

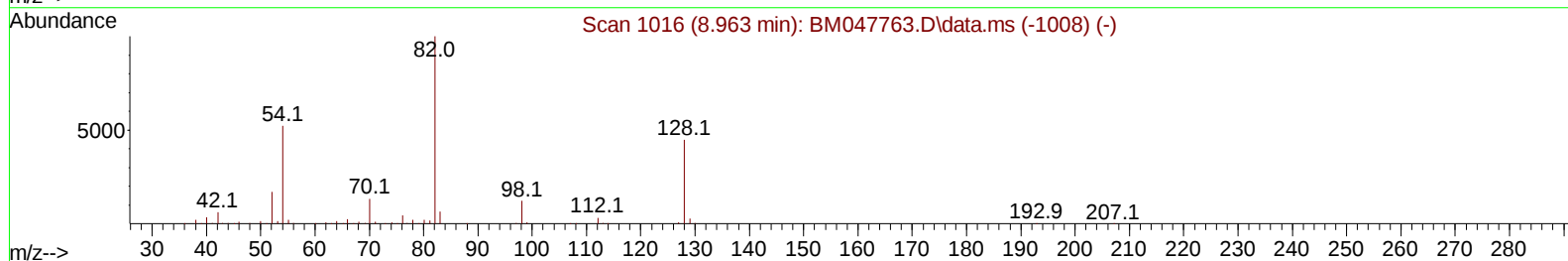
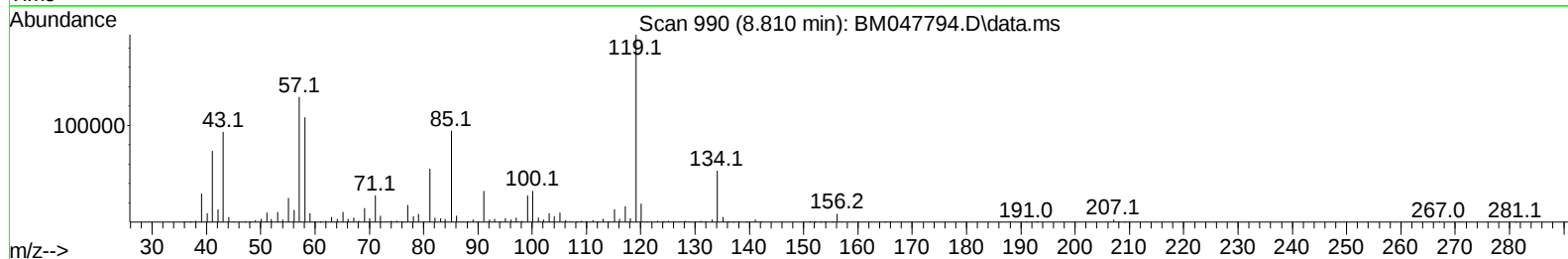
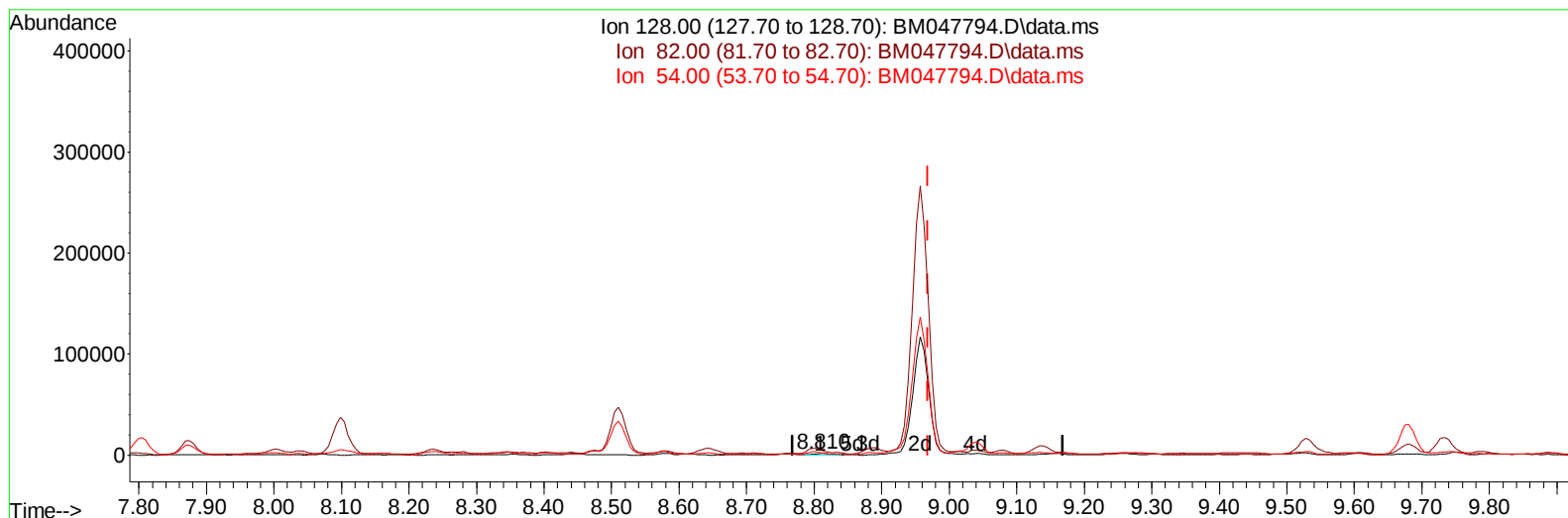
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100224\
Data File : BM047794.D
Acq On : 03 Oct 2024 03:08
Operator : RC/JU
Sample : P4020-03
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DDCD9

Manual IntegrationsAPPROVED

Quant Time: Oct 03 03:41:19 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Sep 28 02:30:36 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 10/03/2024
Supervised By :mohammad ahmed 10/04/2024



TIC: BM047794.D\data.ms

(21) Nitrobenzene-d5 (S)

8.810min (-0.159) 0.26 ng/u1

response 2461

Ion	Exp%	Act%
128.00	100.00	100.00
82.00	266.90	307.20
54.00	152.80	170.52
0.00	0.00	0.00

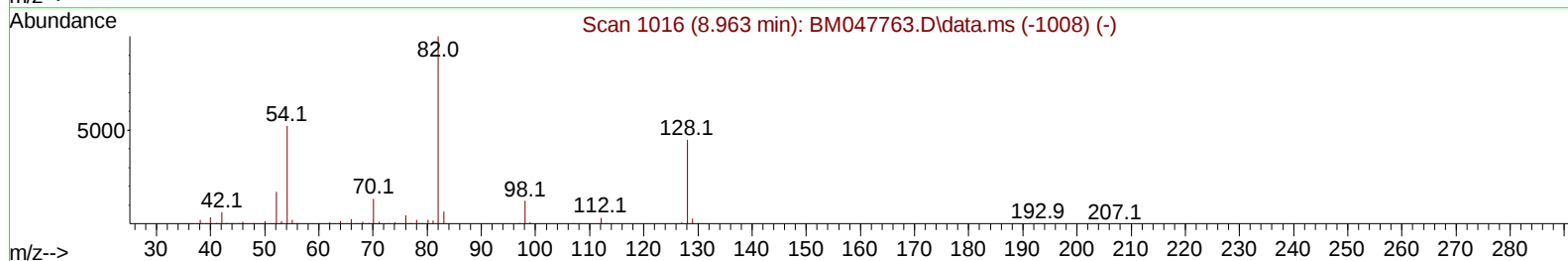
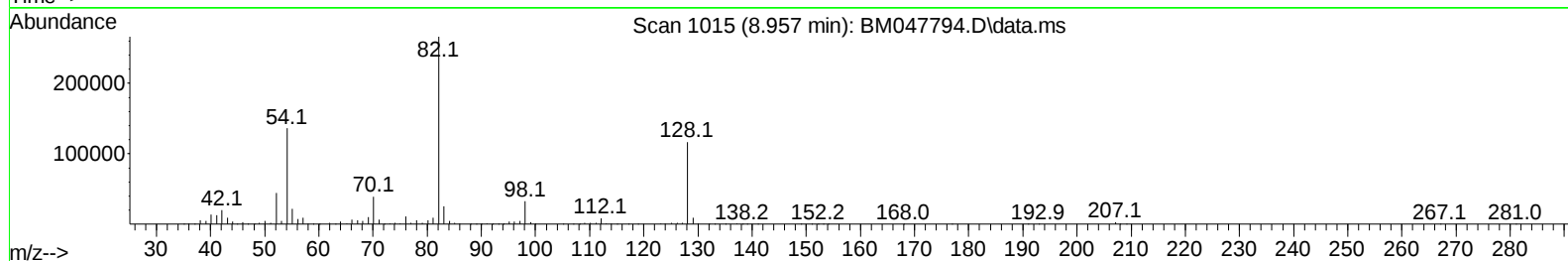
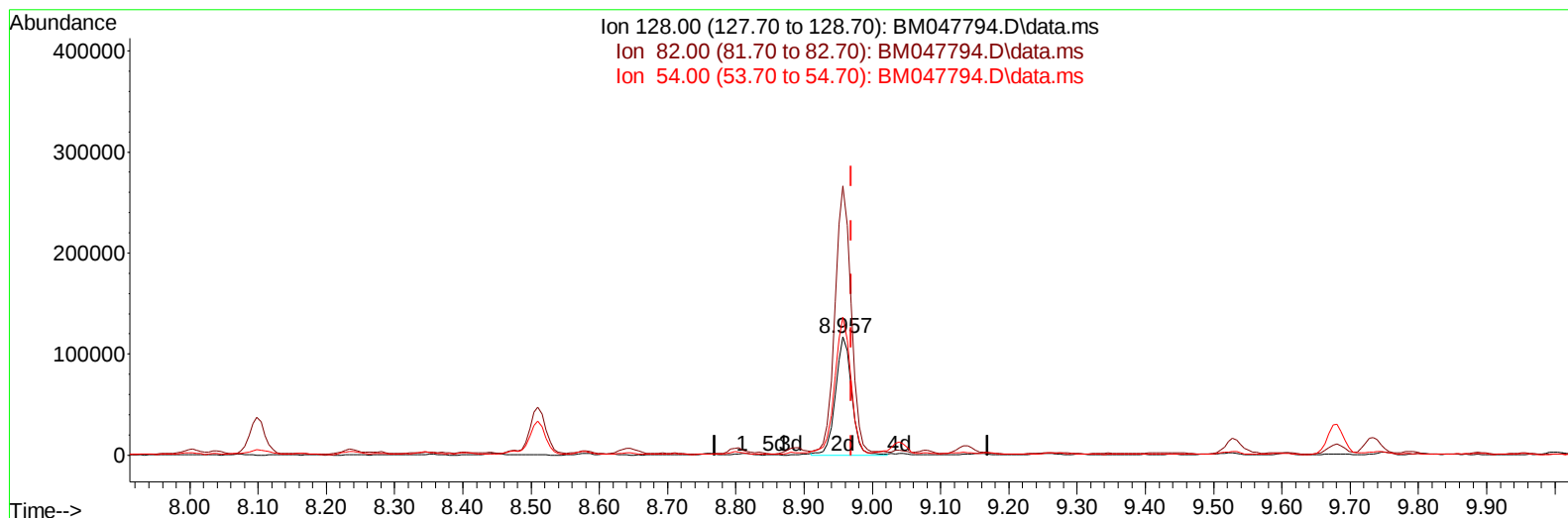
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(21) Nitrobenzene-d5 (S)

8.957min (-0.012) 20.43 ng/ul m

response 191837

Ion	Exp%	Act%
128.00	100.00	100.00
82.00	266.90	227.88
54.00	152.80	116.68#
0.00	0.00	0.00

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 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 10/03/2024
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Quant Time: Oct 03 03:42:19 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Sep 28 02:30:36 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.804	152	325614	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.598	136	1179428	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.439	164	908789	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.186	188	1728022	20.000	ng/ul	0.00
79) Chrysene-d12	21.403	240	1683871	20.000	ng/ul	0.00
88) Perylene-d12	24.403	264	1926889	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.252	96	26270	2.607	ng/uL	0.00
4) Pyridine-d5	3.669	84	329252	11.160	ng/ul	0.00
7) Phenol-d5	6.969	99	490435	15.928	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.134	67	288438	13.202	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.340	132	403893	18.030	ng/ul	0.00
15) 4-Methylphenol-d8	8.510	113	387007	16.906	ng/ul	0.00
21) Nitrobenzene-d5	8.957	128	191837m	20.428	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.681	143	205549	22.841	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.222	165	391008	21.751	ng/ul	0.00
31) 4-Chloroaniline-d4	10.734	131	366048	13.171	ng/ul	0.00
46) Dimethylphthalate-d6	13.851	166	1214058	18.426	ng/ul	0.00
49) Acenaphthylene-d8	14.133	160	1412101	18.076	ng/ul	0.00
54) 4-Nitrophenol-d4	14.633	143	187525	14.295	ng/ul	0.00
60) Fluorene-d10	15.433	176	1057554	18.531	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.539	200	128560	12.713	ng/ul	-0.01
73) Anthracene-d10	17.280	188	1536203	18.080	ng/ul	0.00
81) Pyrene-d10	19.562	212	1913256	20.020	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.197	264	1882642	18.831	ng/ul	-0.01
Target Compounds						
					Qvalue	
30) Naphthalene	10.645	128	135974	2.004	ng/ul	98
36) 2-Methylnaphthalene	12.263	142	251483	5.834	ng/ul	99
37) 1-Methylnaphthalene	12.480	142	159048	3.627	ng/ul	99
58) 2,3,4,6-Tetrachlorophenol	15.069	232	2021069	136.982	ng/ul#	93
71) Pentachlorophenol	16.880	266	17045207	1293.185	ng/ul	99
72) Phenanthrene	17.227	178	313542	3.094	ng/ul	98
80) Fluoranthene	19.227	202	140562	1.249	ng/ul#	90
82) Pyrene	19.592	202	140169	1.119	ng/ul	99

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

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