

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP102219\
Quantitation Report (QT Reviewed)

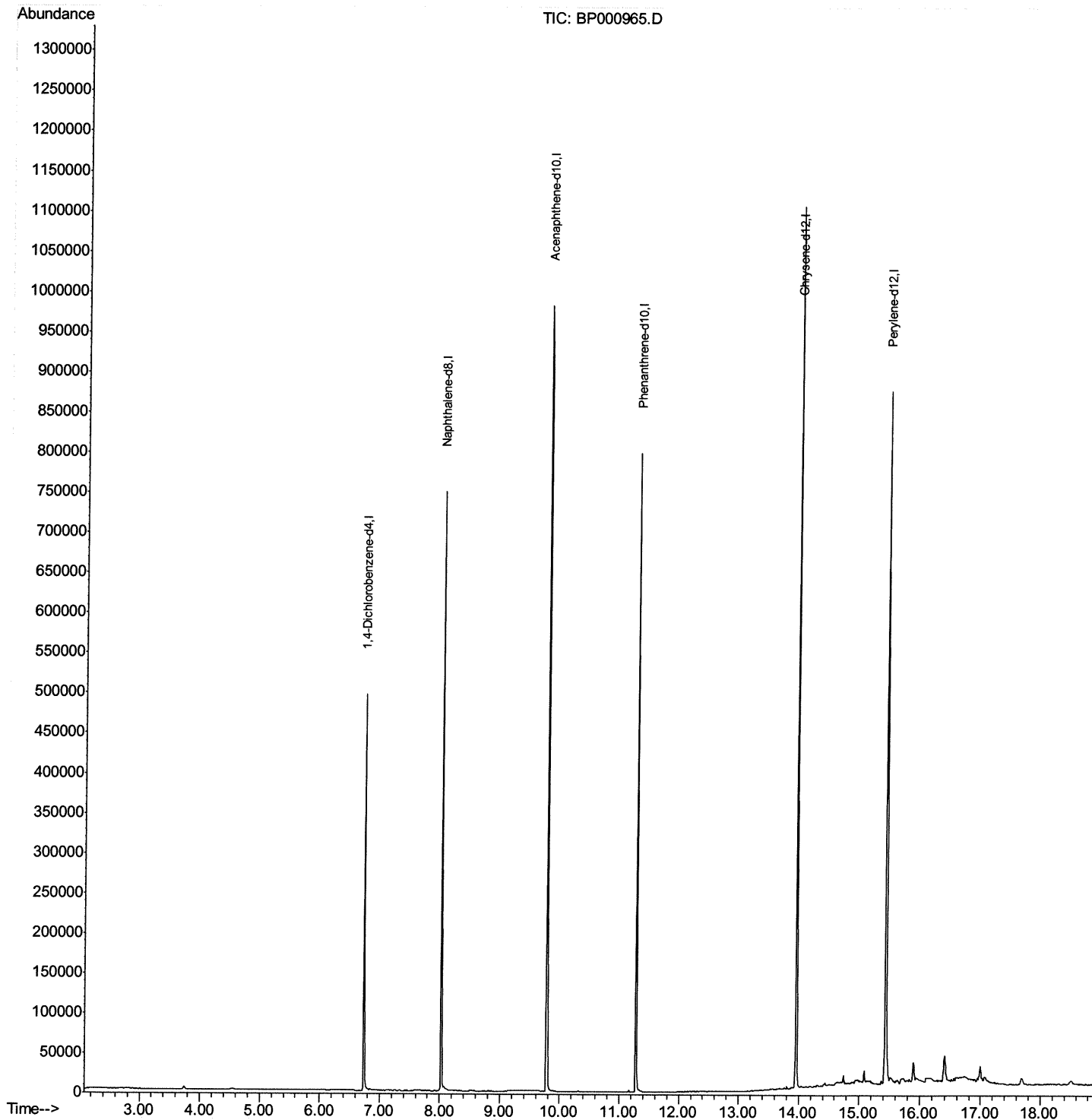
Data File : BP000965.D
Acq On : 23 Oct 2019 01:07
Operator : JU
Sample : K5449-10
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SVOC-GPC2-BLANK

Manual Integrations
APPROVED

mohammad
10/24/2019 8:08:08 AM

Quant Time: Oct 23 07:07:13 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP100919MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Oct 23 05:34:10 2019
Response via : Initial Calibration



Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP102219\
Quantitation Report (Qedit)

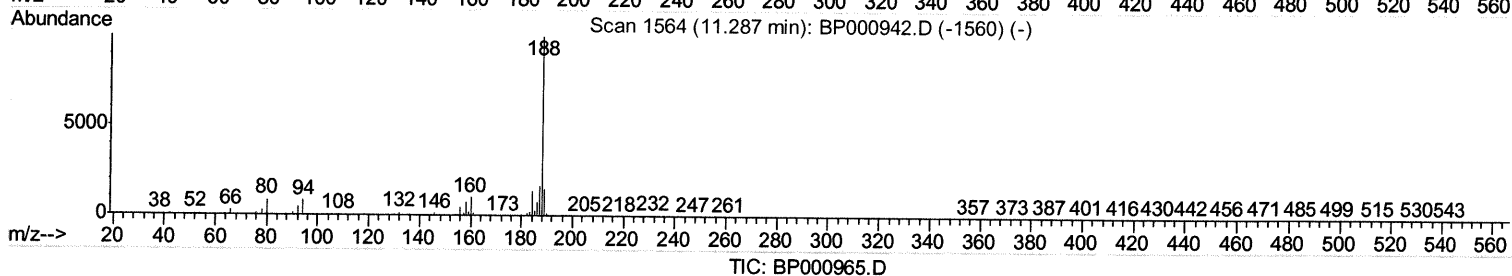
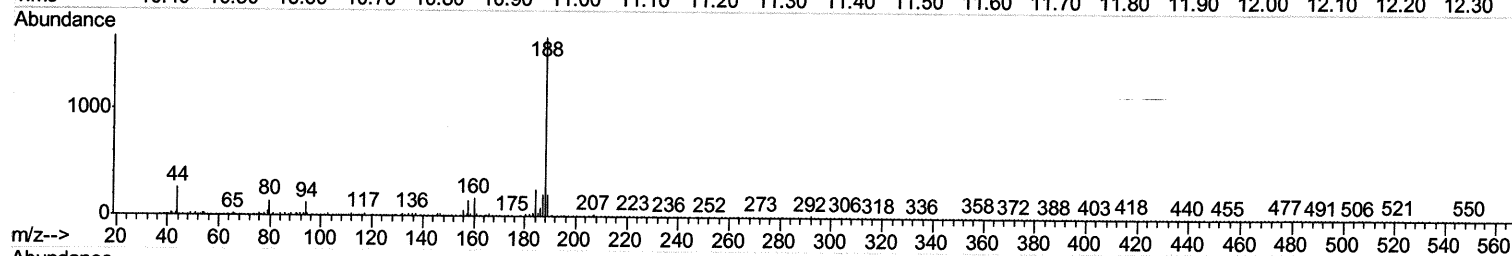
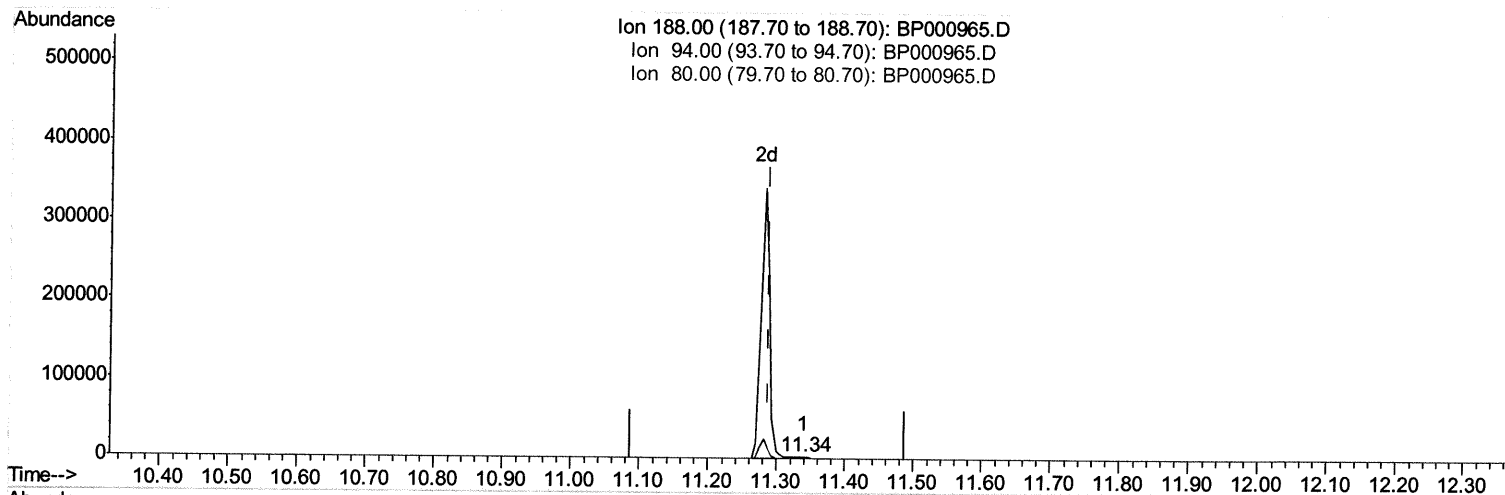
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Quant Time: Oct 23 05:40:18 2019
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TIC: BP000965.D

(62) Phenanthrene-d10 (I)

11.340min (+0.053) 20.00ng/ul

response 1447

Ion	Exp%	Act%
188.00	100	100
94.00	9.10	9.06
80.00	9.30	8.28
0.00	0.00	0.00

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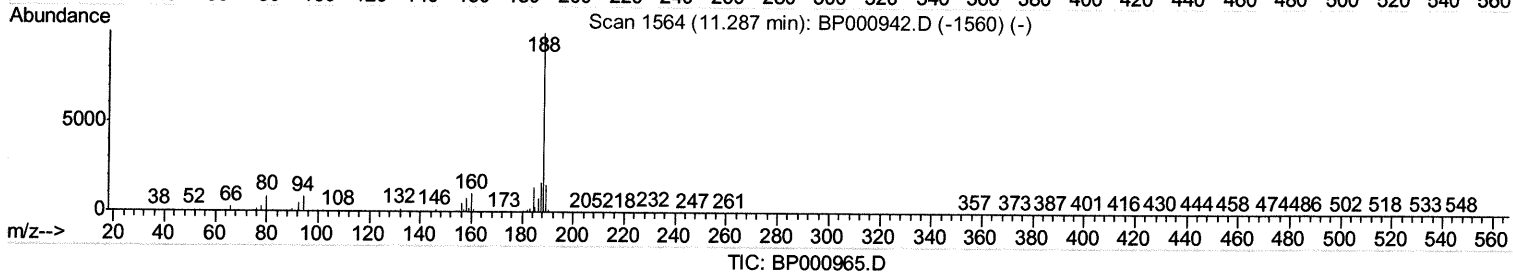
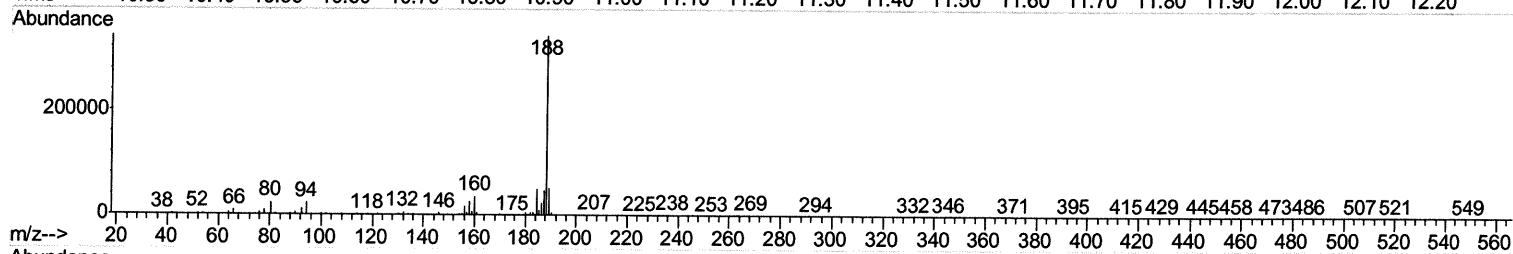
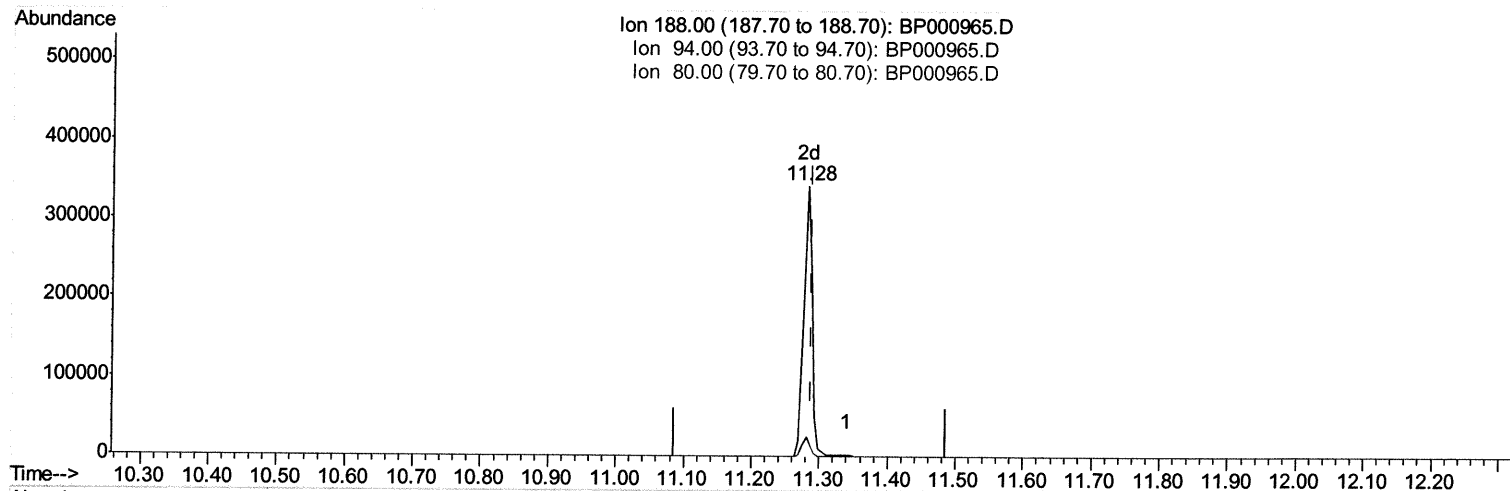
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Sample : K5449-10
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SVOC-GPC2-BLANK

Manual Integrations
APPROVED

mohammad
10/24/2019 8:08:08 AM

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QLast Update : Wed Oct 23 05:34:10 2019
Response via : Initial Calibration



(62) Phenanthrene-d10 (I)

11.281min (-0.006) 20.00ng/ul m *24 10.23.19*

response 286624

Ion	Exp%	Act%
188.00	100	100
94.00	9.10	6.63#
80.00	9.30	7.17#
0.00	0.00	0.00

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Operator : JU
Sample : K5449-10
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SVOC-GPC2-BLANK

Manual Integrations
APPROVED

mohammad
10/24/2019 8:08:08 AM

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.74	152	84082	20.00	ng/ul	0.00
18) Naphthalene-d8	8.02	136	329061	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.78	164	210109	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.28	188	286624m	20.00	ng/ul	0.00
78) Chrysene-d12	13.96	240	381352	20.00	ng/ul	0.00
86) Perylene-d12	15.43	264	416135	20.00	ng/ul	-0.01

7 10-23-19

System Monitoring Compounds

3) 1,4-Dioxane-d8	0.00	96	0d	0.00	ng/uL
5) Phenol-d5	0.00	99	0d	0.00	ng/ul
7) Bis-(2-Chloroethyl) ether-d	0.00	67	0d	0.00	ng/ul
9) 2-Chlorophenol-d4	0.00	132	0d	0.00	ng/ul
13) 4-Methylphenol-d8	0.00	113	0d	0.00	ng/ul
19) Nitrobenzene-d5	0.00	128	0d	0.00	ng/ul
22) 2-Nitrophenol-d4	0.00	143	0d	0.00	ng/ul
26) 2,4-Dichlorophenol-d3	0.00	165	0d	0.00	ng/ul
29) 4-Chloroaniline-d4	0.00	131	0d	0.00	ng/ul
44) Dimethylphthalate-d6	0.00	166	0d	0.00	ng/ul
47) Acenaphthylene-d8	0.00	160	0d	0.00	ng/ul
52) 4-Nitrophenol-d4	0.00	143	0d	0.00	ng/ul
58) Fluorene-d10	0.00	176	0d	0.00	ng/ul
63) 4,6-Dinitro-2-methylphenol	0.00	200	0d	0.00	ng/ul
71) Anthracene-d10	0.00	188	0d	0.00	ng/ul
79) Pyrene-d10	0.00	212	0d	0.00	ng/ul
90) Benzo(a)pyrene-d12	0.00	264	0d	0.00	ng/ul

Target Compounds Qvalue

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