

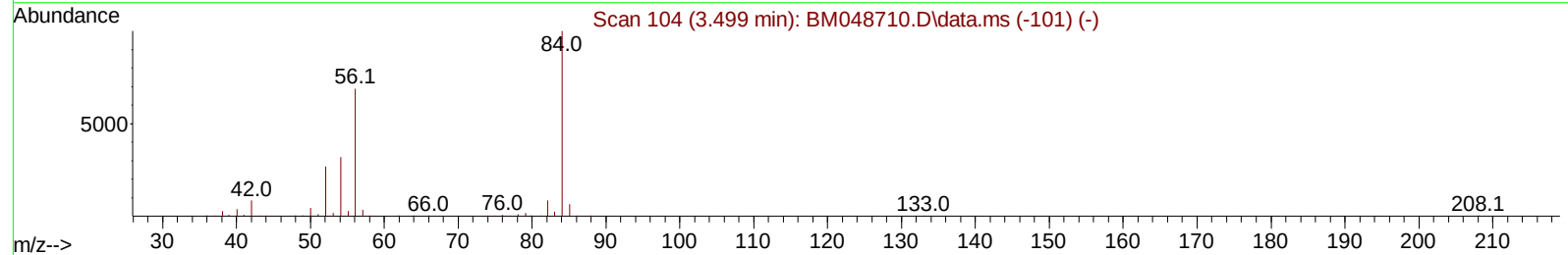
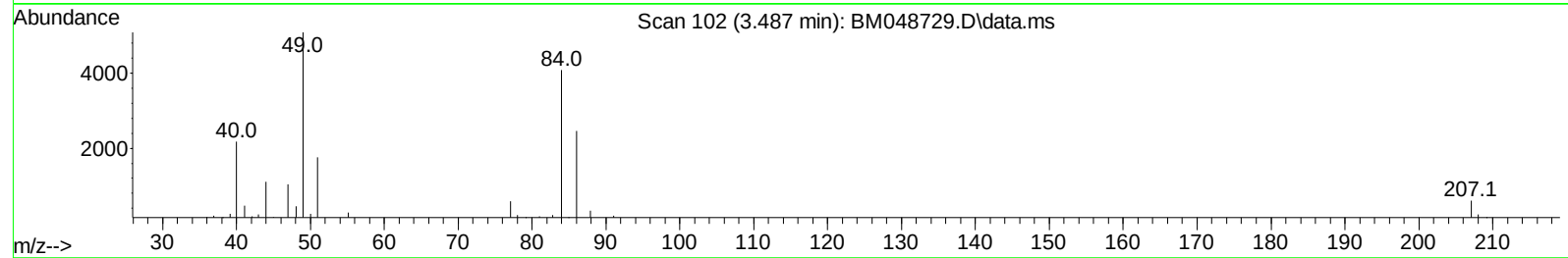
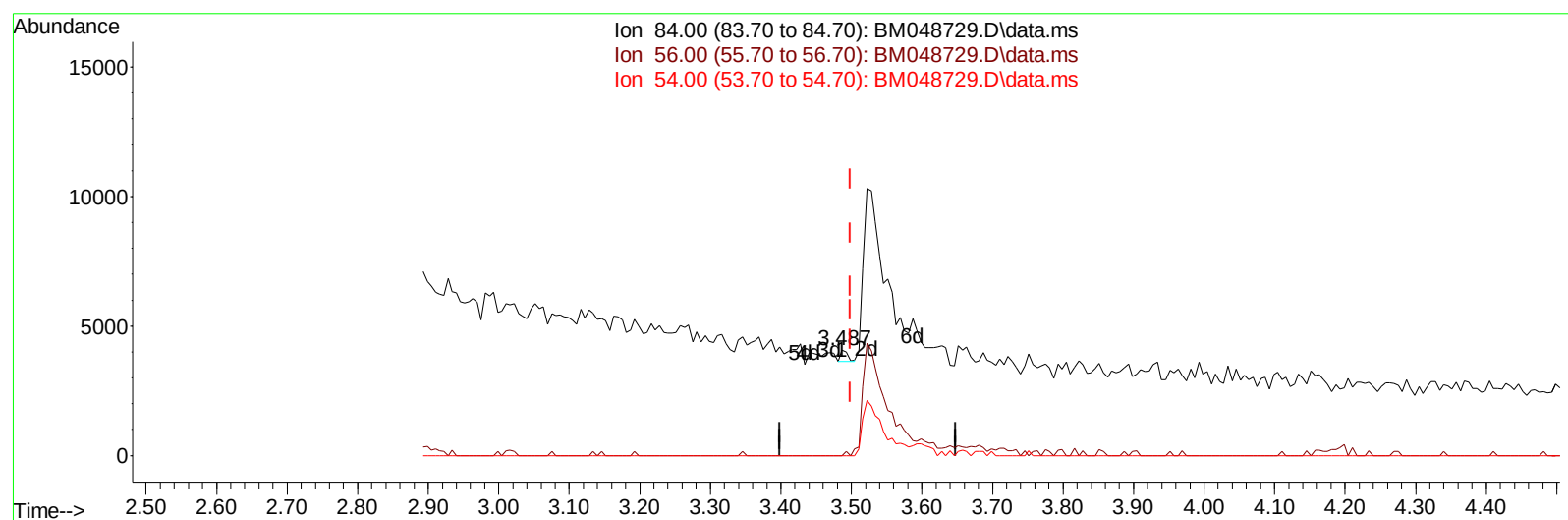
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111524\
Data File : BM048729.D
Acq On : 16 Nov 2024 05:02
Operator : RC/JU
Sample : P4751-21
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BH9G2

Manual IntegrationsAPPROVED

Quant Time: Nov 16 05:36:22 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM110724.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Nov 15 16:55:33 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/18/2024
Supervised By :mohammad ahmed 11/18/2024



TIC: BM048729.D\data.ms

(4) Pyridine-d5 (S)

3.487min (-0.012) 0.03 ng/u1

response 306

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.80	0.00#
54.00	29.50	0.00#
0.00	0.00	0.00

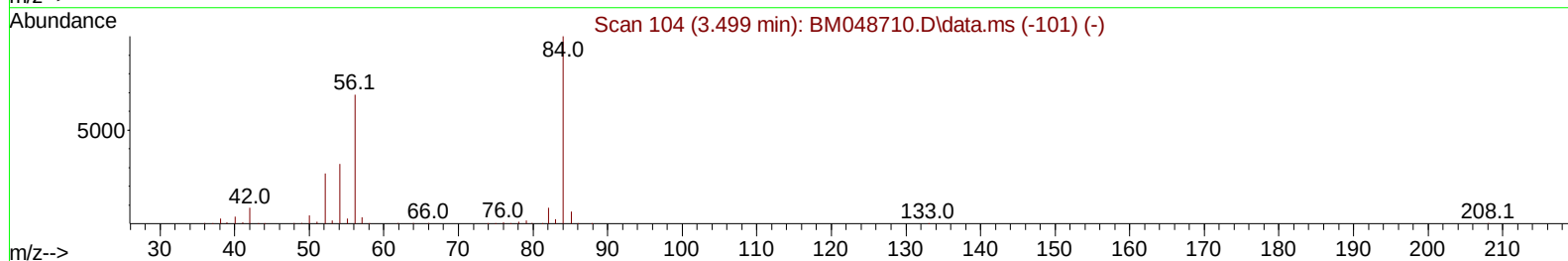
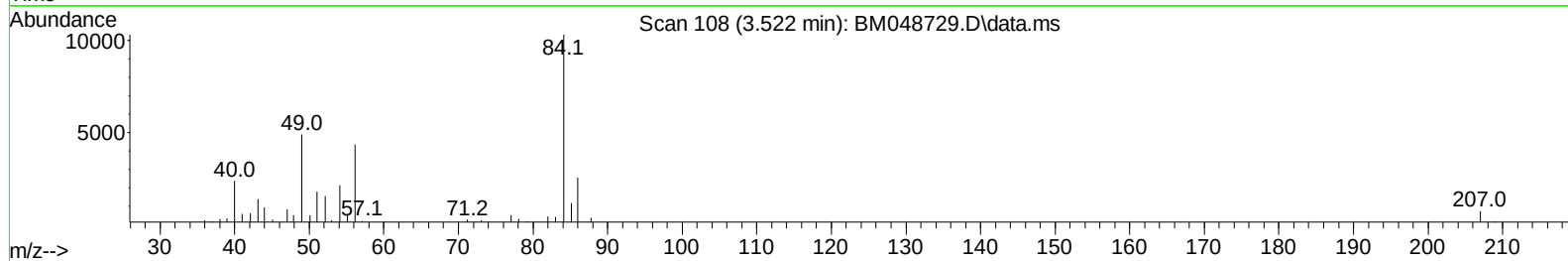
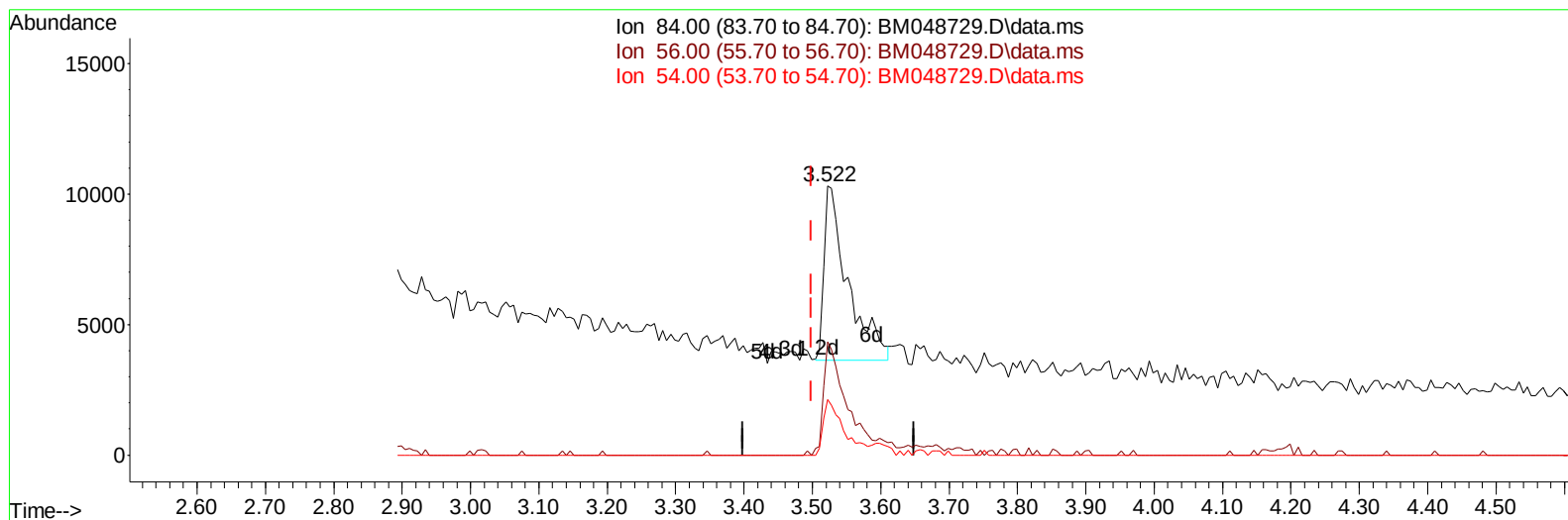
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TIC: BM048729.D\data.ms

(4) Pyridine-d5 (S)

3.522min (+ 0.024) 1.76 ng/ul m

response 16049

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.80	42.08#
54.00	29.50	20.76#
0.00	0.00	0.00

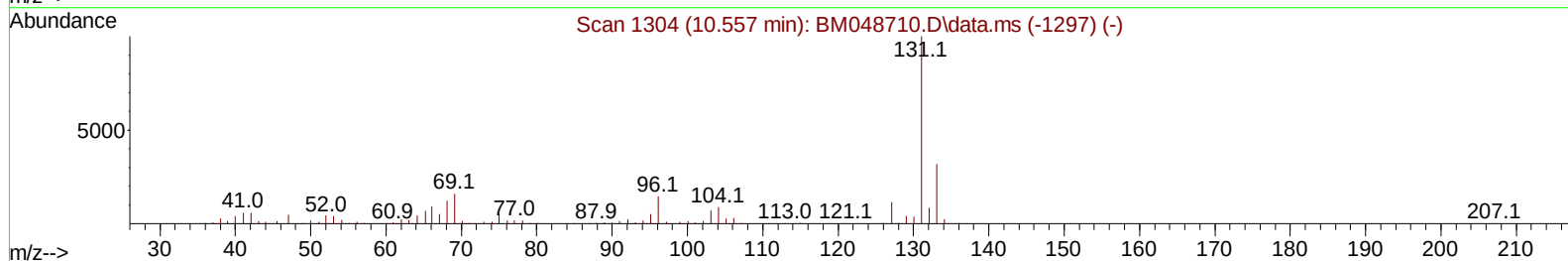
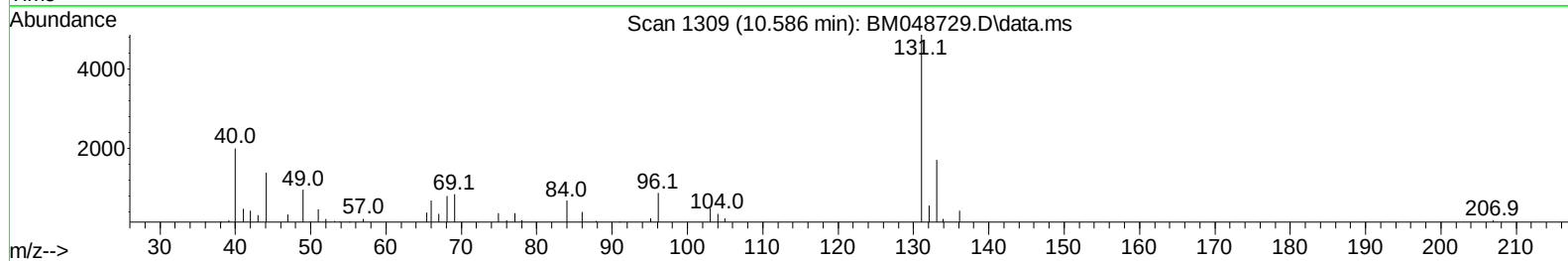
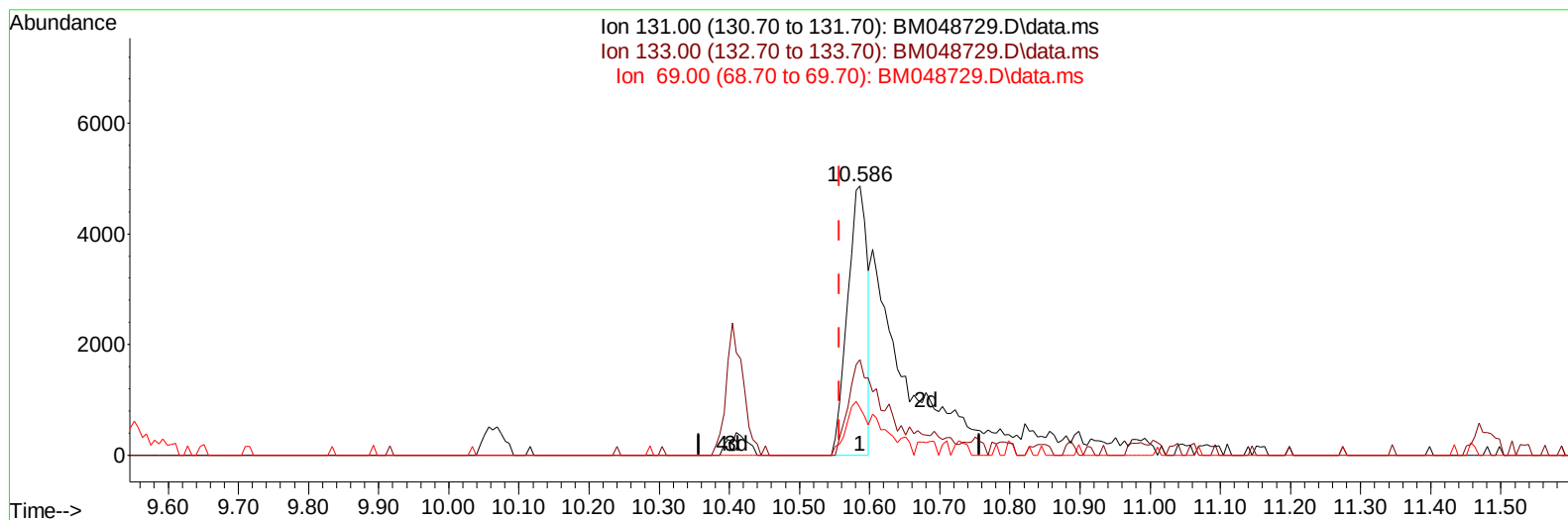
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Data File : BM048729.D
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Operator : RC/JU
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ALS Vial : 22 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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TIC: BM048729.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.586min (+ 0.029) 0.89 ng/u1

response 9470

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	31.70	35.39
69.00	15.00	17.50
0.00	0.00	0.00

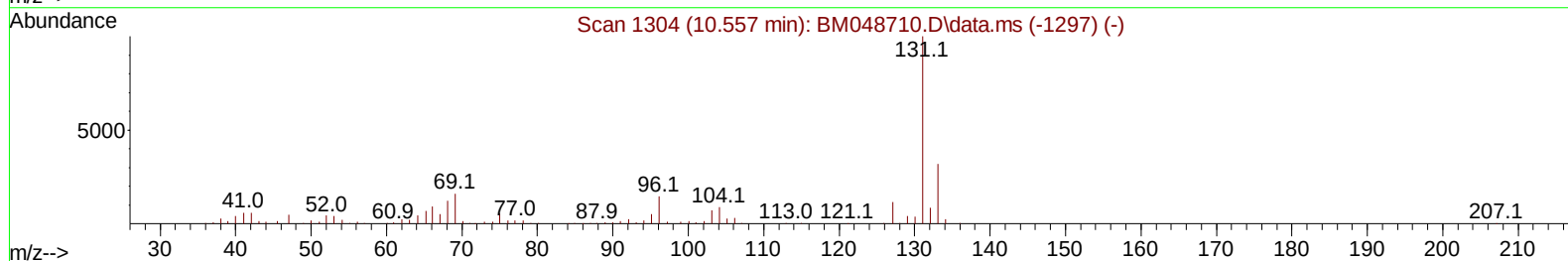
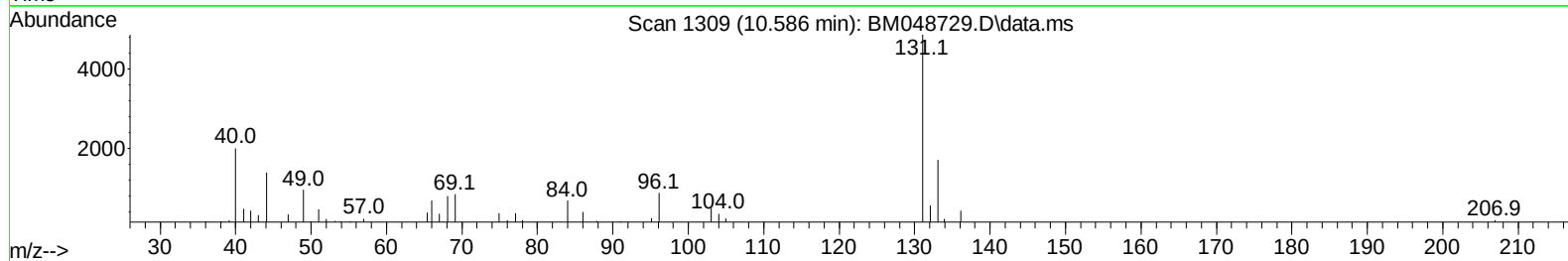
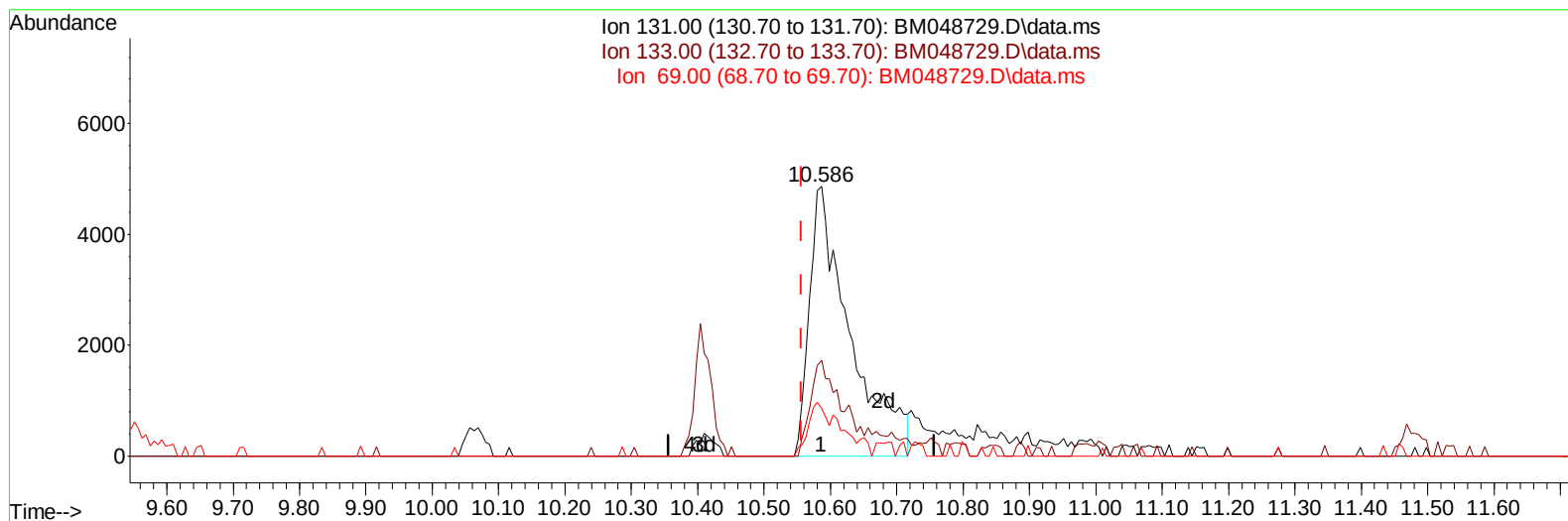
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111524\
Data File : BM048729.D
Acq On : 16 Nov 2024 05:02
Operator : RC/JU
Sample : P4751-21
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_M
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Manual IntegrationsAPPROVED

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TIC: BM048729.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.586min (+ 0.029) 1.94 ng/ul m

response 20522

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	31.70	35.39
69.00	15.00	17.50
0.00	0.00	0.00

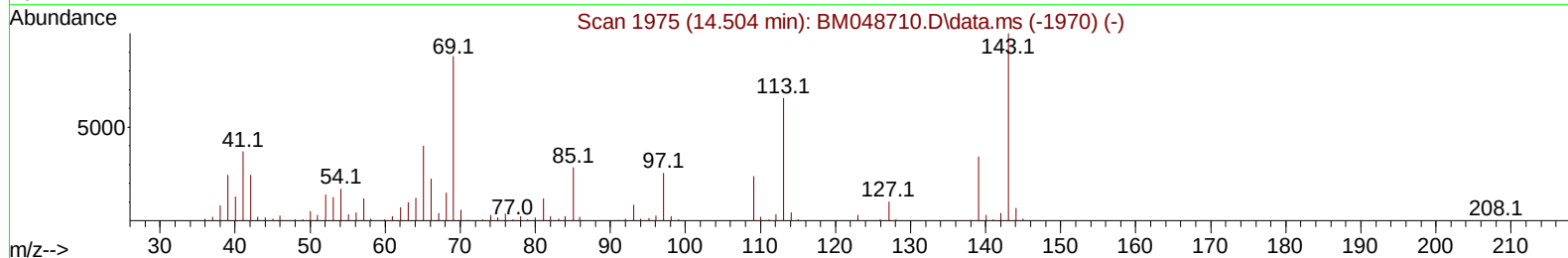
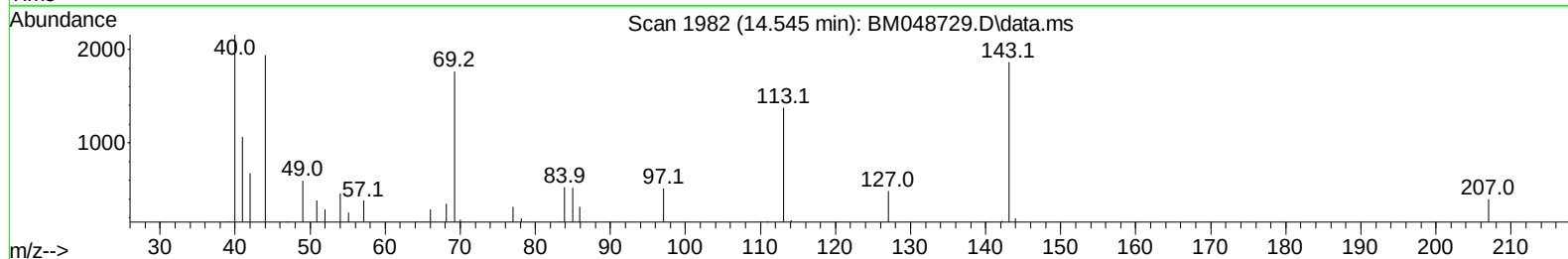
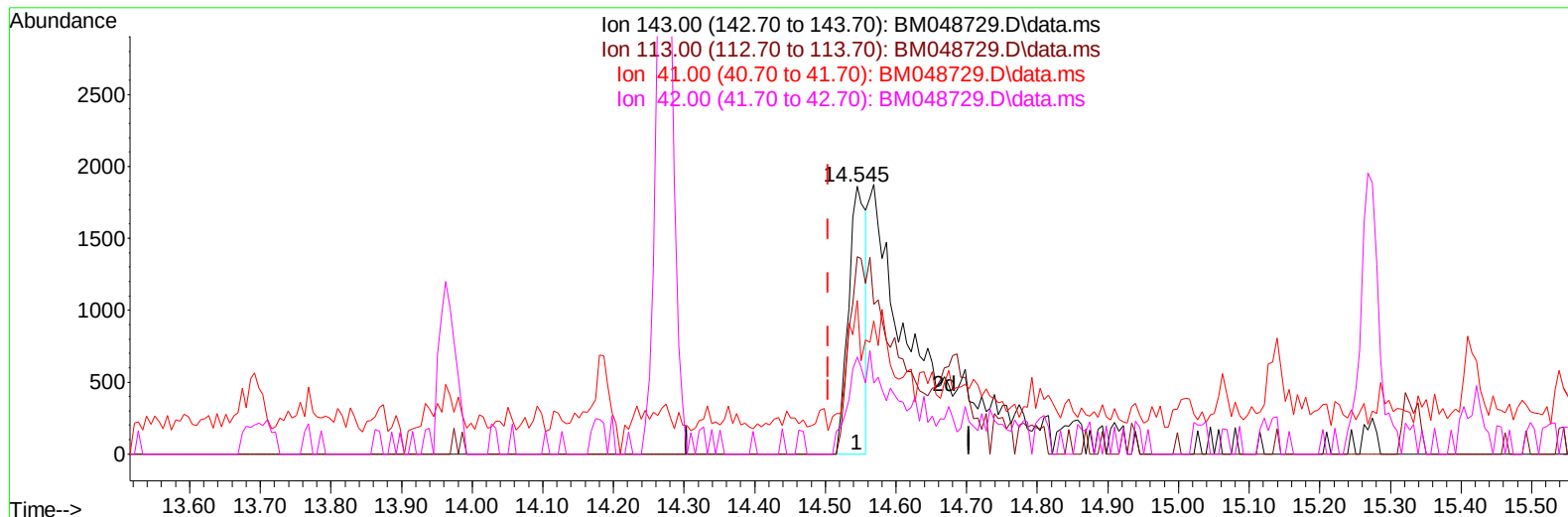
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111524\
Data File : BM048729.D
Acq On : 16 Nov 2024 05:02
Operator : RC/JU
Sample : P4751-21
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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TIC: BM048729.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.545min (+ 0.041) 0.87 ng/ul

response 3147

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	66.50	73.71
41.00	36.20	57.19#
42.00	24.50	36.21#

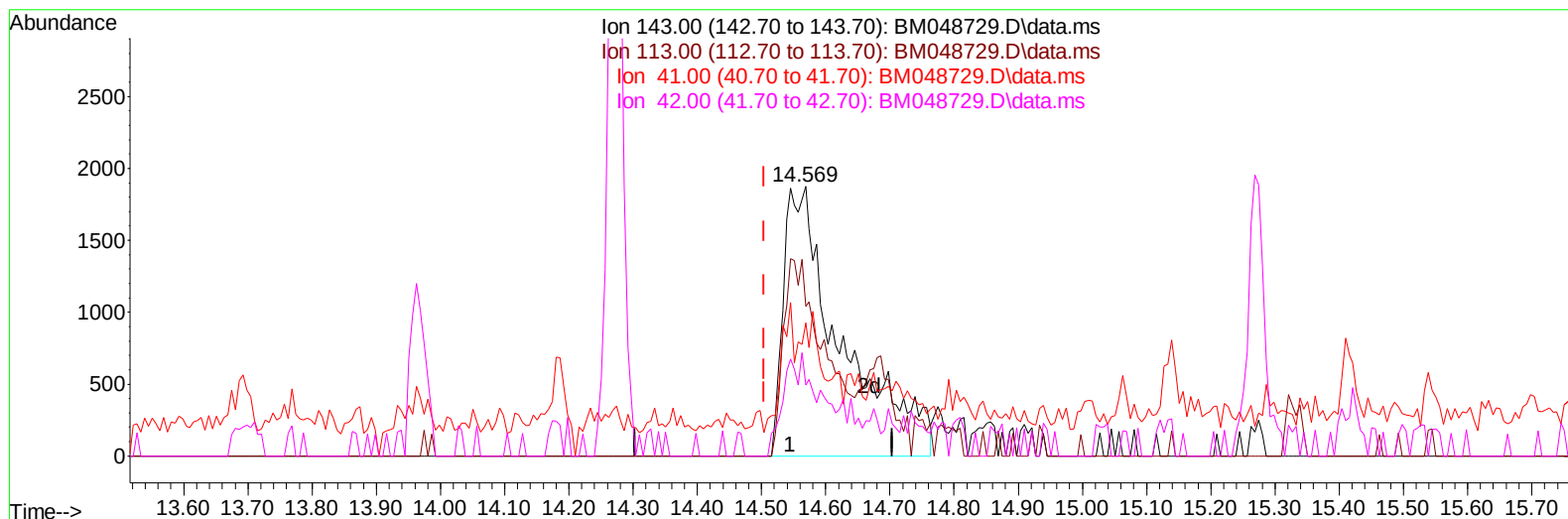
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111524\
Data File : BM048729.D
Acq On : 16 Nov 2024 05:02
Operator : RC/JU
Sample : P4751-21
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BH9G2

Manual IntegrationsAPPROVED

Quant Time: Nov 16 05:36:22 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM110724.MA.M
Quant Title : SVOA CALIBRATION
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Reviewed By :Yogesh Patel 11/18/2024
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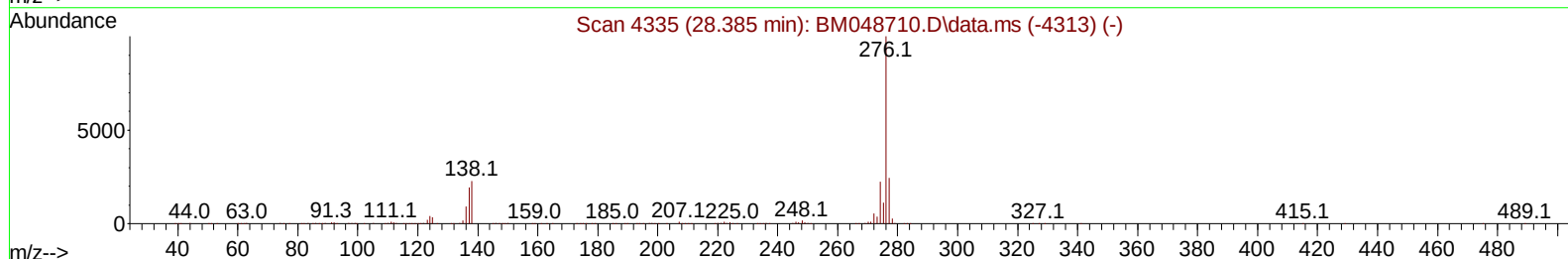
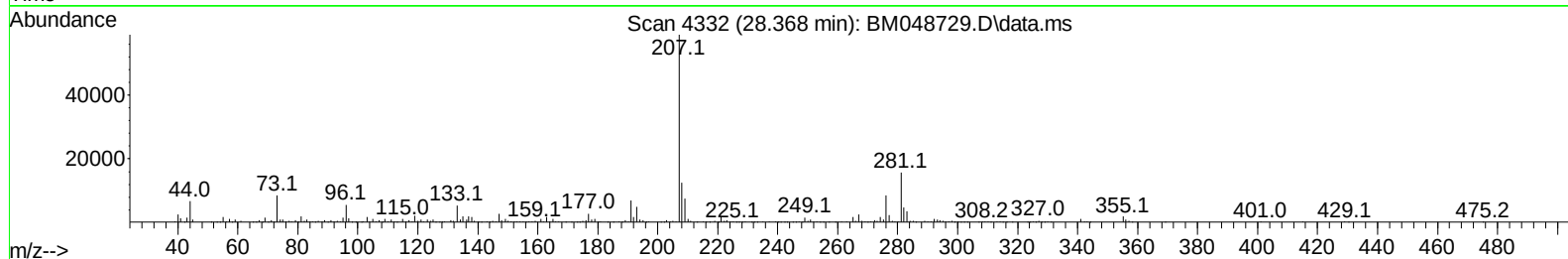
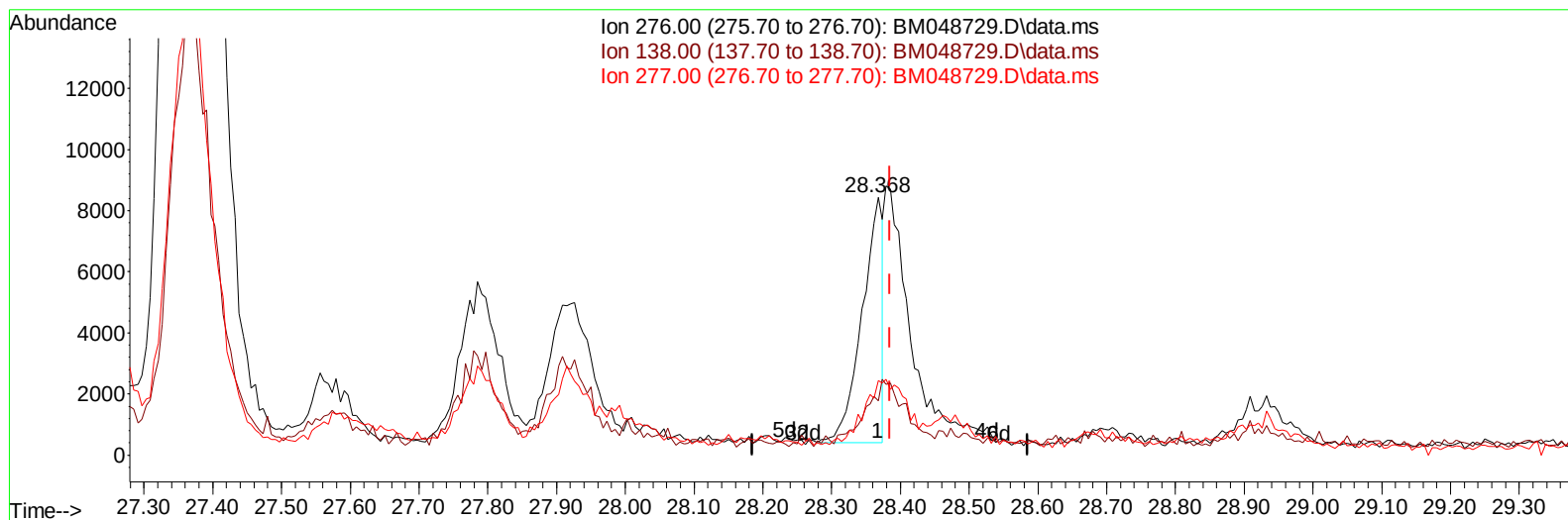
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Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
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Quant Time: Nov 16 05:37:40 2024
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Quant Title : SVOA CALIBRATION
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TIC: BM048729.D\data.ms

(96) Benzo(g,h,i)perylene

28.368min (-0.018) 0.43 ng/u1

response 16925

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	22.00	21.18
277.00	23.30	28.85#
0.00	0.00	0.00

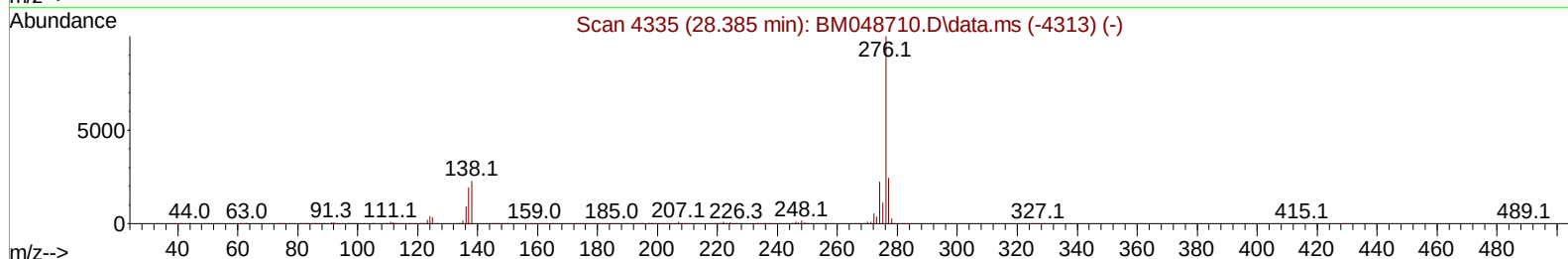
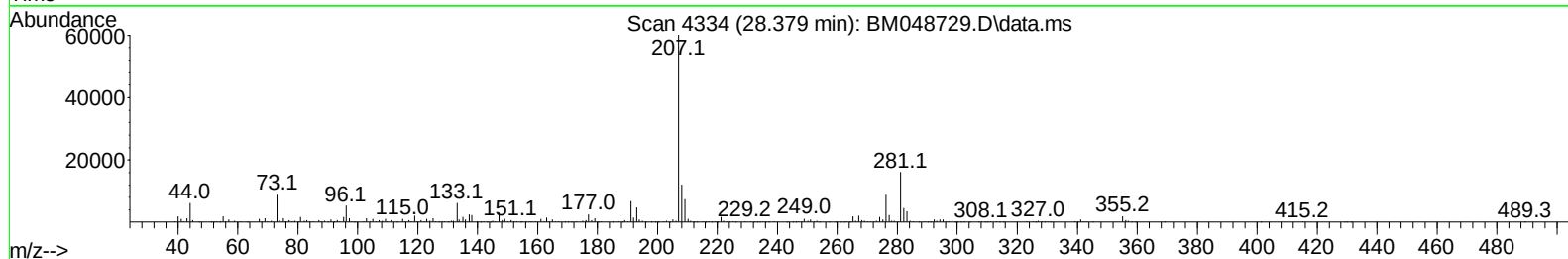
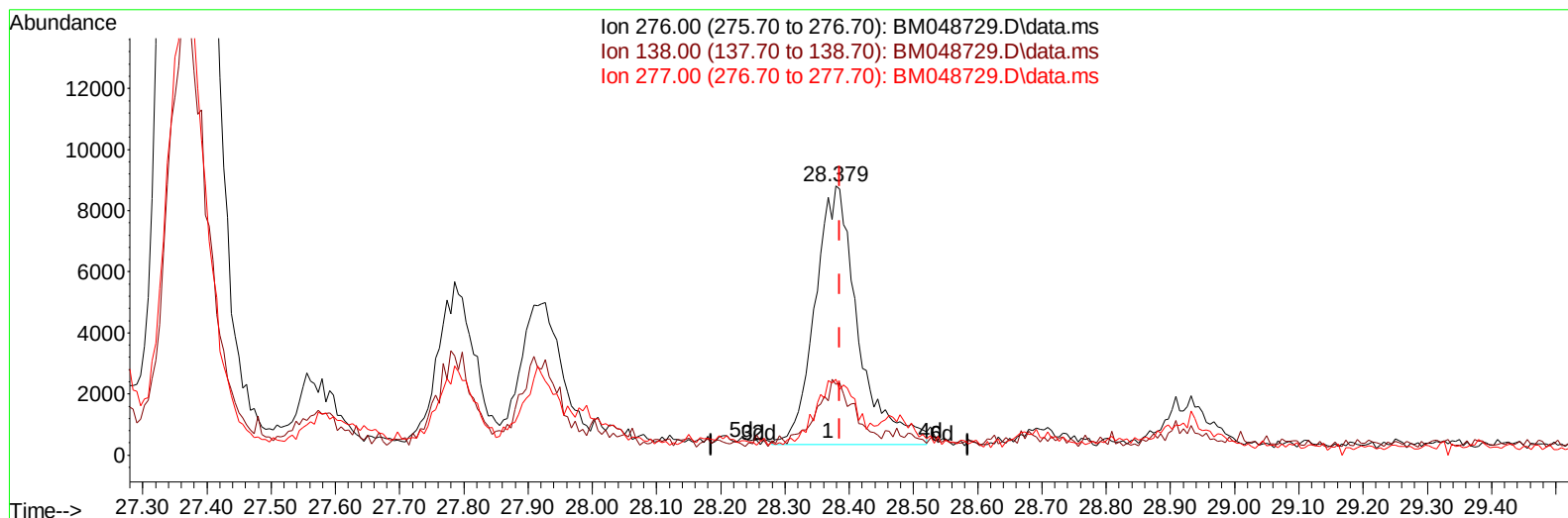
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111524\
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Operator : RC/JU
Sample : P4751-21
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
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Manual IntegrationsAPPROVED

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TIC: BM048729.D\data.ms

(96) Benzo(g,h,i)perylene

28.379min (-0.006) 1.01 ng/ul m

response 39543

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	22.00	26.42#
277.00	23.30	28.25#
0.00	0.00	0.00

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 Operator : RC/JU
 Sample : P4751-21
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BH9G2

Manual IntegrationsAPPROVED

Quant Time: Nov 16 05:38:42 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM110724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 15 16:55:33 2024
 Response via : Initial Calibration

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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.622	152		131645	20.000	ng/ul	0.00
20) Naphthalene-d8	10.404	136		520476	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.274	164		305329	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.021	188		612422	20.000	ng/ul	0.00
79) Chrysene-d12	21.245	240		601183	20.000	ng/ul	0.00
88) Perylene-d12	24.133	264		743349	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.087	96		4256	1.289	ng/uL	0.00
4) Pyridine-d5	3.522	84		16049m	1.763	ng/ul	0.02
7) Phenol-d5	6.810	99		59380	5.925	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.963	67		39119	6.499	ng/ul	0.00
11) 2-Chlorophenol-d4	7.157	132		54716	6.591	ng/ul	0.00
15) 4-Methylphenol-d8	8.351	113		29929	3.798	ng/ul	0.01
21) Nitrobenzene-d5	8.787	128		22328	5.919	ng/ul	0.00
24) 2-Nitrophenol-d4	9.504	143		25169	6.273	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.063	165		45511	6.106	ng/ul	0.02
31) 4-Chloroaniline-d4	10.586	131		20522m	1.938	ng/ul	0.03
46) Dimethylphthalate-d6	13.692	166		154284	7.916	ng/ul	0.00
49) Acenaphthylene-d8	13.963	160		171140	7.447	ng/ul	0.00
54) 4-Nitrophenol-d4	14.569	143		11709m	3.234	ng/ul	0.06
60) Fluorene-d10	15.274	176		137199	8.064	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.416	200		9963	3.234	ng/ul	0.00
73) Anthracene-d10	17.121	188		199800	7.710	ng/ul	0.00
81) Pyrene-d10	19.415	212		204564	6.780	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.933	264		158354	4.559	ng/ul	0.00
Target Compounds							
						Qvalue	
8) Phenol	6.840	94		12604	1.199	ng/ul	96
16) Acetophenone	8.451	105		68258	5.557	ng/ul	99
52) Acenaphthene	14.339	153		25186	1.412	ng/ul	96
61) Fluorene	15.327	166		28139	1.445	ng/ul	94
72) Phenanthrene	17.062	178		543159	17.883	ng/ul	99
74) Anthracene	17.157	178		111510	3.658	ng/ul	99
77) Carbazole	17.433	167		51336	1.870	ng/ul	99
80) Fluoranthene	19.086	202		1192244	34.057	ng/ul	99
82) Pyrene	19.445	202		764573	20.142	ng/ul	99
85) Benzo(a)anthracene	21.227	228		590566	15.583	ng/ul	99
87) Chrysene	21.286	228		572678	15.725	ng/ul	97
90) Benzo(b)fluoranthene	23.227	252		789671	19.298	ng/ul	98
91) Benzo(k)fluoranthene	23.280	252		286114	7.013	ng/ul	97
93) Benzo(a)pyrene	23.991	252		263418	6.847	ng/ul	97
94) Indeno(1,2,3-cd)pyrene	27.362	276		253140	5.063	ng/ul	96
95) Dibenzo(a,h)anthracene	27.397	278		99984	2.448	ng/ul#	89
96) Benzo(g,h,i)perylene	28.379	276		39543m	1.009	ng/ul	

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

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