Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112921\

Data File : BG051271.D

Acq On : 29 Nov 2021 14:16

Operator : CG/JU

Sample : SSTDCCC020EC

Misc :

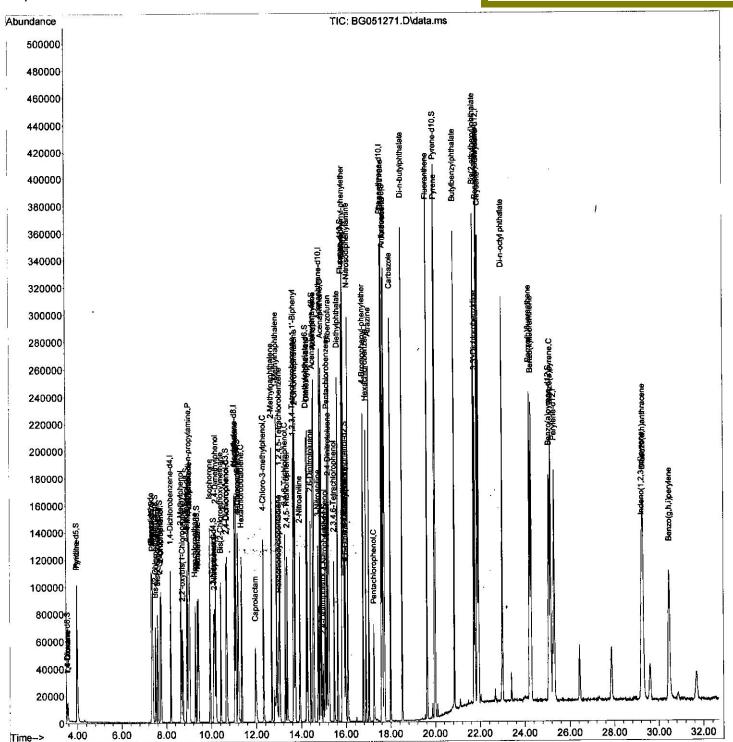
ALS Vial : 2 Sample Multiplier: 1

Quant Time: Nov 29 14:59:35 2021

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

Quant Title : SVOA CALIBRATION QLast Update : Wed Nov 24 06:04:50 2021 Response via : Initial Calibration Instrument:
BNA_G
LabSampleId:
SSTDCCC020EC

Manual IntegrationsAPPROVED



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112921\

Data File : BG051271.D

Acq On : 29 Nov 2021 14:16

Operator : CG/JU

Sample : SSTDCCC020EC

Misc

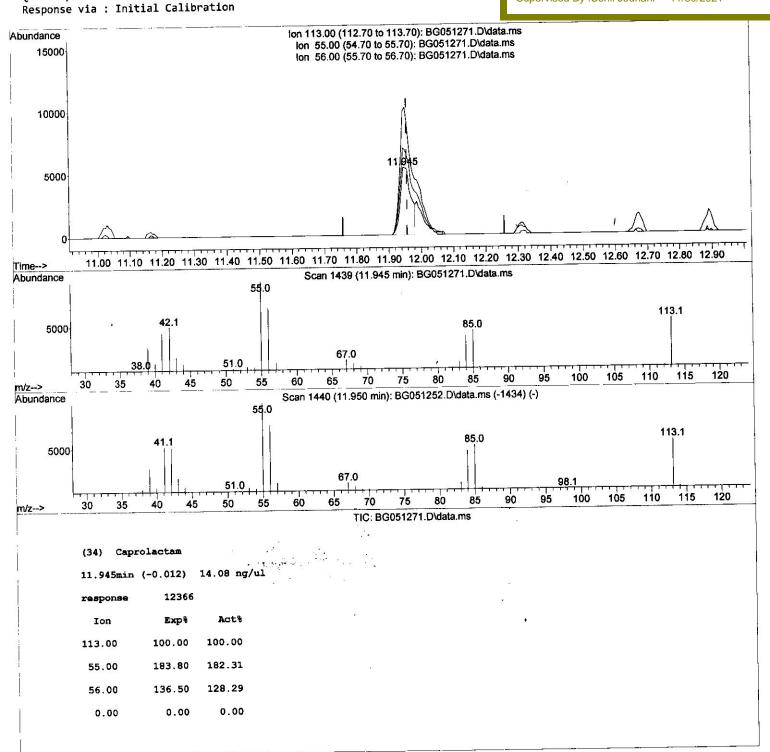
ALS Vial : 2 Sample Multiplier: 1

Quant Time: Nov 29 14:59:35 2021

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

Quant Title : SVOA CALIBRATION QLast Update : Wed Nov 24 06:04:50 2021 Instrument:
BNA_G
LabSampleId:
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Manual IntegrationsAPPROVED



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Acq On : 29 Nov 2021 14:16

Operator : CG/JU

Sample : SSTDCCC020EC

Misc

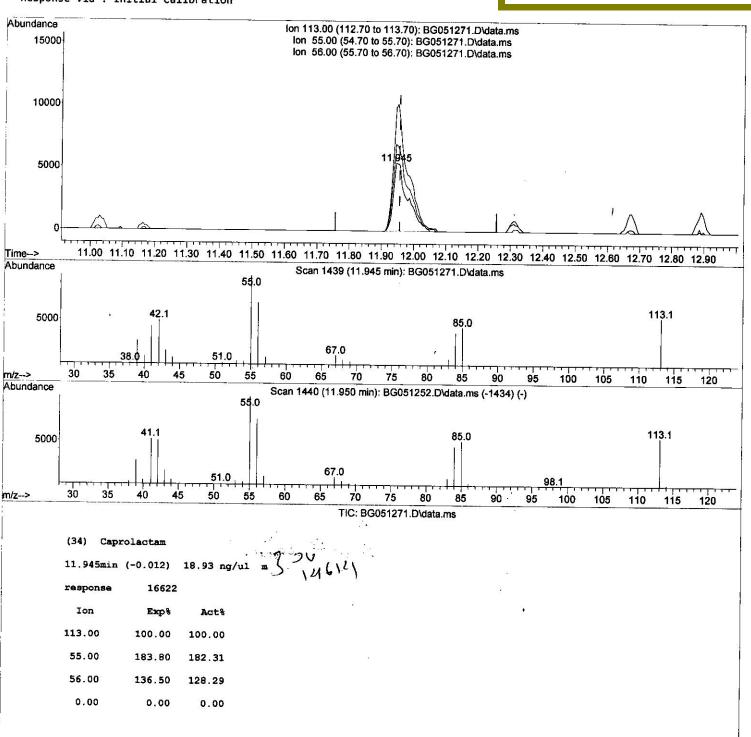
ALS Vial : 2 Sample Multiplier: 1

Quant Time: Nov 29 14:59:35 2021

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

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Manual IntegrationsAPPROVED



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Misc

ALS Vial : 2 Sample Multiplier: 1

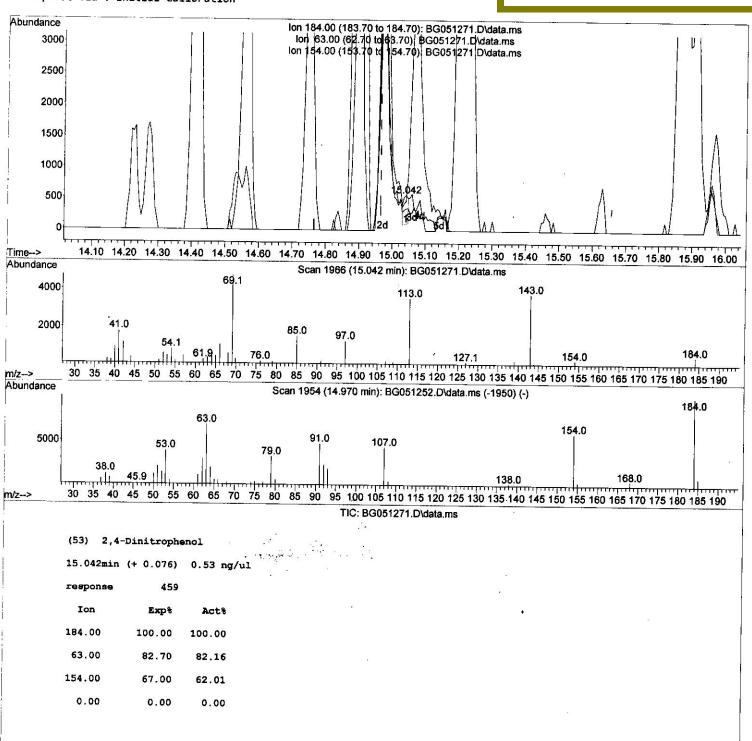
Quant Time: Nov 29 14:59:35 2021

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

Quant Title : SVOA CALIBRATION
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Response via : Initial Calibration

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Manual IntegrationsAPPROVED



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112921\

Data File : BG051271.D

Acq On : 29 Nov 2021 14:16

Operator : CG/JU Sample : SSTDCCC020EC

Misc :

ALS Vial : 2 Sample Multiplier: 1

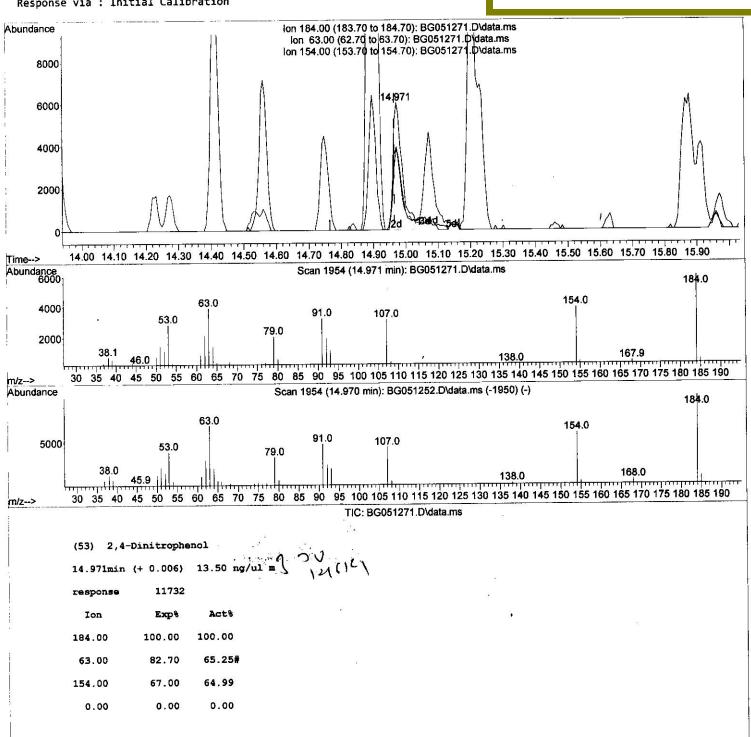
Quant Time: Nov 29 14:59:35 2021

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

Quant Title : SVOA CALIBRATION
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Response via : Initial Calibration

Instrument: BNA_G LabSampleId: SSTDCCC020EC

Manual IntegrationsAPPROVED



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

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112921\

Data File : BG051271.D

Acq On : 29 Nov 2021 14:16

Operator : CG/JU

Sample : SSTDCCC020EC

Compound

Misc

ALS Vial : 2 Sample Multiplier: 1

Quant Time: Nov 29 14:59:35 2021

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

Quant Title : SVOA CALIBRATION QLast Update : Wed Nov 24 06:04:50 2021 Response via : Initial Calibration Instrument: BNA_G LabSampleId: SSTDCCC020EC

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/30/2021 Supervised By :Sohil Jodhani 11/30/2021

	rnal Standards			D4055	20.000	1-1	6 65
	1,4-Dichlorobenzene-d4	8.203	152	31963	20.000	2000 200 50	0.00
September 1	Naphthalene-d8	11.029	136	140439	20.000		0.00
	Acenaphthene-d10	14.836	164	96136	20.000		0.00
	Phenanthrene-d10	17.586	188	214044	20.000		0.00
,	Chrysene-d12	21.887	240	183473	20.000	4000 - 100	0.00
88)	Perylene-d12	25.300	264	184916	20,000	ng/ul	0.02
	em Monitoring Compounds			een.	7 400	was foot	0.00
	1,4-Dioxane-d8	3.538	96	6604	7.180		0.00
	Pyridine-d5	3.967	84	48210	17.862	_	-0.01
	Phenol-d5	7.357	99	55913	17.699		0.00
	Bis-(2-Chloroethyl)eth	7.515	67	36633	18.464		0.00
	2-Chlorophenol-d4	7.727	132	42224	18.562		0.00
	4-Methylphenol-d8	8.914	113	45256	17.753		0.00
	Nitrobenzene-d5	9.378	128	21911	18.482		0.00
	2-Nitrophenol-d4	10.106	143	24977	18.677		0.00
	2,4-Dichlorophenol-d3	10.653	165	42431	18.701	2011	0.00
	4-Chloroaniline-d4	11.170	131	60004	18.074		0.00
46)	Dimethylphthalate-d6	14.225	166	136500	18.453		0.00
	Acenaphthylene-d8	14.531	160	175724	18.839		0.00
	4-Nitrophenol-d4	15.054	143	1769 9	14.782		0.00
50)	Fluorene-d10	15.823	176	122972	18.461		0.00
55)	4,6-Dinitro-2-methylph	15.958	200	19829	15.013		0.00
73)	Anthracene-d10	17.686	188	189908	18.551		0.00
81)	Pyrene-d10	19.965	212	210054	18.921		0.00
2)	Benzo(a)pyrene-d12	25.065	264	180142	18.241	ng/ul	0.02
	et Compounds					(Carlo)	ralue
	1,4-Dioxane	3.573	88	6985		ng/uL#	
	Pyridine	3.990	79	50623	18.025		98
6)	Benzaldehyde	7.333	77	42604	21.177	The second secon	99
	Phenol	7.386	94	57423	17.547		100
10)	Bis(2-Chloroethyl)ether	7.609	93	44956	18.158	110000000000000000000000000000000000000	96
12)	2-Chlorophenol	7.762	128	42017	18.126		92
	2-Methylphenol	8.649	108	43761			96
	2,2'-oxybis(1-Chloropr	8.726	45	67024	18.760	_	99
16)	Acetophenone	9.031	-	70568	17.897		97
	N-Nitroso-di-n-propyla	9.002	70	42073	18.568	•	96
	4-Methylphenol	8.978	108	47114	18.075		95
19)	Hexachloroethane	9.284	117	17091	17.455	35.50	98-
	Nitrobenzene	9.419	77	59631	19.183		99
23)	Isophorone	9.936	82	111815		ng/ul	100
25)	2-Nitrophenol	10.142	139	25345		ng/ul	98
	2,4-Dimethylphenol	10.189	107	54465		ng/ul	99
	Bis(2-Chloroethoxy)met	10.418	93	62576		ng/ul	100
	2,4-Dichlorophenol	10.682	162	41560	18.607	ng/ul	95
	Naphthalene	11.082	128	140670	18.409	ng/ul	98
	4-Chloroaniline	11.193	127	61550	18.467	ng/ul	97
	Hexachlorobutadiene	11.340	225	28114 1		ng/ul	98
	Caprolactam	11.945	113	16622m \		ng/ul	
341	Capiotaciali						

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Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112921\

Data File : BG051271.D

Acq On : 29 Nov 2021 14:16

Operator : CG/JU Sample : SSTDCCC020EC

Misc

ALS Vial : 2 Sample Multiplier: 1

Quant Time: Nov 29 14:59:35 2021

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

Quant Title : SVOA CALIBRATION QLast Update : Wed Nov 24 06:04:50 2021 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc Units Dev	(Min)
36) 2-Methylnaphthalene	12.674	142	97417	10 743 (-1	
37) 1-Methylnaphthalene	12.891	142	98895	18.742 ng/ul 18.494 ng/ul	99
39) 1,2,4,5-Tetrachloroben	13.032		55347	18.338 ng/ul	99
40) Hexachlorocyclopentadiene		237	18104	14.841 ng/ul	96 94
41) 2,4,6-Trichlorophenol	13.279	196	35258	18.616 ng/ul	98
42) 2,4,5-Trichlorophenol	13.361	196	35432	17.865 ng/ul	98
43) 1,1'-Biphenyl	13.667	154	132862	18.503 ng/ul	97
44) 2-Chloronaphthalene	13.720	162	104728	18.335 ng/ul	98
45) 2-Nitroaniline	13.925	65	38383	19.417 ng/ul	91
47) Dimethylphthalate	14.272	163	136301	18.204 ng/ul	100
48) 2,6-Dinitrotoluene	14.413	165	29261	18.605 ng/ul	93
50) Acenaphthylene	14.560	152	172104	18.675 ng/ul	98
51) 3-Nitroaniline	14.748	138	31337	20.157 ng/ul	92
52) Acenaphthene	14.901	153	112366		97
53) 2,4-Dinitrophenol	14.971	184	11732m	\ 13.495 ng/ul	
55) 4-Nitrophenol	15.071	109	21025	20.242 ng/ul	95
56) Dibenzofuran	15.230	168	161478	18.420 ng/ul	99
57) 2,4-Dinitrotoluene	15.206	165	41065	18.281 ng/ul#	97
58) 2,3,4,6-Tetrachlorophenol	15.465	232	25493	16.368 ng/ul	98
59) Diethylphthalate	15.629	149	144190	18.346 ng/ul	98
61) Fluorene	15.882	166	131420	18.716 ng/ul	99
62) 4-Chlorophenyl-phenyle	15.864	204	68393	18.073 ng/ul	99
63) 4-Nitroaniline	15.911	138	31890	21.079 ng/ul	93
66) 4,6-Dinitro-2-methylph	15.976	198	19158	15.040 ng/ul#	96
67) N-Nitrosodiphenylamine	16.082	169	116781	19.058 ng/ul	97
68) 4-Bromophenyl-phenylether	16.757	248	42361	18.466 ng/ul	93
69) Hexachlorobenzene	16.887	284	43335	18.526 ng/ul	98
70) Atrazine	17.022	200	47958	18.622 ng/ul	97
71) Pentachlorophenol	17.245	266	14437	13.928 ng/ul	96
72) Phenanthrene	17.627	178	219852	18.603 ng/ul	99
74) Anthracene	17.721	178	219999	18.744 ng/ul	97
75) 1,2,3,4-Tetrachloroben	13.638	216	57957	18.564 ng/uL	96
76) Pentachlorobenzene	15.153	250	54094	18.595 ng/uL	98
77) Carbazole	17.991	167	193354	18.767 ng/ul	98
78) Di-n-butylphthalate	18.514	149	251811	18.956 ng/ul	99
80) Fluoranthene	19.631	202	257142	18.859 ng/ul	98
82) Pyrene	19.995	202	255895	19.186 ng/ul	97
83) Butylbenzylphthalate84) 3,3'-Dichlorobenzidine			107254	19.342 ng/ul	95
85) Benzo(a)anthracene	21.775	252	81946	19.183 ng/ul	99
86) Ris(2-o+bylboxyl)=b+b-	21.869	228	230519	18.524 ng/ul	98
86) Bis(2-ethylhexyl)phtha87) Chrysene	21.728	149	153073	19.184 ng/ul	98
	21.940	228	220122	18.413 ng/ul	98
<pre>89) Di-n-octyl phthalate 90) Benzo(b)fluoranthene</pre>	22.991	149	258555	19.300 ng/ul	100
91) Benzo(k)fluoranthene	24.208	252	230783	18.493 ng/ul	98
93) Benzo(a)pyrene	24.278	252	210040	17.936 ng/ul	98
94) Indeno(1,2,3-cd)pyrene	25.136	252	220241	18.499 ng/ul	97
95) Dibenzo(a,h)anthracene	29.219	276	240536	18.055 ng/ul	98
96) Benzo(g,h,i)perylene	29.272	278	204078	18.056 ng/ul	96
	30.459	276	193307	17.246 ng/ul	97

Manual IntegrationsAPPROVED

Instrument: BNA_G **LabSampleld**: SSTDCCC020EC

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