(QT Reviewed) Quantitation Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\

Data File : BG051353.D

: 6 Dec 2021 16:59 Acq On

Operator : CG/JU : SSTDCCC020 Sample

Misc

Sample Multiplier: 1 : 10 ALS Vial

Quant Time: Dec 06 17:48:26 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

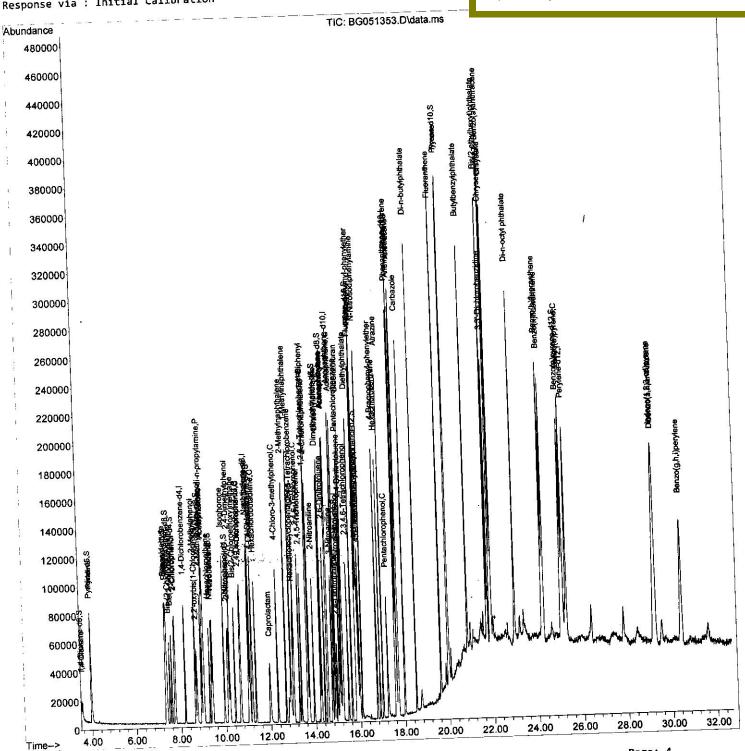
Quant Title : SVOA CALIBRATION

QLast Update : Fri Dec 03 15:23:09 2021

Response via : Initial Calibration

Instrument: BNA_G **LabSampleId**: STDCCC020

Manual IntegrationsAPPROVED



SFAM-EPA-BG112321.M Mon Dec 06 17:52:48 2021

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\

Data File : BG051353.D

Acq On : 6 Dec 2021 16:59

Operator : CG/JU Sample : SSTDCCC020

Misc

ALS Vial : 10 Sample Multiplier: 1

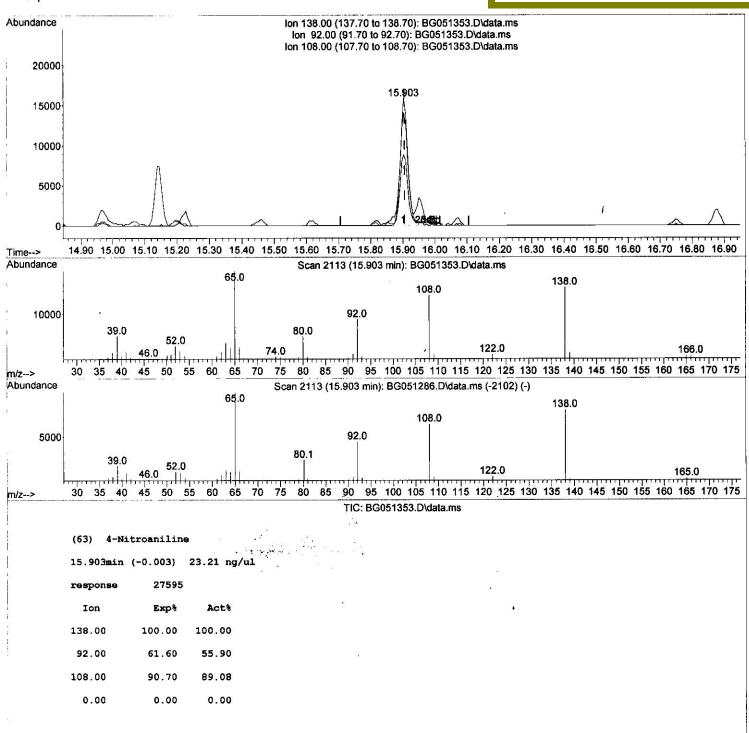
Quant Time: Dec 06 17:48:26 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

Quant Title : SVOA CALIBRATION
QLast Update : Fri Dec 03 15:23:09 2021
Response via : Initial Calibration

Instrument :
BNA_G
LabSampleId :
SSTDCCC020

Manual IntegrationsAPPROVED



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\

Data File : BG051353.D

Acq On : 6 Dec 2021 16:59

Operator : CG/JU Sample : SSTDCCC020

Misc

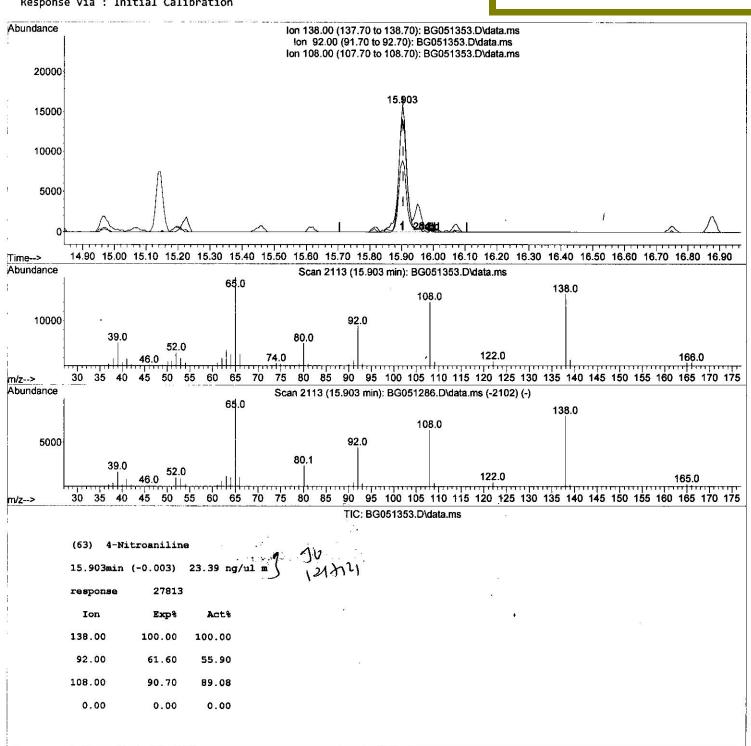
ALS Vial : 10 Sample Multiplier: 1

Quant Time: Dec 06 17:48:26 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

Quant Title : SVOA CALIBRATION QLast Update : Fri Dec 03 15:23:09 2021 Response via : Initial Calibration Instrument :
BNA_G
LabSampleId :
SSTDCCC020

Manual IntegrationsAPPROVED



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\

Data File : BG051353.D

: 6 Dec 2021 16:59 Acq On

; CG/JU Operator : SSTDCCC020 Sample

Misc

Sample Multiplier: 1 ALS Vial : 10

Quant Time: Dec 06 17:48:26 2021

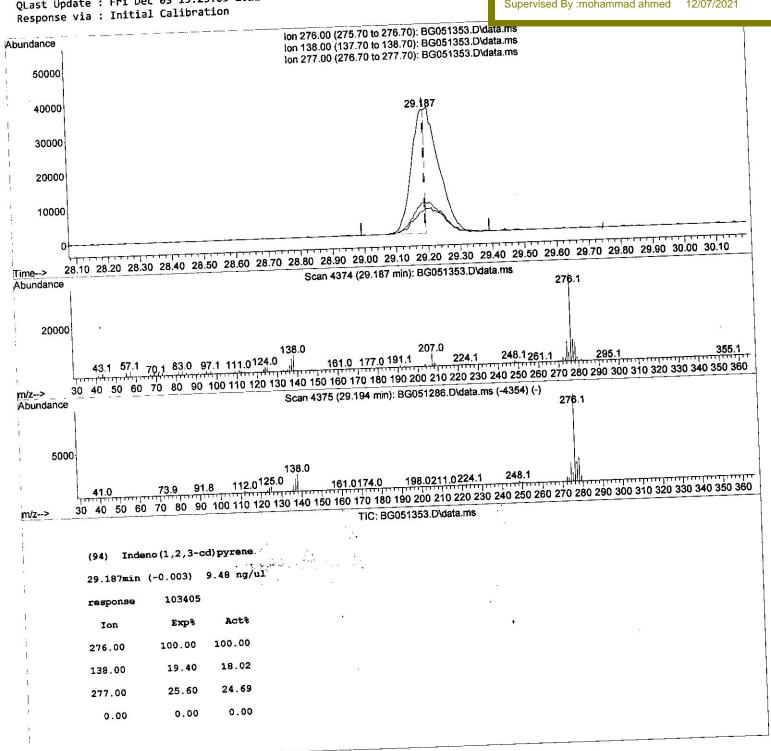
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

Quant Title : SVOA CALIBRATION

QLast Update : Fri Dec 03 15:23:09 2021

Instrument: BNA_G LabSampleId : SSTDCCC020

Manual IntegrationsAPPROVED



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\

Data File: BG051353.D

Acq On : 6 Dec 2021 16:59

Operator : CG/JU Sample : SSTDCCC020

Misc

ALS Vial : 10 Sample Multiplier: 1

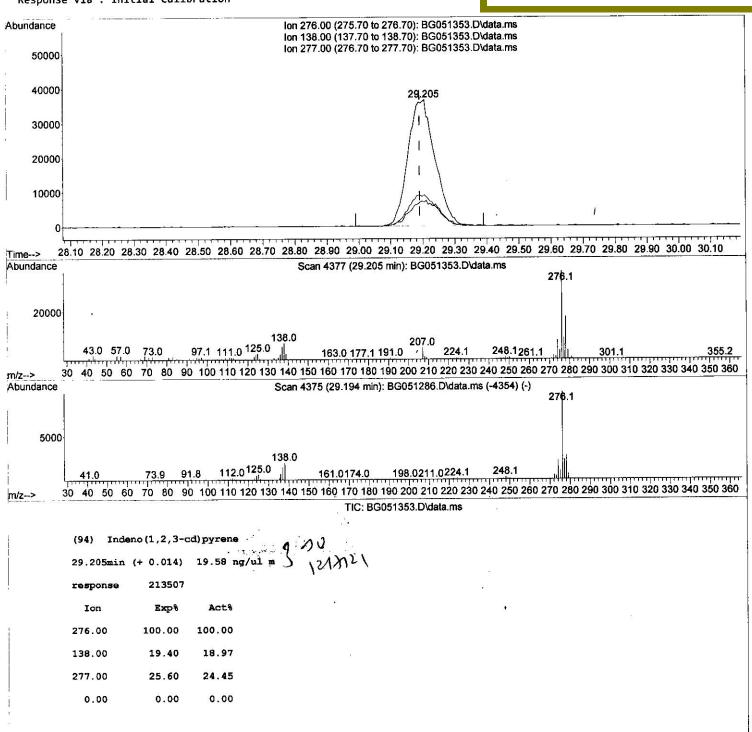
Quant Time: Dec 06 17:48:26 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

Quant Title : 5VOA CALIBRATION

QLast Update : Fri Dec 03 15:23:09 2021 Response via : Initial Calibration Instrument:
BNA_G
LabSampleId:
SSTDCCC020

Manual IntegrationsAPPROVED



Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\

Data File : BG051353.D

Acq On : 6 Dec 2021 16:59

Operator : CG/JU Sample : SSTDCCC020

Misc :

ALS Vial : 10 Sample Multiplier: 1

Quant Time: Dec 06 17:48:26 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

Quant Title : SVOA CALIBRATION QLast Update : Fri Dec 03 15:23:09 2021 Response via : Initial Calibration Instrument:
BNA_G
LabSampleId:
SSTDCCC020

Manual IntegrationsAPPROVED

Compound	R.T. (lon	Response	Conc Units Dev	Min)
Tabana Chandande					
Internal Standards 1) 1,4-Dichlorobenzene-d4	8.194	152	23950	20.000 ng/ul	0.00
20) Naphthalene-d8	11.020	136	106399	20.000 ng/ul	0.00
38) Acenaphthene-d10	14.827	164	75551	20.000 ng/ul	0.00
64) Phenanthrene-d10	17.577	188	175821	20.000 ng/ul	0.00
79) Chrysene-d12	21.878	240	150962	20.000 ng/ul	0.00
88) Perylene-d12	25.274	264	151373	20.000 ng/ul	0.00
88) Perylene-uiz					
System Monitoring Compounds				*** ******* • ***	
3) 1,4-Dioxane-d8	3.529	96	5104	7.406 ng/uL	-0.02
4) Pyridine-d5	3.958	84	35179	17.395 ng/ul	-0.02
7) Phenol-d5	7.354	99	43900	18.546 ng/ul	0.00
Bis-(2-Chloroethyl)eth	7.507	67	27991	18.829 ng/ul	0.00
11) 2-Chlorophenol-d4	7.724	132	32938	19.324 ng/ul	0.00
<pre>15) 4-Methylphenol-d8</pre>	8.911	113	36601	19.161 ng/ul	0.00
21) Nitrobenzene-d5	9.375	128	17617	19.614 ng/ul	0.00
24) 2-Nitrophenol-d4	10.098	143	20166	19.904 ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.644	165	34052	19.809 ng/ul	0.00
31) 4-Chloroaniline-d4	11.161	131	49572	19.708 ng/ul	0.00
46) Dimethylphthalate-d6	14.216	166	113972	19.606 ng/ul	0.00
49) Acenaphthylene-d8	14.522	160	142688	19.465 ng/ul	0.00
54) 4-Nitrophenol-d4	15.051	143	17824	18.942 ng/ul	0.00
60) Fluorene-d10	15.814	176	101596	19.408 ng/ul	0.00
65) 4,6-Dinitro-2-methylph	15.955	200	19160	17.660 ng/ul	0.00
73) Anthracene-d10	17.677	188	167902	19.967 ng/ul	0.00
81) Pyrene-d10	19.957	212	187323	20.508 ng/ul	0.00
92) Benzo(a)pyrene-d12	25.039	264	155074	19.182 ng/ul	0.00
Target Compounds				Qv	alue
2) 1,4-Dioxane	3.564	88	5626	7.238 ng/uL	90
5) Pyridine	3.981	79	38614	18.349 ng/ul	99
6) Benzaldehyde	7.324	77	33346	22.121 ng/ul	96
8) Phenol	7.383	94	46035	18.773 ng/ul	97
10) Bis(2-Chloroethyl)ether	7.601	93	34952	18.841 ng/ul	94
12) 2-Chlorophenol	7.759	128	33577	19.331 ng/ul	92
13) 2-Methylphenol	8.641	108	33771	18.489 ng/ul	97
14) 2,2'-oxybis(1-Chloropr	8.711	45	51722	19.321 ng/ul	98
16) Acetophenone	9.022		56620	19.164 ng/ul	96
17) N-Nitroso-di-n-propyla	8.993	70	34033	20.045 ng/ul	98
18) 4-Methylphenol	8.975	108	37559	19.230 ng/ul	94
19) Hexachloroethane	9.275	117	13615	18.558 ng/ul	93
22) Nitrobenzene	9.410	77	47199	20.041 ng/ul	98
23) Isophorone	9.927	82	92229	20.157 ng/ul	100
25) 2-Nitrophenol	10.127	139	20588	19.618 ng/ul	95
26) 2,4-Dimethylphenol	10.180	107	42884	19.987 ng/ul	96
27) Bis(2-Chloroethoxy)met	10.409	93	49405	19.559 ng/ul	96
29) 2,4-Dichlorophenol	10.673	162	33169	19.602 ng/ul	96
30) Naphthalene	11.067	128		19.439 ng/ul	97
32) 4-Chloroaniline	11.185	127		19.893 ng/ul	97
33) Hexachlorobutadiene	11.331			19.318 ng/ul	98
34) Caprolactam	11.966	113		21.086 ng/ul	97
35) 4-Chloro-3-methylphenol	12.301	107		20.011 ng/ul	95
,,,				MAKENE WINDOWS TRANS AND ASSESSED.	

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Operator : CG/JU Sample : SSTDCCC020

Misc :

ALS Vial : 10 Sample Multiplier: 1

Quant Time: Dec 06 17:48:26 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

Quant Title : SVOA CALIBRATION

QLast Update : Fri Dec 03 15:23:09 2021 Response via : Initial Calibration Instrument:
BNA_G
LabSampleId:
SSTDCCC020

Manual IntegrationsAPPROVED

	Compound	R.T.	QIon	Response	Conc Units Dev(M	1in)
261	2-Methylnaphthalene	12.665	142	78288	19.881 ng/ul	99
	1-Methylnaphthalene	12.883	142	80288	19.818 ng/ul	99
	1,2,4,5-Tetrachloroben	13.024	216	44374	18.709 ng/ul	97
	Hexachlorocyclopentadiene	12.988	237	21025	21.931 ng/ul#	97
	2,4,6-Trichlorophenol	13.270	196	29006	19.488 ng/ul	97
	2,4,5-Trichlorophenol	13.359	HEROLINE.	30716	19.707 ng/ul	95
	1,1'-Biphenyl	13.658	154	107657	19.078 ng/ul	97
	2-Chloronaphthalene	13.711	162	84259	18.771 ng/ul	100
	2-Nitroaniline	13.917		31790	20.463 ng/ul	90
2000	Dimethylphthalate	14.263	163	116358	19.775 ng/ul	100
	2,6-Dinitrotoluene	14.404	165	24130	19.523 ng/ul	92
100000	Acenaphthylene	14.551	152	139264	19.229 ng/ul	98
	3-Nitroaniline	14.745	138	25469	20.847 ng/ul	98
	Acenaphthene	14.892	153	91280	19.111 ng/ul	97
	2,4-Dinitrophenol	14.963		15531	22.733 ng/ul	89
2000000	4-Nitrophenol	15.068		18716	22.928 ng/ul	94
	Dibenzofuran	15.221		133374	19.360 ng/ul	99
000000000000000000000000000000000000000	2,4-Dinitrotoluene	15.198		36206	20.509 ng/ul#	92
	2,3,4,6-Tetrachlorophenol	15.456		24386	19.924 ng/ul	96
	Diethylphthalate	15.615	149	126403	20.465 ng/ul	98
	Fluorene	15.873		108383	19.641 ng/ul	100
155	4-Chlorophenyl-phenyle	15.856		57017	19.173 ng/ul	94
	4-Nitroaniline	15.903		27813m	23.394 ng/ul	
	4,6-Dinitro-2-methylph	15.967		17924	17.130 ng/ul	98
	N-Nitrosodiphenylamine	16.073		98145	19.499 ng/ul	95
	4-Bromophenyl-phenylether	16.749		35663	18.926 ng/ul	95
	Hexachlorobenzene	16.878		37745	19.644 ng/ul	97
tanana Sana	Atrazine	17.013		44076	20.836 ng/ul	99
Liver Co.	Pentachlorophenol	17.236		18675	21.934 ng/ul	92
100000000000000000000000000000000000000	Phenanthrene	17.618		190032	19.575 ng/ul	99
	Anthracene	17.712		193329	20.052 ng/ul	98
	1,2,3,4-Tetrachloroben	13.635		47356	18.466 ng/uL	98
200	Pentachlorobenzene	15.145		44478	18.614 ng/uL	98
	Carbazole	17.983		176333	20.836 ng/ul	99
	Di-n-butylphthalate	18.505		228233	20.916 ng/ul	99
	Fluoranthene	19.622		234133	- 12 Villa	95
	Pyrene	19.986	. 202	227127	20.696 ng/ul	95
	Butylbenzylphthalate	20.844		90398	19.813 ng/ul	94
50000000000	3,3'-Dichlorobenzidine	21.760		71163	20.247 ng/ul	99
0.00	Benzo(a)anthracene	21.854		198941	19.430 ng/ul	98
1000000	Bis(2-ethylhexyl)phtha	21.713		127883	19.479 ng/ul	98
87.0	Chrysene	21.925	228	187717	19.084 ng/ul	99
	Di-n-octyl phthalate	22.977	149	214299	19.541 ng/ul	100
	Benzo(b)fluoranthene	24.187		193745	18.966 ng/ul	97
	Benzo(k)fluoranthene	24.258		184846	/19.282 ng/ul	99
	Benzo(a)pyrene	25.115		190151	/ 19.511 ng/ul	97
30.00	Indeno(1,2,3-cd)pyrene	29.205		213507m	19.577 ng/ul	
	Dibenzo(a,h)anthracene	29.246		181862	, 19.656 ng/ul	97
	Benzo(g,h,i)perylene	30.415		179097	19.518 ng/ul	95

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