Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112421\

Data File: BN017571.D

Acq On : 23 Nov 2021 17:57

Operator : CG/JU Sample : PB140870BS

Misc

ALS Vial : 12 Sample Multiplier: 1

Quant Time: Nov 24 00:58:24 2021

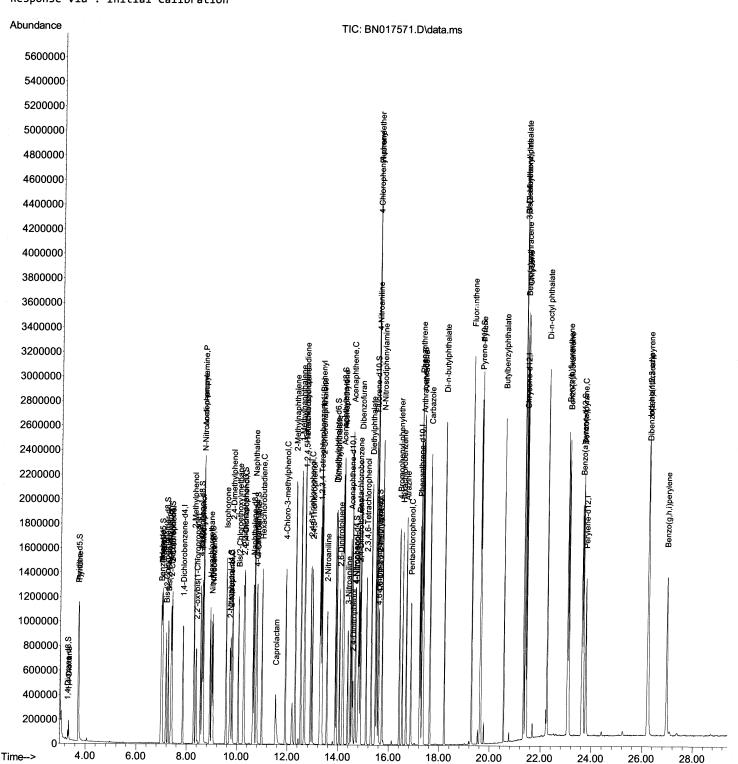
Quant Title : SVOA CALIBRATION

QLast Update : Mon Nov 22 16:16:36 2021 Response via : Initial Calibration



Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/24/2021 Supervised By :mohammad ahmed 11/26/2021



SFAM-EPA-BN112221.M Wed Nov 24 01:16:17 2021

Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112421\

Data File : BN017571.D

Acq On : 23 Nov 2021 17:57

Operator : CG/JU Sample : PB140870BS

Misc

ALS Vial : 12 Sample Multiplier: 1

Quant Time: Nov 24 00:58:24 2021

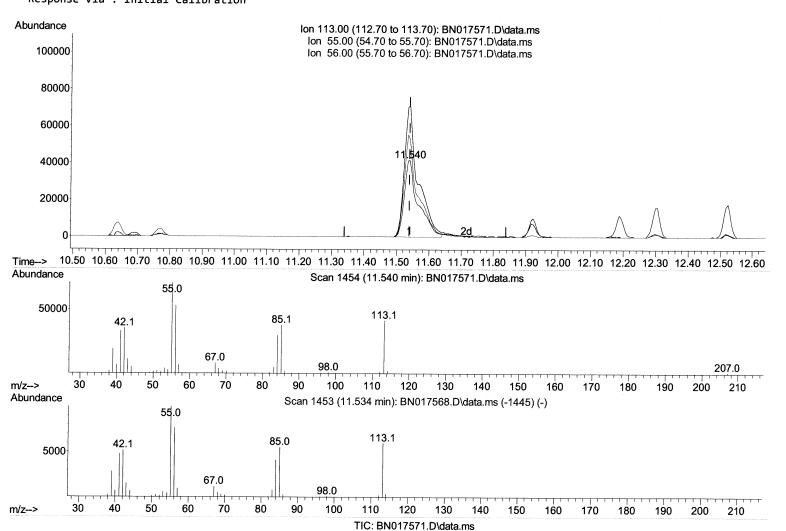
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN112221.M

Quant Title : SVOA CALIBRATION QLast Update : Mon Nov 22 16:16:36 2021 Response via : Initial Calibration



Manual IntegrationsAPPROVED

Reviewed By: Jagrut Upadhyay 11/24/2021 Supervised By: mohammad ahmed 11/26/2021



(34) Caprolactam

11.540min (+ 0.000) 26.85 ng/ul

response	125533	
Ion	Ежр%	Act%
113.00	100.00	100.00
55.00	176.80	170.33
56.00	129.90	128.01
0.00	0.00	0.00

Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112421\

Data File : BN017571.D

Acq On : 23 Nov 2021 17:57

Operator : CG/JU Sample : PB140870BS

Misc

ALS Vial : 12 Sample Multiplier: 1

Quant Time: Nov 24 00:58:24 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN112221.M

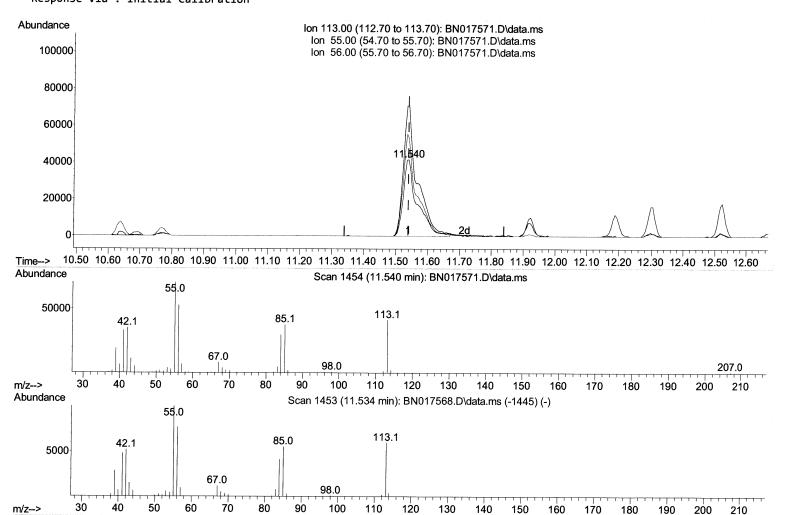
Quant Title : SVOA CALIBRATION

QLast Update : Mon Nov 22 16:16:36 2021 Response via : Initial Calibration

Instrument:
BNA_N
ClientSampleId:
SLCS870

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/24/2021 Supervised By :mohammad ahmed 11/26/2021



TIC: BN017571.D\data.ms

(34) Caprolactam

11.540min (+ 0.000) 27.22 ng/ul m (1/30/2174

response	127275	
Ion	Ехр%	Act%
113.00	100.00	100.00
55.00	176.80	170.33
56.00	129.90	128.01
0.00	0.00	0 00

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112421\

Data File : BN017571.D

Acq On : 23 Nov 2021 17:57

Operator : CG/JU Sample : PB140870BS

Misc

ALS Vial : 12 Sample Multiplier: 1

Quant Time: Nov 24 00:58:24 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN112221.M

Quant Title : SVOA CALIBRATION
QLast Update : Mon Nov 22 16:16:36 2021

Response via : Initial Calibration

Instrument : BNA_N

ClientSampleId :

SLCS870

Manual IntegrationsAPPROVED

Reviewed By: Jagrut Upadhyay 11/24/2021 Supervised By: mohammad ahmed 11/26/2021

Compound	R.T.	QIon	Response	Conc Ui	nits Dev	v(Min)	
Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.846	152	254076	20 000	ng/ul	0.00	
20) Naphthalene-d8	10.640				ng/ul	0.00	
38) Acenaphthene-d10	14.475			20.000	ng/ul	0.00	
64) Phenanthrene-d10	17.222				ng/ul	0.00	
79) Chrysene-d12	21.410				ng/ul	0.00	
88) Perylene-d12	23.774				ng/ul	0.00	
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.305	96	36685	5.160	ng/uL	0.00	
4) Pyridine-d5	3.705	84	482890	25.885	ng/ul	0.00	
7) Phenol-d5	7.005	99	641553		ng/ul	0.00	
<pre>9) Bis-(2-Chloroethyl)eth</pre>	7.175	67	392231	27.061	ng/ul	0.00	
11) 2-Chlorophenol-d4	7.375	132	496888	28.372	ng/ul	0.00	
15) 4-Methylphenol-d8	8.546	113	487930	26.818	ng/ul	0.00	
21) Nitrobenzene-d5	8.999	128	238019	30.326	ng/ul	0.00	
24) 2-Nitrophenol-d4	9.716	143	240636	30.989	ng/ul	0.00	
28) 2,4-Dichlorophenol-d3	10.257	165	445145	29.118	ng/ul	0.00	
31) 4-Chloroaniline-d4	10.769	131	558340	25.434	ng/ul	0.00	
46) Dimethylphthalate-d6	13.887	166	1162639	28.011		0.00	
49) Acenaphthylene-d8	14.169	160	1558188	29.239	_	0.00	
54) 4-Nitrophenol-d4	14.663	143	199228	28.365		0.00	
60) Fluorene-d10	15.469	176	968880	27.675		0.00	
65) 4,6-Dinitro-2-methylph	15.581	200	149323	27.937		0.00	
73) Anthracene-d10	17.322	188	1392750	29.118		0.00	
81) Pyrene-d1092) Benzo(a)pyrene-d12	19.616 23.621	212 264	1468979 1338446	26.870 29.231		0.00 0.00	
Target Compounds					_		
2) 1,4-Dioxane	3.340	00	74467	10 600	-	alue	
5) Pyridine	3.728	88 79	74467	10.608	-	97	
6) Benzaldehyde	6.975	73 77	496419 367781	26.442		90	
8) Phenol	7.034	94	664146	29.974	_	91	
10) Bis(2-Chloroethyl)ether	7.263	93	530500	28.351 27.618		97 100	
12) 2-Chlorophenol	7.405	128	514216	28.493		100 96	
13) 2-Methylphenol	8.281	108	488356	27.678		90 97	
14) 2,2'-oxybis(1-Chloropr	8.375	45	705295	27.236		97 97	
16) Acetophenone	8.657	105	745403	27.396		94	
17) N-Nitroso-di-n-propyla	8.652	70	389524	27.468		94	
18) 4-Methylphenol	8.610	108	516530	27.327		99	
19) Hexachloroethane	8.928	117	217147	28.477		98	
22) Nitrobenzene	9.040	77	588636	29.676		96	
23) Isophorone	9.563	82	1073998	28.929		97	
25) 2-Nitrophenol	9.751	139	260799	30.448	_	91	
26) 2,4-Dimethylphenol	9.816	107	557367	28.137		93	
<pre>27) Bis(2-Chloroethoxy)met</pre>	10.051	93	676094	27.979		100	
29) 2,4-Dichlorophenol	10.287	162	449753	29.175		97	
30) Naphthalene	10.687	128	1573970	28.398	-	100	
32) 4-Chloroaniline	10.793	127	565645	25.395		100	
33) Hexachlorobutadiene	10.987	225	291520	28.750		99	
34) Caprolactam	11.540	113	127275m >	27.224			70
35) 4-Chloro-3-methylphenol	11.922	107	481658	27.772		95	<i>0</i> 4

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112421\

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Operator : CG/JU Sample : PB140870BS

Misc

ALS Vial : 12 Sample Multiplier: 1

Quant Time: Nov 24 00:58:24 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN112221.M

Quant Title : SVOA CALIBRATION QLast Update : Mon Nov 22 16:16:36 2021 Response via : Initial Calibration Instrument : BNA_N ClientSampleId : SLCS870

Manual IntegrationsAPPROVED

Reviewed By: Jagrut Upadhyay 11/24/2021 Supervised By: mohammad ahmed 11/26/2021

Compound	R.T.	QIon	Response	Conc Units Dev(Min)
36) 2-Methylnaphthalene	12.298	142	1016290	27.475 ng/ul	100
37) 1-Methylnaphthalene	12.522	142	1022936	27.311 ng/ul	99
39) 1,2,4,5-Tetrachloroben	12.675	216	497755	29.397 ng/ul	98
40) Hexachlorocyclopentadiene	12.663	237	300746	27.693 ng/ul	93
41) 2,4,6-Trichlorophenol	12.910	196	318293	30.770 ng/ul	98
42) 2,4,5-Trichlorophenol	12.975	196	342871	30.397 ng/ul	99
43) 1,1'-Biphenyl	13.316	154	1316712	29.245 ng/ul	100
44) 2-Chloronaphthalene	13.351	162	987626	28.848 ng/ul	95
45) 2-Nitroaniline	13.551	65	290790	30.688 ng/ul	96
47) Dimethylphthalate	13.934	163	1151761	27.995 ng/ul	99
48) 2,6-Dinitrotoluene	14.051	165	238464	30.884 ng/ul	93
50) Acenaphthylene	14.198	152	1554024	28.613 ng/ul	98
51) 3-Nitroaniline	14.375	138	220036	26.917 ng/ul	94
52) Acenaphthene	14.539	153	999065	28.530 ng/ul	99
53) 2,4-Dinitrophenol	14.581	184	101886	26.250 ng/ul	99
55) 4-Nitrophenol	14.681	109	170011	27.752 ng/ul#	81
56) Dibenzofuran	14.875	168	1422244	28.292 ng/ul	97
57) 2,4-Dinitrotoluene	14.834	165	324451	29.654 ng/ul	93
58) 2,3,4,6-Tetrachlorophenol	15.104	232	242423	29.048 ng/ul	99
59) Diethylphthalate	15.304	149	1158414	28.114 ng/ul	99
61) Fluorene	15.528	166	1046299	27.727 ng/ul	97
62) 4-Chlorophenyl-phenyle	15.522	204	513698	27.895 ng/ul	92
63) 4-Nitroaniline	15.534	138	212080	30.771 ng/ul	88
66) 4,6-Dinitro-2-methylph	15.598	198	155151	28.912 ng/ul#	89
67) N-Nitrosodiphenylamine	15.734	169	916888	29.557 ng/ul	97
68) 4-Bromophenyl-phenylether	16.416	248	294659	28.999 ng/ul#	91
69) Hexachlorobenzene	16.533	284	317159	29.289 ng/ul#	90
70) Atrazine	16.686	200	306480	27.976 ng/ul	99
71) Pentachlorophenol	16.875	266	178908	29.056 ng/ul	97
72) Phenanthrene	17.263	178	1647218	29.336 ng/ul	99
74) Anthracene	17.357	178	1611830	28.947 ng/ul	100
75) 1,2,3,4-Tetrachloroben	13.281	216	509865	31.600 ng/uL	98
76) Pentachlorobenzene	14.804	250	414879	28.278 ng/uL	98
77) Carbazole	17.622	167	1447866	30.297 ng/ul	99
78) Di-n-butylphthalate	18.198	149	1791996	30.857 ng/ul	99
80) Fluoranthene 82) Pyrene	19.280	202	1799285	27.360 ng/ul	99
83) Butylbenzylphthalate	19.645	202	1775193	26.964 ng/ul	98
84) 3,3'-Dichlorobenzidine	20.551	149	768977	31.994 ng/ul	94
85) Benzo(a)anthracene	21.327	252 228	481224	29.375 ng/ul	99
86) Bis(2-ethylhexyl)phtha	21.392 21.333	228 149	1629713	28.568 ng/ul	98 97
			1149687	33.050 ng/ul#	
87) Chrysene 89) Di-n-octyl phthalate	21.445	228	1597265	28.494 ng/ul	100
90) Benzo(b)fluoranthene	22.251 23.057	149 252	1976485 1672188	29.880 ng/ul 27.767 ng/ul	100 98
91) Benzo(k)fluoranthene	23.104	252	1579648	27.787 ng/ul 27.780 ng/ul	100
93) Benzo(a)pyrene	23.668	252	1627522	28.464 ng/ul	
94) Indeno(1,2,3-cd)pyrene	26.221	276	1822089	30.690 ng/ul	100 100
95) Dibenzo(a,h)anthracene	26.239	278	1585923	31.450 ng/ul	98
96) Benzo(g,h,i)perylene	26.962	276	1589948	31.333 ng/ul	97
(8),-/k					

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