Data File : BM033361.D

Acq On : 09 Dec 2021 16:48

Operator : CG/JU Sample : M4960-07

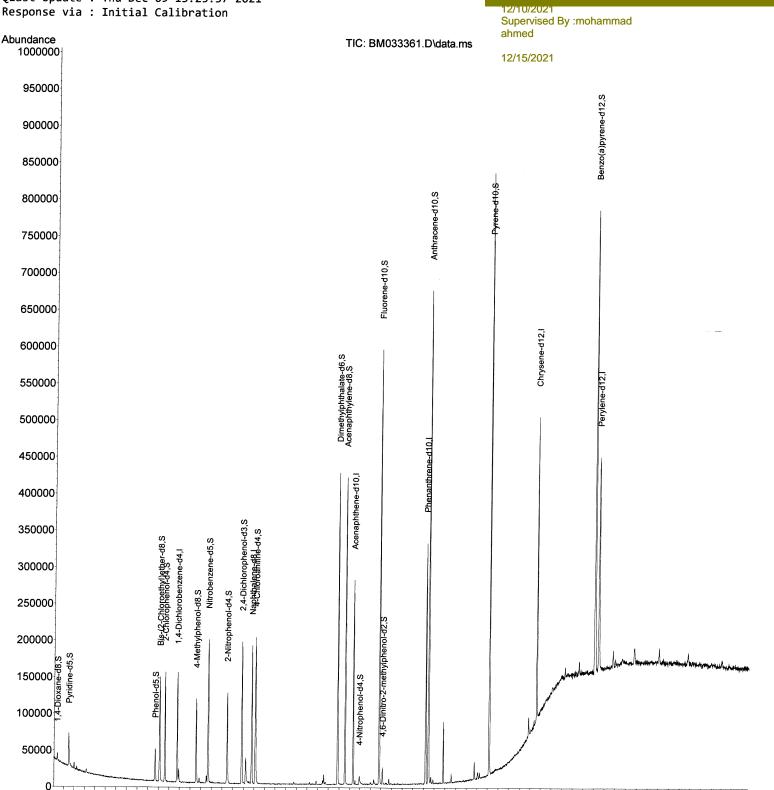
Misc

ALS Vial : 14 Sample Multiplier: 1

Quant Time: Dec 10 01:14:43 2021

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM120921.M

Quant Title : SVOA CALIBRATION
QLast Update : Thu Dec 09 13:25:37 2021
Response via : Initial Calibration



Instrument : BNA_M

BGKR7

ClientSampleId :

Upadhyay

Reviewed By :Jagrut

Manual IntegrationsAPPROVED

6.00

8.00

4.00

Time-->

10.00

12.00

14.00

16.00

18.00

20.00

22.00

24.00

26.00

Page: 2

28.00

Data File : BM033361.D

Acq On : 09 Dec 2021 16:48

Operator : CG/JU Sample : M4960-07

Misc

ALS Vial : 14 Sample Multiplier: 1

Quant Time: Dec 10 01:14:43 2021

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM120921.M

Quant Title : SVOA CALIBRATION

QLast Update : Thu Dec 09 13:25:37 2021 Response via : Initial Calibration

Instrument:
BNA_M
ClientSampleId:
BGKR7

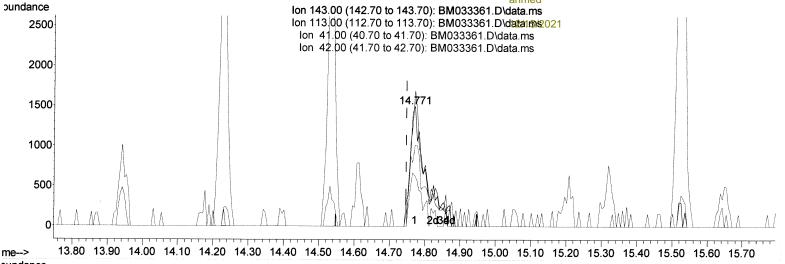
Manual IntegrationsAPPROVED

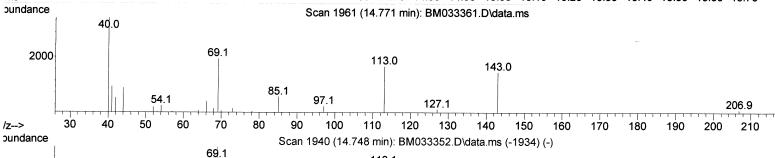
Reviewed By :Jagrut Upadhyay

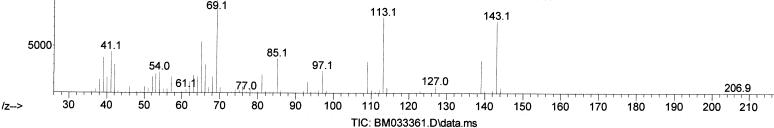
12/10/2021

Supervised By :mohammad

ahme







(54) 4-Nitrophenol-d4 (S)

14.771min (+ 0.024) 2.35 ng/ul

response	2830			
Ion	Ехр%	Act%		
143.00	100.00	100.00		
113.00	105.00	112.97		
41.00	57.20	67.51		
42.00	39.50	42.11		

Data File : BM033361.D

: 09 Dec 2021 16:48 Acq On

Operator : CG/JU Sample : M4960-07

Misc

<u>oundance</u>

ALS Vial : 14 Sample Multiplier: 1

Quant Time: Dec 10 01:14:43 2021

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM120921.M

Quant Title : SVOA CALIBRATION

QLast Update : Thu Dec 09 13:25:37 2021 Response via : Initial Calibration

Upadhyay

12/10/2021 Supervised By:mohammad

Reviewed By :Jagrut

Manual Integrations APPROVED

ahmed

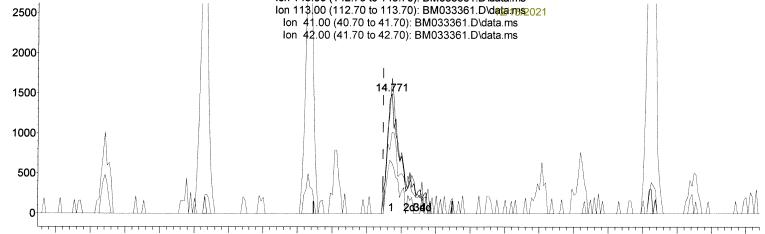
Instrument:

ClientSampleId :

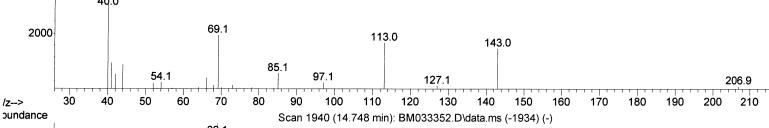
BNA_M

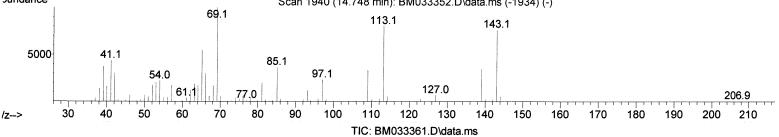
BGKR7





13.80 13.90 14.00 14.10 14.20 14.30 14.40 14.50 14.60 14.70 14.80 14.90 15.00 15.10 15.20 15.30 15.40 15.50 15.60 15.70 15.80 me--> **oundance** Scan 1961 (14.771 min): BM033361.D\data.ms 40.0 69.1





4-Nitrophenol-d4 (S)

14.771min (+ 0.024) 3.30 ng/ul m

response	3981	Herry.		
Ion	Exp%	Act%		
143.00	100.00	100.00		
113.00	105.00	112.97		
41.00	57.20	67.51		
42.00	39.50	42.11		

Data File : BM033361.D

: 09 Dec 2021 16:48 Acq On

Operator : CG/JU Sample : M4960-07

۱isc

ALS Vial : 14 Sample Multiplier: 1

Quant Time: Dec 10 01:14:43 2021

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM120921.M

Quant Title : SVOA CALIBRATION

QLast Update : Thu Dec 09 13:25:37 2021 Response via : Initial Calibration

Instrument: BNA_M

ClientSampleId:

BGKR7

Manual Integrations APPROVED

Reviewed By :Jagrut

Upadhyay

12/10/2021

Supervised By:mohammad

ahmed

Compound	R.T.	QIon	Response	Conc Units Dev(M	in) ^{ahmed}
Internal Standards					12/15/2021
1) 1,4-Dichlorobenzene-d4	7.913	152	36093	20.000 ng/ul	0.00
20) Naphthalene-d8	10.701		140703	20.000 ng/ul	0.00
38) Acenaphthene-d10	14.536	164	88818	20.000 ng/ul	0.00
64) Phenanthrene-d10	17.271	188	185763	20.000 ng/ul	0.00
79) Chrysene-d12	21.436	240	187141	20.000 ng/ul	0.00
88) Perylene-d12	23.759	264	199436	20.000 ng/ul	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.366	96	4651	4.842 ng/uL (0.00
4) Pyridine-d5	3.796	84	24571	•	0.01
7) Phenol-d5	7.084	99	20213	5.914 ng/ul (0.00
9) Bis-(2-Chloroethyl)eth	7.243	67	66458	29.702 ng/ul (0.00
<pre>11) 2-Chlorophenol-d4</pre>	7.443	132	55970	23.390 ng/ul (0.00
<pre>15) 4-Methylphenol-d8</pre>	8.619	113	37111	13.873 ng/ul (0.00
21) Nitrobenzene-d5	9.072	128	36594	32.053 ng/ul (0.00
24) 2-Nitrophenol-d4	9.789	143	29041		0.00
28) 2,4-Dichlorophenol-d3	10.331	165	59520	26.888 ng/ul (0.00

31) 4-Chloroaniline-d4 10.848 131 91662 27.929 ng/ul 0.00 46) Dimethylphthalate-d6 13.942 166 243590 0.00 36.706 ng/ul 3981m 3.304 ng/ul 0.00 210435 35.451 ng/ul 0.02 JU(1/26/2) 49) Acenaphthylene-d8 14.230 160 280317 54) 4-Nitrophenol-d4 14.771 143

60) Fluorene-d10 15.524 176 210435 35.451 ng/ul 65) 4,6-Dinitro-2-methylph... 15.648 200 4031 3.595 ng/ul 0.00 73) Anthracene-d10 17.371 188 360444 39.261 ng/ul 0.00

81) Pyrene-d10 19.653 212 423165 40.458 ng/ul 0.00 92) Benzo(a)pyrene-d12 23.612 264 413362 38.239 ng/ul 0.00

Target Compounds

Qvalue

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