

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD\_R\Data\PR103019\  
Data File : PR042648.D  
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch  
Acq On : 31 Oct 2019 03:31  
Operator : SM\AJ  
Sample : K5538-01  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

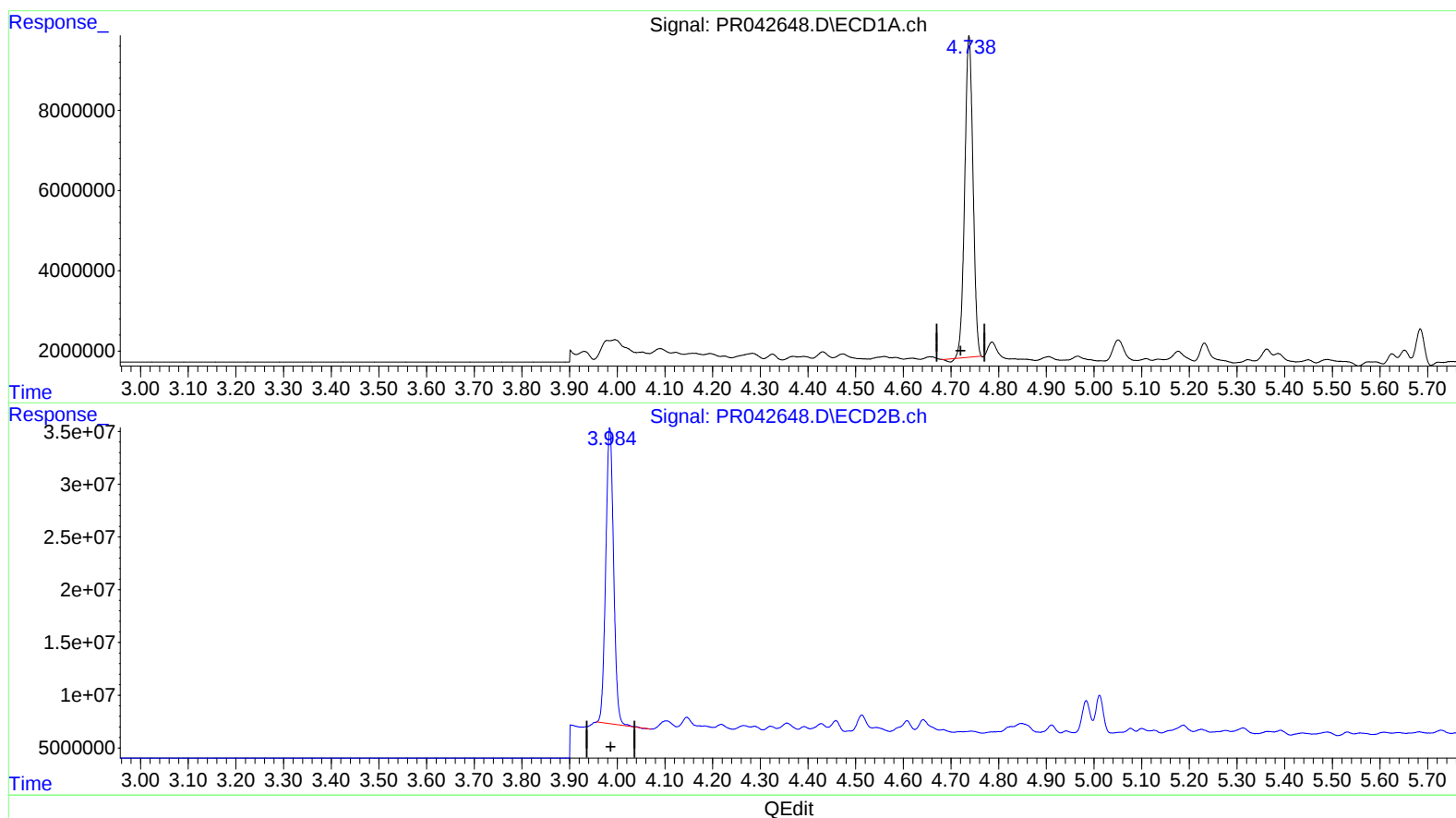
Instrument :  
ECD\_R  
ClientSampleId :  
BF6M1

Manual Integrations  
APPROVED

Ankita  
10/31/2019 1:07:43 PM

Integration File signal 1: autoint1.e  
Integration File signal 2: autoint2.e  
Quant Time: Oct 31 09:04:00 2019  
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD\_R\Method\PR100319CLP.M  
Quant Title : GC EXTRACTABLES  
QLast Update : Fri Oct 04 06:59:35 2019  
Response via : Initial Calibration  
Integrator: ChemStation

Volume Inj. : 1 µl  
Signal #1 Phase : ZB MR1 Signal #2 Phase: ZB MR2  
Signal #1 Info : 30Mx0.32mmx 0.5µm Signal #2 Info : 30M x 0.32mm x 0.25µm



(1) Tetrachloro-m-xylene (SA)

4.738min 23.595 ng/ml

response 95985474

(1) Tetrachloro-m-xylene #2 (SA)

3.984min 20.987 ng/ml

response 320908421

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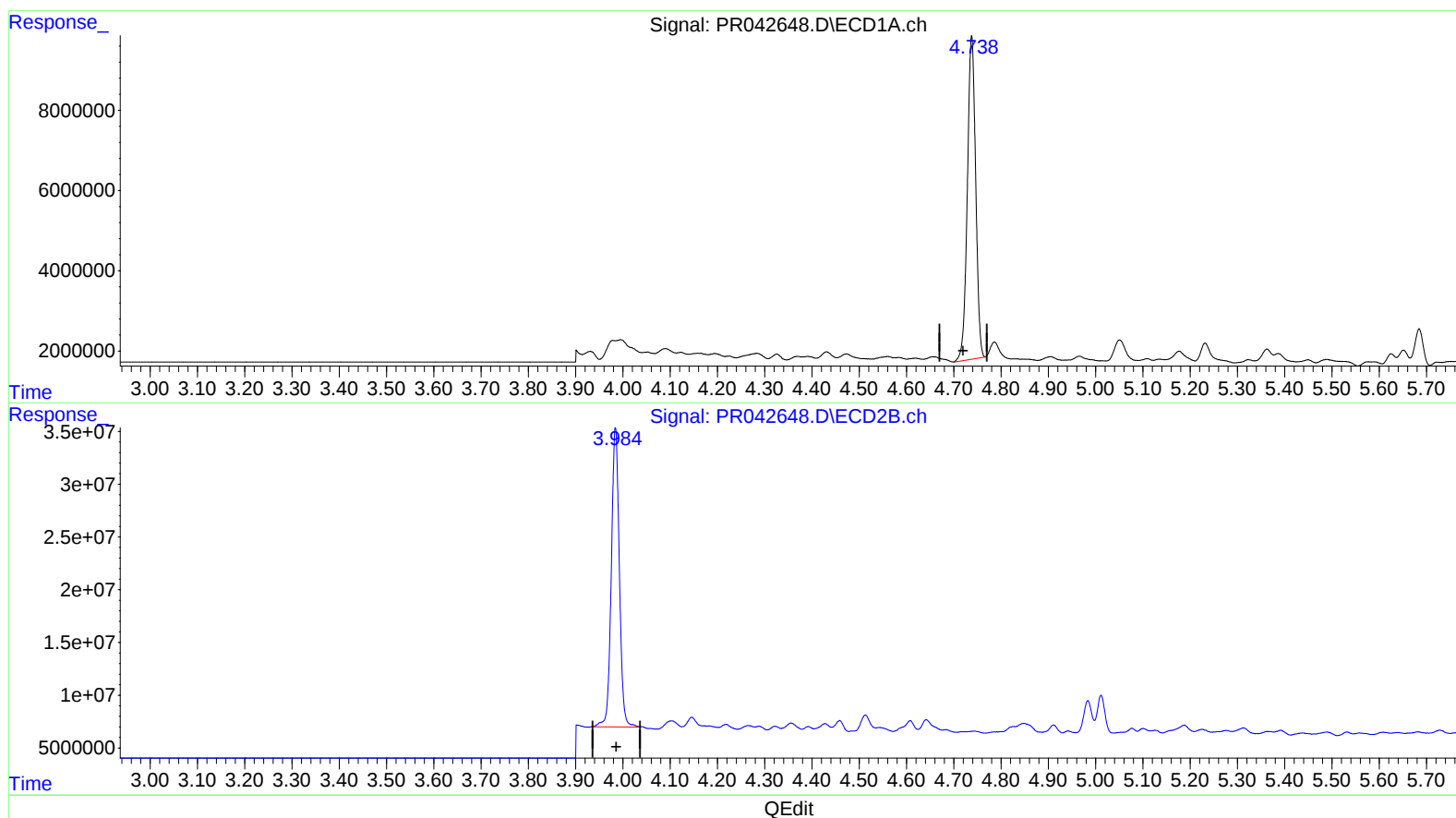
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(1) Tetrachloro-m-xylene (SA)

4.738min 24.211 ng/ml m

response 98488084

(1) Tetrachloro-m-xylene #2 (SA)

3.984min 21.916 ng/ml m

response 335122099

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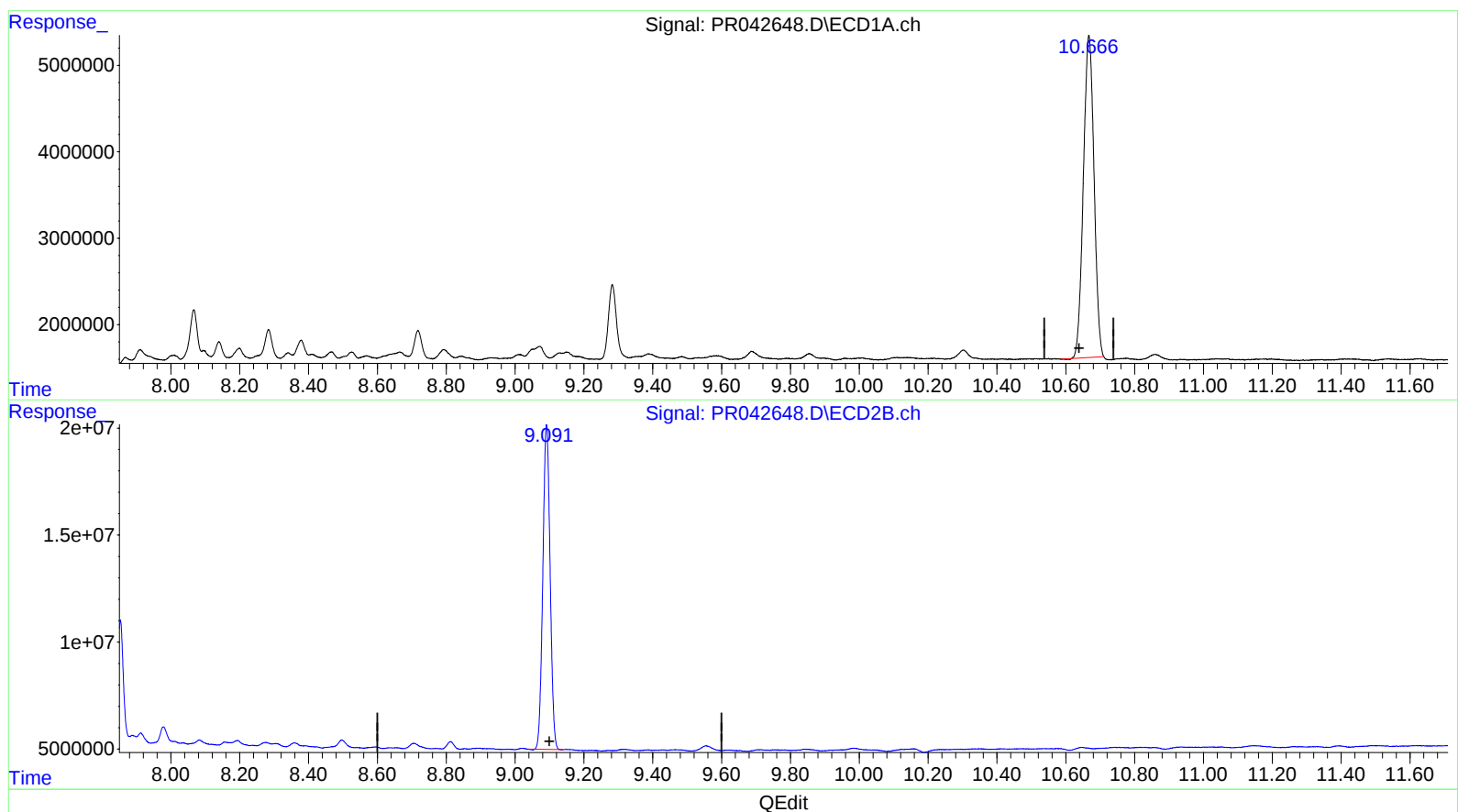
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(2) Decachlorobiphenyl (SA)

10.667min 23.238 ng/ml

response 79848911

(2) Decachlorobiphenyl #2 (SA)

9.091min 22.407 ng/ml

response 222150351

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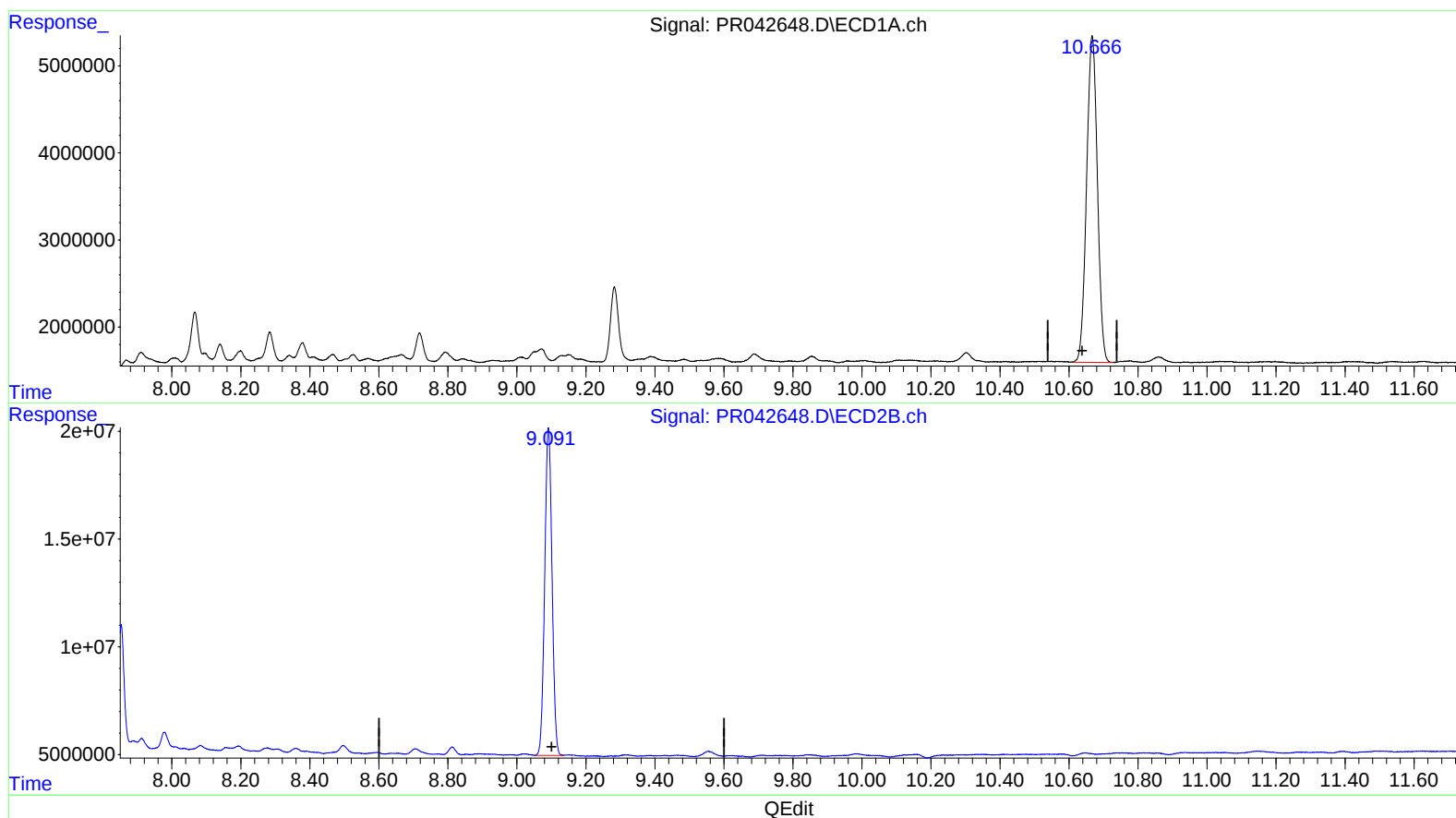
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(2) Decachlorobiphenyl (SA)

10.666min 23.672 ng/ml m

response 81340724

(2) Decachlorobiphenyl #2 (SA)

9.091min 22.519 ng/ml m

response 223260564

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| Compound | RT#1 | RT#2 | Resp#1 | Resp#2 | ng/ml | ng/ml |
|----------|------|------|--------|--------|-------|-------|
|----------|------|------|--------|--------|-------|-------|

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System Monitoring Compounds

|                    |        |       |          |         |         |         |
|--------------------|--------|-------|----------|---------|---------|---------|
| 1) SA Tetrachlo... | 4.738  | 3.984 | 98488084 | 335.1E6 | 24.211m | 21.916m |
| 2) SA Decachlor... | 10.666 | 9.091 | 81340724 | 223.3E6 | 23.672m | 22.519m |

Target Compounds  
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(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

J A-J  
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