

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060524\
 Data File : BN031804.D
 Acq On : 05 Jun 2024 17:45
 Operator : MA/JU
 Sample : P2591-13DL 5X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :

BNA_N

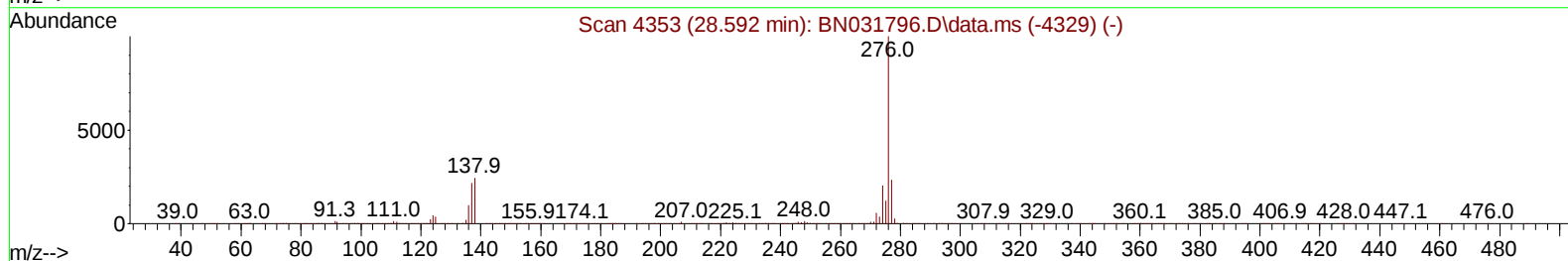
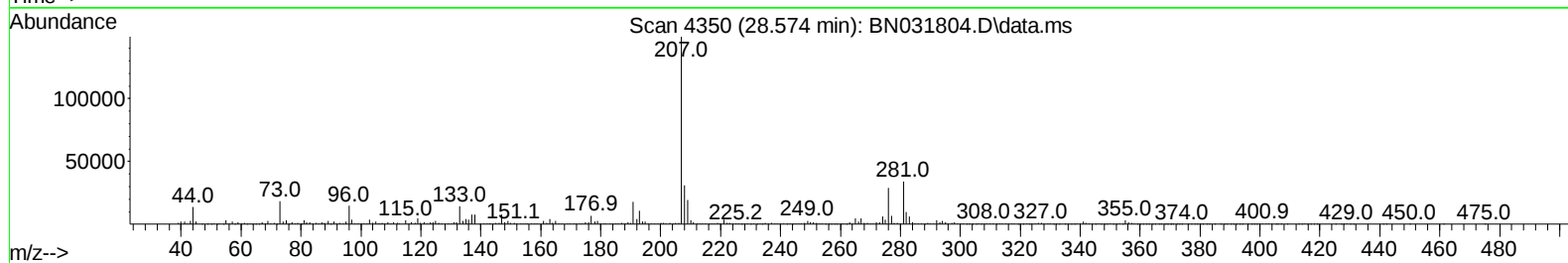
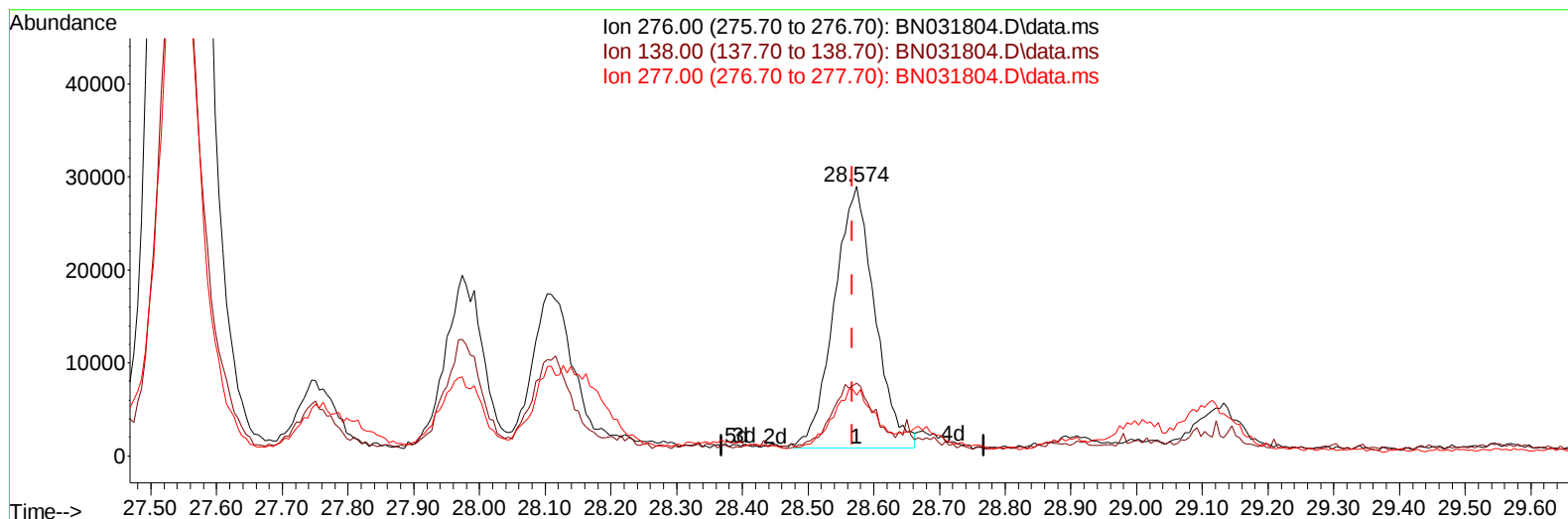
ClientSampleId :

BHBQ4DL

Manual IntegrationsAPPROVED

Quant Time: Jun 05 18:18:22 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed May 29 15:13:07 2024
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/06/2024
 Supervised By :mohammad ahmed 06/07/2024



TIC: BN031804.D\data.ms

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

28.574min (+ 0.006) 0.96 ng/u1

response 121935

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	26.10	26.97
277.00	22.50	22.60
0.00	0.00	0.00

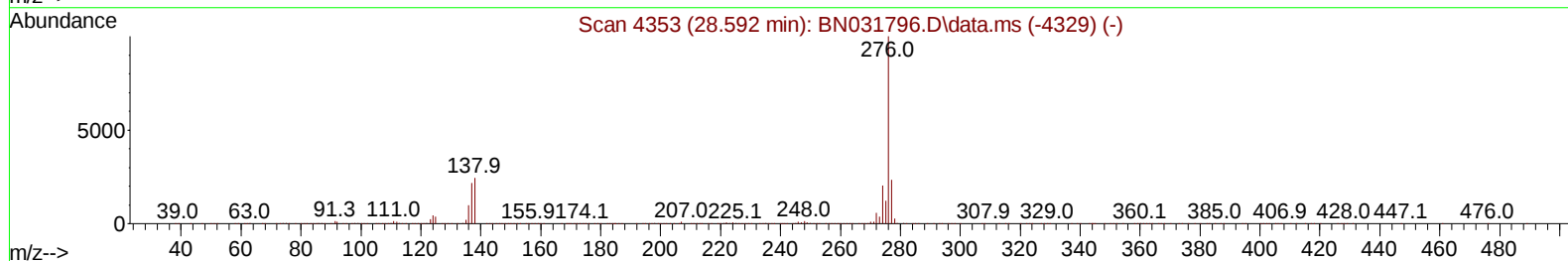
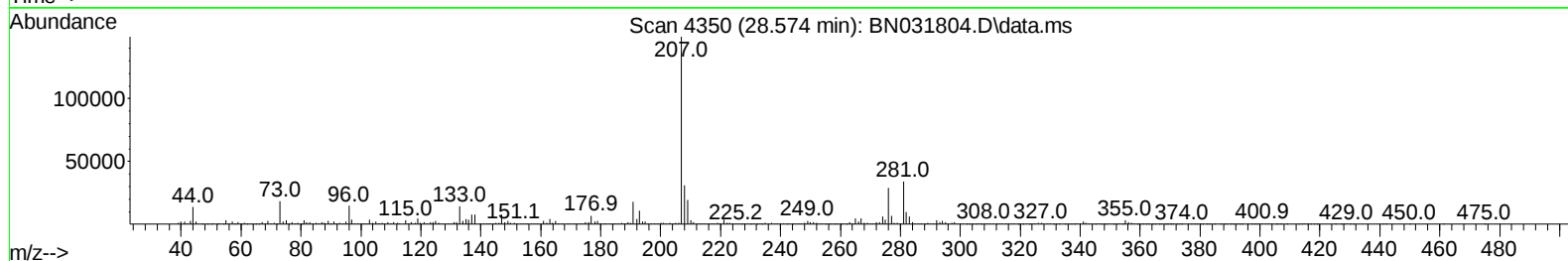
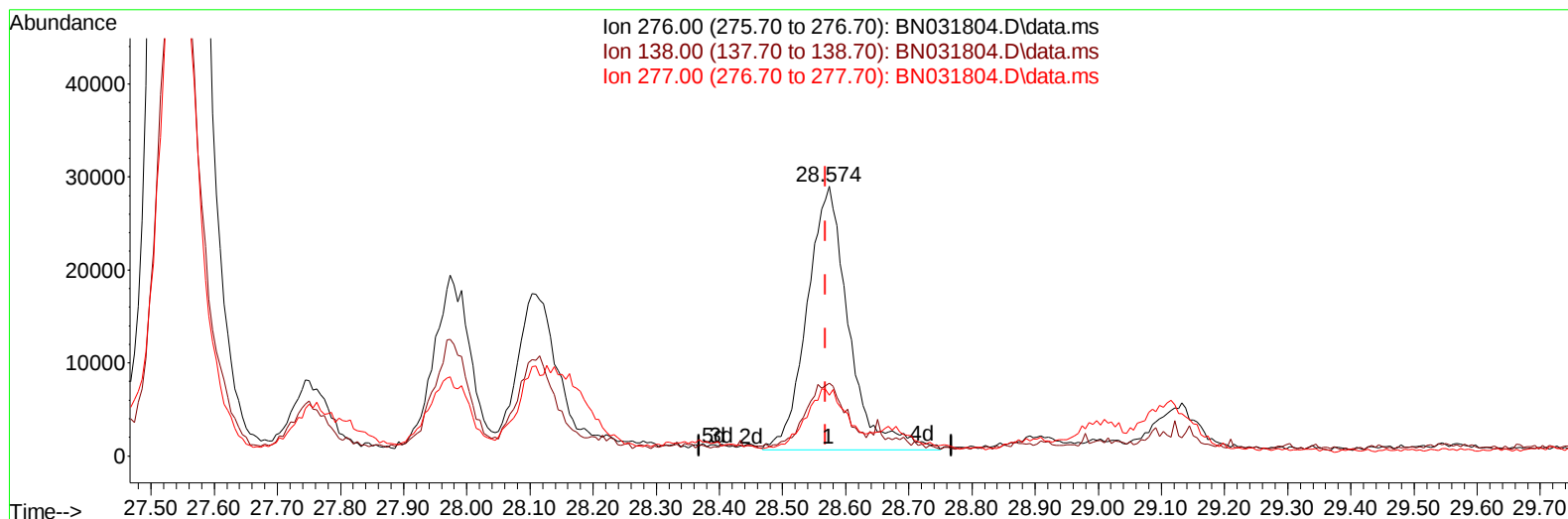
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060524\
Data File : BN031804.D
Acq On : 05 Jun 2024 17:45
Operator : MA/JU
Sample : P2591-13DL 5X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BHBQ4DL

Manual IntegrationsAPPROVED

Quant Time: Jun 05 18:18:22 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed May 29 15:13:07 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/06/2024
Supervised By :mohammad ahmed 06/07/2024



TIC: BN031804.D\data.ms

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

28.574min (+ 0.006) 1.02 ng/u1 m

response 129990

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	26.10	26.97
277.00	22.50	22.60
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060524\
 Data File : BN031804.D
 Acq On : 05 Jun 2024 17:45
 Operator : MA/JU
 Sample : P2591-13DL 5X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BHBQ4DL

Manual IntegrationsAPPROVED

Quant Time: Jun 06 01:06:10 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed May 29 15:13:07 2024
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/06/2024
 Supervised By :mohammad ahmed 06/07/2024

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

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.675	152		297956	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.457	136		1383910	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.316	164		940365	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.069	188		2014641	20.000	ng/ul	0.00
79) Chrysene-d12	21.304	240		2060521	20.000	ng/ul	0.00
88) Perylene-d12	24.239	264		2487456	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.187	96		7663	1.064	ng/uL	0.00
4) Pyridine-d5	3.599	84		59204	2.920	ng/ul	0.00
7) Phenol-d5	6.846	99		184719	7.232	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.016	67		98837	6.615	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.204	132		140156	7.191	ng/ul	-0.01
15) 4-Methylphenol-d8	8.375	113		146679	7.045	ng/ul	-0.01
21) Nitrobenzene-d5	8.828	128		70831	6.694	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.546	143		73811	6.753	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.081	165		143152	7.015	ng/ul	0.00
31) 4-Chloroaniline-d4	10.598	131		152868	5.046	ng/ul	-0.01
46) Dimethylphthalate-d6	13.734	166		473343	7.225	ng/ul	-0.01
49) Acenaphthylene-d8	14.010	160		545998	7.248	ng/ul	0.00
54) 4-Nitrophenol-d4	14.528	143		79708	6.077	ng/ul	0.00
60) Fluorene-d10	15.310	176		425009	7.667	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	0.000	200		0d	0.000	ng/ul	
73) Anthracene-d10	17.163	188		672998	7.803	ng/ul	0.00
81) Pyrene-d10	19.463	212		763540	6.925	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.033	264		693177	5.836	ng/ul	0.00
Target Compounds							
						Qvalue	
30) Naphthalene	10.504	128		153869	2.148	ng/ul	99
36) 2-Methylnaphthalene	12.122	142		49144	1.004	ng/ul	92
52) Acenaphthene	14.381	153		96173	1.707	ng/ul	99
56) Dibenzofuran	14.716	168		174816	2.284	ng/ul	94
61) Fluorene	15.369	166		116189	1.885	ng/ul	98
72) Phenanthrene	17.110	178		3299019	32.145	ng/ul	100
74) Anthracene	17.198	178		662791	6.374	ng/ul	99
77) Carbazole	17.469	167		471472	5.317	ng/ul	98
80) Fluoranthene	19.127	202		5218592	39.602	ng/ul	99
82) Pyrene	19.492	202		3868823	28.615	ng/ul	99
85) Benzo(a)anthracene	21.286	228		2603537	18.877	ng/ul	98
87) Chrysene	21.345	228		2331191	18.096	ng/ul	96
90) Benzo(b)fluoranthene	23.315	252		3239150	21.786	ng/ul	98
91) Benzo(k)fluoranthene	23.368	252		1040133	7.129	ng/ul	99
93) Benzo(a)pyrene	24.104	252		1408556	10.115	ng/ul	97
94) Indeno(1,2,3-cd)pyrene	27.545	276		982108	6.120	ng/ul	94
95) Dibenzo(a,h)anthracene	27.580	278		342262	2.613	ng/ul#	91
96) Benzo(g,h,i)perylene	28.574	276		129990m	1.019	ng/ul	

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

