

Instrument :
BNA_E
ClientSampleId :
C0AA5DL

Manual Integrations

Sohil
8/16/2017 6:43:06 PM

[illegible]

Quantitation Report (Qedit)

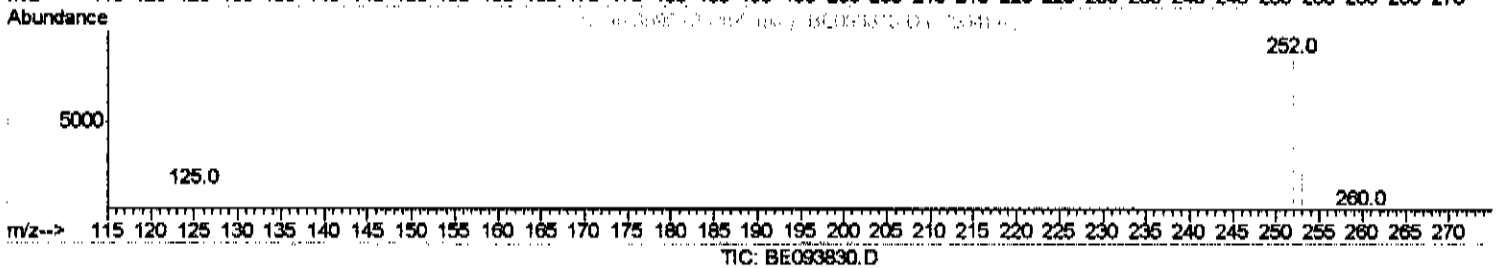
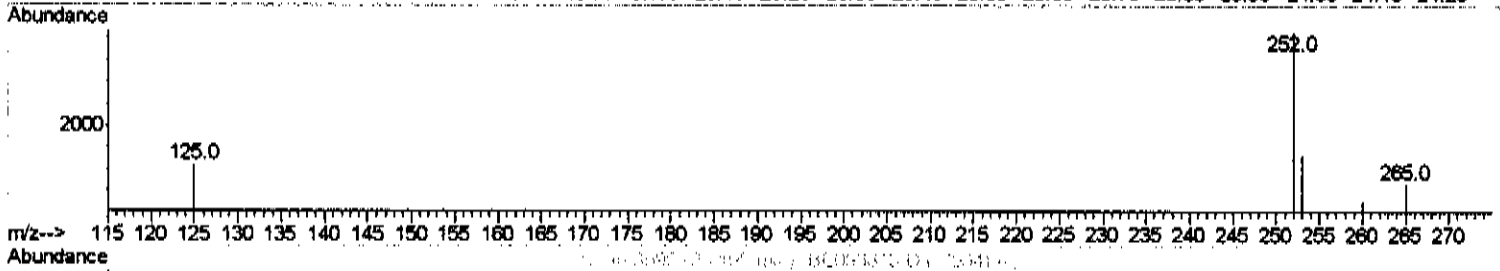
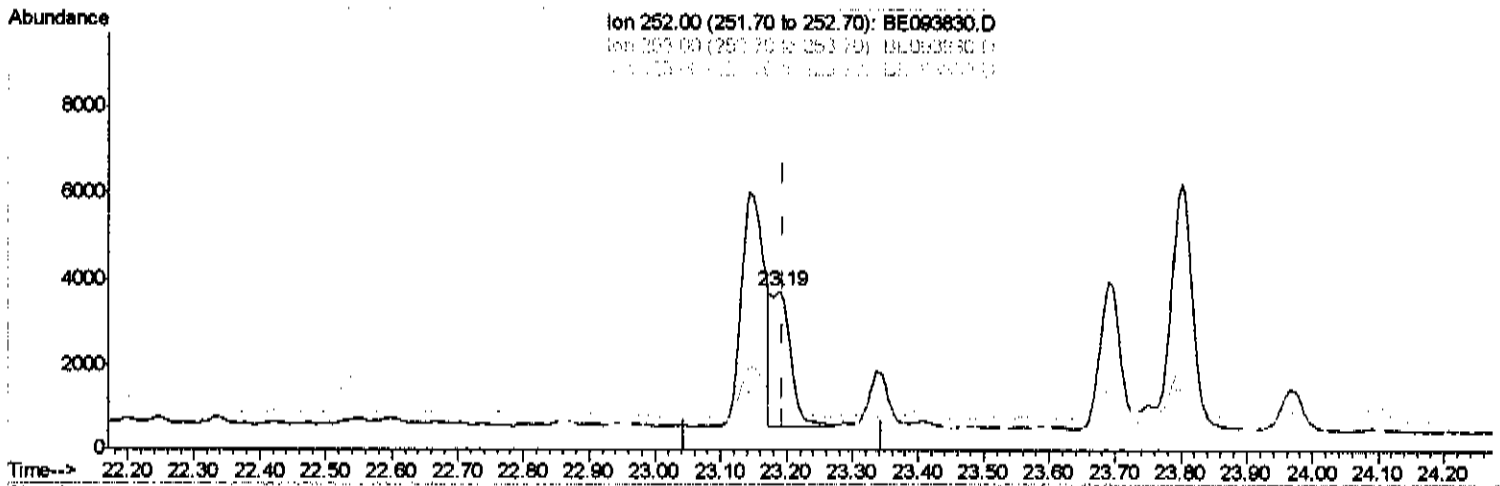
Data Path : Z:\HPCHEM1\BNA E\DATA\BE081517\
 Data File : BE093830.D
 Acq On : 16 Aug 2017 5:05
 Operator : SJ/JU
 Sample : I4529-11DL 5X
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleID :
 C0AA5DL

Manual Integrations
 APPROVED

Sohil
 8/16/2017 6:43:06 PM

Quant Time: Aug 16 07:39:14 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE072417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 16 07:30:49 2017
 Response via : Initial Calibration



(22) Benzo(k)fluoranthene

23.190min (-0.005) 0.12ng/ui m

response 6567

Ion	Exp%	Act%
252.00	100	100
253.00	30.60	38.51%
125.00	17.10	33.95%
0.00	0.00	0.00

Quantitation Report (Qedit)

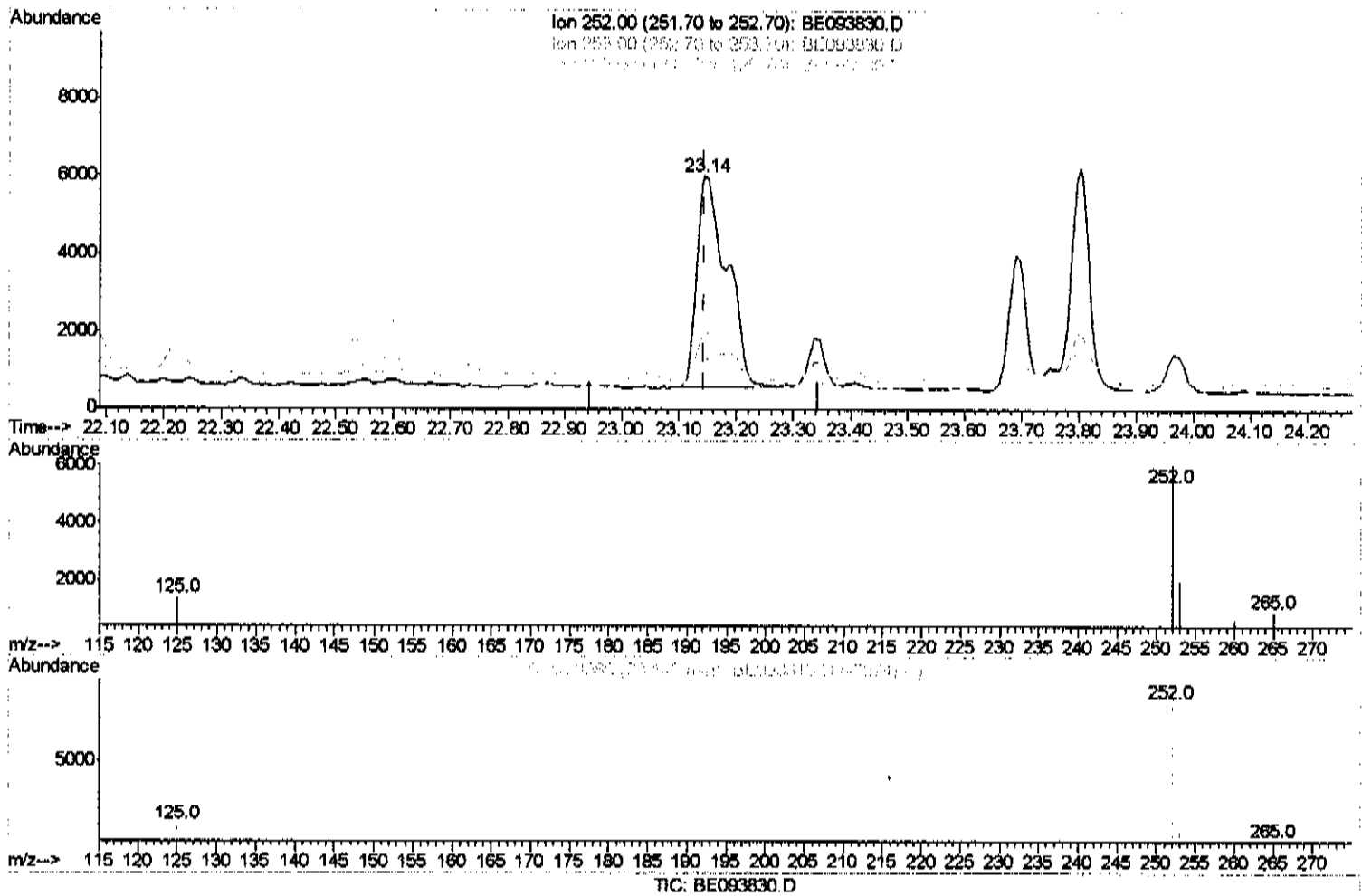
Data Path : Z:\HPCHEM1\BNA E\DATA\BE081517\
 Data File : BE093830.D
 Acq On : 16 Aug 2017 5:05
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 Sample : I4529-11DL 5X
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Instrument :
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 ClientSampled :
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Manual Integrations
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(21) Benzo(b)fluoranthene

23.144min (-0.001) 0.37ng/ul

response 19930

Ion	Exp%	Act%
252.00	100	100
253.00	32.10	33.41
125.00	17.50	22.99
0.00	0.00	0.00

Quantitation Report (Qedit)

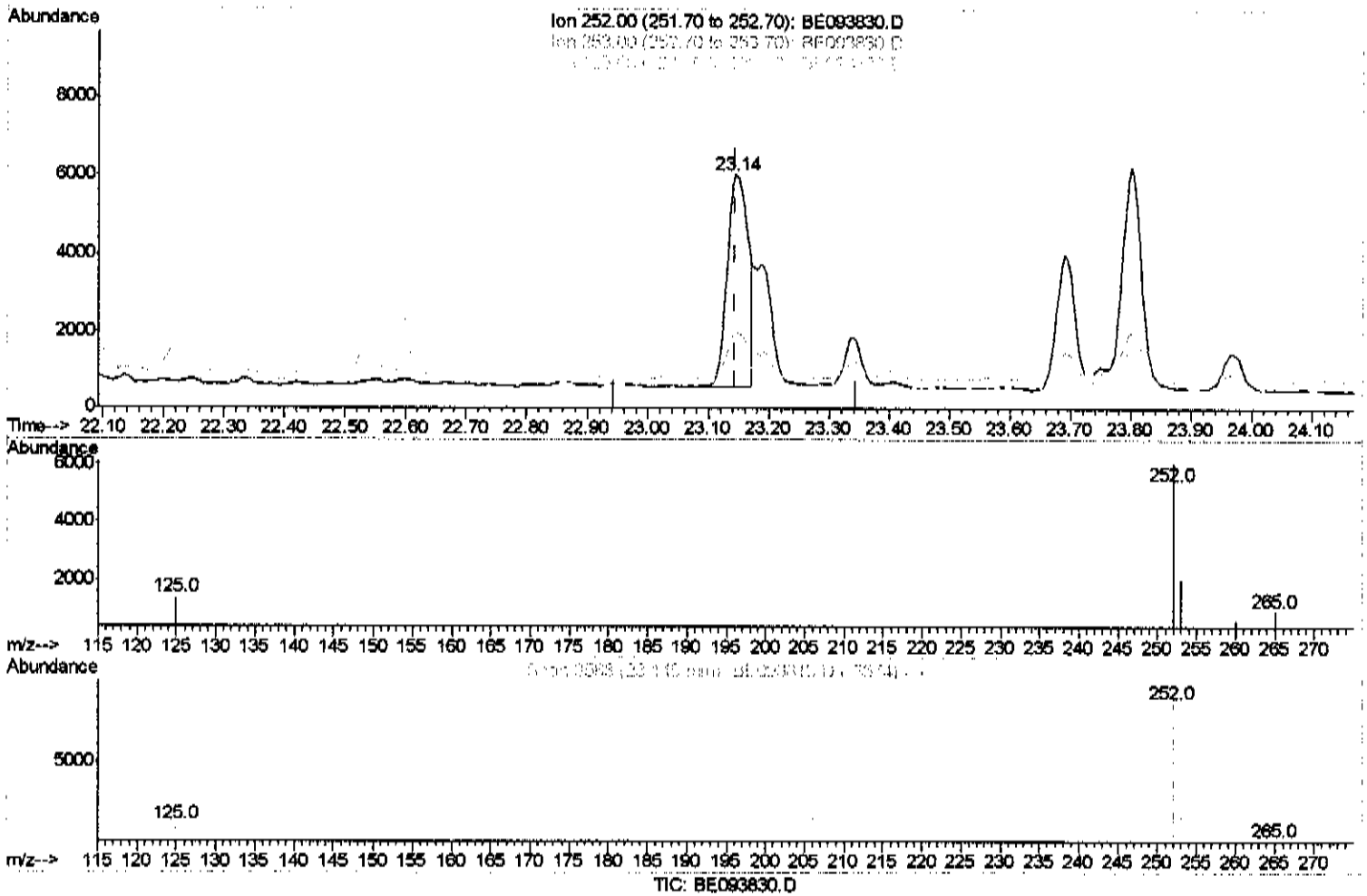
Data Path : Z:\HPCHEM1\BNA E\DATA\BE081517\
 Data File : BE093830.D
 Acq On : 16 Aug 2017 5:05
 Operator : SJ/JU
 Sample : I4529-11DL 5X
 Misc :
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Manual Integrations
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 Response via : Initial Calibration



(21) Benzo(b)fluoranthene

23.144min (-0.001) 0.25ng/ul m

response 13310

Ion	Exp%	Act%
252.00	100	100
253.00	32.10	33.41
125.00	17.50	22.99
0.00	0.00	0.00

Quantitation Report (Qedit)

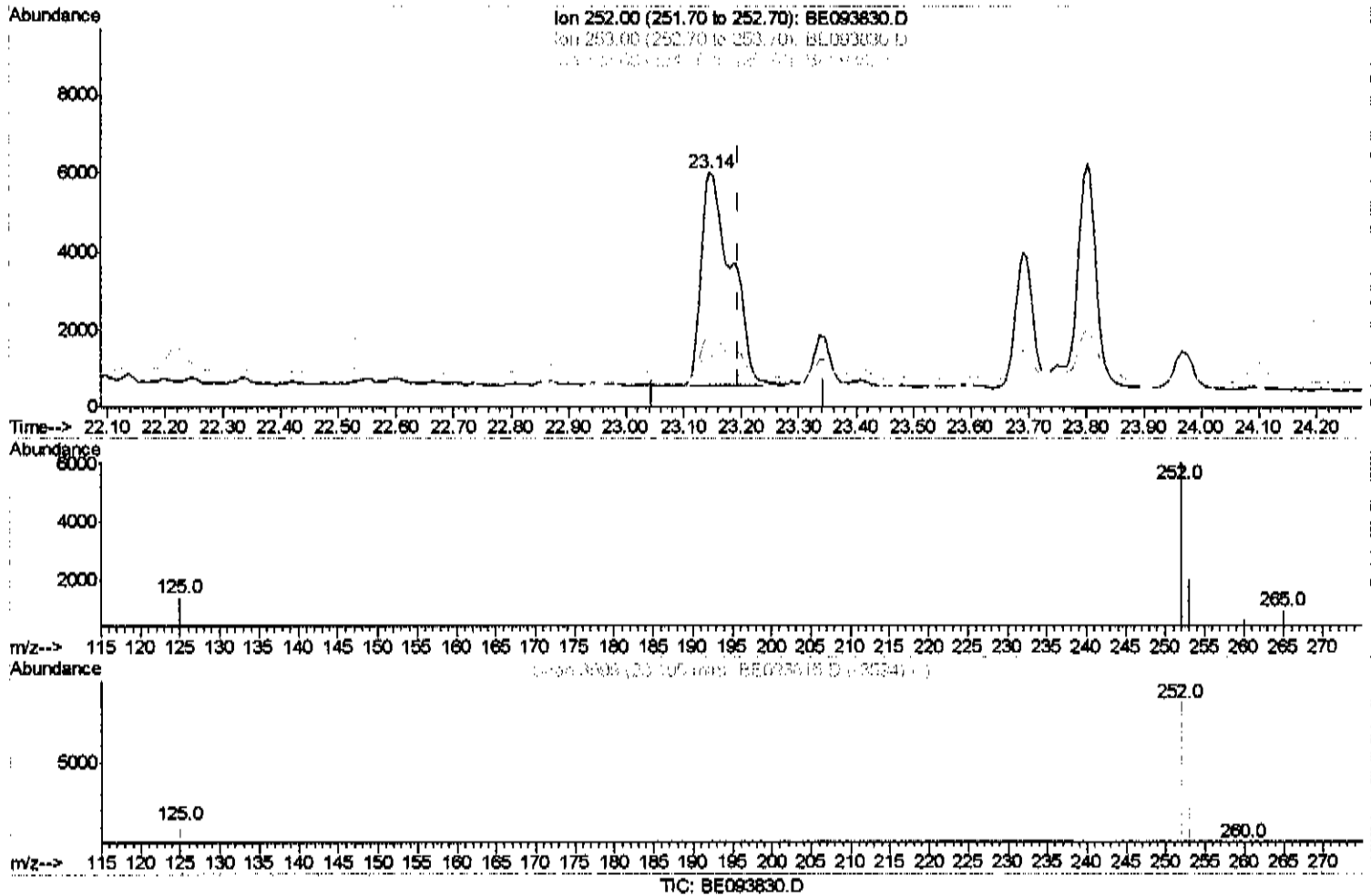
Data Path : Z:\HPCHEM1\BNA E\DATA\BE081517\
 Data File : BE093830.D
 Acq On : 16 Aug 2017 5:05
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 Sample : I4529-11DL 5X
 Misc :
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Instrument :
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Manual Integrations
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Quant Time: Aug 16 07:39:14 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE072417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 16 07:30:49 2017
 Response via : Initial Calibration



(22) Benzo(k)fluoranthene
 23.144min (-0.051) 0.38ng/ul
 response 18930

Ion	Exp%	Act%
252.00	100	100
253.00	30.60	33.41
125.00	17.10	22.99
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA E\DATA\BE081517\
 Data File : BE093830.D
 Acq On : 16 Aug 2017 5:05
 Operator : SJ/JU
 Sample : I4529-11DL 5X
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampled :
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Manual Integrations
 APPROVED

Sohil
 8/16/2017 6:43:06 PM

Quant Time: Aug 16 08:44:06 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE072417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 16 07:30:49 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	2363	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	12562	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	8197	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	15988	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	16977	0.40	ng/ul	0.00
20) Perylene-d12	23.91	264	17238	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.28	152	1825	0.09	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	5255	0.10	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
3) Naphthalene	10.75	128	25019	0.819	ng/ul#	93
5) 2-Methylnaphthalene	12.35	142	31561	1.416	ng/ul	100
7) Acenaphthylene	14.24	152	5244	0.153	ng/ul#	65
8) Acenaphthene	14.58	153	7585	0.281	ng/ul	95
9) Fluorene	15.57	166	39407	1.156	ng/ul	92
12) Phenanthrene	17.29	178	93915	2.157	ng/ul#	92
13) Anthracene	17.38	178	23218	0.547	ng/ul#	91
15) Fluoranthene	19.30	202	49573	0.910	ng/ul	83
17) Pyrene	19.66	202	58805	1.187	ng/ul#	83
18) Benzo(a)anthracene	21.39	228	18097	0.368	ng/ul#	81
19) Chrysene	21.45	228	18740	0.397	ng/ul#	82
21) Benzo(b)fluoranthene	23.14	252	13310m}	0.250	ng/ul	96
22) Benzo(k)fluoranthene	23.19	252	6567m}	0.125	ng/ul	
23) Benzo(a)pyrene	23.80	252	12482	0.246	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.55	276	5571	0.099	ng/ul#	83
26) Benzo(a,h,i)perylene	27.37	276	4302	0.091	ng/ul#	82

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