Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120221\

Data File : BG051343.D

Acq On : 4 Dec 2021 00:54

Operator : CG/JU Sample : SSTDCCC020EC

Misc

ALS Vial : 48 Sample Multiplier: 1

Quant Time: Dec 04 02:28:16 2021

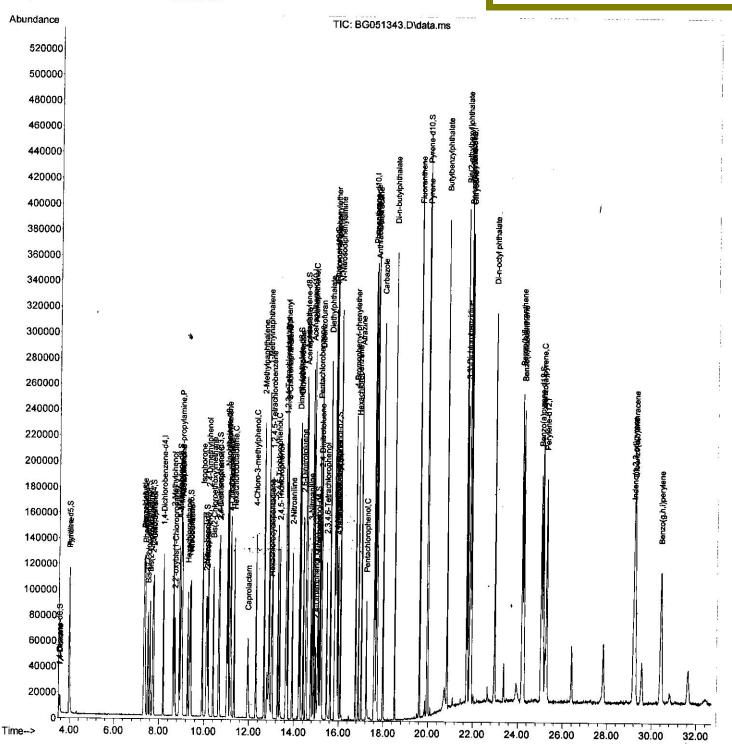
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

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



Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/06/2021 Supervised By :mohammad ahmed 12/07/2021



SFAM-EPA-BG112321.M Sat Dec 04 02:31:48 2021

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Acq On 4 Dec 2021 00:54

Operator : CG/JU Sample : SSTDCCC020EC

Misc

ALS Vial : 48 Sample Multiplier: 1

Quant Time: Dec 04 02:28:16 2021

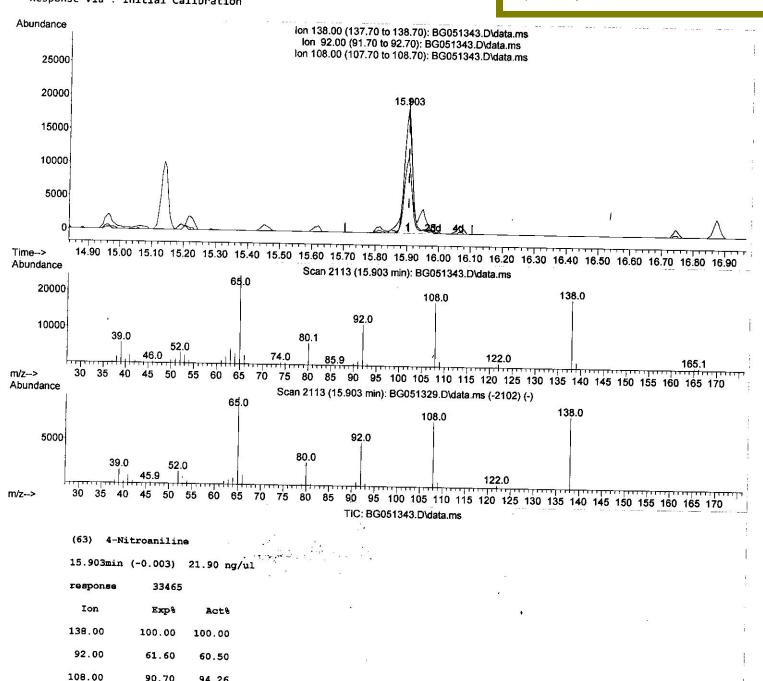
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

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

Instrument: BNA_G LabSampleId: STDCCC020E

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/06/2021 Supervised By :mohammad ahmed 12/07/2021



0.00

90.70

0.00

94.26

0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120221\

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Acq On : 4 Dec 2021 00:54

Operator : CG/JU

Sample : SSTDCCC020EC

Misc

ALS Vial : 48 Sample Multiplier: 1

Quant Time: Dec 04 02:28:16 2021

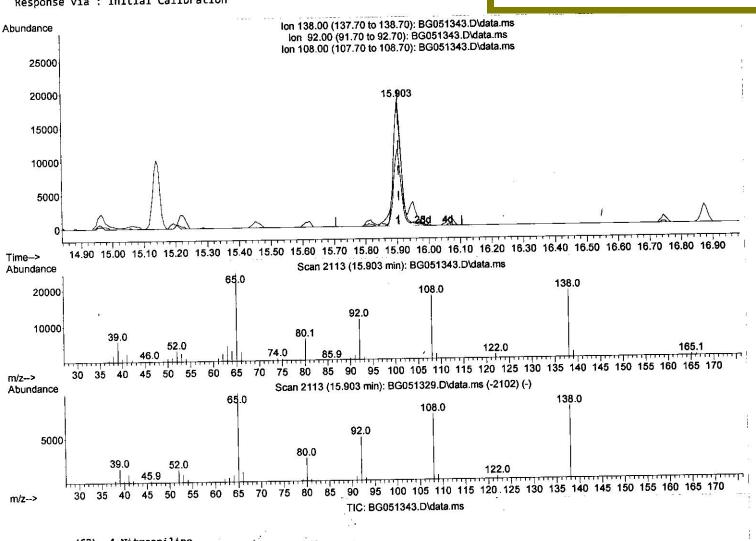
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

Quant Title : SVOA CALIBRATION

QLast Update : Fri Dec 03 15:23:09 2021 Response via : Initial Calibration **Instrument :** BNA_G **LabSampleld :** SSTDCCC020EC

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/06/2021 Supervised By :mohammad ahmed 12/07/2021



(63) 4-Nitroaniline

15.903min (-0.003) 22.00 ng/ul m \ \2/3/2\

response	33609	
Ion	Exp%	Act*
138.00	100.00	100.00
92.00	61.60	60.50
108.00	90.70	94.26
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120221\

Data File : BG051343.D : 4 Dec 2021 00:54 Acq On

: CG/JU Operator

Sample : SSTDCCC020EC

Misc

ALS Vial : 48 Sample Multiplier: 1

Quant Time: Dec 04 02:28:16 2021

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

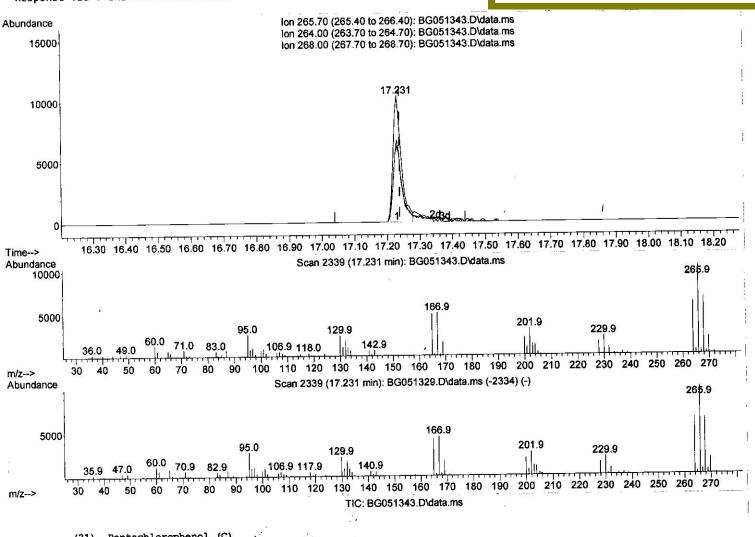
Quant Title : SVOA CALIBRATION

QLast Update : Fri Dec 03 15:23:09 2021 Response via : Initial Calibration

Instrument: BNA_G LabSampleId: STDCCC020EC

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/06/2021 Supervised By:mohammad ahmed 12/07/2021



(71) Pentachlorophenol (C)

17.231min (-0.009) 17.82 ng/ul

response	18281	
Ion	Exp8	Act*
265.70	100.00	100.00
264.00	67.90	59.92
268.00	63.80	64.76
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120221\

Data File : BG051343.D

: 4 Dec 2021 00:54 Acq On

Operator : CG/JU

: SSTDCCC020EC Sample

Misc

Sample Multiplier: 1 ALS Vial : 48

Quant Time: Dec 04 02:28:16 2021

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

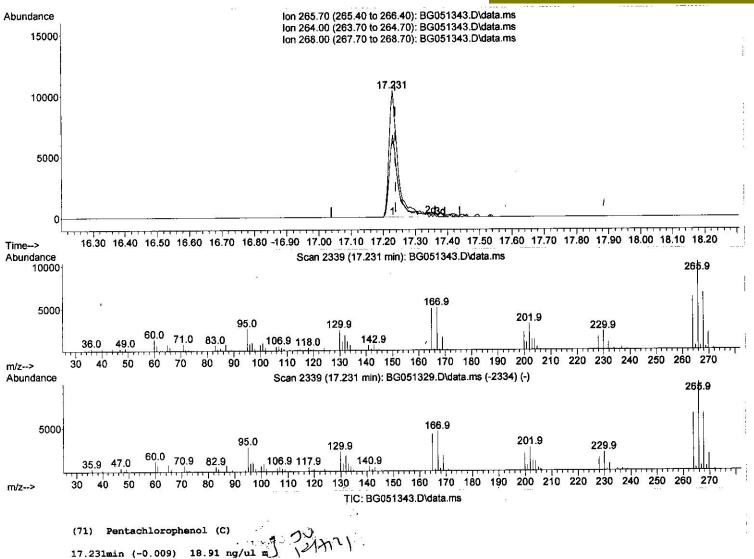
Quant Title : SVOA CALIBRATION

QLast Update : Fri Dec 03 15:23:09 2021 Response via : Initial Calibration

Instrument: BNA_G L<u>ab</u>Sampleld : STDCCC020E0

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/06/2021 Supervised By:mohammad ahmed 12/07/2021



18.91 ng/ul m

response	19394	
Ion	Exp%	Act*
265.70	100.00	100.00
264.00	67.90	59.92
268.00	63.80	64.76
0.00	0.00	0.00

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Data File : BG051343.D

Acq On : 4 Dec 2021 00:54

Operator : CG/JU

Sample : SSTDCCC020EC

Misc

ALS Vial : 48 Sample Multiplier: 1

Quant Time: Dec 04 02:28:16 2021

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

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QLast Update : Fri Dec 03 15:23:09 2021
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Instrument: BNA_G **LabSampleld:** SSTDCCC020EC

Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 12/06/2021 Supervised By :mohammad ahmed 12/07/2021

Compound	R.T.	QIon	Response	Conc Units Dev	(Min)
Tabana 1 Standard					
Internal Standards	0 100	450	24500		
1) 1,4-Dichlorobenzene-d4	8.188	152	34608	20.000 ng/ul	-0.01
<pre>20) Naphthalene-d8 38) Acenaphthene-d10</pre>	11.020		150736	20.000 ng/ul	0.00
[11] [12] [13] [13] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] - [14] -	14.822		97090	20.000 ng/ul	0.00
64) Phenanthrene-d10 79) Chrysene-d12	17.572		211828	20.000 ng/ul	0.00
88) Perylene-d12	21.872		184149	20.000 ng/ul	0.00
88) Perylene-ulz	25.274	264	184684	20.000 ng/ul	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.529	96	7523	7.554 ng/uL	-0.01
4) Pyridine-d5	3.964	84	54147	18.529 ng/ul	-0.01
7) Phenol-d5	7.354	99	63470	18.556 ng/ul	0.00
<pre>9) Bis-(2-Chloroethyl)eth</pre>	7.507	67	40635	18.916 ng/ul	0.00
11) 2-Chlorophenol-d4	7.724	132	46692	18.957 ng/ul	0.00
15) 4-Methylphenol-d8	8.905	113	50875	18.432 ng/ul	0.00
21) Nitrobenzene-d5	9.369	128	24642	19.366 ng/ul	0.00
24) 2-Nitrophenol-d4	10.092	143	28969	20.182 ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.644	165	47904	19.670 ng/ul	0.00
31) 4-Chloroaniline-d4	11.156	131	67186	18.854 ng/ul	0.00
46) Dimethylphthalate-d6	14.217	166	142375	19.058 ng/ul	0.00
49) Acenaphthylene-d8	14.522	160	189295	20.095 ng/ul	0.00
54) 4-Nitrophenol-d4	15.051	143	18798		
60) Fluorene-d10	15.815	176	128235	15.545 ng/ul 19.062 ng/ul	0.00
65) 4,6-Dinitro-2-methylph	15.950	200	-20344	15.564 ng/ul	0.00 0.00
73) Anthracene-d10	17.671	188	197605	19.505 ng/ul	0.00
81) Pyrene-d10	19.951	212	220295	19.771 ng/ul	0.00
92) Benzo(a)pyrene-d12	25.033	264	189517	19.7/1 mg/ul	
32) Denzo(d)pyrene dzz	20.000	204	103317	19.214 lig/u1	0.00
Target Compounds				Ωv	alue
2) 1,4-Dioxane	3.570	88	7444	6.628 ng/uL	88
5) Pyridine	3.982	79	57190	18.807 ng/ul	98
Benzaldehyde	7.325	77	47324	21.726 ng/ul	92
8) Phenol	7.384	94	67165	18.955 ng/ul	99
10) Bis(2-Chloroethyl)ether	7.601	93	51827	19.333 ng/ul	91
12) 2-Chlorophenol	7.754	128	47880	19.076 ng/ul	99
13) 2-Methylphenol	8.641	108	49128	18.614 ng/ul	97
14) 2,2'-oxybis(1-Chloropr	8.706		75417	19.496 ng/ul	97
16) Acetophenone	9.023		78817	18.461 ng/ul	98
17) N-Nitroso-di-n-propyla	8.993	70	46571	18.982 ng/ul	98
18) 4-Methylphenol	8.970	108	53419	18.928 ng/ul	98
19) Hexachloroethane	9.275	117	20085	18.945 ng/ul	91
22) Nitrobenzene	9.411	77	66998	20.081 ng/ul	97
23) Isophorone	9.928	82	126155	19.462 ng/ul	98
25) 2-Nitrophenol	10.127	139	28759	19.344 ng/ul	99
26) 2,4-Dimethylphenol	10.180	107	60010	19.742 ng/ul	99
<pre>27) Bis(2-Chloroethoxy)met</pre>	10.404	93	70124	19.596 ng/ul	98
29) 2,4-Dichlorophenol	10.674	162	47086	19.641 ng/ul	99
30) Naphthalene	11.067	128	158316	19.302 ng/ul	98
32) 4-Chloroaniline	11.179	127	67977	19.002 ng/ul	98
33) Hexachlorobutadiene	11.332	225	30658	18.541 ng/ul	96
34) Caprolactam	11.943	113	17773	18.858 ng/ul	93
35) 4-Chloro-3-methylphenol	12.301	107	56183	19.510 ng/ul	97
±				/-	###

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Compound	R.T.	QIon	Response	Conc Units Dev	(Min)
36) 2-Methylnaphthalene	12.660	142	107443	19.259 ng/ul	97
37) 1-Methylnaphthalene	12.877	142	108544	18.912 ng/ul	99
39) 1,2,4,5-Tetrachloroben	13.024	216	60595	19.880 ng/ul	98
40) Hexachlorocyclopentadiene	12.989	237	22677	18.407 ng/ul	97
41) 2,4,6-Trichlorophenol	13.271	196	37439	19.573 ng/ul	98
42) 2,4,5-Trichlorophenol	13.353	196	36910	18.427 ng/ul	98
43) 1,1 -Biphenyl	13.659	154	144133	19.876 ng/ul	98
44) 2-Chloronaphthalene	13.706	162	114484	19.846 ng/ul	99
45) 2-Nitroaniline	13.917	65	41845	20.960 ng/ul	93
47) Dimethylphthalate	14.264	163	143199	18.937 ng/ul	100
48) 2,6-Dinitrotoluene	14.405	165	31092	19.575 ng/ul	97
50) Acenaphthylene	14.552	152	182513	19.610 ng/ul	99
51) 3-Nitroaniline	14.740	138	32686	20.819 ng/ul	90
52) Acenaphthene	14.887	153	120619	19.651 ng/ul	96
53) 2,4-Dinitrophenol	14.963	184	15687	17.868 ng/ul#	86
55) 4-Nitrophenol	15.063	109	17124	16.324 ng/ul	90
56) Dibenzofuran	15.221	168	170415	19.249 ng/ul	96
57) 2,4-Dinitrotoluene	15.198	165	42717	18.829 ng/ul#	98
58) 2,3,4,6-Tetrachlorophenol	15.451	232	27045	17.194 ng/ul	97
59) Diethylphthalate	15.615	149	152695	19.238 ng/ul	99
61) Fluorene '	15.868	166	136294	19.219 ng/ul	98
62) 4-Chlorophenyl-phenyle	15.850	204	71509		94
63) 4-Nitroaniline	15.903	138	33609m (21.997 ng/ul	
66) 4,6-Dinitro-2-methylph	15.962	198	20085	15.933 ng/ul#	97
67) N-Nitrosodiphenylamine	16.067	169	121461	20.029 ng/ul	98
68) 4-Bromophenyl-phenylether	16.749	248	44001	19.381 ng/ul	92
69) Hexachlorobenzene	16.878	284	44881	19.387 ng/ul	97
70) Atrazine	17.008	200	51008	20.014 ng/ul	96
71) Pentachlorophenol	17.231	266	19394m	18.906 ng/ul	
72) Phenanthrene	17.619	178	229383	19.612 ng/ul	99
74) Anthracene	17.707	178	228598	19.680 ng/ul	98
75) 1,2,3,4-Tetrachloroben	13.629	216	62619	20.267 ng/uL	96
76) Pentachlorobenzene	15.139	250	58013	20.151 ng/uL	98
77) Carbazole	17.983	167	205804	20.185 ng/ul	99
78) Di-n-butylphthalate	18.500	149	262080	19.935 ng/ul	99
80) Fluoranthene	19.616	202		20.267 ng/ul	97
82) Pyrene	19.981	202	268419	20.051 ng/ul	97
83) Butylbenzylphthalate	20.838			19.874 ng/ul	98
84) 3,3'-Dichlorobenzidine	21.761	252	82525	19.248 ng/ul	98
85) Benzo(a)anthracene	21.855	228	246182	19.710 ng/ul	99
86) Bis(2-ethylhexyl)phtha	21.714	149	160229	20.007 ng/ul	99
87) Chrysene	21.925	228	232950	19.414 ng/ul	99
89) Di-n-octyl phthalate	22.971	149	272445	20.363 ng/ul	100
90) Benzo(b)fluoranthene	24.187	252	237384	19.046 ng/ul	99
91) Benzo(k)fluoranthene	24.252	252	227395	19.442 ng/ul	99
93) Benzo(a)pyrene 94) Indeno(1,2,3-cd)pyrene	25.110	252	229791	19.325 ng/ul	97
95) Dibenzo(a,h)anthracene	29.187	276	255840	19.227 ng/ul	99
96) Benzo(g,h,i)perylene	29.240	278	217201	19.241 ng/ul	97
	30.421	276	215136	19.217 ng/ul	98

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/06/2021 Supervised By :mohammad ahmed 12/07/2021

12441

Instrument : BNA_G **LabSampleld :** SSTDCCC020EC

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