

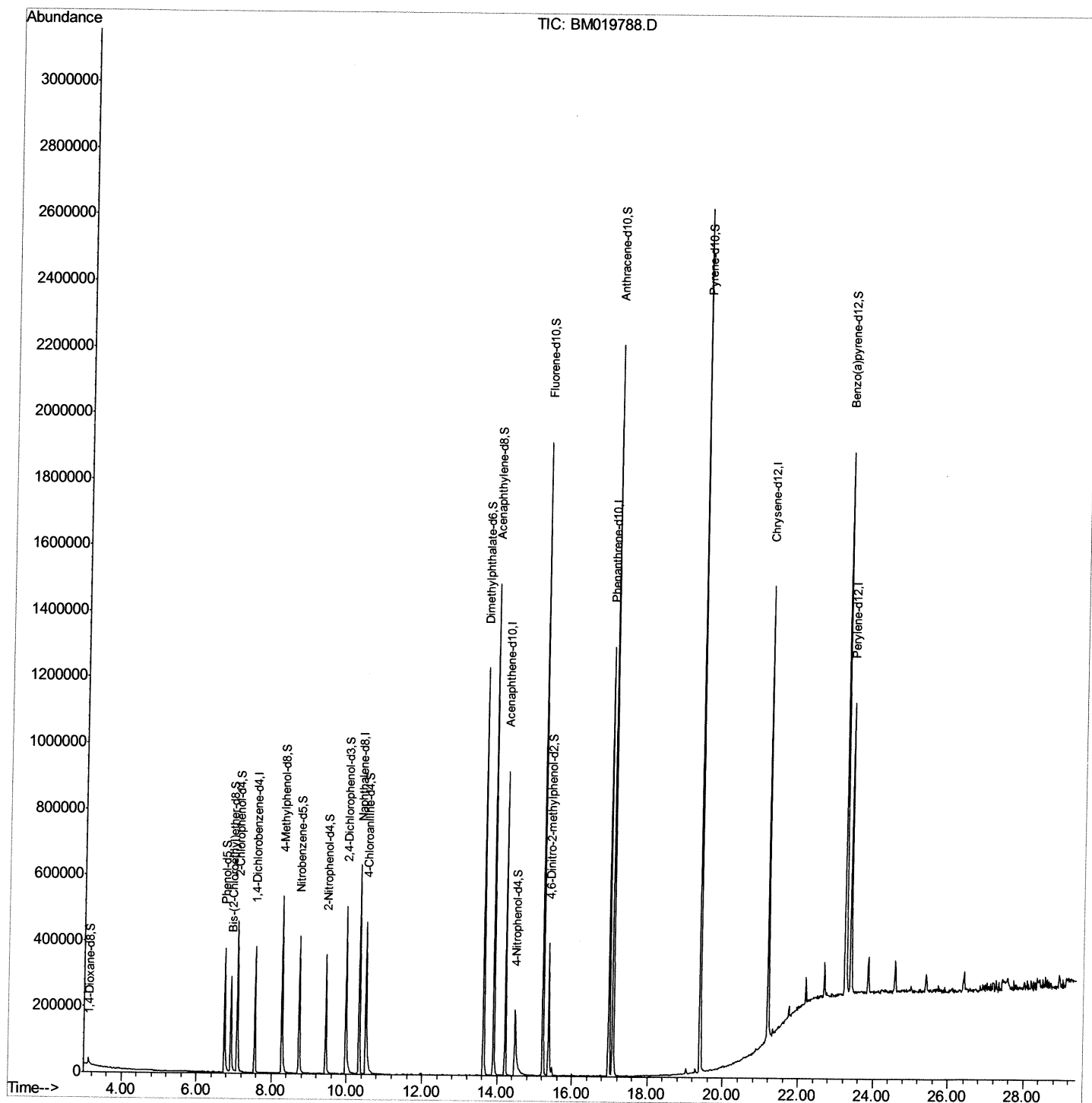
Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM040819\
Data File : BM019788.D
Acq On : 08 Apr 2019 11:03
Operator : JU/SJ
Sample : PB118630BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SBLK30

Manual Integrations
APPROVED

Sohil
4/10/2019 6:16:21 PM

Quant Time: Apr 09 06:38:40 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Apr 06 00:13:11 2019
Response via : Initial Calibration



Quantitation Report (Qedit)

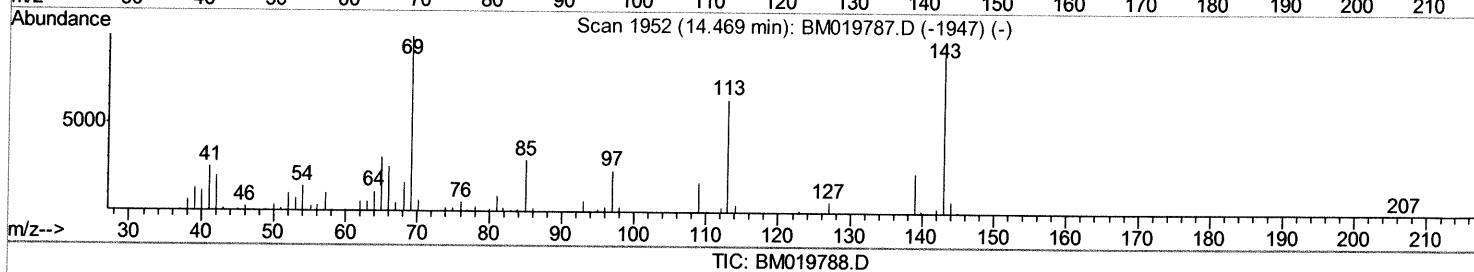
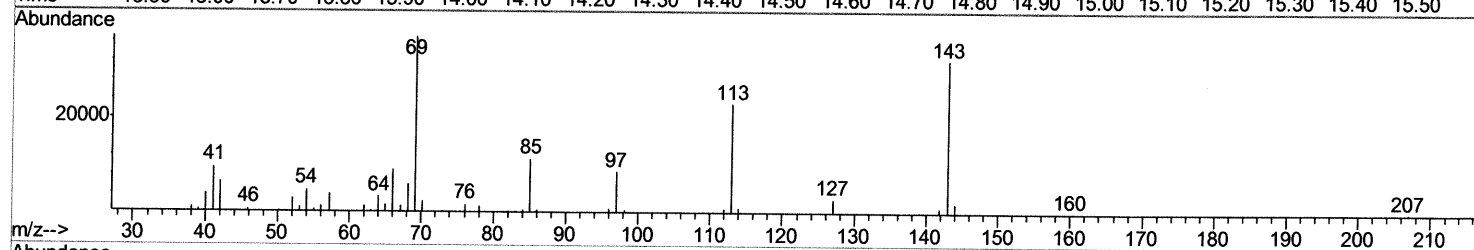
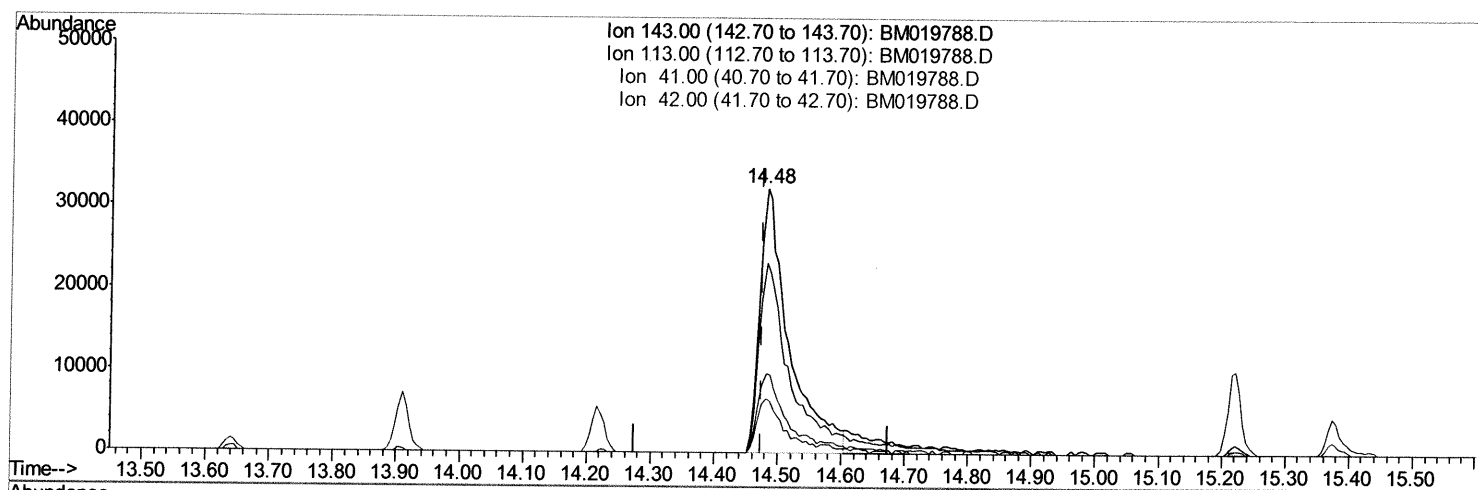
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(52) 4-Nitrophenol-d4 (S)

14.480min (+0.006) 26.66ng/ul

response 102210

Ion	Exp%	Act%
143.00	100	100
113.00	77.10	71.73
41.00	32.60	30.10
42.00	22.30	20.67

Quantitation Report (Qedit)

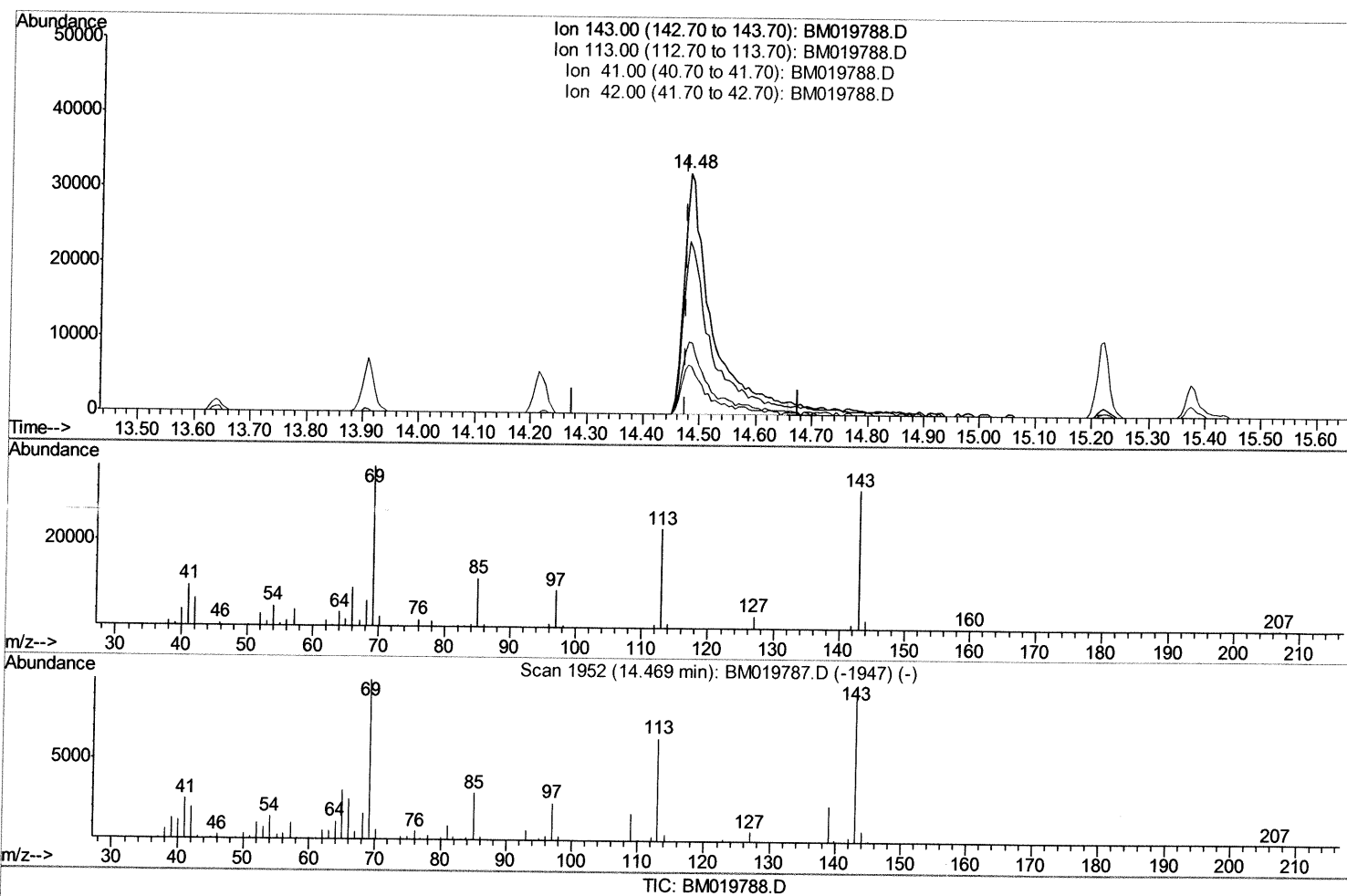
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Manual Integrations
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Response via : Initial Calibration



(52) 4-Nitrophenol-d4 (S)

14.480min (+0.006) 28.38ng/ul m> SJ 4/16/19

response 108816

Ion	Exp%	Act%
143.00	100	100
113.00	77.10	71.73
41.00	32.60	30.10
42.00	22.30	20.67

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM040819\
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 Sample : PB118630BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK30

Manual Integrations
 APPROVED

Sohil
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Quant Time: Apr 09 06:38:40 2019
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 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Apr 06 00:13:11 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	105010	20.00	ng/ul	0.00
18) Naphthalene-d8	10.33	136	500521	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.22	164	303119	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.98	188	694099	20.00	ng/ul	0.00
78) Chrysene-d12	21.19	240	651694	20.00	ng/ul	0.00
86) Perylene-d12	23.39	264	595822	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.12	96	13343	4.97	ng/uL	0.00
5) Phenol-d5	6.75	99	272054	29.09	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.91	67	171633	27.42	ng/ul	0.00
9) 2-Chlorophenol-d4	7.09	132	214628	30.28	ng/ul	0.00
13) 4-Methylphenol-d8	8.27	113	234124	33.35	ng/ul	0.00
19) Nitrobenzene-d5	8.73	128	109706	30.73	ng/ul	0.00
22) 2-Nitrophenol-d4	9.44	143	124801	31.41	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.97	165	222401	29.66	ng/ul	0.00
29) 4-Chloroaniline-d4	10.50	131	335196	37.23	ng/ul	0.00
44) Dimethylphthalate-d6	13.64	166	843527	34.49	ng/ul	0.00
47) Acenaphthylene-d8	13.91	160	1003974	32.83	ng/ul	0.00
52) 4-Nitrophenol-d4	14.48	143	108816m>	28.38	ng/ul	0.00
58) Fluorene-d10	15.22	176	721679	34.25	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.38	200	102123	22.25	ng/ul	0.00
71) Anthracene-d10	17.08	188	1200057	36.03	ng/ul	0.00
79) Pyrene-d10	19.39	212	1283873	42.65	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.25	264	1063694	33.68	ng/ul	0.00

Target Compounds Qvalue

53 41 16 11

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