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

Data File : BN017568.D

Acq On : 23 Nov 2021 16:09

Operator : CG/JU Sample : SSTDCCC020

Misc

ALS Vial : 9 Sample Multiplier: 1

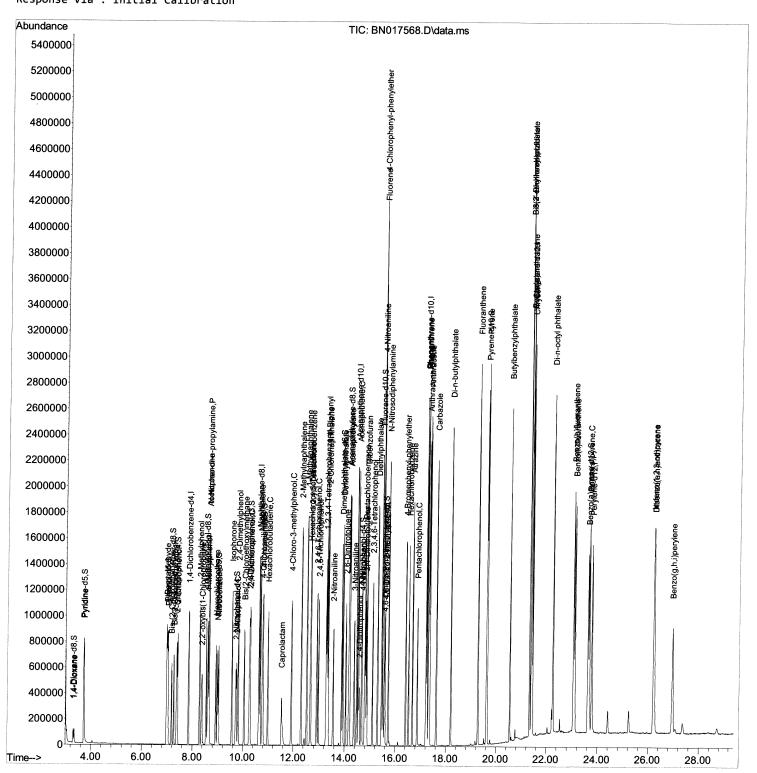
Quant Time: Nov 23 17:20:22 2021

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 LabSampleId : SSTDCCC020

Manual IntegrationsAPPROVED



Quantitation Report (Qedit)

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

Data File : BN017568.D

Acq On : 23 Nov 2021 16:09

Operator : CG/JU Sample : SSTDCCC020

Misc

ALS Vial : 9 Sample Multiplier: 1

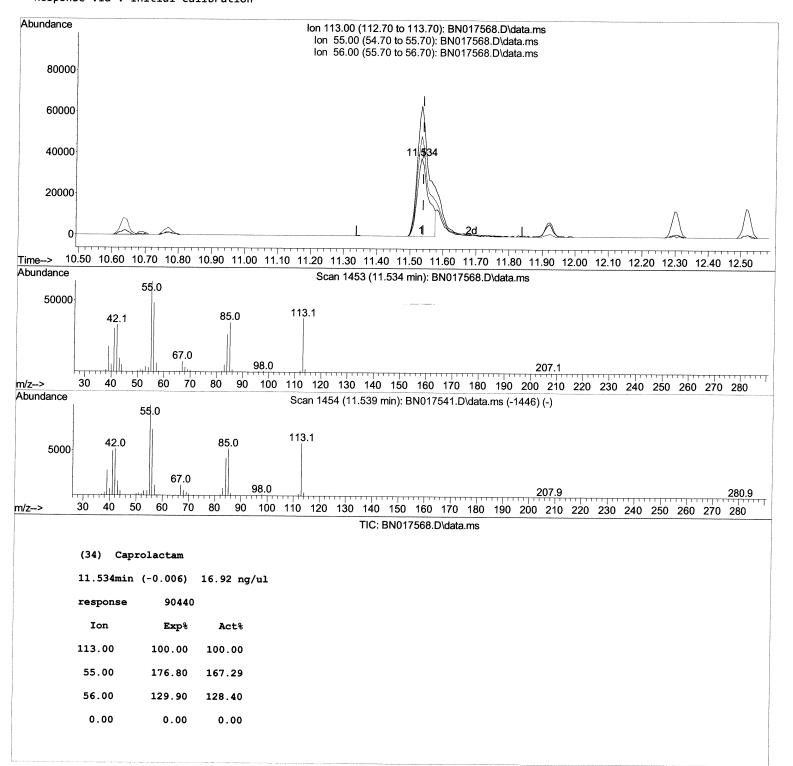
Quant Time: Nov 23 17:20:22 2021

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
LabSampleId :
SSTDCCC020

Manual IntegrationsAPPROVED



Quantitation Report (Qedit)

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

Data File : BN017568.D

Acq On : 23 Nov 2021 16:09

Operator : CG/JU Sample : SSTDCCC020

Misc

ALS Vial : 9 Sample Multiplier: 1

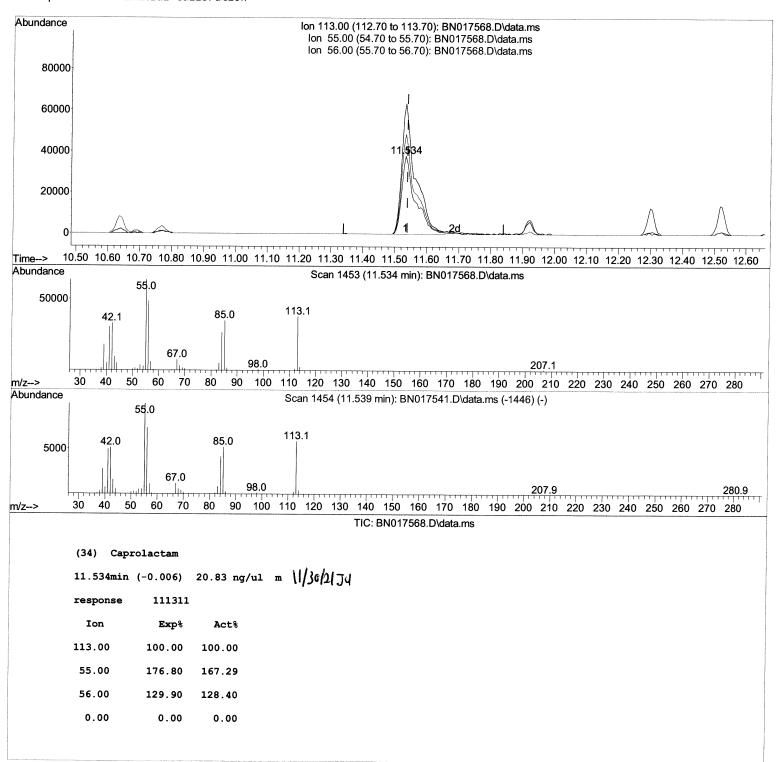
Quant Time: Nov 23 17:20:22 2021

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
LabSampleId :
SSTDCCC020

Manual Integrations APPROVED



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

Data File : BN017568.D

Acq On : 23 Nov 2021 16:09

Operator : CG/JU Sample : SSTDCCC020

Misc

ALS Vial : 9 Sample Multiplier: 1

Quant Time: Nov 23 17:20:22 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 **LabSampleld** : SSTDCCC020

Manual IntegrationsAPPROVED

Compound				Conc Units Dev	
Internal Standards					
 1,4-Dichlorobenzene-d4 	7.846	152	269623	20.000 ng/ul	0.00
20) Naphthalene-d8	10.640			20.000 ng/ul	0.00
38) Acenaphthene-d10	14.475			20.000 ng/ul	0.00
64) Phenanthrene-d10	17.222	188		20.000 ng/ul	0.00
79) Chrysene-d12	21.410			20.000 ng/ul	0.00
88) Perylene-d12	23.774	264	985420	20.000 ng/ul	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.305	96	51869	6.875 ng/uL	0.00
4) Pyridine-d5	3.705	84	349950	17.677 ng/ul	0.00
7) Phenol-d5	7.005	99	449871	18.328 ng/ul	0.00
<pre>9) Bis-(2-Chloroethyl)eth</pre>	7.169	67	277645	18.051 ng/ul	0.00
11) 2-Chlorophenol-d4	7.375	132	344090	18.514 ng/ul	0.00
15) 4-Methylphenol-d8	8.546	113	358554	18.571 ng/ul	0.00
21) Nitrobenzene-d5	8.993	128	173238	19.306 ng/ul	0.00
24) 2-Nitrophenol-d4	9.722	143	179088	20.172 ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.257	165	335642	19.203 ng/ul	0.00
31) 4-Chloroaniline-d4	10.769	131	492239	19.613 ng/ul	0.00
46) Dimethylphthalate-d6	13.887	166	996971	18.221 ng/ul	0.00
49) Acenaphthylene-d8	14.169	160	1310717	18.658 ng/ul	
54) 4-Nitrophenol-d4	14.663	143	180982	19.547 ng/ul	0.00 0.00
60) Fluorene-d10	15.469	176	857772	18.587 ng/ul	
65) 4,6-Dinitro-2-methylph	15.581	200	136041	17.418 ng/ul	0.00 0.00
73) Anthracene-d10	17.322	188	1307305	18.705 ng/ul	0.00
81) Pyrene-d10	19.616	212	1426477	19.258 ng/ul	0.00
92) Benzo(a)pyrene-d12	23.621	264	975310	18.826 ng/ul	0.00
Target Compounds				Qva	luo
2) 1,4-Dioxane	3.340	88	52589	7.060 ng/uL	98
5) Pyridine	3.728	79	357075	17.923 ng/ul	91
6) Benzaldehyde	6.981	77	296344	22.759 ng/ul	93
8) Phenol	7.034	94	454555	18.285 ng/ul	96
10) Bis(2-Chloroethyl)ether	7.263	93	363322	17.824 ng/ul	99
12) 2-Chlorophenol	7.405	128	352995	18.432 ng/ul	95
13) 2-Methylphenol	8.281	108	343481	18.432 ng/ul	93 97
14) 2,2'-oxybis(1-Chloropr	8.369	45	486965	17.720 ng/ul	98
16) Acetophenone	8.657	105	555643	19.244 ng/ul	93
17) N-Nitroso-di-n-propyla	8.652	70	288377	19.163 ng/ul	94
18) 4-Methylphenol	8.610	108	374048	18.648 ng/ul	99
19) Hexachloroethane	8.928	117	147472	18.225 ng/ul	96
22) Nitrobenzene	9.034	77	420346	18.536 ng/ul	95
23) Isophorone	9.563	82	811397	19.116 ng/ul	97
25) 2-Nitrophenol	9.751	139	192806	19.688 ng/ul#	92
26) 2,4-Dimethylphenol	9.816	107	425225	18.776 ng/ul#	96
27) Bis(2-Chloroethoxy)met	10.051	93	499721	18.088 ng/ul	99
29) 2,4-Dichlorophenol	10.287	162	333794	18.939 ng/ul	99 97
30) Naphthalene	10.687	128	1152675	18.190 ng/ul	99
32) 4-Chloroaniline	10.793	127	491936	19.318 ng/ul	99
33) Hexachlorobutadiene	10.793	225	211082	18.208 ng/ul	99
34) Caprolactam	11.534	113	111311m >	20.825 ng/ul >	11/30/21 74
35) 4-Chloro-3-methylphenol	11.922	107	385332	19.433 ng/ul	97

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

Data File : BN017568.D

Acq On : 23 Nov 2021 16:09

Operator : CG/JU Sample : SSTDCCC020

Misc

ALS Vial : 9 Sample Multiplier: 1

Quant Time: Nov 23 17:20:22 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
LabSampleId:
SSTDCCC020

Manual IntegrationsAPPROVED

Compound	R.T.	QIon	Response	Conc Units Dev(Min)
36) 2-Methylnaphthalene	12.298	142	784003	18.538 ng/ul	99
37) 1-Methylnaphthalene	12.522	142	798948	18.657 ng/ul	97
39) 1,2,4,5-Tetrachloroben	12.675	216	393875	17.646 ng/ul	96
40) Hexachlorocyclopentadiene	12.657	237	229504	16.031 ng/ul	98
41) 2,4,6-Trichlorophenol	12.910	196	259824	19.054 ng/ul	99
42) 2,4,5-Trichlorophenol	12.981	196	275509	18.529 ng/ul	98
43) 1,1'-Biphenyl	13.310	154	1059238	17.847 ng/ul	98
44) 2-Chloronaphthalene	13.351	162	802060	17.772 ng/ul	96
45) 2-Nitroaniline	13.545	65	247157	19.787 ng/ul	88
47) Dimethylphthalate	13.934	163	1012692	18.672 ng/ul	99
48) 2,6-Dinitrotoluene	14.045	165	202192	19.865 ng/ul	88
50) Acenaphthylene	14.198	152	1311766	18.322 ng/ul	100
51) 3-Nitroaniline	14.375	138	221622	20.567 ng/ul	91
52) Acenaphthene	14.539	153	841762	18.235 ng/ul	97
53) 2,4-Dinitrophenol	14.581	184	87363	17.075 ng/ul	97
55) 4-Nitrophenol	14.681	109	156482	19.378 ng/ul	83
56) Dibenzofuran	14.875	168	1180118	17.808 ng/ul	96
57) 2,4-Dinitrotoluene	14.834	165	294285	20.404 ng/ul	96
58) 2,3,4,6-Tetrachlorophenol	15.098	232	216249	19.656 ng/ul	96
59) Diethylphthalate	15.298	149	1029365	18.952 ng/ul	99
61) Fluorene	15.528	166	919297	18.480 ng/ul	100
62) 4-Chlorophenyl-phenyle	15.522	204	445449	18.350 ng/ul	93
63) 4-Nitroaniline	15.534	138	216349	23.812 ng/ul#	86
66) 4,6-Dinitro-2-methylph	15.598	198	139477	17.788 ng/ul#	90
67) N-Nitrosodiphenylamine	15.733	169	836608	18.457 ng/ul	98
68) 4-Bromophenyl-phenylether	16.416	248	266154	17.926 ng/ul	94
69) Hexachlorobenzene	16.533	284	279433	17.660 ng/ul#	85
70) Atrazine	16.681	200	308171	19.251 ng/ul	97
71) Pentachlorophenol	16.875	266	167434	18.610 ng/ul	99
72) Phenanthrene	17.263	178	1526661	18.607 ng/ul	98
74) Anthracene	17.351	178	1526733	18.765 ng/ul	100
75) 1,2,3,4-Tetrachloroben	13.275	216	406016	17.221 ng/uL	95
76) Pentachlorobenzene	14.804	250	373254	17.411 ng/uL	93
77) Carbazole	17.622	167	1386723	19.859 ng/ul	99
78) Di-n-butylphthalate 80) Fluoranthene	18.198	149	1723416	20.309 ng/ul	99
82) Pyrene	19.280	202	1702052	19.102 ng/ul	100
	19.645	202	1702080	19.081 ng/ul	98
83) Butylbenzylphthalate84) 3,3'-Dichlorobenzidine	20.551	149	738481	22.677 ng/ul	94
85) Benzo(a)anthracene	21.327 21.392	252 228	462695	20.845 ng/ul	99
86) Bis(2-ethylhexyl)phtha	21.332	226 149	1449994	18.760 ng/ul	99
87) Chrysene	21.445	228	1063215	22.558 ng/ul# 18.326 ng/ul	99
89) Di-n-octyl phthalate	22.251	228 149	1391835 1727715	_	98
90) Benzo(b)fluoranthene	23.057	252	1272549	23.085 ng/ul 18.676 ng/ul	100 97
91) Benzo(k)fluoranthene	23.104	252	1140179	17.722 ng/ul	95
93) Benzo(a)pyrene	23.668	252	1172222	18.120 ng/ul	
94) Indeno(1,2,3-cd)pyrene	26.215	276	1216523	18.110 ng/ul	97 100
95) Dibenzo(a,h)anthracene	26.233	278	1016261	17.812 ng/ul	97
96) Benzo(g,h,i)perylene	26.962	276	1010201	17.512 ng/ul 17.592 ng/ul	95
(6),,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,					

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