

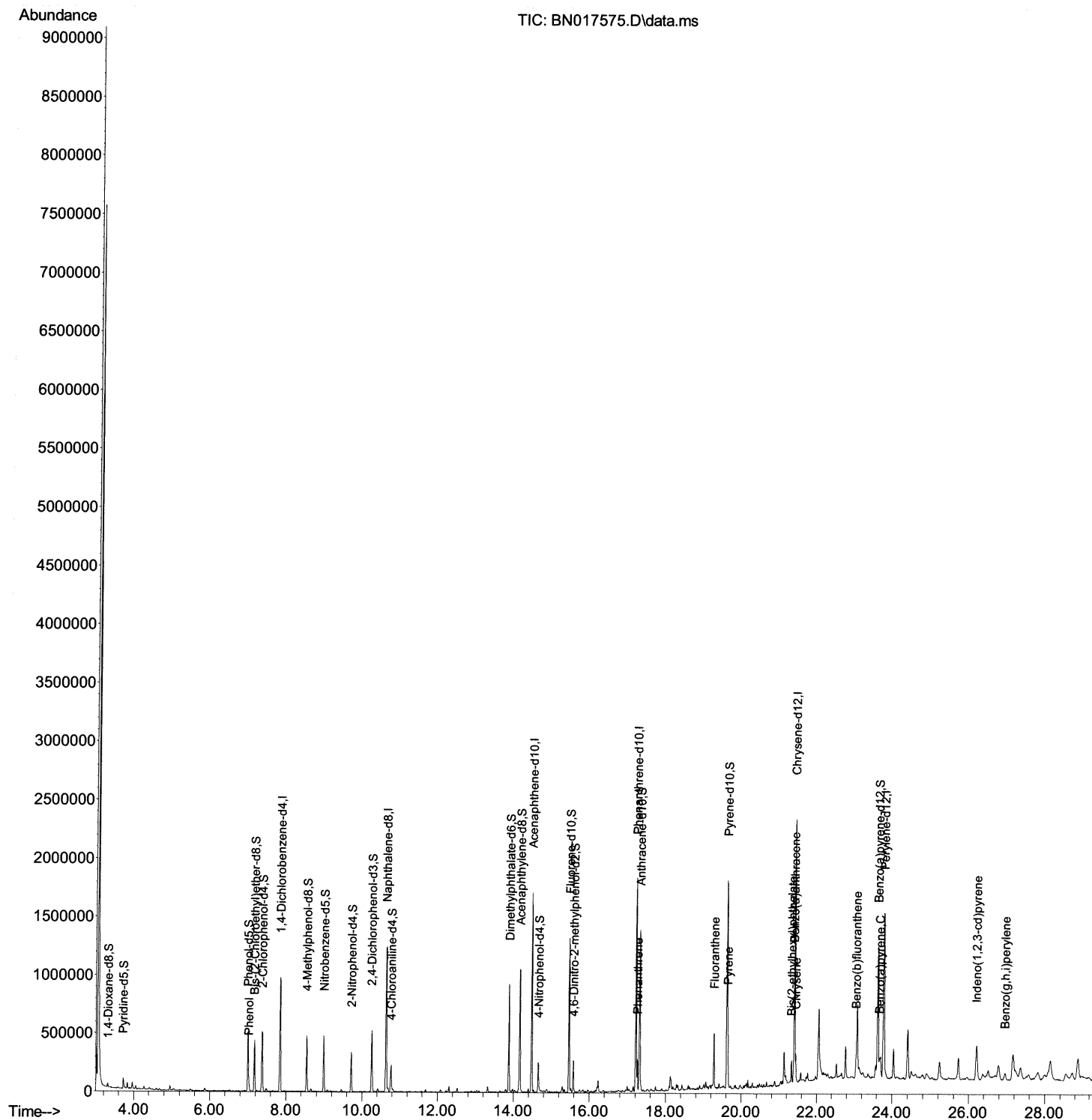
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112421\
Data File : BN017575.D
Acq On : 23 Nov 2021 20:21
Operator : CG/JU
Sample : M4702-10
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBLP5

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/24/2021
Supervised By :mohammad ahmed 11/26/2021

Quant Time: Nov 24 00:59:31 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN112221.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Nov 22 16:16:36 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

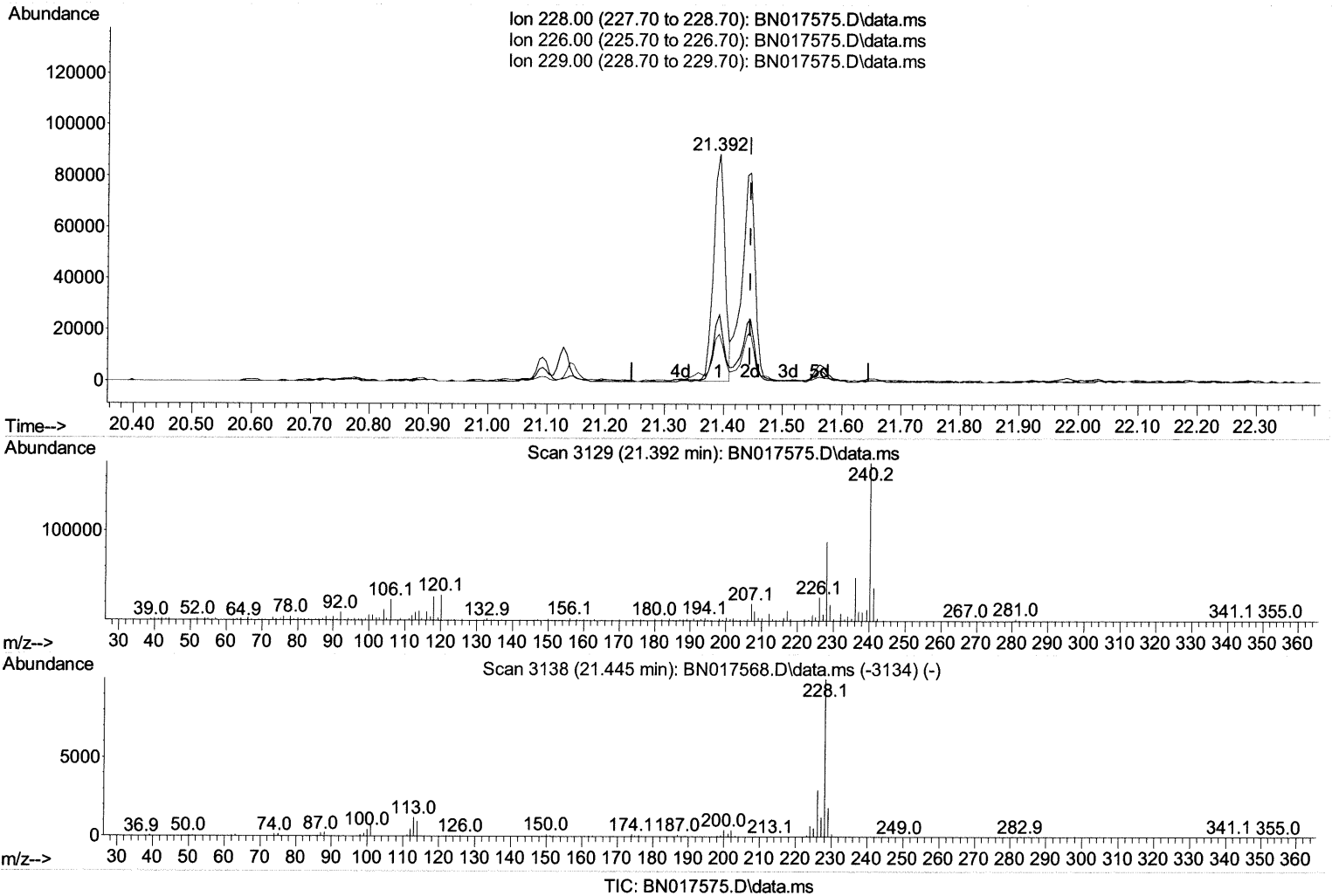
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(87) Chrysene

21.392min (-0.053) 1.88 ng/ul

response 119550

Ion	Exp%	Act%
228.00	100.00	100.00
226.00	30.10	29.63
229.00	19.50	20.84
0.00	0.00	0.00

Quantitation Report (Qedit)

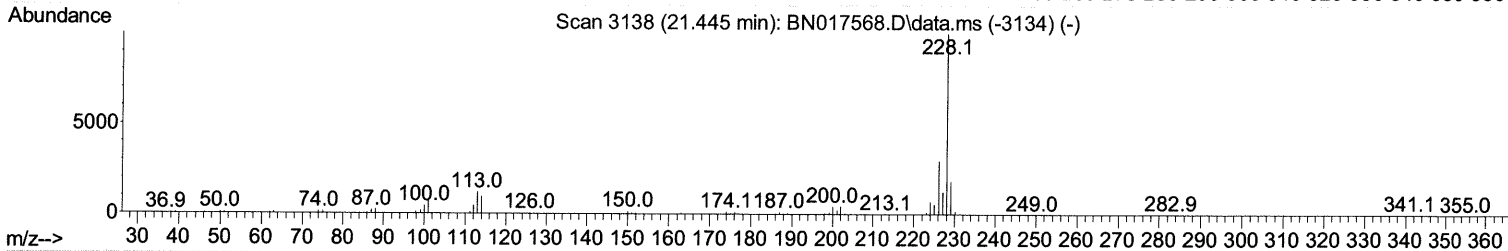
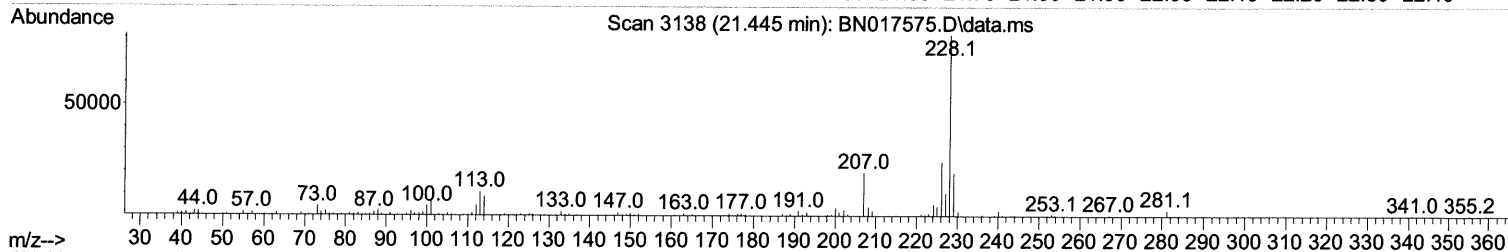
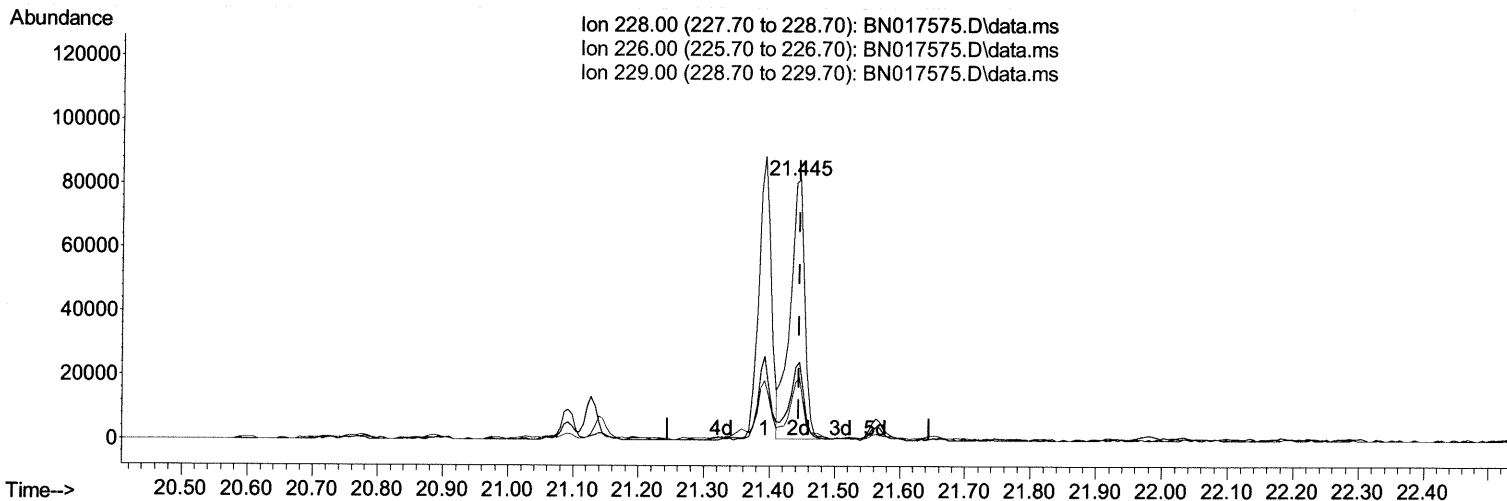
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TIC: BN017575.D\data.ms

(87) Chrysene

21.445min (+ 0.000) 2.02 ng/ul m 11/30/21 JU

response 128365

Ion	Exp%	Act%
228.00	100.00	100.00
226.00	30.10	29.96
229.00	19.50	23.91#
0.00	0.00	0.00

Quantitation Report (Qedit)

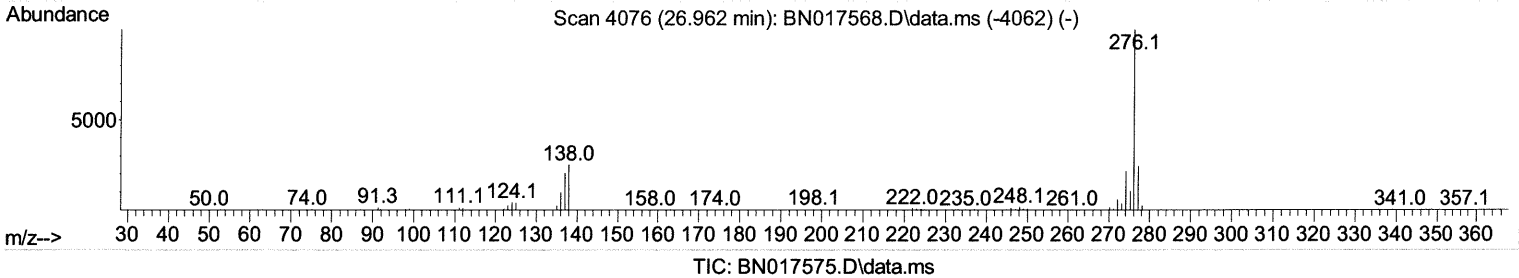
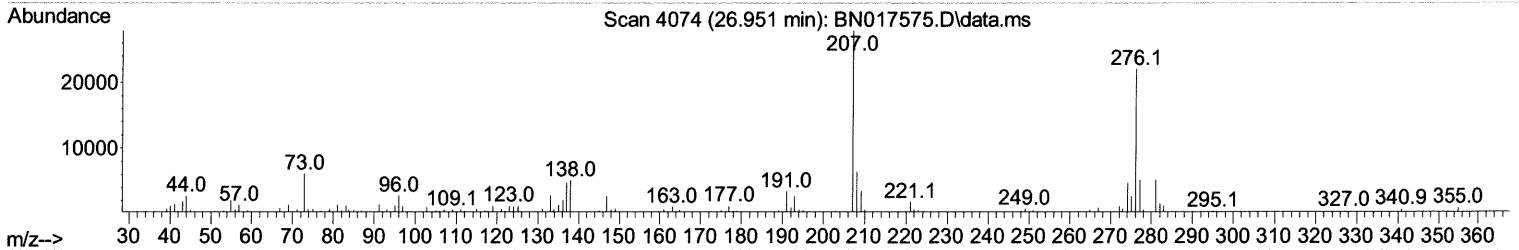
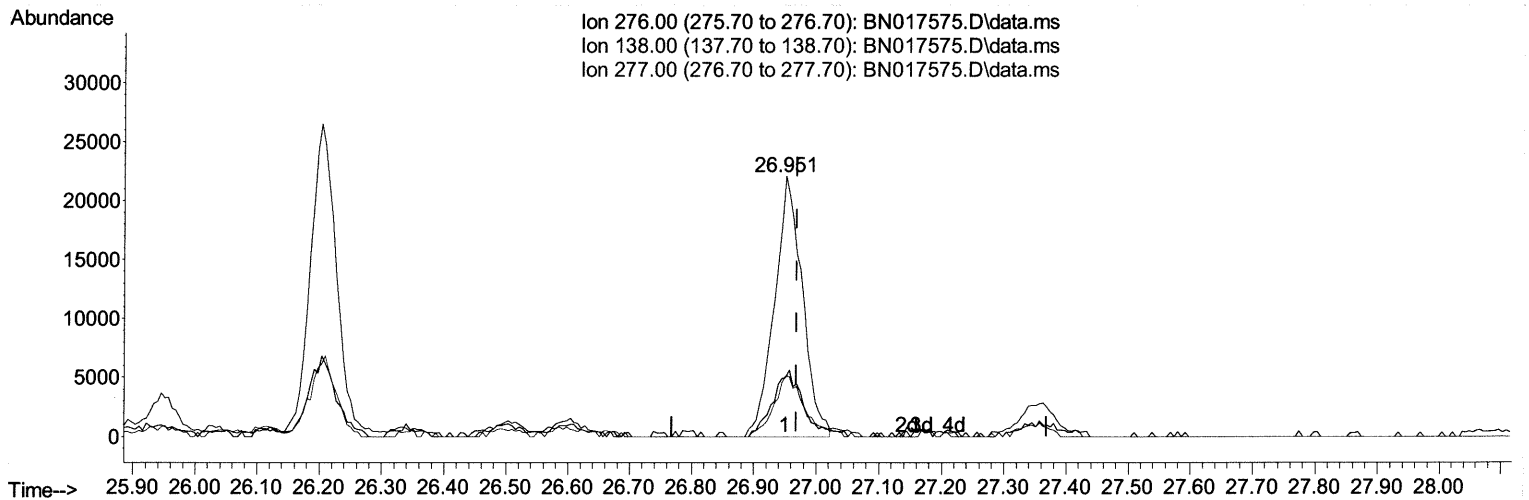
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(96) Benzo(g,h,i)perylene

26.951min (-0.017) 1.19 ng/ul

response 68180

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	22.80	22.82
277.00	21.70	23.26
0.00	0.00	0.00

Quantitation Report (Qedit)

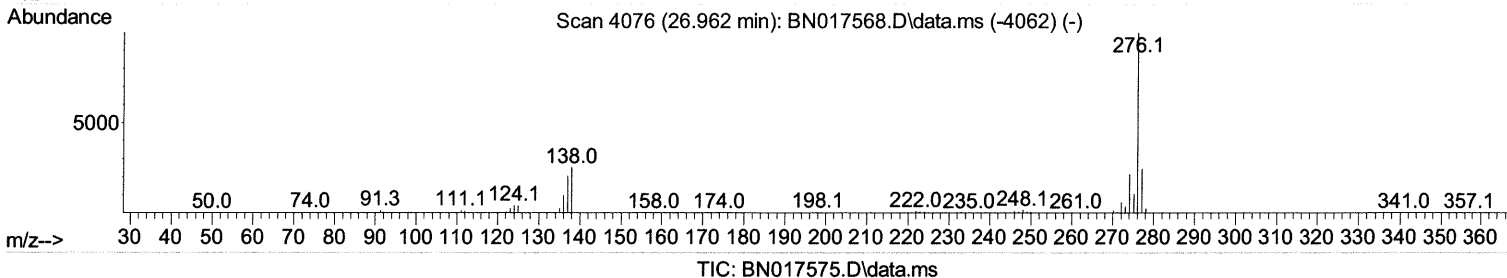
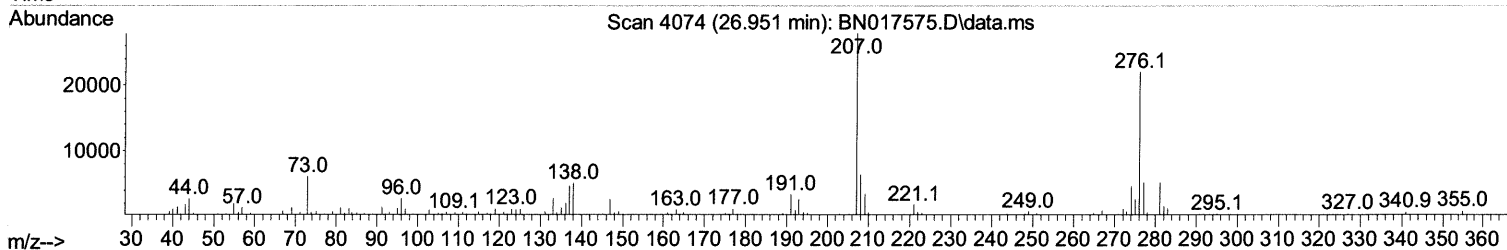
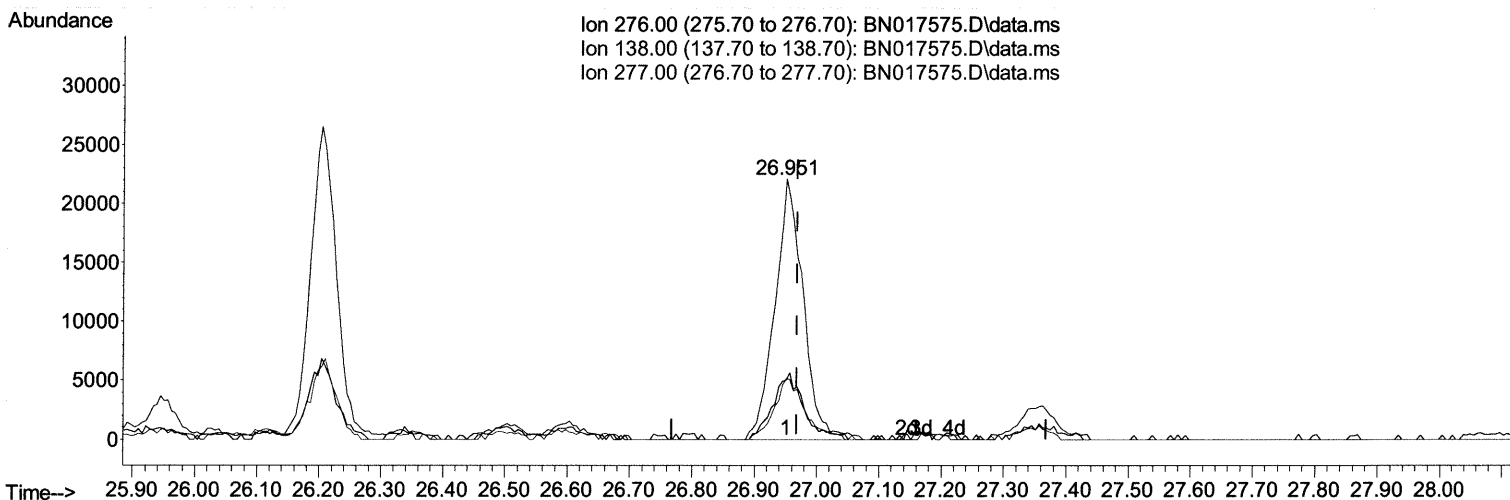
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(96) Benzo(g,h,i)perylene

26.951min (-0.017) 1.21 ng/ul m 11/30/21 JU

response 69592

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	22.80	22.82
277.00	21.70	23.26
0.00	0.00	0.00

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.840	152	257530	20.000	ng/ul	0.00
20) Naphthalene-d8	10.640	136	1020592	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.475	164	560543	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.222	188	1064183	20.000	ng/ul	0.00
79) Chrysene-d12	21.404	240	977906	20.000	ng/ul	0.00
88) Perylene-d12	23.774	264	985204	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.311	96	17228	2.391	ng/uL	0.00
4) Pyridine-d5	3.716	84	44368	2.346	ng/ul	0.00
7) Phenol-d5	7.005	99	283791	12.105	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.169	67	195928	13.336	ng/ul	0.00
11) 2-Chlorophenol-d4	7.375	132	229997	12.956	ng/ul	0.00
15) 4-Methylphenol-d8	8.540	113	175036	9.492	ng/ul	-0.01
21) Nitrobenzene-d5	8.993	128	105052	13.587	ng/ul	0.00
24) 2-Nitrophenol-d4	9.716	143	100542	13.143	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.251	165	189937	12.612	ng/ul	0.00
31) 4-Chloroaniline-d4	10.769	131	135602	6.270	ng/ul	0.00
46) Dimethylphthalate-d6	13.887	166	595306	14.607	ng/ul	0.00
49) Acenaphthylene-d8	14.169	160	745814	14.253	ng/ul	0.00
54) 4-Nitrophenol-d4	14.657	143	63182	9.161	ng/ul	0.00
60) Fluorene-d10	15.469	176	514330	14.962	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.581	200	53272	9.638	ng/ul	0.00
73) Anthracene-d10	17.316	188	746980	15.102	ng/ul	0.00
81) Pyrene-d10	19.610	212	896237	14.437	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.615	264	783803	15.133	ng/ul	-0.01
Target Compounds						
8) Phenol	7.028	94	33032	1.391	ng/ul	94
72) Phenanthrene	17.257	178	157135	2.706	ng/ul	99
80) Fluoranthene	19.280	202	268436	3.595	ng/ul	98
82) Pyrene	19.639	202	263290	3.522	ng/ul	99
85) Benzo(a)anthracene	21.392	228	118190	1.825	ng/ul	96
86) Bis(2-ethylhexyl)phtha...	21.333	149	61549	1.558	ng/ul	97
87) Chrysene	21.445	228	128365m >	2.017	ng/ul >	11/30/21JU
90) Benzo(b)fluoranthene	23.051	252	158085	2.321	ng/ul	97
93) Benzo(a)pyrene	23.662	252	106418	1.645	ng/ul	99
94) Indeno(1,2,3-cd)pyrene	26.203	276	77490	1.154	ng/ul	98
96) Benzo(g,h,i)perylene	26.951	276	69592m >	1.212	ng/ul >	11/30/21JU

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