Data File : BM033352.D

Acq On : 09 Dec 2021 10:52

Operator : CG/JU Sample : SSTD02013

Misc

ALS Vial : 5 Sample Multiplier: 1

Quant Time: Dec 09 13:01:52 2021

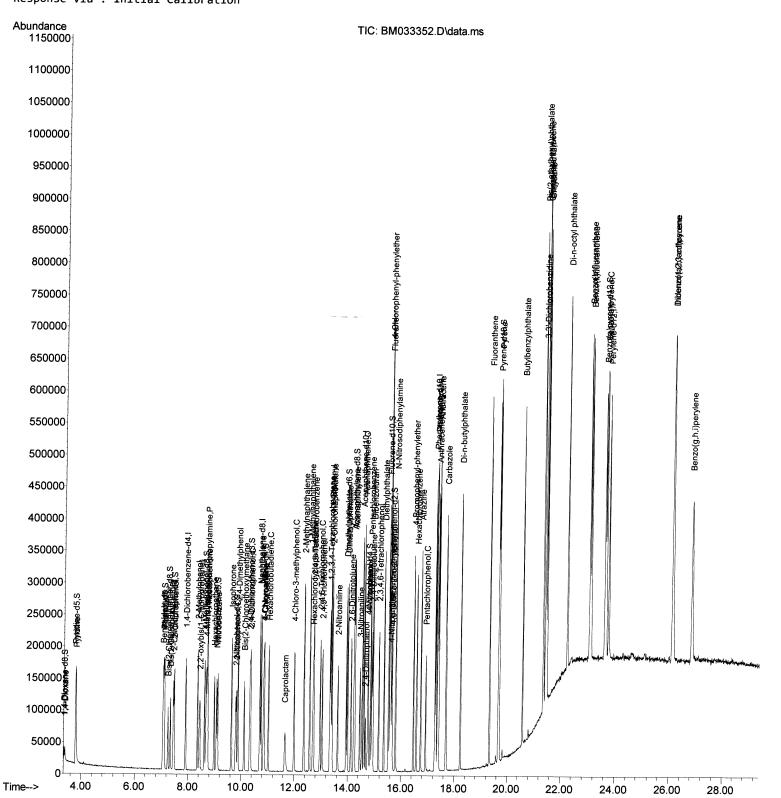
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SFAM-EPA-BM120921.M

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

Instrument :
BNA\_M
ClientSampleId :
SSTD020013

### **Manual IntegrationsAPPROVED**

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



Data File : BM033352.D

Acq On : 09 Dec 2021 10:52

Operator : CG/JU Sample : SSTD02013

Misc

ALS Vial : 5 Sample Multiplier: 1

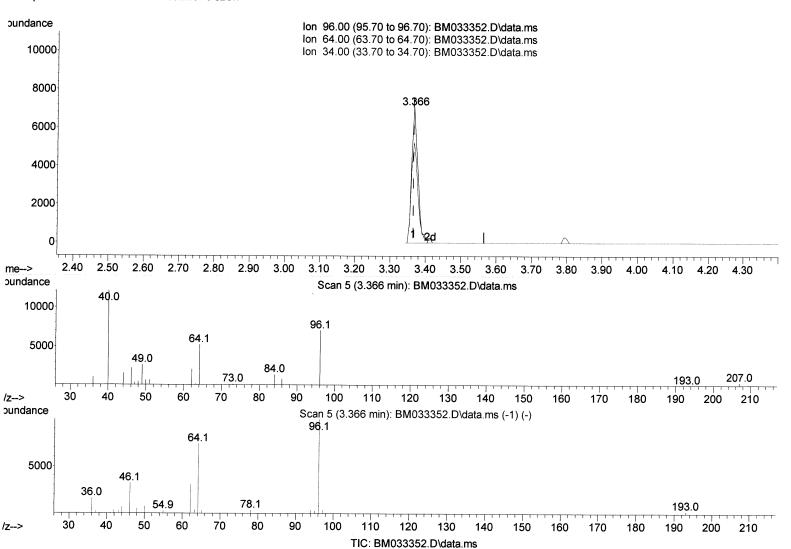
Quant Time: Dec 09 13:01:52 2021

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:
SSTD020013

#### **Manual IntegrationsAPPROVED**

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

3.366min (0.000) 6.86 ng/uL

response	7693				
Ion	Exp%	Act%			
96.00	100.00	100.00			
64.00	74.20	74.17			
34.00	0.00	0.00			
0.00	0.00	0.00			

Data File : BM033352.D

Acq On : 09 Dec 2021 10:52

Operator : CG/JU Sample : SSTD02013

Misc

ALS Vial : 5 Sample Multiplier: 1

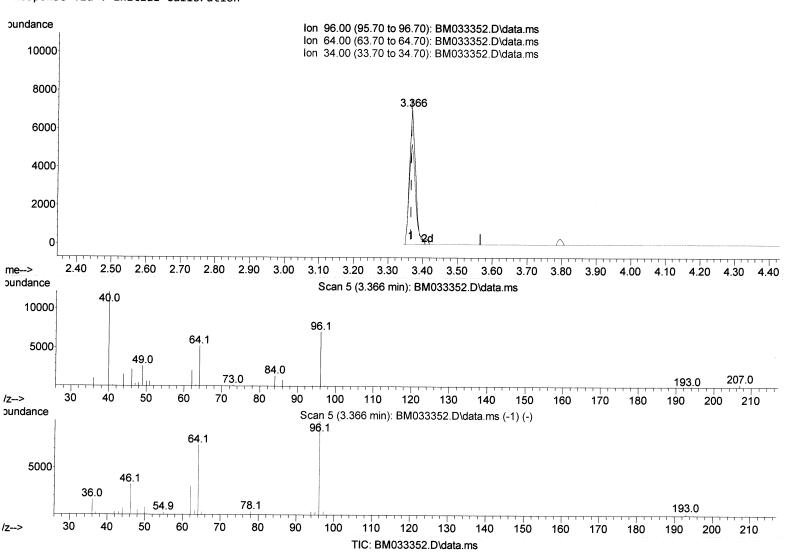
Quant Time: Dec 09 13:01:52 2021

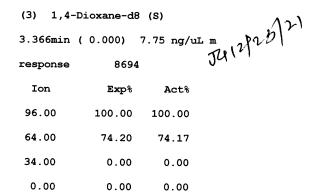
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SFAM-EPA-BM120921.M

Quant Title : SVOA CALIBRATION QLast Update : Thu Dec 09 13:01:40 2021 Response via : Initial Calibration Instrument : BNA\_M ClientSampleId : SSTD020013

#### **Manual IntegrationsAPPROVED**

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





Data File : BM033352.D

Acq On : 09 Dec 2021 10:52

Dperator : CG/JU
Sample : SSTD02013

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ALS Vial : 5 Sample Multiplier: 1

Quant Time: Dec 09 13:01:52 2021

 ${\tt Quant\ Method: Z:\SVOASRV\ HPCHEM1\ BNA\_M\ METHODS\ SFAM-EPA-BM120921.M}$ 

Quant Title : SVOA CALIBRATION

QLast Update : Thu Dec 09 13:01:40 2021 Response via : Initial Calibration Instrument : BNA\_M ClientSampleId : SSTD020013

## **Manual IntegrationsAPPROVED**

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

Compound	R.T.	QIon	Response	Conc Units Dev	v(Min)
Internal Standards					
1) 1,4-Dichlorobenzene-d4	7.913	152	43237	20.000 ng/ul	0.00
20) Naphthalene-d8	10.707		178084	20.000 ng/ul	0.00
38) Acenaphthene-d10	14.536	164	119509	20.000 ng/ul	0.00
64) Phenanthrene-d10	17.277	188	259377	20.000 ng/ul	0.00
79) Chrysene-d12	21.436		280357	20.000 ng/ul	0.00
88) Perylene-d12	23.759		280809	20.000 ng/ul	0.00
System Monitoring Compounds					>0.007412/25/21 0.00
3) 1,4-Dioxane-d8	3.366	96	8694m >	> 7.752 ng/uL	50 00 TU 12/25/
4) Pyridine-d5	3.784	84	58096	18.805 ng/ul	9 99
7) Phenol-d5	7.084	99	68758	18.800 ng/ul	0.00
<pre>9) Bis-(2-Chloroethyl)eth</pre>		67	46405	20.004 ng/ul	0.00
11) 2-Chlorophenol-d4	7.448	132	51526	18.601 ng/ul	0.00
15) 4-Methylphenol-d8	8.619	113	53992	19.022 ng/ul	0.00
21) Nitrobenzene-d5	9.072	128	26010	20.343 ng/ul	0.00
24) 2-Nitrophenol-d4	9.795	143	27069	21.139 ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.330	165	53475	18.236 ng/ul	0.00
31) 4-Chloroaniline-d4	10.848	131	72677	18.637 ng/ul	0.00
46) Dimethylphthalate-d6	13.942	166	165694	18.902 ng/ul	0.00
49) Acenaphthylene-d8	14.230	160	209167	18.522 ng/ul	0.00
54) 4-Nitrophenol-d4	14.748	143	26239	18.376 ng/ul	0.00
60) Fluorene-d10	15.524	176	150202	19.101 ng/ul	0.00
65) 4,6-Dinitro-2-methylph	15.642	200	25657	21.048 ng/ul	0.00
73) Anthracene-d10	17.371	188	244257	19.536 ng/ul	0.00
81) Pyrene-d10	19.659	212	285949	17.260 ng/ul	0.00
92) Benzo(a)pyrene-d12	23.612	264	282168	18.723 ng/ul	0.00
Target Compounds				Ova	alue
2) 1,4-Dioxane	3.401	88	9556	8.392 ng/uL	100
5) Pyridine	3.807	79	58109	18.433 ng/ul	100
6) Benzaldehyde	7.054	77	52728	25.297 ng/ul	100
8) Phenol	7.107	94	72772	19.981 ng/ul	100
<pre>10) Bis(2-Chloroethyl)ether</pre>	7.337	93	55870	19.311 ng/ul	100
12) 2-Chlorophenol	7.478	128	53187	18.603 ng/ul	100
<pre>13) 2-Methylphenol</pre>	8.354	108	51559	18.480 ng/ul	100
<pre>14) 2,2'-oxybis(1-Chloropr</pre>	8.436	45	98259	21.990 ng/ul	100
16) Acetophenone	8.736	105	90502	20.201 ng/ul	100
17) N-Nitroso-di-n-propyla	8.713	70	51230	21.224 ng/ul	100
<pre>18) 4-Methylphenol</pre>	8.683	108	56637	19.365 ng/ul	100
19) Hexachloroethane	8.983	117	27150	20.845 ng/ul	100
22) Nitrobenzene	9.113	77	78989	22.137 ng/ul	100
23) Isophorone	9.636	82	132452	20.309 ng/ul	100
25) 2-Nitrophenol	9.825	139	29620	21.823 ng/ul	100
<pre>26) 2,4-Dimethylphenol</pre>	9.883	107	70678	19.819 ng/ul	100
<pre>27) Bis(2-Chloroethoxy)met</pre>	10.119	93	75913	19.560 ng/ul	100
<pre>29) 2,4-Dichlorophenol</pre>	10.360	162	52337	18.406 ng/ul	100
30) Naphthalene	10.754	128	180587	19.070 ng/ul	100
32) 4-Chloroaniline	10.872	127	72935	18.514 ng/ul	100
33) Hexachlorobutadiene	11.036	225	37718	17.029 ng/ul	100
34) Caprolactam	11.654	113	15944	19.504 ng/ul	100
35) 4-Chloro-3-methylphenol	11.989	107	61631	19.904 ng/ul	100

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ALS Vial : 5 Sample Multiplier: 1

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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 : SSTD020013

## **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)
36) 2-Methylnaphthalene	12.360	142	125139	19.072 ng/ul	100
37) 1-Methylnaphthalene	12.583	142	128479	19.182 ng/ul	100
39) 1,2,4,5-Tetrachloroben	12.724	216	67820	17.057 ng/ul	100
40) Hexachlorocyclopentadiene	12.701	237	41738	14.918 ng/ul	100
41) 2,4,6-Trichlorophenol	12.971	196	40859	17.449 ng/ul	100
42) 2,4,5-Trichlorophenol	13.048	196	42496	16.924 ng/ul	100
43) 1,1'-Biphenyl	13.371	154	173074	18.587 ng/ul	100
44) 2-Chloronaphthalene	13.413	162	132292	18.423 ng/ul	100
45) 2-Nitroaniline	13.624	65	47010	24.838 ng/ul	100
47) Dimethylphthalate	13.989	163	162152	18.979 ng/ul	100
48) 2,6-Dinitrotoluene	14.118	165	31725	21.703 ng/ul	100
50) Acenaphthylene	14.260	152	216917	18.929 ng/ul	100
51) 3-Nitroaniline	14.448	138	31934	21.682 ng/ul	100
52) Acenaphthene	14.601	153	142991	19.128 ng/ul	100
53) 2,4-Dinitrophenol	14.654	184	14181	19.007 ng/ul	100
55) 4-Nitrophenol	14.760	109	29404	20.162 ng/ul	100
56) Dibenzofuran	14.936	168	208498	19.029 ng/ul	100
57) 2,4-Dinitrotoluene	14.901	165	47789	23.771 ng/ul	100
58) 2,3,4,6-Tetrachlorophenol	15.160	232	37046	18.075 ng/ul	100
59) Diethylphthalate	15.348	149	170722	19.951 ng/ul	100
61) Fluorene	15.583	166	171554	19.699 ng/ul	100
62) 4-Chlorophenyl-phenyle	15.571	204	84628	18.618 ng/ul	100
63) 4-Nitroaniline	15.612	138	34884	24.087 ng/ul	100
66) 4,6-Dinitro-2-methylph	15.660	198	26465	21.662 ng/ul	100
67) N-Nitrosodiphenylamine	15.789	169	142770	18.715 ng/ul	100
68) 4-Bromophenyl-phenylether	16.465	248	48930	16.964 ng/ul	100
69) Hexachlorobenzene	16.577	284	55756	16.873 ng/ul	100
70) Atrazine	16.736	200	54371	18.445 ng/ul	100
71) Pentachlorophenol	16.924	266	27345	14.316 ng/ul	100
72) Phenanthrene	17.318	178	282664	19.674 ng/ul	100
74) Anthracene	17.406	178	287791	19.974 ng/ul	100
75) 1,2,3,4-Tetrachloroben	13.336	216	69637	16.345 ng/uL	100
76) Pentachlorobenzene	14.848	250	69254	16.581 ng/uL	100
77) Carbazole	17.683	167	251733	19.667 ng/ul	100
78) Di-n-butylphthalate	18.230	149	287299	20.527 ng/ul	100
80) Fluoranthene	19.324	202	335155	17.268 ng/ul	100
82) Pyrene	19.683	202	353117	17.918 ng/ul	100
83) Butylbenzylphthalate	20.571	149	131917	19.592 ng/ul	100
84) 3,3'-Dichlorobenzidine	21.353	252	119115	18.859 ng/ul	100
85) Benzo(a)anthracene	21.418	228	337152	18.593 ng/ul	100
86) Bis(2-ethylhexyl)phtha	21.341	149	190013	19.910 ng/ul	100
87) Chrysene	21.471	228	337233	19.044 ng/ul	100
89) Di-n-octyl phthalate	22.236	149	340369	19.893 ng/ul	100
90) Benzo(b)fluoranthene 91) Benzo(k)fluoranthene	23.053	252	352246	18.625 ng/ul	100
93) Benzo(k)fluoranthene 93) Benzo(a)pyrene	23.100	252	331507	19.135 ng/ul	100
94) Indeno(1,2,3-cd)pyrene	23.653	252	344835	19.293 ng/ul	100
95) Dibenzo(a,h)anthracene	26.135	276	369691	18.574 ng/ul	100
96) Benzo(g,h,i)perylene	26.141 26.859	278 276	323553 317626	19.001 ng/ul	100
	20.000	276	317626	18.386 ng/ul	100
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<sup>(#) =</sup> qualifier out of range (m) = manual integration (+) = signals summed