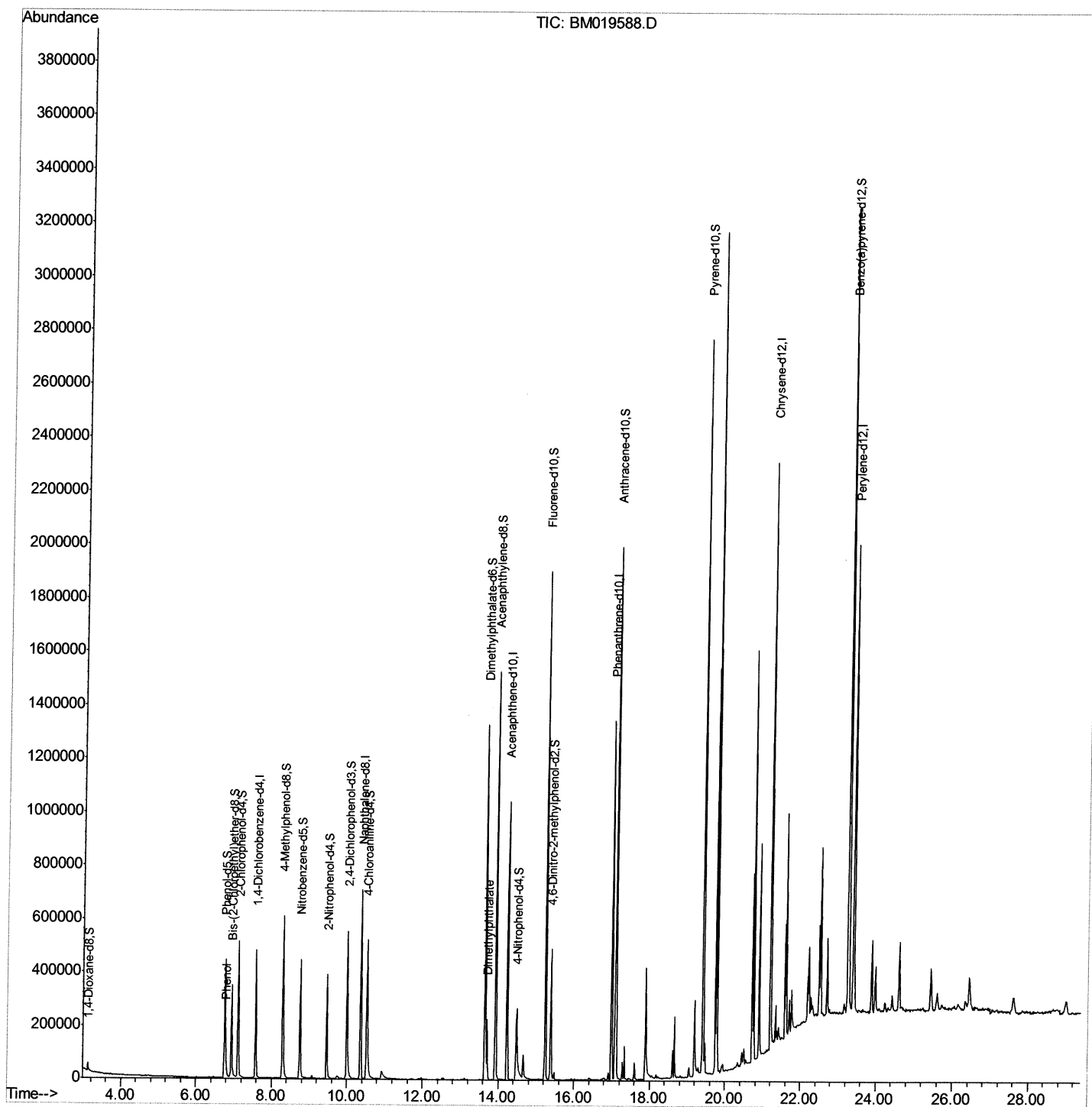
Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032919\
Data File : BM019588.D
Acq On : 29 Mar 2019 20:13
Operator : JU/SJ
Sample : K2085-13
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BF7T8

Manual Integrations
APPROVED

Sohil
4/1/2019 5:49:06 PM

Quant Time: Mar 29 23:41:15 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Mar 29 14:33:17 2019
Response via : Initial Calibration



Quantitation Report (Qedit)

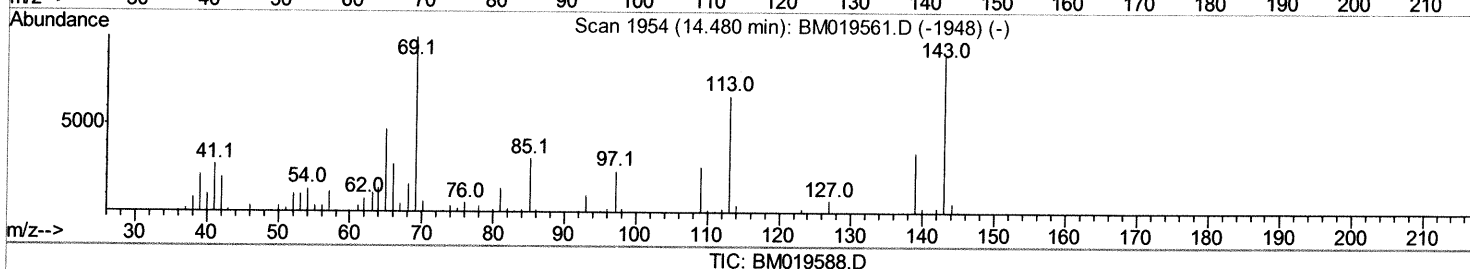
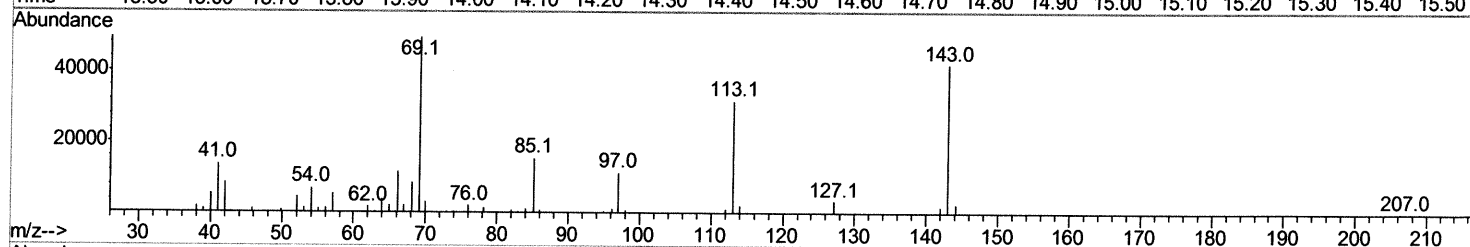
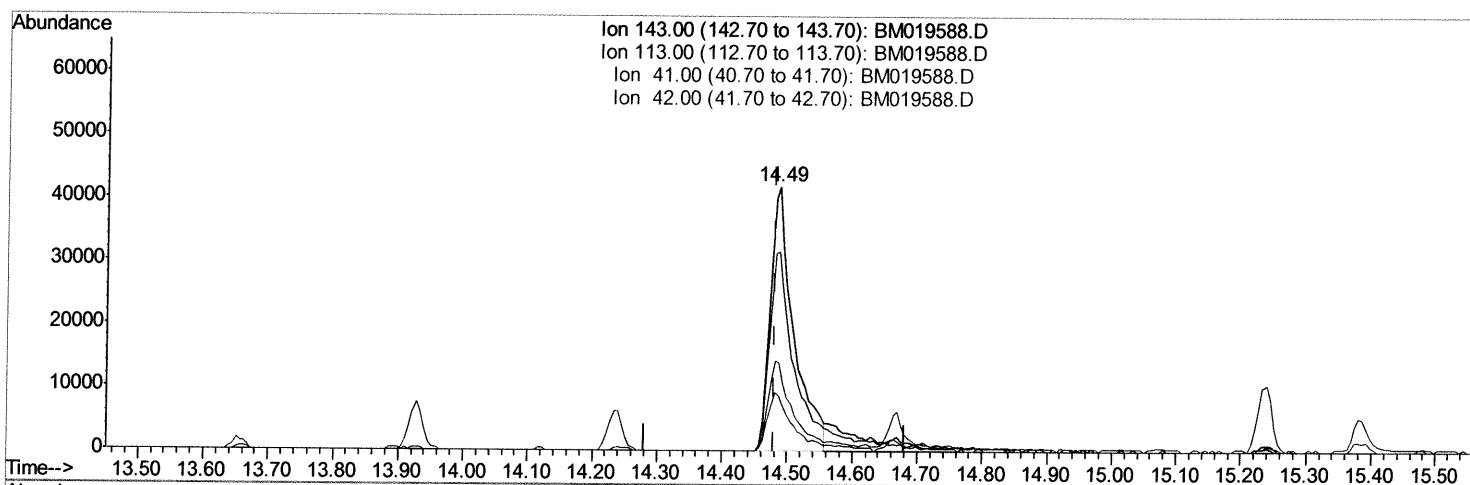
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(52) 4-Nitrophenol-d4 (S)

14.486min (+0.006) 24.44ng/ul

response 104012

Ion	Exp%	Act%
143.00	100	100
113.00	77.10	75.49
41.00	32.60	32.99
42.00	22.30	21.13

Quantitation Report (Qedit)

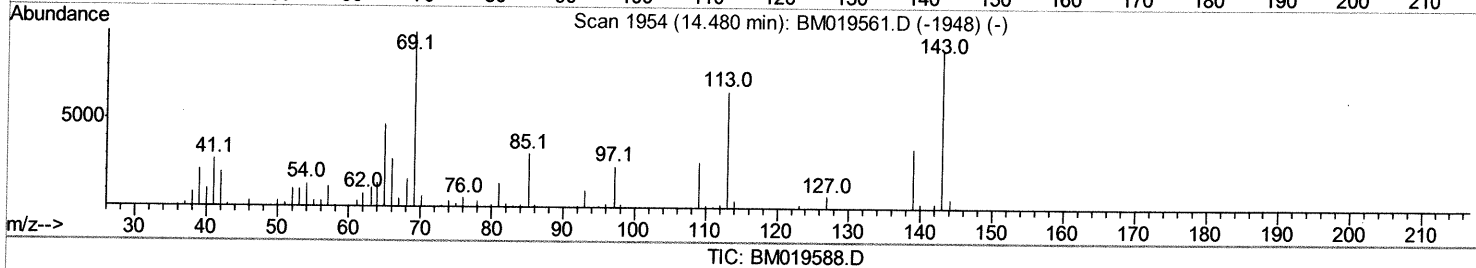
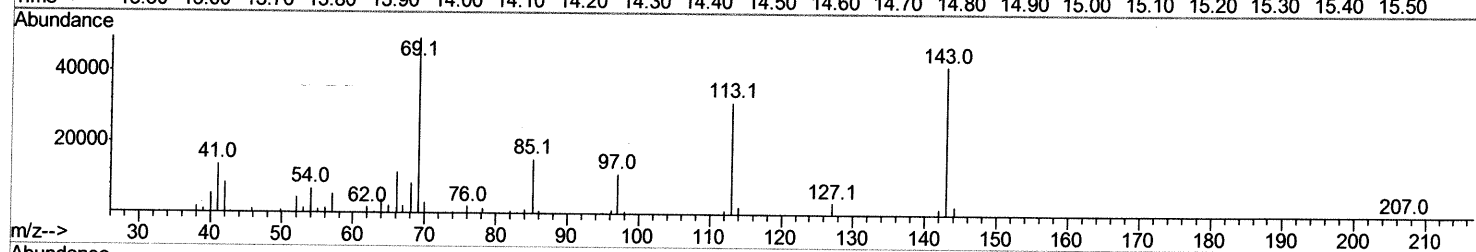
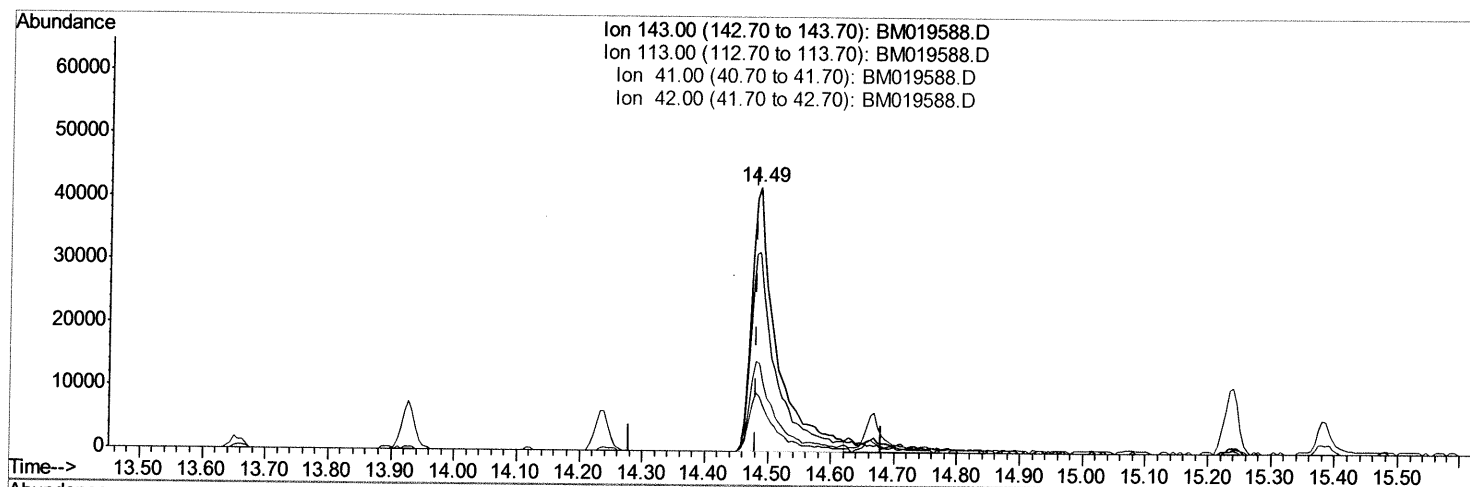
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Manual Integrations
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Response via : Initial Calibration



(52) 4-Nitrophenol-d4 (S)

14.486min (+0.006) 27.18ng/ul m> SD 4/2/19

response 115708

Ion	Exp%	Act%
143.00	100	100
113.00	77.10	75.49
41.00	32.60	32.99
42.00	22.30	21.13

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032919\
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 Acq On : 29 Mar 2019 20:13
 Operator : JU/SJ
 Sample : K2085-13
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BF7T8

Manual Integrations
 APPROVED

Sohil
 4/1/2019 5:49:06 PM

Quant Time: Mar 29 23:41:15 2019
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 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Mar 29 14:33:17 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	129892	20.00	ng/ul	0.00
18) Naphthalene-d8	10.36	136	566785	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.23	164	336523	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.00	188	718977	20.00	ng/ul	0.00
78) Chrysene-d12	21.20	240	917639	20.00	ng/ul	0.00
86) Perylene-d12	23.41	264	1142089	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.13	96	21376	6.44	ng/uL	0.00
5) Phenol-d5	6.76	99	307705	26.60	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.93	67	197161	25.46	ng/ul	0.00
9) 2-Chlorophenol-d4	7.11	132	239340	27.30	ng/ul	0.00
13) 4-Methylphenol-d8	8.29	113	257112	29.61	ng/ul	0.00
19) Nitrobenzene-d5	8.75	128	110312	27.28	ng/ul	0.00
22) 2-Nitrophenol-d4	9.46	143	129707	28.82	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.99	165	229681	27.05	ng/ul	0.00
29) 4-Chloroaniline-d4	10.52	131	341671	33.51	ng/ul	0.00
44) Dimethylphthalate-d6	13.66	166	836433	30.80	ng/ul	0.00
47) Acenaphthylene-d8	13.93	160	1017156	29.96	ng/ul	0.00
52) 4-Nitrophenol-d4	14.49	143	115708m	27.18	ng/ul	0.00
58) Fluorene-d10	15.24	176	709294	30.32	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.39	200	109441	23.02	ng/ul	0.00
71) Anthracene-d10	17.10	188	1089803	31.59	ng/ul	0.00
79) Pyrene-d10	19.40	212	1235715	29.15	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.27	264	1768501	29.21	ng/ul	0.00

SD 4/2/19

Target Compounds

				Ovalue	
6) Phenol	6.79	94	26381	2.186 ng/ul#	91
45) Dimethylphthalate	13.70	163	145600	5.344 ng/ul#	98

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