

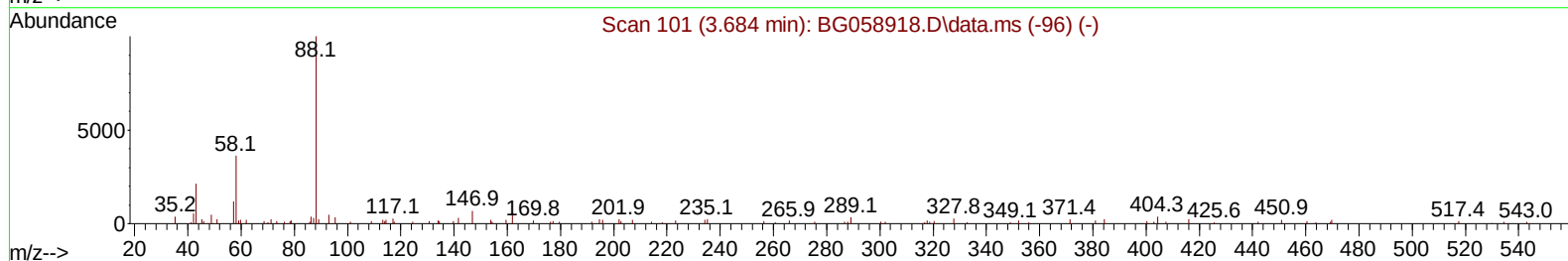
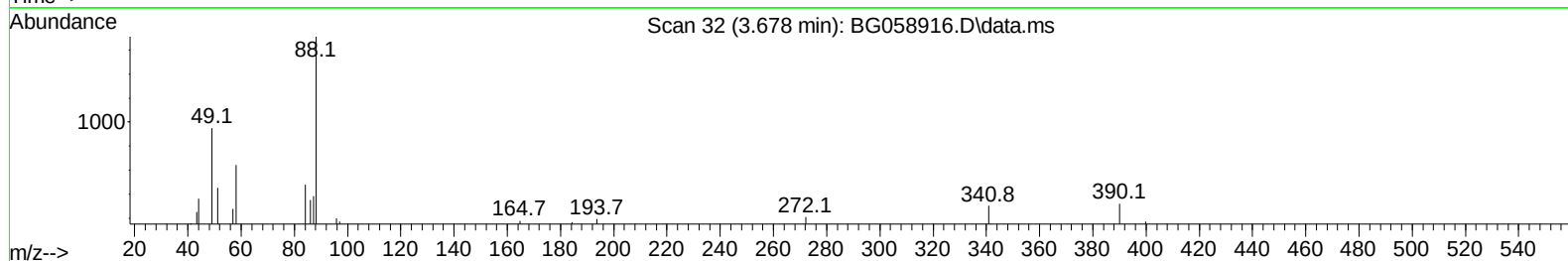
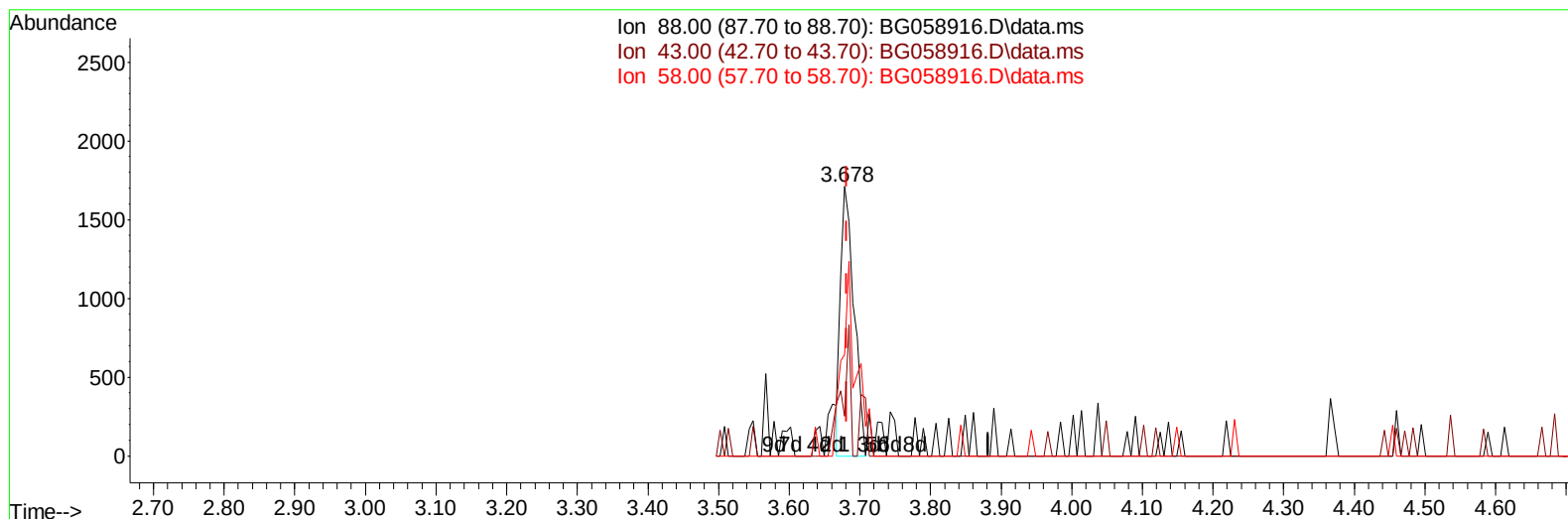
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090823\
 Data File : BG058916.D
 Acq On : 8 Sep 2023 10:41
 Operator : MA/JU
 Sample : SST00501
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SST005449

Manual IntegrationsAPPROVED

Quant Time: Sep 08 23:25:50 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090823.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Sep 08 22:56:23 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 09/09/2023
 Supervised By :mohammad ahmed 09/10/2023



TIC: BG058916.D\data.ms

(2) 1,4-Dioxane

3.678min (-0.003) 1.84 ng/uL

response 2262

Ion	Exp%	Act%
88.00	100.00	100.00
43.00	24.80	14.67#
58.00	62.40	37.64#
0.00	0.00	0.00

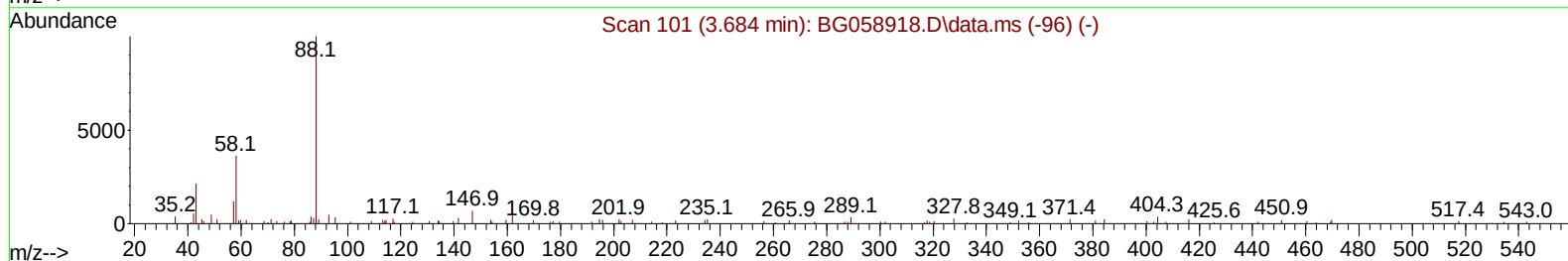
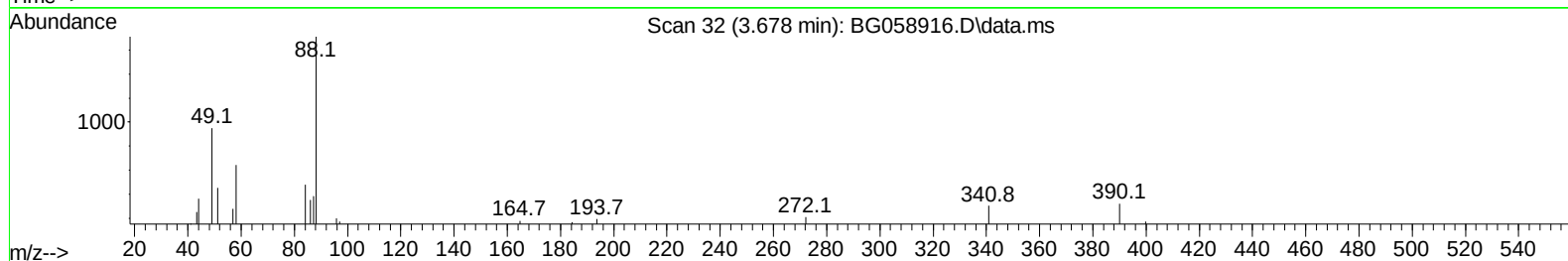
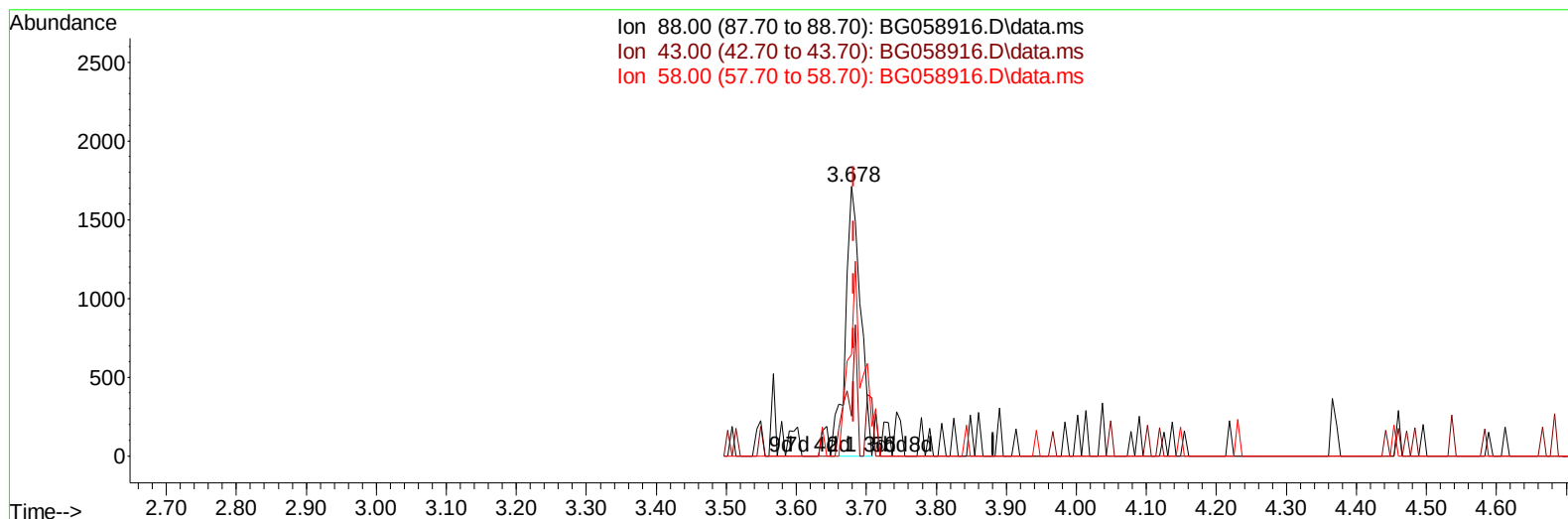
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TIC: BG058916.D\data.ms

(2) 1,4-Dioxane

3.678min (-0.003) 2.10 ng/uL m

response 2583

Ion	Exp%	Act%
88.00	100.00	100.00
43.00	24.80	14.67#
58.00	62.40	37.64#
0.00	0.00	0.00

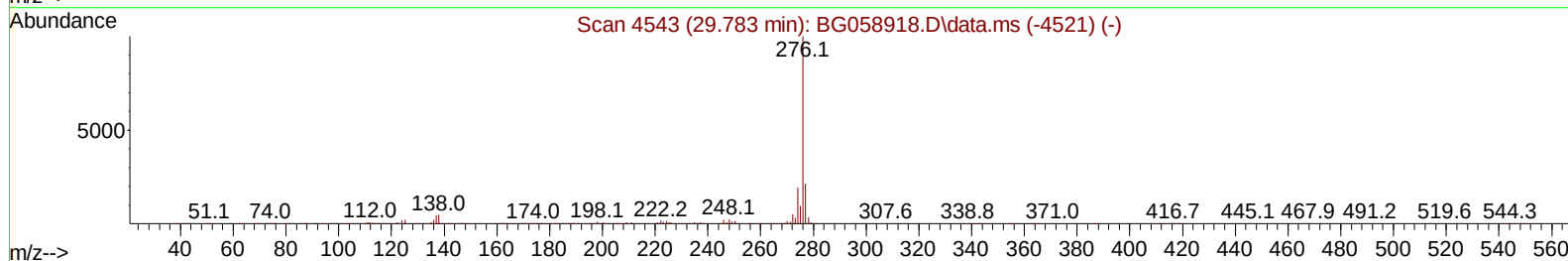
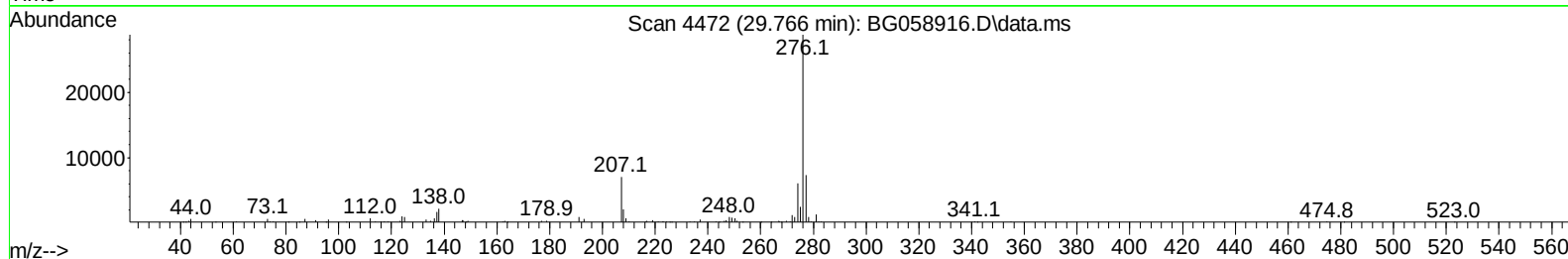
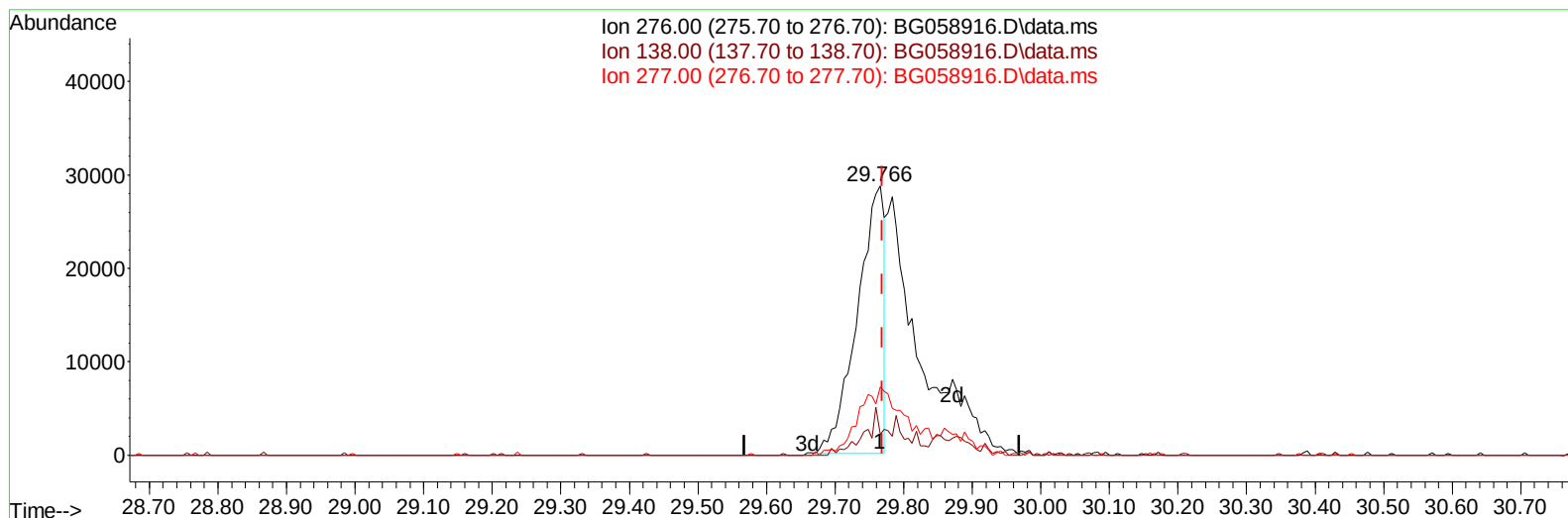
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090823\
 Data File : BG058916.D
 Acq On : 8 Sep 2023 10:41
 Operator : MA/JU
 Sample : SST00501
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
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Quant Time: Sep 08 23:35:20 2023
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TIC: BG058916.D\data.ms

(94) Indeno(1,2,3-cd)pyrene

29.766min (-0.003) 2.10 ng/u1

response 78363

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	7.70	7.62
277.00	22.20	25.34
0.00	0.00	0.00

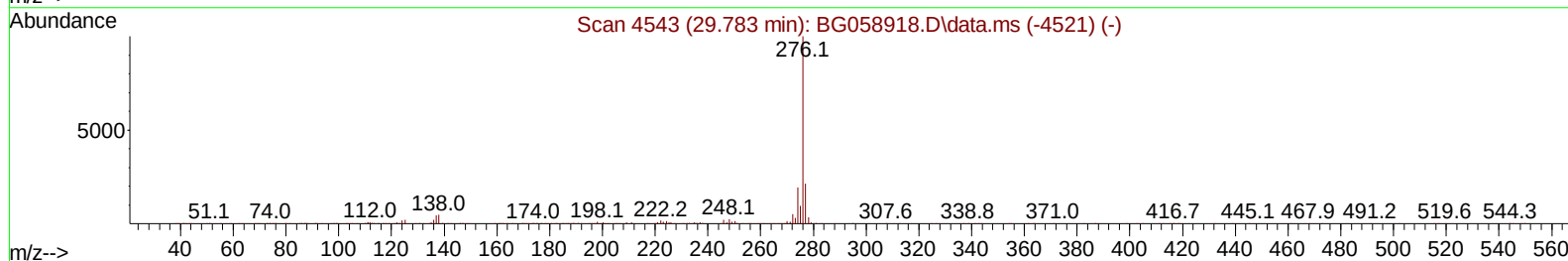
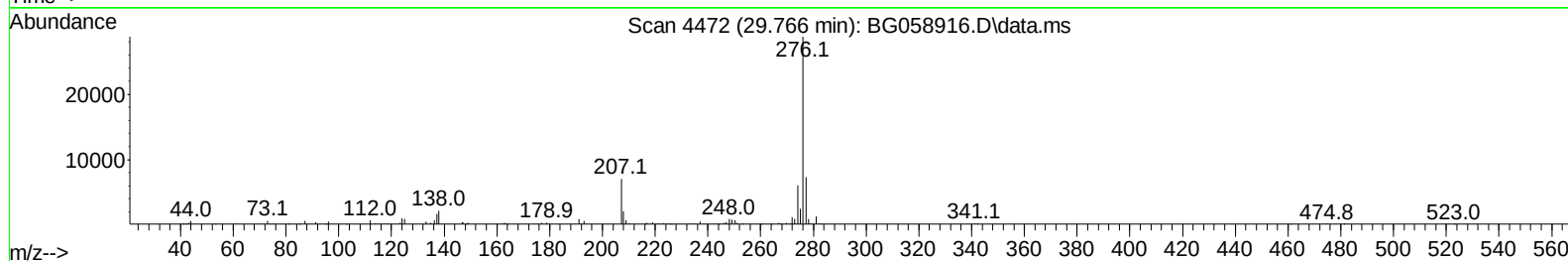
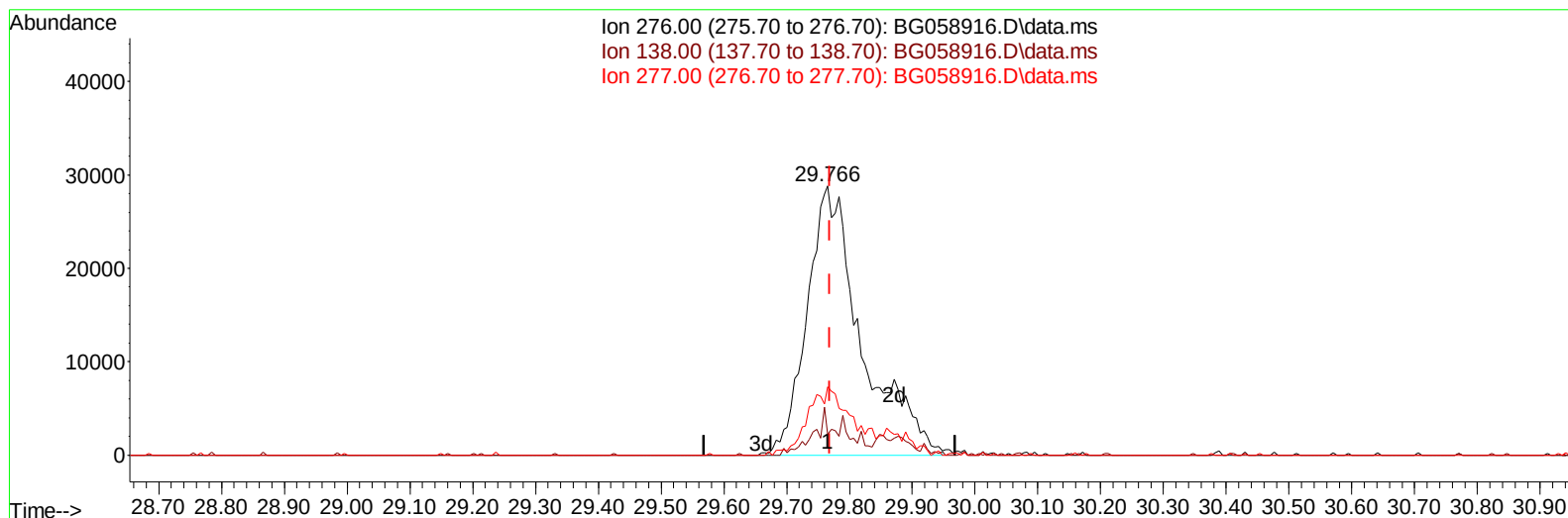
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090823\
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Instrument :
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TIC: BG058916.D\data.ms

(94) Indeno(1,2,3-cd)pyrene

29.766min (-0.003) 4.65 ng/ul m

response 173078

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	7.70	7.62
277.00	22.20	25.34
0.00	0.00	0.00

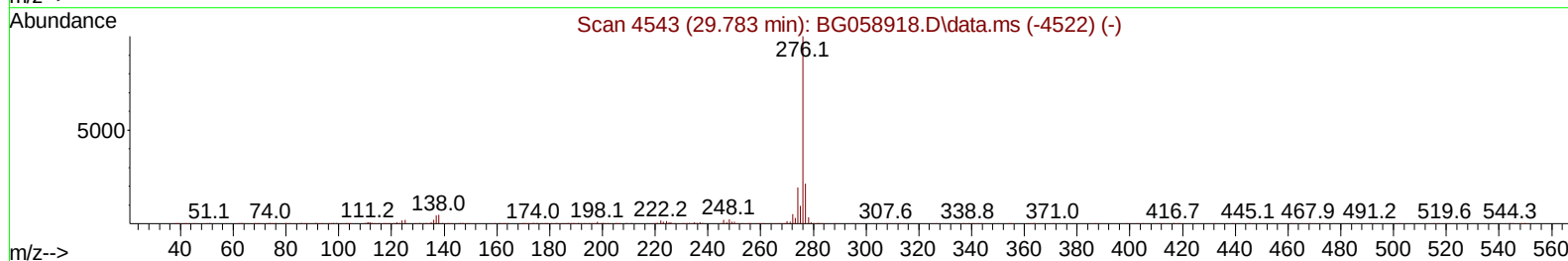
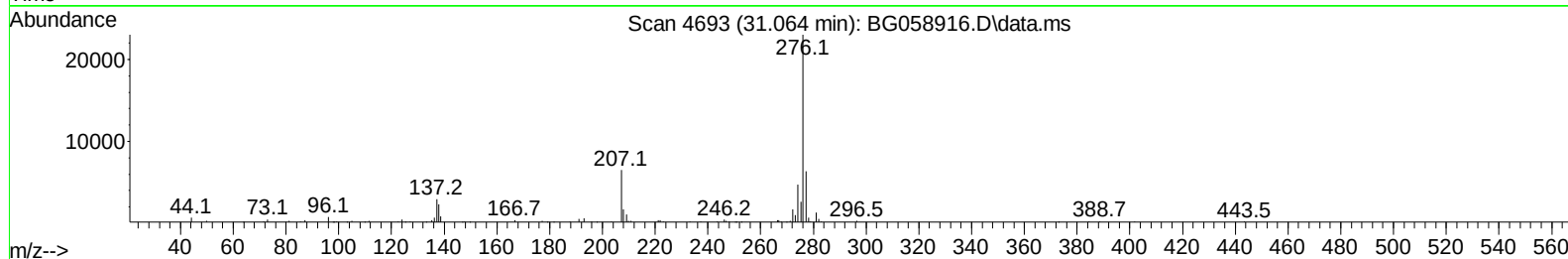
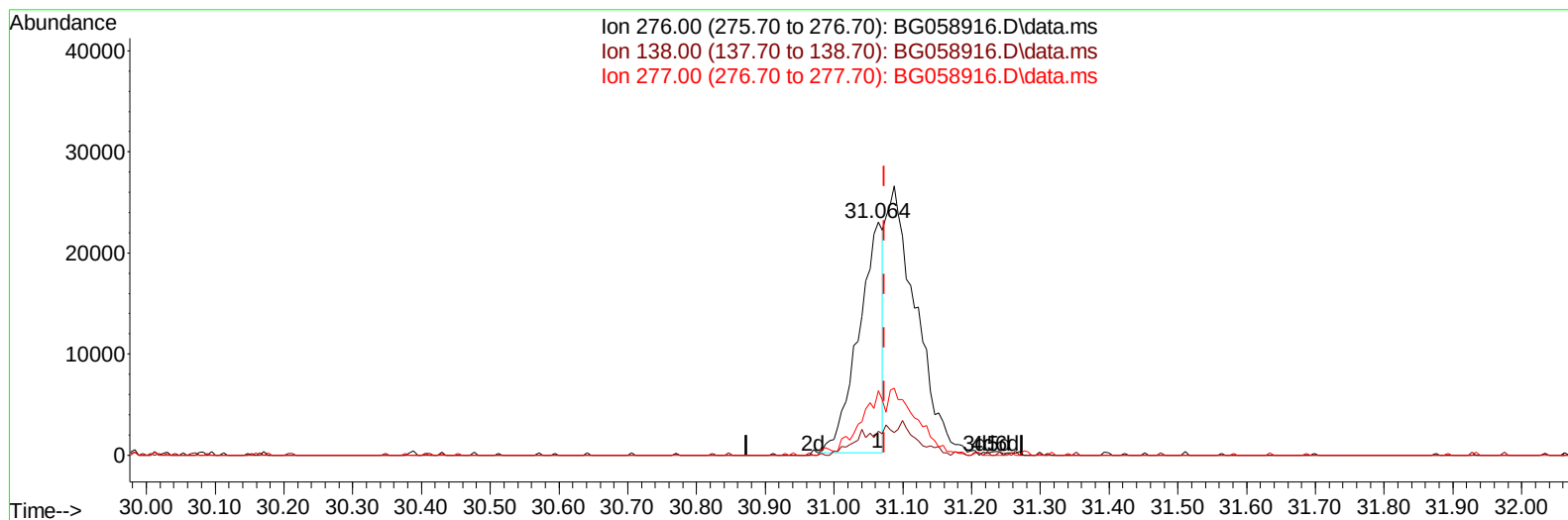
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Data File : BG058916.D
Acq On : 8 Sep 2023 10:41
Operator : MA/JU
Sample : SSTD00501
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SSTD005449

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 09/09/2023
Supervised By :mohammad ahmed 09/10/2023

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QLast Update : Fri Sep 08 22:56:23 2023
Response via : Initial Calibration



TIC: BG058916.D\data.ms

(96) Benzo(g,h,i)perylene**31.064min (-0.009) 1.89 ng/u1****response 55950**

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	10.90	13.89#
277.00	24.60	27.73
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090823\
 Data File : BG058916.D
 Acq On : 8 Sep 2023 10:41
 Operator : MA/JU
 Sample : SST00501
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :

BNA_G

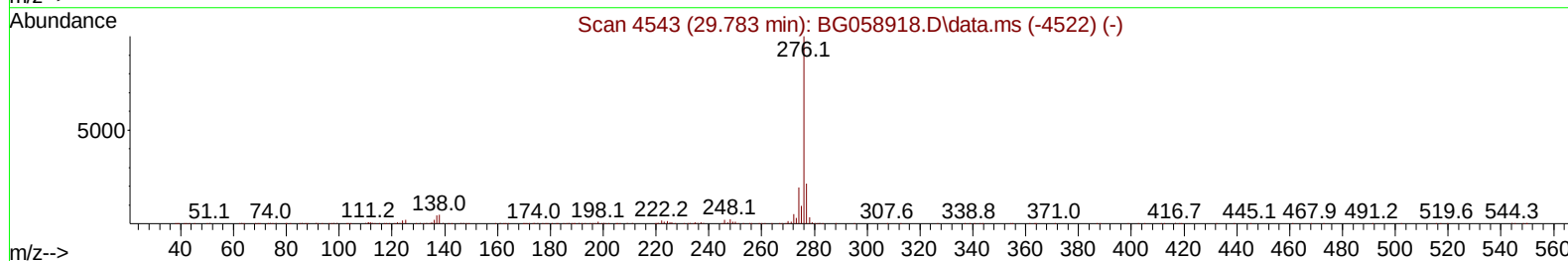
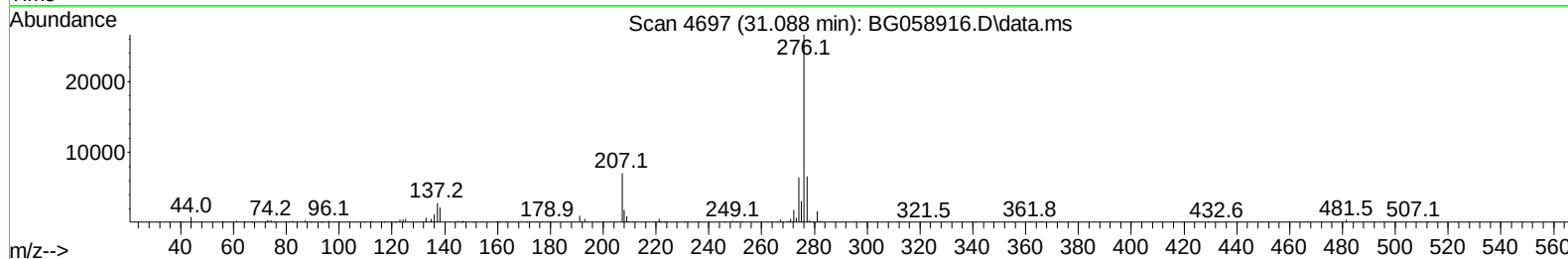
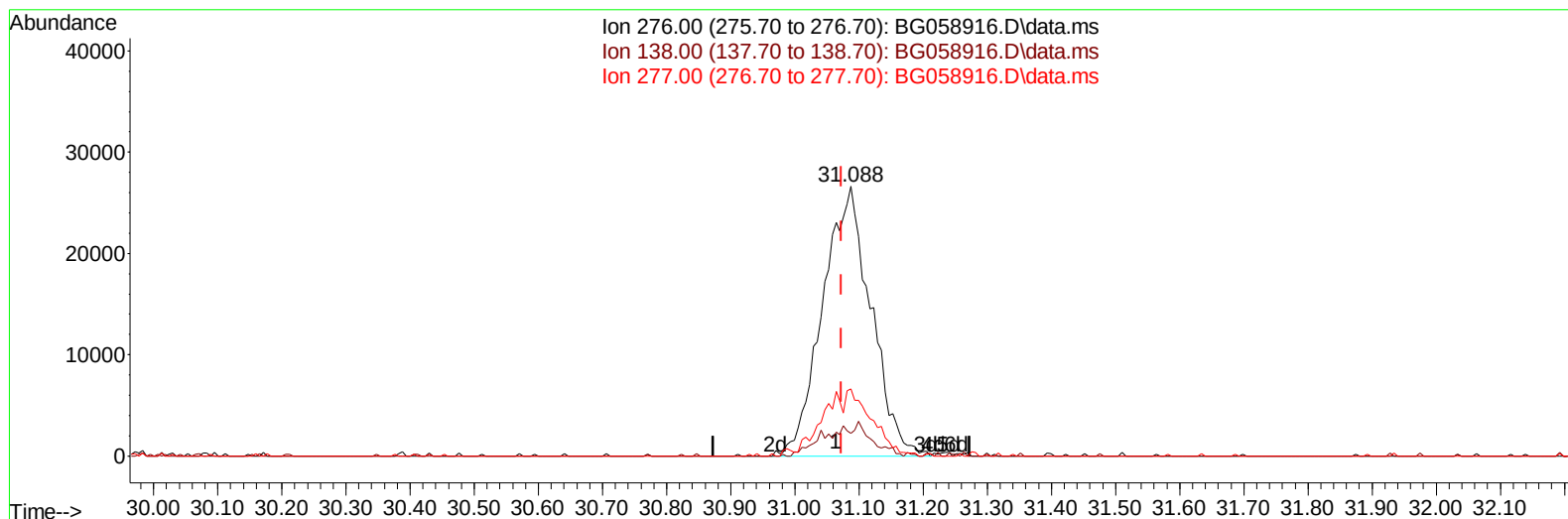
ClientSampleId :

SSTD005449

Manual IntegrationsAPPROVED

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TIC: BG058916.D\data.ms

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

31.088min (+ 0.014) 4.70 ng/ul m

response 139101

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	10.90	8.37#
277.00	24.60	24.78
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

Quant Time: Sep 08 23:36:58 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.332	152	47067	20.000	ng/ul	0.00
20) Naphthalene-d8	11.181	136	206678	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.965	164	188913	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.715	188	544441	20.000	ng/ul	0.00
79) Chrysene-d12	22.045	240	465308	20.000	ng/ul	0.00
88) Perylene-d12	25.623	264	529750	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.637	96	2148	2.197	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	0.000	99	0d	0.000	ng/ul	
9) Bis-(2-Chloroethyl)eth...	0.000	67	0d	0.000	ng/ul	
11) 2-Chlorophenol-d4	7.844	132	11737	4.215	ng/ul	0.00
15) 4-Methylphenol-d8	0.000	113	0d	0.000	ng/ul	
21) Nitrobenzene-d5	9.530	128	4339	3.564	ng/ul	0.00
24) 2-Nitrophenol-d4	10.253	143	4694	3.520	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.770	165	15915	4.113	ng/ul	0.00
31) 4-Chloroaniline-d4	0.000	131	0d	0.000	ng/ul	
46) Dimethylphthalate-d6	14.360	166	70203	4.814	ng/ul	0.00
49) Acenaphthylene-d8	14.660	160	73901	4.688	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
60) Fluorene-d10	15.946	176	65889	4.794	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	0.000	200	0d	0.000	ng/ul	
73) Anthracene-d10	17.809	188	122039	4.927	ng/ul	0.00
81) Pyrene-d10	20.083	212	129737	4.808	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.371	264	123245	4.723	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.678	88	2583m	2.096	ng/uL	
12) 2-Chlorophenol	7.885	128	12591	4.485	ng/ul#	81
17) N-Nitroso-di-n-propyla...	9.143	70	14287	4.234	ng/ul#	86
19) Hexachloroethane	9.419	117	5333	4.654	ng/ul#	79
22) Nitrobenzene	9.572	77	15257	3.664	ng/ul#	85
23) Isophorone	10.083	82	47451	4.568	ng/ul#	92
25) 2-Nitrophenol	10.277	139	4751	3.303	ng/ul#	92
26) 2,4-Dimethylphenol	10.300	107	21039	4.704	ng/ul	87
27) Bis(2-Chloroethoxy)met...	10.570	93	25899	4.814	ng/ul#	81
29) 2,4-Dichlorophenol	10.800	162	14577	3.975	ng/ul	96
30) Naphthalene	11.234	128	51673	4.628	ng/ul#	95
33) Hexachlorobutadiene	11.469	225	16566	4.612	ng/ul	89
35) 4-Chloro-3-methylphenol	12.409	107	17800	4.131	ng/ul#	84
36) 2-Methylnaphthalene	12.809	142	40430	4.783	ng/ul	95
37) 1-Methylnaphthalene	13.026	142	41767	4.831	ng/ul	88
39) 1,2,4,5-Tetrachloroben...	13.150	216	33551	4.760	ng/ul#	95
41) 2,4,6-Trichlorophenol	13.391	196	17599	4.296	ng/ul	94
42) 2,4,5-Trichlorophenol	13.455	196	19007	4.320	ng/ul	85
43) 1,1'-Biphenyl	13.802	154	64620	5.027	ng/ul#	97
44) 2-Chloronaphthalene	13.855	162	50654	4.941	ng/ul	97
45) 2-Nitroaniline	14.060	65	6828	3.012	ng/ul	98
47) Dimethylphthalate	14.407	163	70834	4.777	ng/ul#	93
48) 2,6-Dinitrotoluene	14.548	165	7720	3.668	ng/ul#	79

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
50) Acenaphthylene	14.695	152	85778	4.877	ng/ul	98
52) Acenaphthene	15.030	153	54993	4.802	ng/ul	97
56) Dibenzofuran	15.359	168	86308	5.048	ng/ul	94
57) 2,4-Dinitrotoluene	15.312	165	11733	3.865	ng/ul#	90
58) 2,3,4,6-Tetrachlorophenol	15.559	232	20079	4.327	ng/ul#	96
59) Diethylphthalate	15.753	149	62404	4.683	ng/ul	98
61) Fluorene	16.005	166	72558	4.999	ng/ul#	88
62) 4-Chlorophenyl-phenyle...	15.993	204	41114	4.722	ng/ul	98
67) N-Nitrosodiphenylamine	16.211	169	62757	4.687	ng/ul	99
68) 4-Bromophenyl-phenylether	16.887	248	32184	4.868	ng/ul	92
69) Hexachlorobenzene	16.975	284	36020	4.722	ng/ul	96
72) Phenanthrene	17.756	178	132256	5.005	ng/ul	95
74) Anthracene	17.844	178	132372	4.905	ng/ul	98
75) 1,2,3,4-Tetrachloroben...	13.755	216	34000	4.301	ng/ul#	83
76) Pentachlorobenzene	15.247	250	35086	4.403	ng/ul	90
78) Di-n-butylphthalate	18.637	149	72377	3.931	ng/ul	97
80) Fluoranthene	19.748	202	164017	5.001	ng/ul	97
82) Pyrene	20.112	202	163697	4.983	ng/ul	98
83) Butylbenzylphthalate	20.982	149	21460	3.768	ng/ul	87
85) Benzo(a)anthracene	22.022	228	151747	4.790	ng/ul	98
86) Bis(2-ethylhexyl)phtha...	21.881	149	34265	3.856	ng/ul	91
87) Chrysene	22.092	228	142668	4.807	ng/ul	98
90) Benzo(b)fluoranthene	24.460	252	157111	4.925	ng/ul	99
91) Benzo(k)fluoranthene	24.542	252	154240	4.805	ng/ul	99
93) Benzo(a)pyrene	25.447	252	146758	4.916	ng/ul#	97
94) Indeno(1,2,3-cd)pyrene	29.766	276	173078m	4.649	ng/ul	
95) Dibenzo(a,h)anthracene	29.877	278	143958	4.757	ng/ul#	92
96) Benzo(g,h,i)perylene	31.088	276	139101m	4.700	ng/ul	

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

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