Data File : BM033350.D

Acq On : 09 Dec 2021 09:41

Operator : CG/JU Sample : SSTD00511

Misc

ALS Vial : 3 Sample Multiplier: 1

Quant Time: Dec 09 13:03:44 2021

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

Quant Title : SVOA CALIBRATION

QLast Update : Thu Dec 09 13:01:40 2021

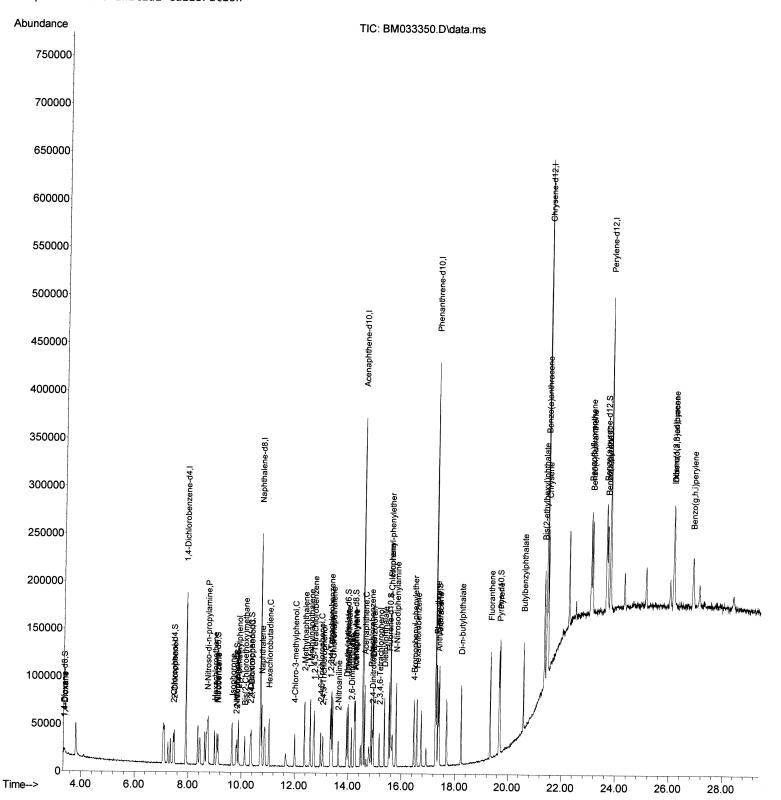
Response via : Initial Calibration



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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/10/2021 Supervised By :mohammad ahmed 12/15/2021



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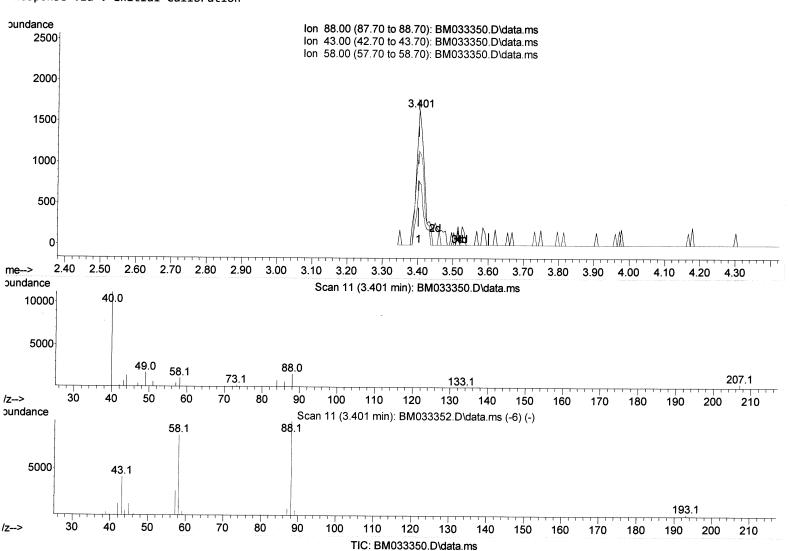
Quant Time: Dec 09 13:03:44 2021

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

Quant Title : SVOA CALIBRATION QLast Update : Thu Dec 09 13:01:40 2021 Response via : Initial Calibration Instrument : BNA_M ClientSampleId : SSTD005011

Manual IntegrationsAPPROVED

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(2) 1,4-Dioxane

3.401min (+ 0.000) 2.04 ng/uL

response	2406		
Ion	Exp%	Act%	
88.00	100.00	100.00	
43.00	45.30	47.09	
58.00	85.60	69.37	
0.00	0.00	0.00	

Data File: BM033350.D

Acq On : 09 Dec 2021 09:41

Operator : CG/JU Sample : SSTD00511

Misc

ALS Vial : 3 Sample Multiplier: 1

Quant Time: Dec 09 13:03:44 2021

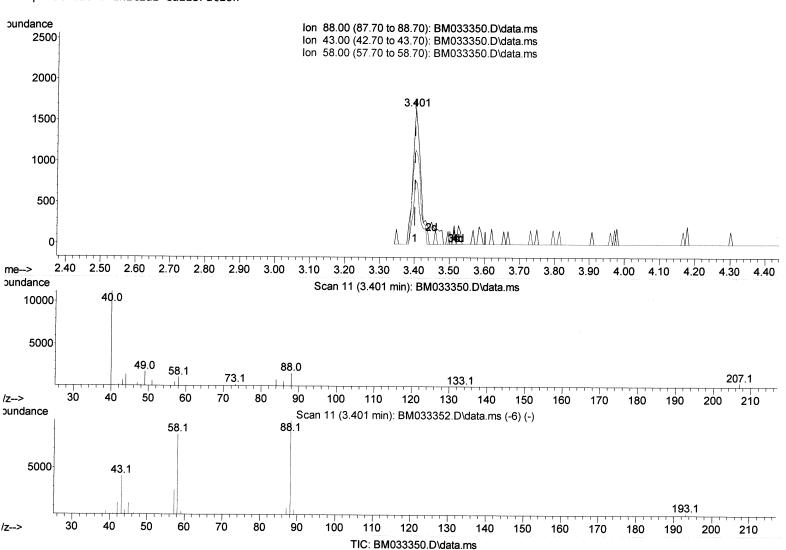
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM120921.M

Quant Title : SVOA CALIBRATION QLast Update : Thu Dec 09 13:01:40 2021 Response via : Initial Calibration



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(2) 1,4-Dioxane

3.401min	(+ 0.000)	2.25 ng/ul	Lm , 2)
response	2650	3	4 124 231 21
Ion	Ехр%	Act%	
88.00	100.00	100.00	
43.00	45.30	47.09	
58.00	85.60	69.37	
0.00	0.00	0.00	

Data File : BM033350.D

Acq On : 09 Dec 2021 09:41

Operator : CG/JU Sample : SSTD00511

Misc

ALS Vial : 3 Sample Multiplier: 1

Quant Time: Dec 09 13:03:44 2021

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

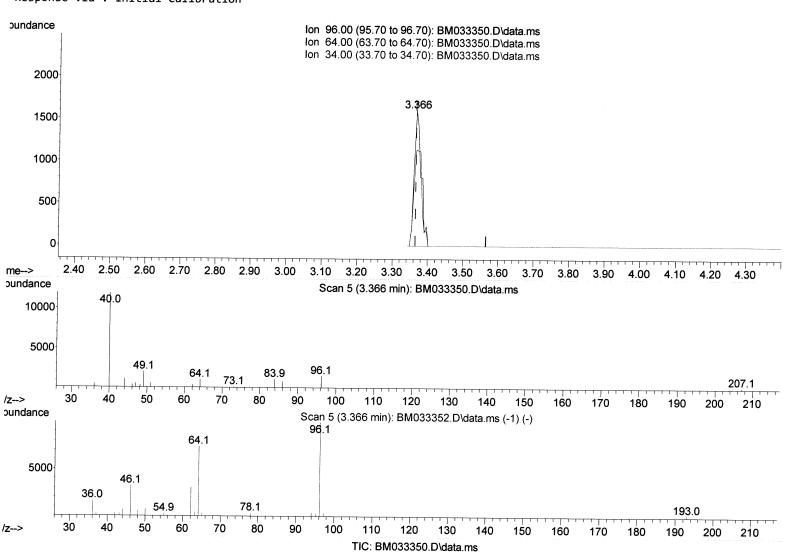
Quant Title : SVOA CALIBRATION

QLast Update : Thu Dec 09 13:01:40 2021 Response via : Initial Calibration



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3) 1,4-Dioxane-d8 (S)

3.366min (0.000) 1.81 ng/uL

response	2102			
Ion	Exp%	Act%		
96.00	100.00	100.00		
64.00	74.20	70.91		
34.00	0.00	0.00		
0.00	0.00	0.00		

Data File : BM033350.D

Acq On : 09 Dec 2021 09:41

Operator : CG/JU Sample : SSTD00511

Misc

ALS Vial : 3 Sample Multiplier: 1

Quant Time: Dec 09 13:03:44 2021

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

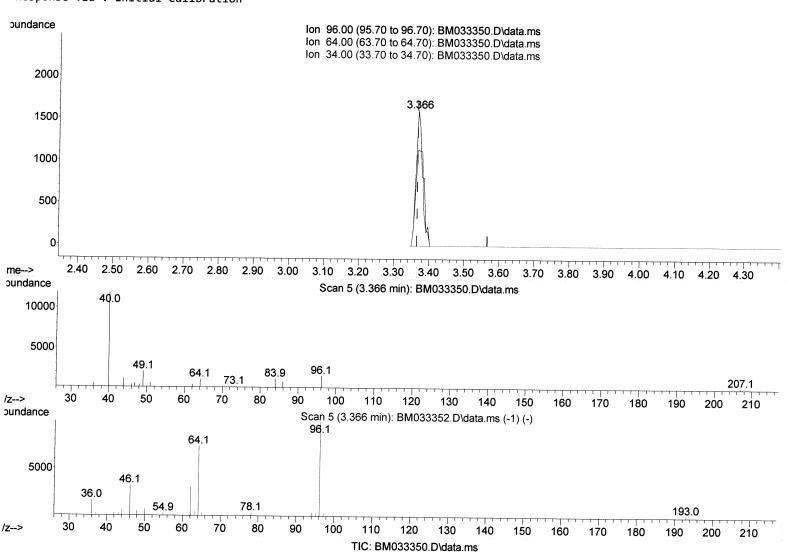
Quant Title : SVOA CALIBRATION

QLast Update : Thu Dec 09 13:01:40 2021 Response via : Initial Calibration



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(3)	1,4-Dioxan	e-d8 (S)
-----	------------	----------

- 14 12 23/21 3.366min (0.000) 1.88 ng/uL m response 2179 Ion Ехр% Act%

96.00	100.00	100.00
64.00	74.20	70.91
34.00	0.00	0.00
0.00	0.00	0.00

Data File: BM033350.D

Acq On : 09 Dec 2021 09:41

Dperator : CG/JU
Sample : SSTD00511

4isc :

ALS Vial : 3 Sample Multiplier: 1

Quant Time: Dec 09 13:03:44 2021

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

Quant Title : SVOA CALIBRATION

Compound

QLast Update : Thu Dec 09 13:01:40 2021
Response via : Initial Calibration

Instrument : BNA_M ClientSampleId : SSTD005011

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			&±0	пеэропэс	Conc on	.cs Dev	((1111)	
Internal	Standards							
1) 1,4-	Dichlorobenzene-d4	7.907	152	44772	20.000	ng/ul	0.00	
20) Naph	nthalene-d8	10.707		181096	20.000		0.00	
38) Acen	naphthene-d10	14.536	164	115636	20.000		0.00	
64) Phen	nanthrene-d10	17.271	188	240197	20.000		0.00	
79) Chry	rsene-d12	21.436	240	235477	20.000		0.00	
88) Pery	lene-d12	23.753	264	236752	20.000		0.00	
System Mo	onitoring Compounds							
	Dioxane-d8	3.366	96	2179m>	1.876	ng/ul	0 00 5 TT	112/23/21
	dine-d5	0.000	84	0d	0.000		0.00	[] -
7) Phen		0.000	99	0d	0.000			
	(2-Chloroethyl)eth		67	0d	0.000			
	_	7.448	132	12066	4.207		0.00	
	thylphenol-d8	0.000	113	9d	0.000	_	0.00	
	obenzene-d5	9.072	128	6537	5.028	_	0.00	
	trophenol-d4	9.795	143	5780	4.439		0.00	
	Dichlorophenol-d3		165	10520	3.528		0.00	
	loroaniline-d4	0.000	131	0d	0.000	-	0.00	
	thylphthalate-d6	13.942	166	38122	4.495		0.00	
	aphthylene-d8	14.230	160	44766	4.097		0.00	
	trophenol-d4	0.000	143	9d	0.000		0.00	
60) Fluo	•	15.524	176	33929	4.459		0.00	
	Dinitro-2-methylph	0.000	200	9d	0.000		0.00	
	racene-d10	17.371	188	50971	4.402		0.00	
81) Pyrei		19.659	212	57270	4.116		0.00	
	o(a)pyrene-d12	23.606	264	53910	4.243		0.00	
						-6,		12/23/21
Target Cor				_			alue	12/23/21
2) 1,4-1		3.401	88	2650m_			> V9 '	(- /
	lorophenol	7.478	128	12550	4.239 r			
	troso-di-n-propyla		70	12253	4.902 r		91	
	chloroethane	8.989	117	6715	4.979 r		85	
22) Nitro		9.113	77	18385	5.067 r		96	
23) Isoph		9.642	82	31081	4.686 r		99	
	trophenol	9.825	139	5982	4.334 r	_	97	
	Dimethylphenol	9.883	107	16525	4.557 r		95	
	2-Chloroethoxy)met		93	17756	4.499 r	g/ul	98	
	Dichlorophenol	10.360	162	11712	4.050 r	-	96	
30) Napht		10.754	128	45325	4.707 r		100	
	chlorobutadiene	11.030	225	9498	4.217 r		97	
	loro-3-methylphenol	11.995	107	12837	4.077 r	g/ul	94	
•	chylnaphthalene	12.366	142	29150	4.369 n	g/ul	96	
	hylnaphthalene	12.583	142	30901	4.537 n	g/ul	100	
	,5-Tetrachloroben	12.724	216	16224	4.217 n		98	
	5-Trichlorophenol	12.977	196	7900	3.487 n		91	
	5-Trichlorophenol	13.054	196	8691	3.577 n		89	
43) 1,1'-		13.371	154	40206	4.463 n	g/ul	99	
	oronaphthalene	13.413	162	30914	4.449 n		99	
	roaniline	13.630	65	8850	4.832 n	g/ul	95	
	hylphthalate	13.989	163	38587	4.668 n		98	
48) 2,6-D	initrotoluene	14.118	165	6488	4.587 n	g/ul#	82	

R.T. QIon Response Conc Units Dev(Min)

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Dperator : CG/JU
Sample : SSTD00511

۹isc :

ALS Vial : 3 Sample Multiplier: 1

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Quant Title : SVOA CALIBRATION

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Manual IntegrationsAPPROVED

Reviewed By: Jagrut Upadhyay 12/10/2021 Supervised By: mohammad ahmed 12/15/2021

Compound	R.T.	QIon	Response	Conc Units Dev(Min)
50) Acenaphthylene	14.260	152	47841	4.315 ng/ul	96
52) Acenaphthene	14.601	153	33128	4.580 ng/ul	97
56) Dibenzofuran	14.930	168	48006	4.528 ng/ul	100
57) 2,4-Dinitrotoluene	14.907	165	8760	4.503 ng/ul#	99
58) 2,3,4,6-Tetrachlorophenol	15.165	232	7061	3.561 ng/ul#	90
59) Diethylphthalate	15.348	149	37635	4.545 ng/ul	97
61) Fluorene	15.577	166	38328	4.548 ng/ul	98
62) 4-Chlorophenyl-phenyle	15.571	204	20263	4.607 ng/ul	89
67) N-Nitrosodiphenylamine	15.789	169	30729	4.350 ng/ul	92
68) 4-Bromophenyl-phenylether	16.471	248	10729	4.017 ng/ul	90
69) Hexachlorobenzene	16.577	284	12908	4.218 ng/ul	99
72) Phenanthrene	17.312	178	62848	4.724 ng/ul	98
74) Anthracene	17.406	178	61732	4.627 ng/ul	99
75) 1,2,3,4-Tetrachloroben	13.336	216	17181	4.355 ng/uL#	86
76) Pentachlorobenzene	14.848	250	16180	4.183 ng/uL	97
78) Di-n-butylphthalate	18.230	149	55563	4.287 ng/ul	98
80) Fluoranthene	19.324	202	68760	4.218 ng/ul	96
82) Pyrene	19.683	202	73128	4.418 ng/ul	99
83) Butylbenzylphthalate	20.571	149	23347	4.128 ng/ul	98
<pre>85) Benzo(a)anthracene</pre>	21.418	228	68525	4.499 ng/ul	98
<pre>86) Bis(2-ethylhexyl)phtha</pre>	21.336	149	34106	4.255 ng/ul#	98
87) Chrysene	21.471	228	70069	4.711 ng/ul	98
90) Benzo(b)fluoranthene	23.053	252	71175	4.464 ng/ul	98
<pre>91) Benzo(k)fluoranthene</pre>	23.100	252	66894	4.580 ng/ul	97
93) Benzo(a)pyrene	23.653	252	68418	4.540 ng/ul	99
94) Indeno(1,2,3-cd)pyrene	26.129	276	76573	4.563 ng/ul	100
95) Dibenzo(a,h)anthracene	26.141	278	66274	4.616 ng/ul	100
96) Benzo(g,h,i)perylene	26.859	276	65301	4.483 ng/ul	96

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