

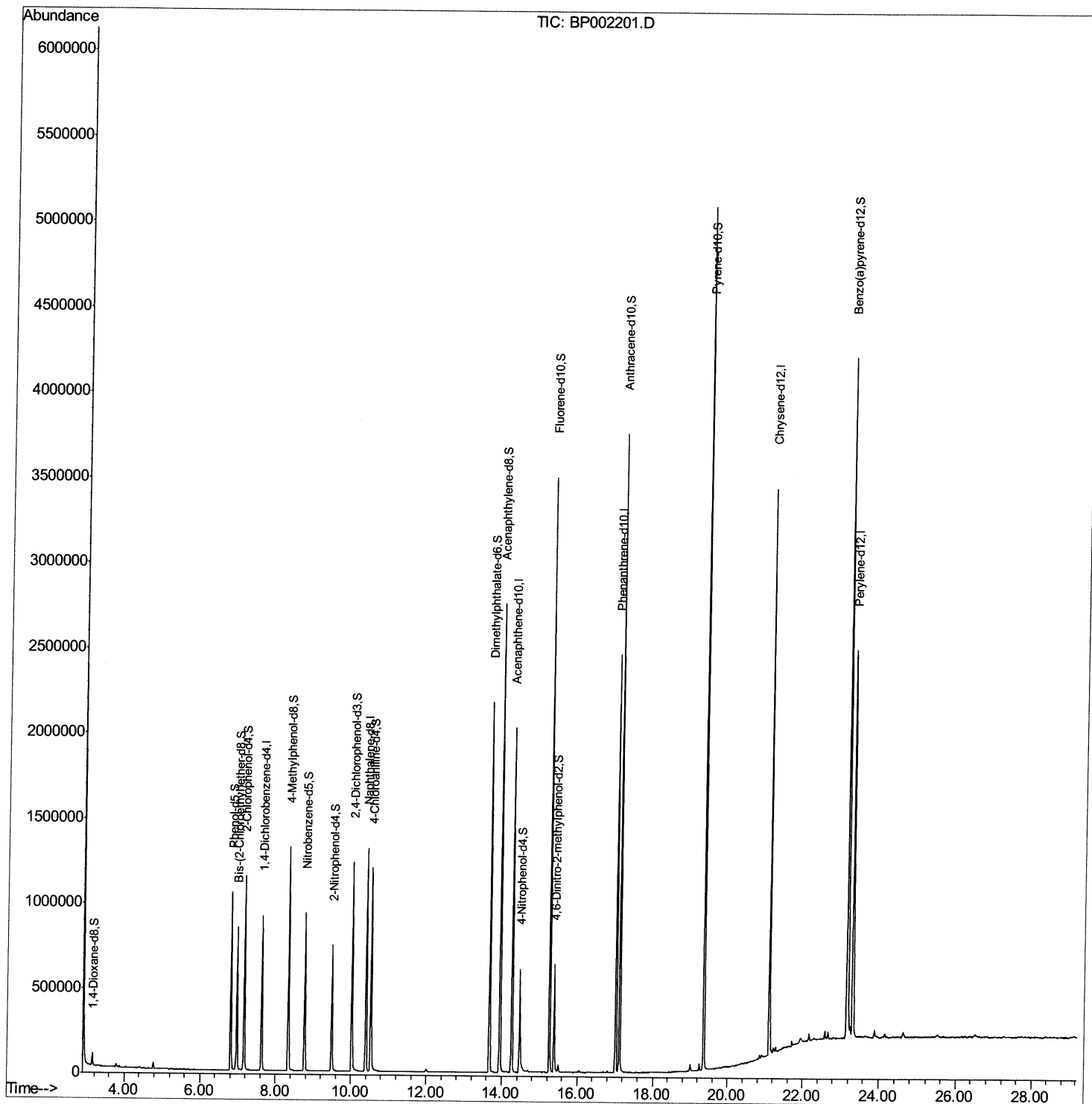
Data Path : Z:\svoasrv\HPCHEM1\BNA P\Data\BP031220\
Data File : BP002201.D
Acq On : 13 Mar 2020 00:57
Operator : CG/JU
Sample : TESTBL-03
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TESTBL-03

Manual Integrations
APPROVED

mohammad
3/13/2020 8:57:34 AM

Quant Time: Mar 13 04:06:20 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Mar 13 02:52:52 2020
Response via : Initial Calibration



Quantitation Report (Qedit)

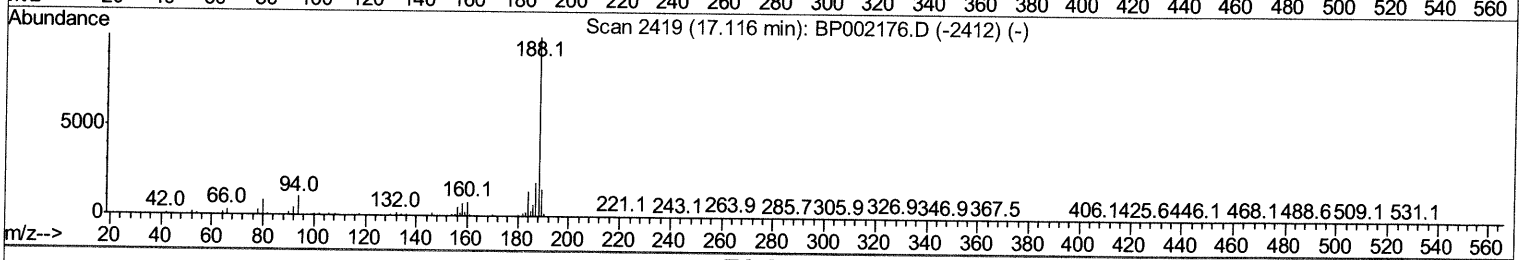
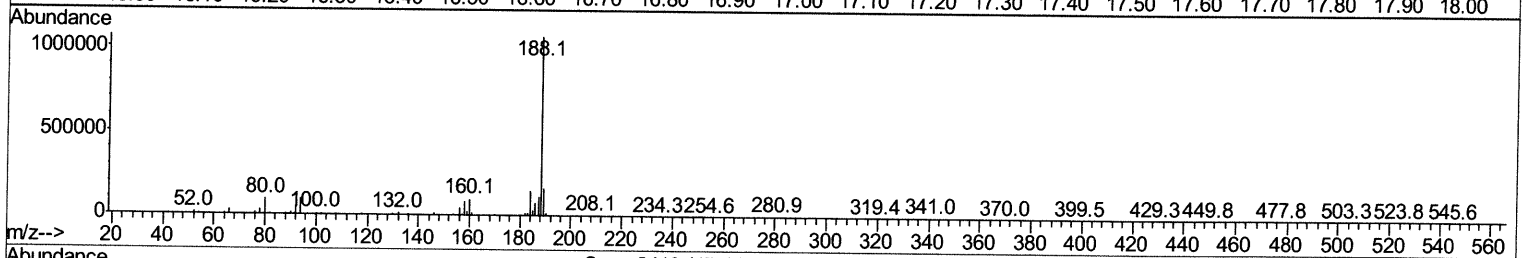
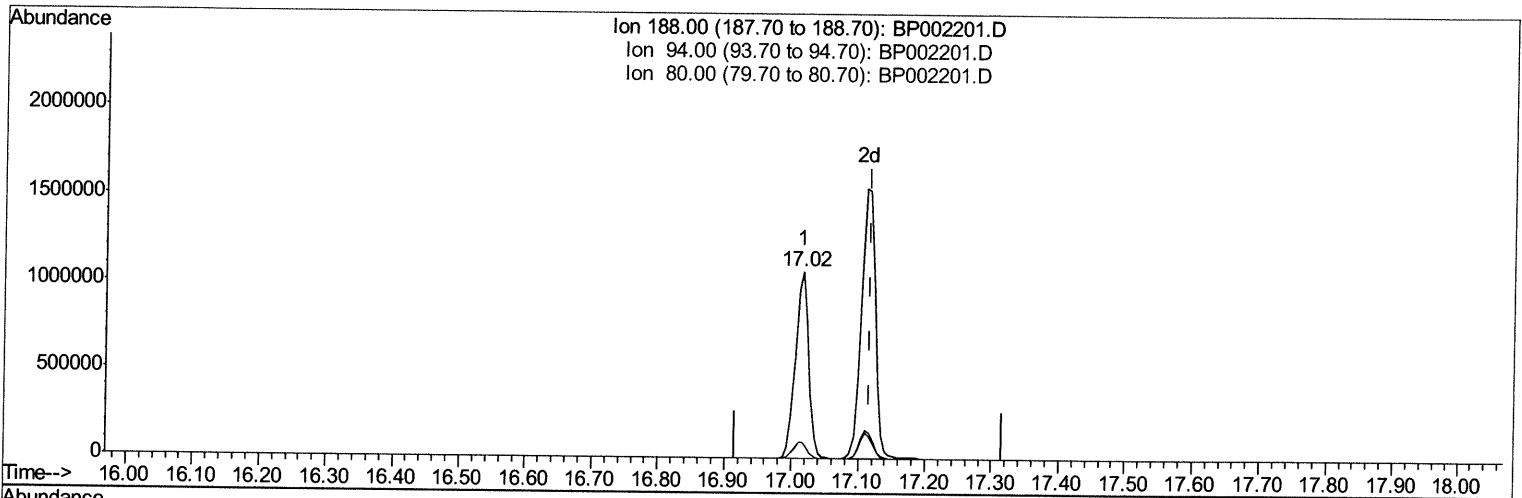
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TIC: BP002201.D

(71) Anthracene-d10 (S)

17.016min (-0.100) 22.25ng/ul

response 1493969

Ion	Exp%	Act%
188.00	100	100
94.00	9.30	8.60
80.00	7.70	8.69
0.00	0.00	0.00

Quantitation Report (Qedit)

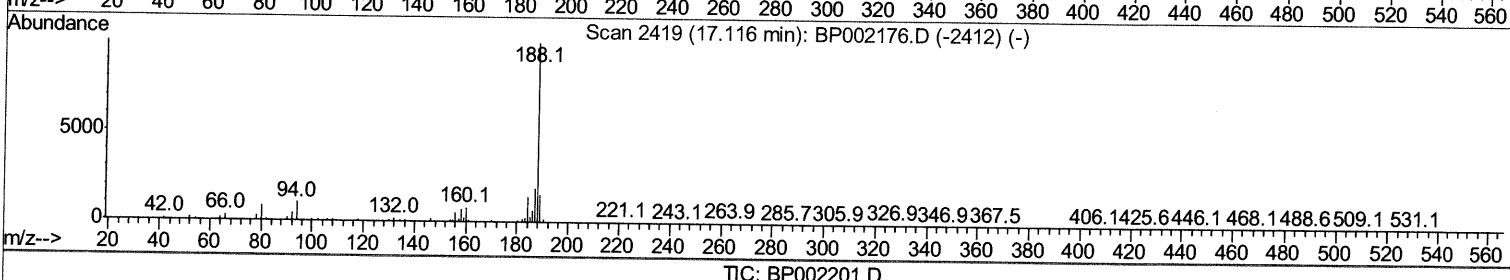
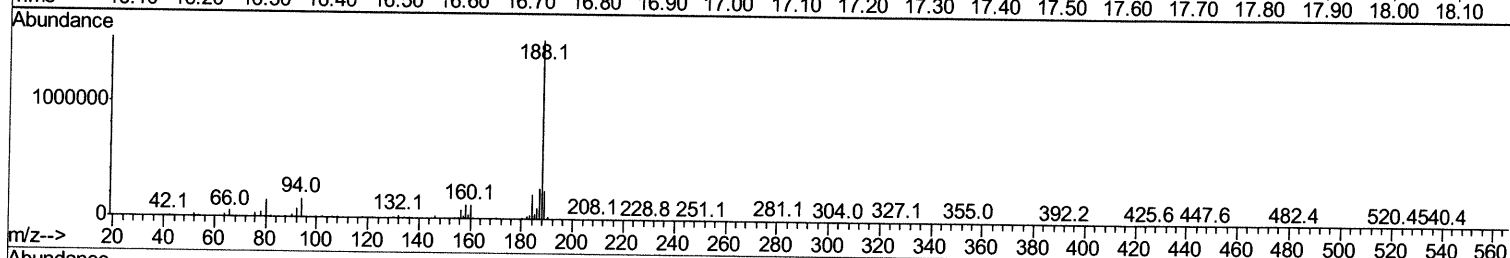
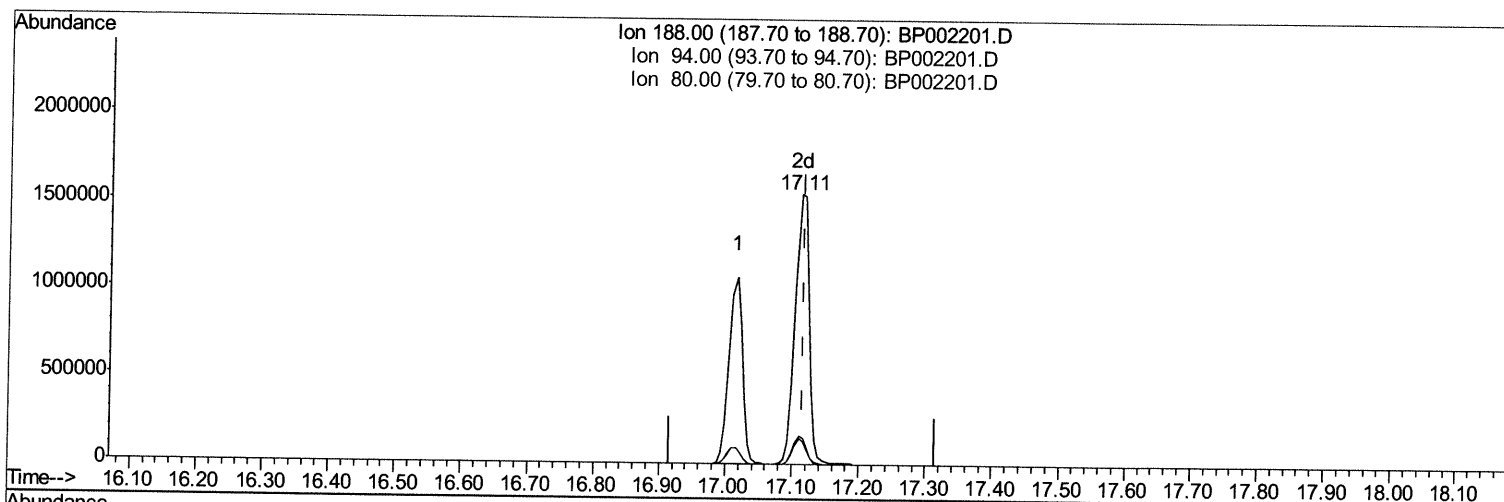
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Manual Integrations
APPROVED

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3/13/2020 8:57:34 AM

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Quant Title : SVOA CALIBRATION
QLast Update : Fri Mar 13 02:52:52 2020
Response via : Initial Calibration



(71) Anthracene-d10 (S)

17.110min (-0.006) 33.57ng/ul m 3403/14/20

response 2254376

Ion	Exp%	Act%
188.00	100	100
94.00	9.30	10.63
80.00	7.70	9.35#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031220\
 Data File : BP002201.D
 Acq On : 13 Mar 2020 00:57
 Operator : CG/JU
 Sample : TESTBL-03
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TESTBL-03

Manual Integrations
 APPROVED

mohammad
 3/13/2020 8:57:34 AM

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 Quant Title : SVOA CALIBRATION
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	271464	20.00	ng/ul	0.00
18) Naphthalene-d8	10.40	136	1121892	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.26	164	696526	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.02	188	1493969	20.00	ng/ul	0.00
78) Chrysene-d12	21.11	240	1462654	20.00	ng/ul	0.00
86) Perylene-d12	23.31	264	1695551	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.15	96	37357	5.92	ng/uL	0.00
5) Phenol-d5	6.81	99	639328	30.87	ng/ul	0.00
7) Bis-(2-Chloroethyl) ether-d	6.96	67	372492	33.97	ng/ul	0.00
9) 2-Chlorophenol-d4	7.16	132	565803	31.58	ng/ul	0.00
13) 4-Methylphenol-d8	8.33	113	515225	30.28	ng/ul	0.00
19) Nitrobenzene-d5	8.77	128	252438	29.29	ng/ul	0.00
22) 2-Nitrophenol-d4	9.49	143	272055	28.06	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.03	165	525977	28.16	ng/ul	0.00
29) 4-Chloroaniline-d4	10.53	131	726531	35.48	ng/ul	0.00
44) Dimethylphthalate-d6	13.68	166	1572764	29.84	ng/ul	0.00
47) Acenaphthylene-d8	13.95	160	1940178	30.78	ng/ul	0.00
52) 4-Nitrophenol-d4	14.47	143	198565	19.70	ng/ul	0.00
58) Fluorene-d10	15.26	176	1443475	31.65	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.39	200	163916	19.90	ng/ul	0.00
71) Anthracene-d10	17.11	188	2254376m>	33.57	ng/ul	0.00
79) Pyrene-d10	19.36	212	2590804	33.84	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.17	264	2875425	33.15	ng/ul	0.00

5u03/14/20

Target Compounds Ovalue

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