Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111721\

Data File : BM033131.D Acq On : 17 Nov 2021 22:12

Operator : CG/JU Sample : PB140762BS

Misc :

ALS Vial : 8 Sample Multiplier: 1

Quant Time: Nov 18 00:36:39 2021

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM111721.M

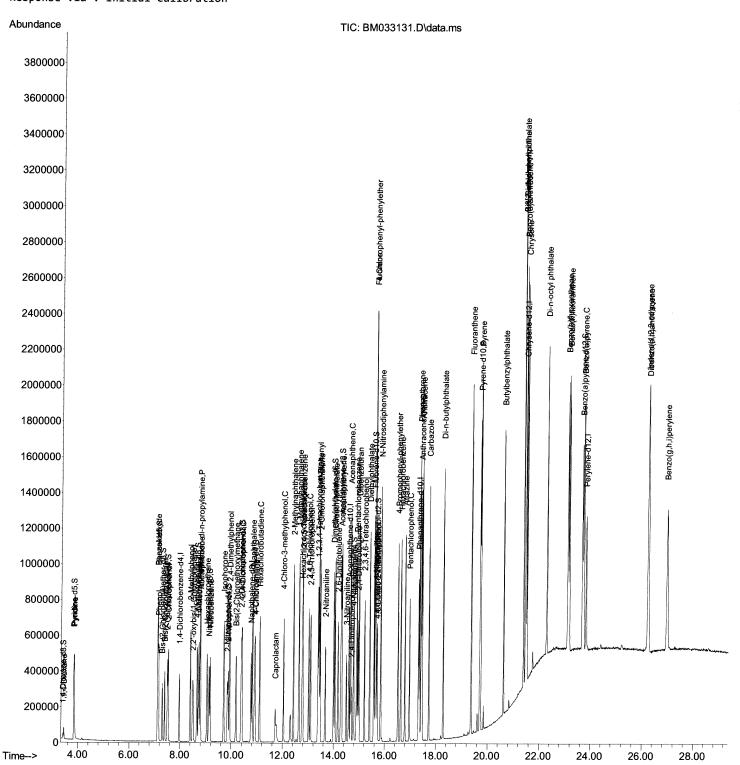
Quant Title : SVOA CALIBRATION

QLast Update : Wed Nov 17 14:14:11 2021 Response via : Initial Calibration



Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/18/2021 Supervised By :mohammad ahmed 11/26/2021



SFAM-EPA-BM111721.M Thu Nov 18 01:11:14 2021

Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111721\

Data File : BM033131.D

Acq On : 17 Nov 2021 22:12

Operator : CG/JU Sample : PB140762BS

Misc

ALS Vial : 8 Sample Multiplier: 1

Quant Time: Nov 18 00:36:39 2021

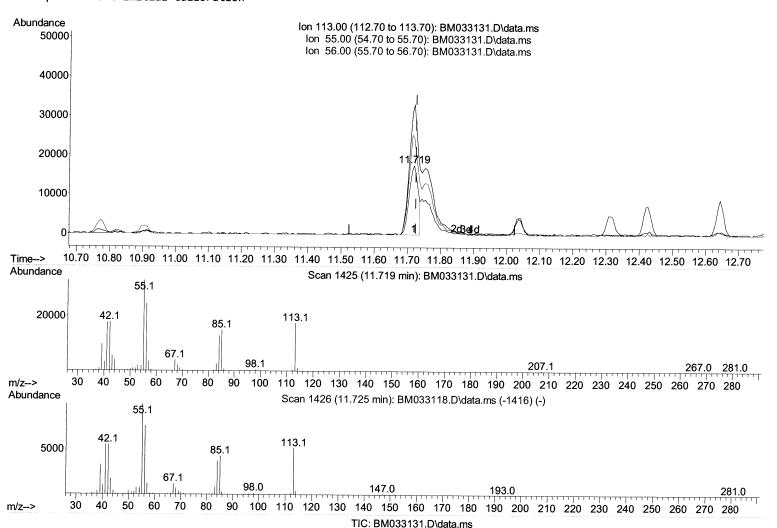
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM111721.M

Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 17 14:14:11 2021
Response via : Initial Calibration



Manual IntegrationsAPPROVED

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(34) Caprolactam

11.719min (-0.006) 18.41 ng/ul

response	33017			
Ion	Ежр%	Act%		
113.00	100.00	100.00		
55.00	196.60	187.74		
56.00	147.80	139.58		
0.00	0.00	0.00		

Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111721\

Data File : BM033131.D

Acq On : 17 Nov 2021 22:12

Operator : CG/JU Sample

: PB140762BS

Misc

ALS Vial : 8 Sample Multiplier: 1

Quant Time: Nov 18 00:36:39 2021

 $\label{thm:lem1} Quant \ \ Method: Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM111721.M$

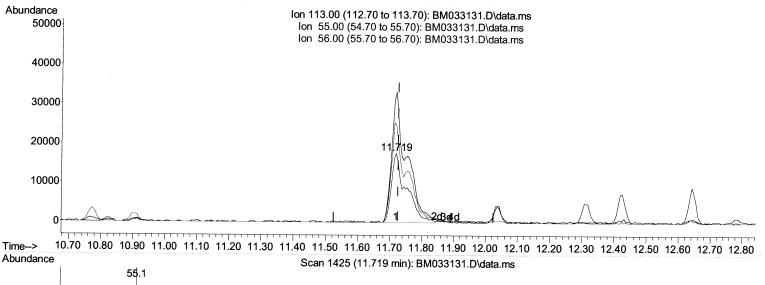
Quant Title : SVOA CALIBRATION

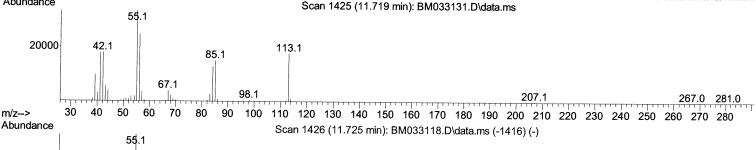
QLast Update : Wed Nov 17 14:14:11 2021 Response via : Initial Calibration

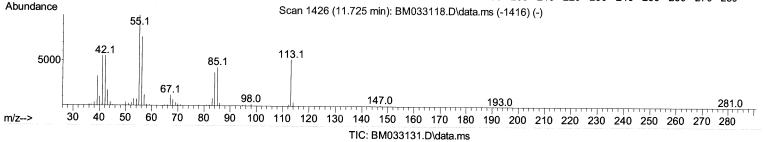


Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/18/2021 Supervised By:mohammad ahmed 11/26/2021







(34) Caprolactam

response	56941		
Ion	Ехр%	Act%	
113.00	100.00	100.00	
55.00	196.60	187.74	
56.00	147.80	139.58	
0.00	0.00	0.00	

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111721\

Data File : BM033131.D

Acq On : 17 Nov 2021 22:12

Operator : CG/JU Sample : PB140762BS

Misc :

ALS Vial : 8 Sample Multiplier: 1

Quant Time: Nov 18 00:36:39 2021

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM111721.M

Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 17 14:14:11 2021
Response via : Initial Calibration

Instrument : BNA_M ClientSampleId : SLCS762

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/18/2021 Supervised By :mohammad ahmed 11/26/2021

Compound	R.T.	QIon	Response	Conc Units Dev	(Min)
Internal Standards					
 1,4-Dichlorobenzene-d4 	7.972	152	96390	20.000 ng/ul	0.00
20) Naphthalene-d8	10.772	136	390790	20.000 ng/ul	0.00
38) Acenaphthene-d10	14.589	164	262603	20.000 ng/ul	0.00
64) Phenanthrene-d10	17.330	188	555913	20.000 ng/ul	0.00
79) Chrysene-d12	21.477	240	543319	20.000 ng/ul	0.00
88) Perylene-d12	23.830	264	533224	20.000 ng/ul	0.00
ystem Monitoring Compounds					
3) 1,4-Dioxane-d8	3.425	96	14316	5.726 ng/uL	0.00
4) Pyridine-d5	3.843	84	189679	27.541 ng/ul	0.00
7) Phenol-d5	7.125	99	237485	29.128 ng/ul	0.00
<pre>9) Bis-(2-Chloroethyl)eth</pre>	7.301	67	149682	28.944 ng/ul	0.00
11) 2-Chlorophenol-d4	7.501	132	183324	29.686 ng/ul	0.00
15) 4-Methylphenol-d8	8.672	113	185044	29.244 ng/ul	0.00
21) Nitrobenzene-d5	9.131	128	88034	31.376 ng/ul	0.00
24) 2-Nitrophenol-d4	9.854	143	90383	32.165 ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.389	165	192475	29.911 ng/ul	0.00
31) 4-Chloroaniline-d4	10.907	131	225098	26.304 ng/ul	0.00
46) Dimethylphthalate-d6	14.001	166	585999	30.423 ng/ul	0.00
49) Acenaphthylene-d8	14.289	160	726339	29.271 ng/ul	0.00
54) 4-Nitrophenol-d4	14.771	143	96407	30.727 ng/ul	0.00
50) Fluorene-d10	15.583	176	515443	29.831 ng/ul	0.00
55) 4,6-Dinitro-2-methylph	15.689	200	83331	31.896 ng/ul	0.00
73) Anthracene-d10	17.424	188	812468	30.319 ng/ul	0.00
31) Pyrene-d10	19.706	212	960187	29.906 ng/ul	0.00
92) Benzo(a)pyrene-d12	23.677	264	883593	30.876 ng/ul	0.00
arget Compounds				0va	lue
2) 1,4-Dioxane	3.460	88	28775	11.335 ng/uL	98
5) Pyridine	3.860	79	200157	28.480 ng/ul	99
Benzaldehyde	7.119	77	153174	32.964 ng/ul	96
8) Phenol	7.154	94	252695	31.123 ng/ul	97
<pre>Description</pre>	7.395	93	199670	30.958 ng/ul	99
.2) 2-Chlorophenol	7.537	128	197795	31.032 ng/ul	98
3) 2-Methylphenol	8.407	108	194133	31.212 ng/ul	99
4) 2,2'-oxybis(1-Chloropr	8.501	45	313197	31.441 ng/ul	98
.6) Acetophenone	8.801	105	312543	31.294 ng/ul	96
7) N-Nitroso-di-n-propyla	8.784	70	172274	32.014 ng/ul	97
<pre>8) 4-Methylphenol</pre>	8.731	108	203672	31.237 ng/ul	99
9) Hexachloroethane	9.054	117	88550	30.496 ng/ul	98
Nitrobenzene	9.178	77	251385	32.105 ng/ul	100
Isophorone	9.707	82	457238	31.949 ng/ul	99
5) 2-Nitrophenol	9.884	139	102191	34.310 ng/ul	97
6) 2,4-Dimethylphenol	9.937	107	241651	30.879 ng/ul	99
7) Bis(2-Chloroethoxy)met	10.178	93	265739	31.202 ng/ul	98
9) 2,4-Dichlorophenol	10.413	162	198211	31.766 ng/ul	97
0) Naphthalene	10.825	128	639551	30.776 ng/ul	99
2) 4-Chloroaniline	10.931	127	235747	27.270 ng/ul	100
3) Hexachlorobutadiene	11.101	225	147430	30.333 ng/ul	98
4) Caprolactam	11.719	113	56941m >	31.742 ng/ul>	11/29/21
5) 4-Chloro-3-methylphenol	12.036	107	220433	32.442 ng/ul	99

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111721\

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Misc :

ALS Vial : 8 Sample Multiplier: 1

Quant Time: Nov 18 00:36:39 2021

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM111721.M

Quant Title : SVOA CALIBRATION

QLast Update : Wed Nov 17 14:14:11 2021 Response via : Initial Calibration

Instrument : BNA_M ClientSampleId : SLCS762

Manual IntegrationsAPPROVED

Reviewed By: Jagrut Upadhyay 11/18/2021 Supervised By: mohammad ahmed 11/26/2021

Compound	R.T.	QIon	Response	Conc Units Dev	(Min)
36) 2-Methylnaphthalene	12.425	142	451284	31.343 ng/ul	 97
37) 1-Methylnaphthalene	12.642			31.015 ng/ul	99
39) 1,2,4,5-Tetrachloroben	12.789			30.198 ng/ul	98
40) Hexachlorocyclopentadiene				27.850 ng/ul	97
41) 2,4,6-Trichlorophenol	13.025			31.404 ng/ul	96
42) 2,4,5-Trichlorophenol	13.095			31.901 ng/ul	99
43) 1,1'-Biphenyl	13.430			30.315 ng/ul	97
44) 2-Chloronaphthalene	13.472		473080	29.982 ng/ul	99
45) 2-Nitroaniline	13.672	65	147045	35.357 ng/ul	96
47) Dimethylphthalate	14.048		603168	32.129 ng/ul	99
48) 2,6-Dinitrotoluene	14.166	165	116070	36.136 ng/ul	93
50) Acenaphthylene	14.319	152	770084	30.582 ng/ul	99
51) 3-Nitroaniline	14.495	138	100424	31.030 ng/ul#	92
52) Acenaphthene	14.654	153	506726	30.848 ng/ul	98
53) 2,4-Dinitrophenol	14.695	184	54169	33.041 ng/ul	92
55) 4-Nitrophenol	14.789	109	102638	32.029 ng/ul	99
56) Dibenzofuran	14.989	168	741886	30.815 ng/ul	
57) 2,4-Dinitrotoluene	14.948	165	165003	_	98 99
58) 2,3,4,6-Tetrachlorophenol	15.207	232	151401	37.352 ng/ul#	
59) Diethylphthalate	15.407	149	618168	33.618 ng/ul	97
61) Fluorene	15.636	166	604266	32.876 ng/ul	98
62) 4-Chlorophenyl-phenyle	15.630	204		31.577 ng/ul	99
63) 4-Nitroaniline	15.654	138	314196 113171	31.457 ng/ul	99
66) 4,6-Dinitro-2-methylph	15.707	198		35.562 ng/ul	96
67) N-Nitrosodiphenylamine	15.842	169	88751 525218	33.894 ng/ul	98
68) 4-Bromophenyl-phenylether	16.518	248		32.123 ng/ul	100
69) Hexachlorobenzene	16.630	284	197326	31.919 ng/ul	96
70) Atrazine	16.789	200	222722	31.448 ng/ul	99
71) Pentachlorophenol	16.789		198167	31.366 ng/ul	99
72) Phenanthrene	17.371	266 178	132228	32.300 ng/ul	98
74) Anthracene	17.460		987125	32.056 ng/ul	99
75) 1,2,3,4-Tetrachloroben	13.395	178 216	985079	31.900 ng/ul	99
76) Pentachlorobenzene	14.907	250	266962	29.237 ng/uL	98
77) Carbazole	17.724	167	262364	29.309 ng/uL	99
78) Di-n-butylphthalate	18.289	149	889491	32.424 ng/ul	100
80) Fluoranthene	19.371	202	1049537	34.987 ng/ul	100
82) Pyrene	19.730	202	1190973	31.663 ng/ul	99
83) Butylbenzylphthalate	20.624	149	1209414	31.667 ng/ul	98
84) 3,3'-Dichlorobenzidine	21.395	252	453539 366290	34.758 ng/ul	96
85) Benzo(a)anthracene	21.395	228	1135625	29.925 ng/ul	99
86) Bis(2-ethylhexyl)phtha	21.403	149		32.316 ng/ul	99
87) Chrysene			657363	35.542 ng/ul	100
89) Di-n-octyl phthalate	21.518	228	1111583	32.391 ng/ul	99
90) Benzo(b)fluoranthene	22.300	149	1098100	33.798 ng/ul	100
91) Benzo(k)fluoranthene	23.112	252	1190620	33.154 ng/ul	100
93) Benzo(a)pyrene	23.165	252	1055974	32.099 ng/ul	99
94) Indeno(1,2,3-cd)pyrene	23.730	252	1118911	32.968 ng/ul	99
95) Dibenzo(a,h)anthracene	26.235	276	1222966	32.358 ng/ul	97
96) Benzo(g,h,i)perylene	26.253 26.977	278	1044690	32.309 ng/ul	98
	20.9//	276	1055918	32.189 ng/ul	98

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