Quantitation Report (Qedit)

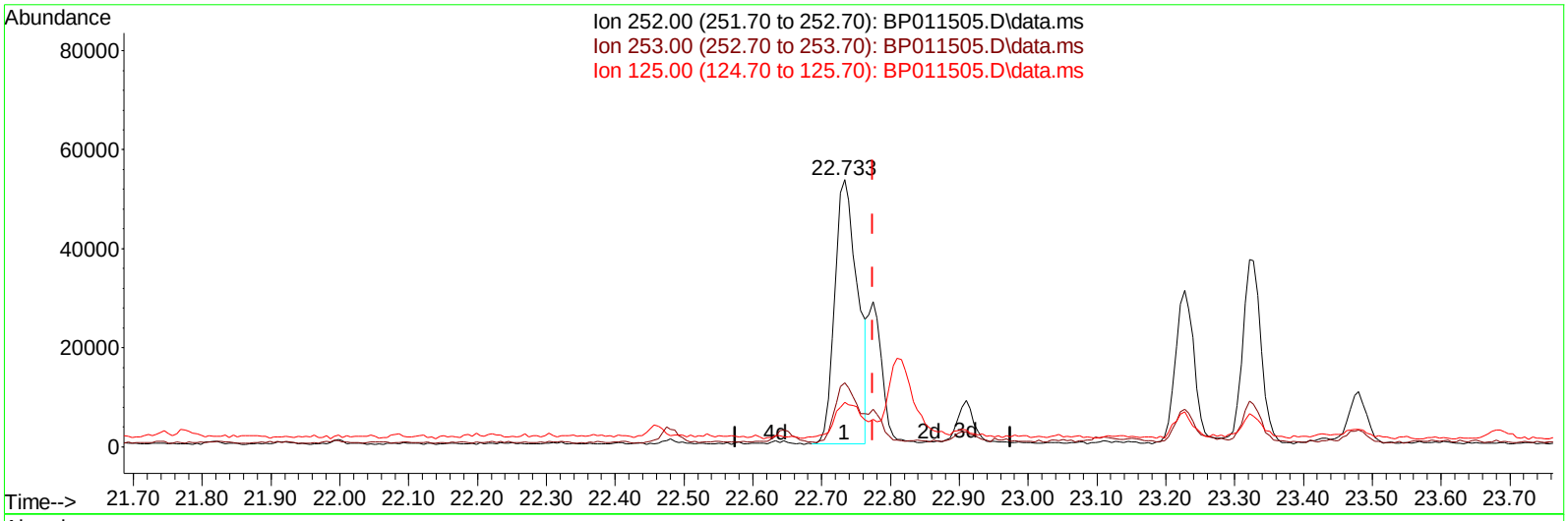
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081922\
Data File : BP011505.D
Acq On : 19 Aug 2022 13:44
Operator : CG/JU
Sample : N4210-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

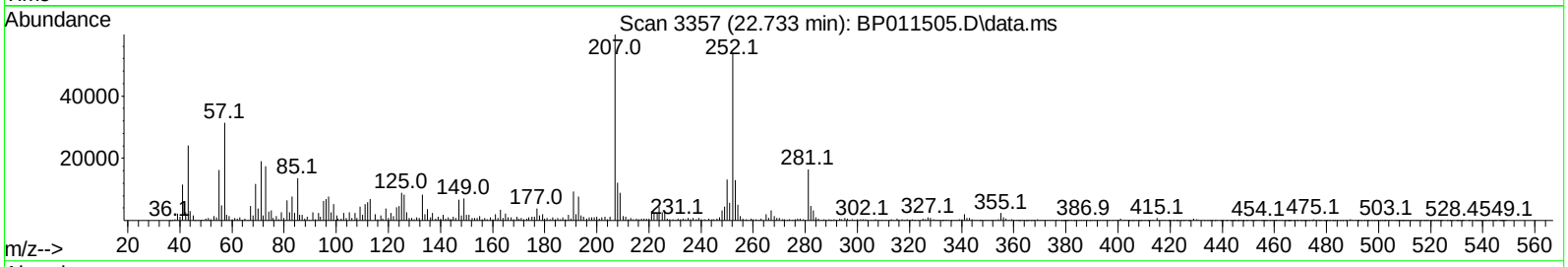
Quant Time: Aug 20 04:35:51 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081622.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Aug 17 01:58:37 2022
Response via : Initial Calibration

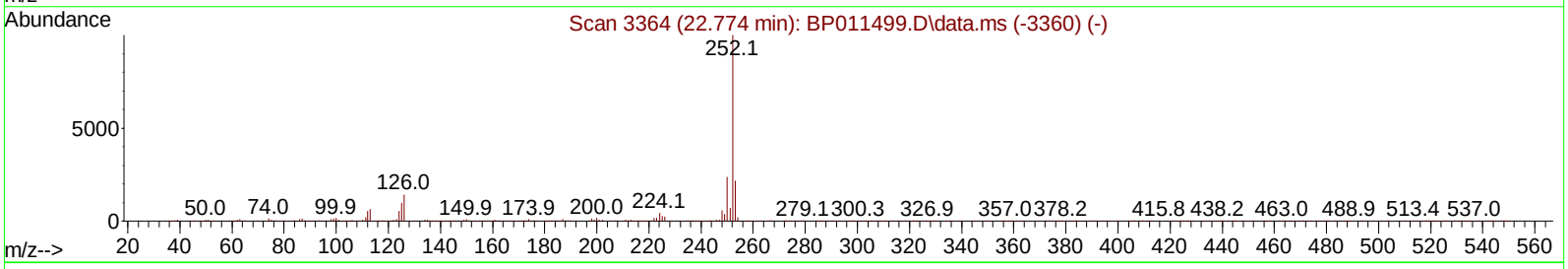
Instrument :
BNA_P
ClientSampleId :
BGC47

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 08/22/2022
Supervised By :Sohil Jodhani 08/22/2022







TIC: BP011505.D\data.ms

(91) Benzo(k)fluoranthene

22.733min (-0.041) 4.95 ng/ul

response 122487

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	22.40	23.97
125.00	12.30	16.70#
0.00	0.00	0.00

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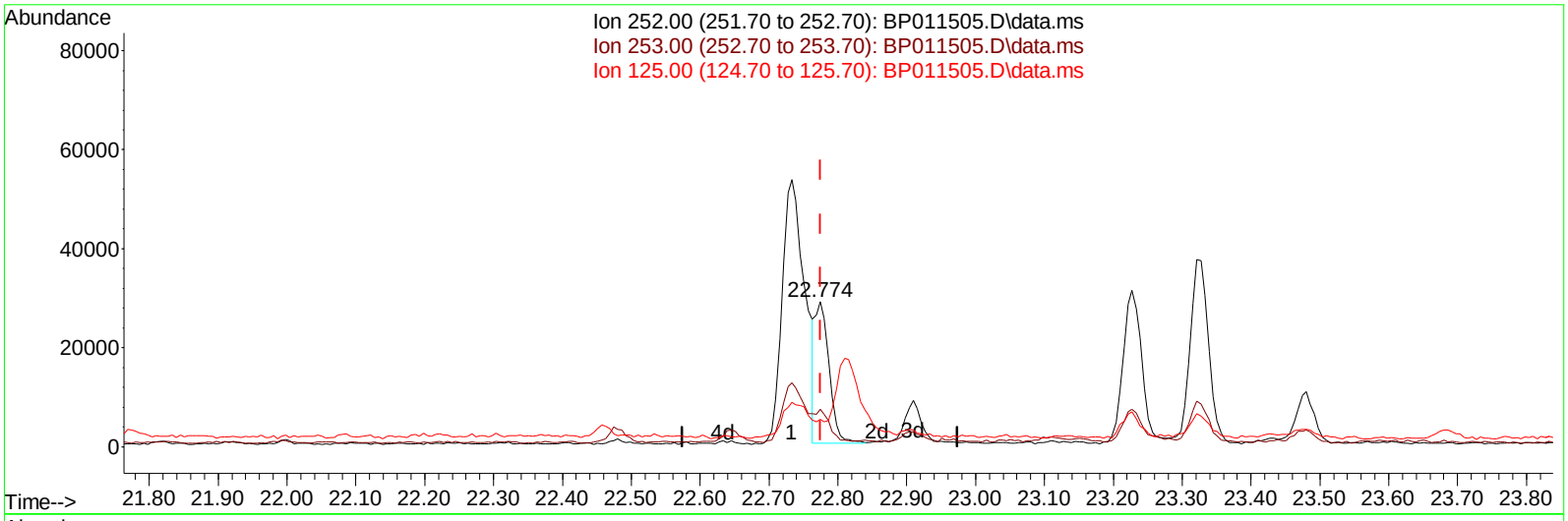

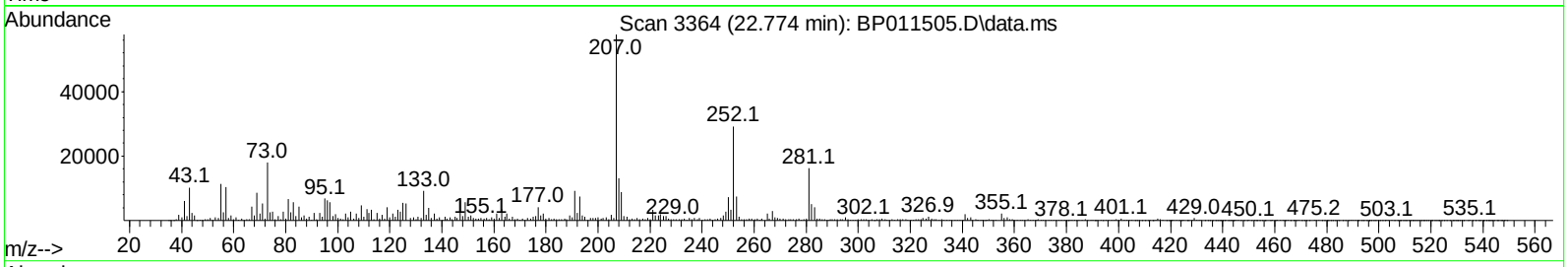

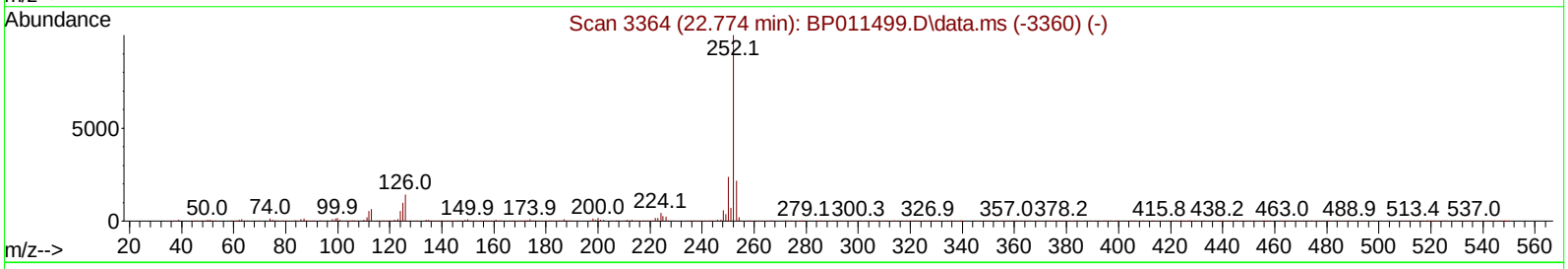

TIC: BP011505.D\data.ms

(91) Benzo(k)fluoranthene

22.774min (-0.000) 1.64 ng/ul m

response 40491

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	22.40	25.71
125.00	12.30	19.07#
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.558	152	96312	20.000	ng/ul	0.00
20) Naphthalene-d8	10.346	136	393328	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.228	164	213654	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.004	188	419548	20.000	ng/ul	0.00
79) Chrysene-d12	21.133	240	393234	20.000	ng/ul	0.00
88) Perylene-d12	23.427	264	416536	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.070	96	8256	2.660	ng/uL	0.00
4) Pyridine-d5	3.487	84	49845	6.273	ng/ul	0.00
7) Phenol-d5	6.746	99	158071	17.860	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.905	67	101264	17.611	ng/ul	0.00
11) 2-Chlorophenol-d4	7.087	132	128176	19.690	ng/ul	0.00
15) 4-Methylphenol-d8	8.275	113	109166	15.410	ng/ul	0.00
21) Nitrobenzene-d5	8.722	128	59946	21.170	ng/ul	0.00
24) 2-Nitrophenol-d4	9.440	143	62536	21.700	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.975	165	106458	20.012	ng/ul	0.00
31) 4-Chloroaniline-d4	10.499	131	145096	16.385	ng/ul	0.00
46) Dimethylphthalate-d6	13.657	166	344729	22.812	ng/ul	0.00
49) Acenaphthylene-d8	13.922	160	375258	22.343	ng/ul	0.00
54) 4-Nitrophenol-d4	14.463	143	47643	16.364	ng/ul	0.00
60) Fluorene-d10	15.234	176	288397	22.685	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.369	200	31921	14.080	ng/ul	0.00
73) Anthracene-d10	17.104	188	458123	24.655	ng/ul	0.00
81) Pyrene-d10	19.369	212	533383	26.809	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.280	264	522275	25.776	ng/ul	0.00
Target Compounds						Qvalue
72) Phenanthrene	17.045	178	105976	4.667	ng/ul	98
80) Fluoranthene	19.039	202	174036	7.113	ng/ul	97
82) Pyrene	19.392	202	133152	5.179	ng/ul	97
85) Benzo(a)anthracene	21.116	228	76254	3.183	ng/ul	97
86) Bis(2-ethylhexyl)phtha...	21.063	149	273555	16.029	ng/ul	99
87) Chrysene	21.169	228	80531	3.422	ng/ul	97
89) Di-n-octyl phthalate	21.963	149	36414	1.088	ng/ul	100
90) Benzo(b)fluoranthene	22.733	252	122487	4.727	ng/ul#	92
91) Benzo(k)fluoranthene	22.774	252	40491m	1.638	ng/ul	
93) Benzo(a)pyrene	23.321	252	68519	2.709	ng/ul#	90
94) Indeno(1,2,3-cd)pyrene	25.786	276	62690	2.222	ng/ul	99
96) Benzo(g,h,i)perylene	26.515	276	53477	2.228	ng/ul	96

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

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