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

Data File : BN017566.D

Acq On : 23 Nov 2021 14:57

Operator : CG/JU Sample : M4702-03MS

Misc

ALS Vial : 7 Sample Multiplier: 1

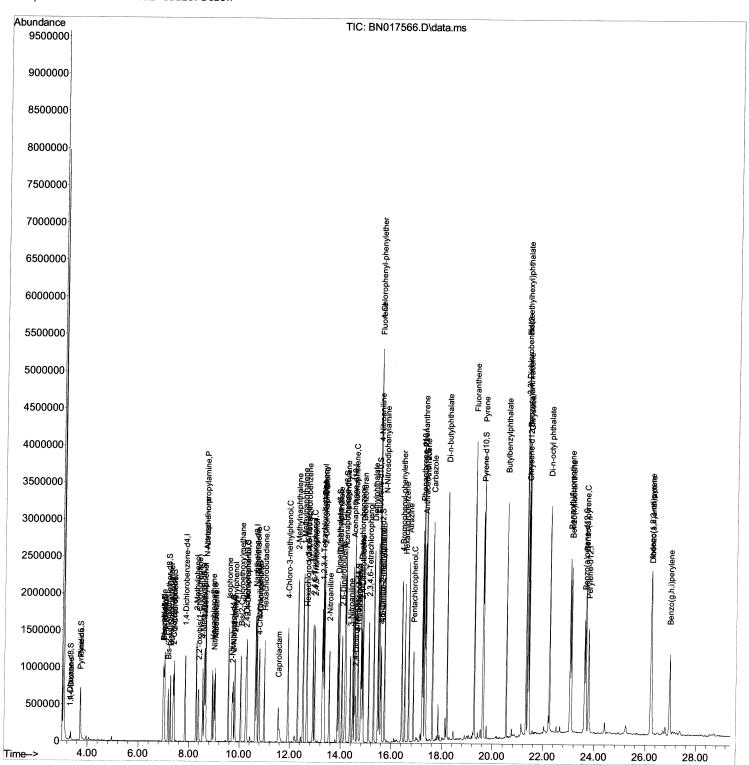
Quant Time: Nov 23 17:12:32 2021

Quant Title : SVOA CALIBRATION

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



Manual IntegrationsAPPROVED



Quantitation Report (Qedit)

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

Data File : BN017566.D

Acq On : 23 Nov 2021 14:57

Operator : CG/JU Sample : M4702-03MS

Misc

ALS Vial : 7 Sample Multiplier: 1

Quant Time: Nov 23 17:12:32 2021

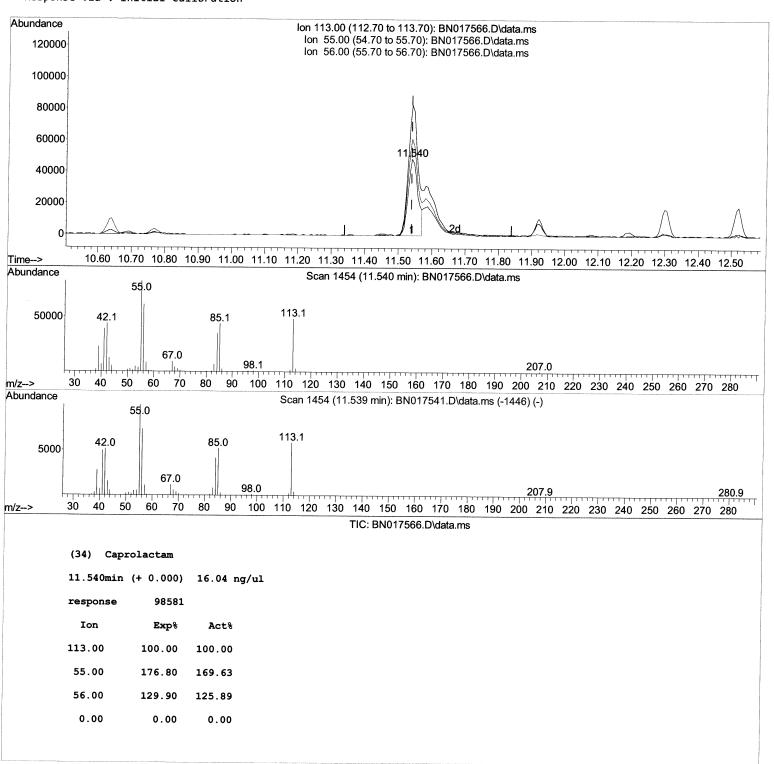
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN112221.M

Quant Title : SVOA CALIBRATION

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



Manual IntegrationsAPPROVED



Quantitation Report (Qedit)

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

Data File: BN017566.D

Acq On : 23 Nov 2021 14:57

Operator : CG/JU Sample : M4702-03MS

Misc

ALS Vial : 7 Sample Multiplier: 1

Quant Time: Nov 23 17:12:32 2021

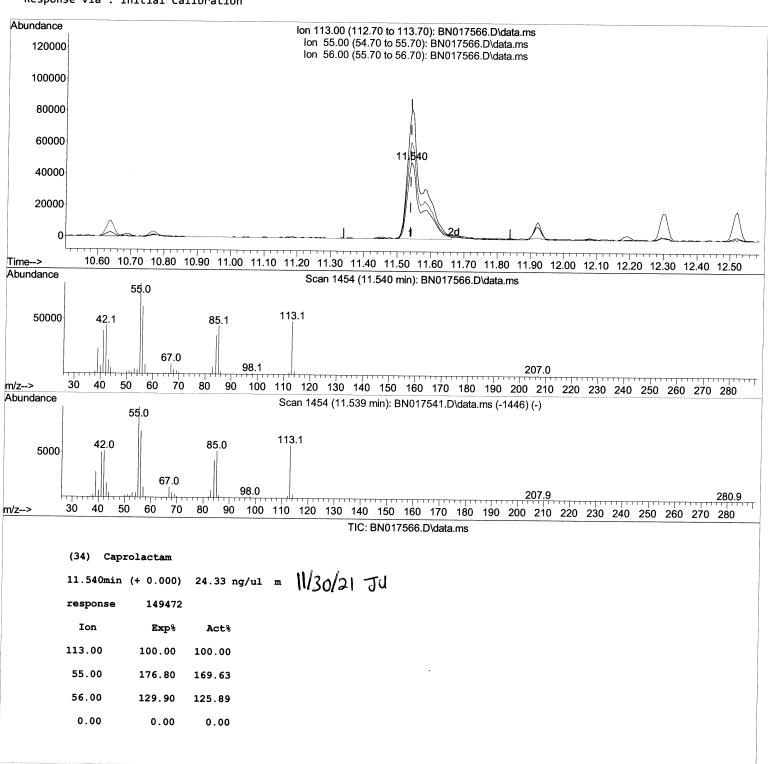
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN112221.M

Quant Title : SVOA CALIBRATION

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



Manual Integrations APPROVED



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

Data File : BN017566.D

Acq On : 23 Nov 2021 14:57 Operator : CG/JU Sample : M4702-03MS

Misc

ALS Vial : 7 Sample Multiplier: 1

Quant Time: Nov 23 17:12:32 2021

 $\label{thm:lem1_BNA_N\ethods\SFAM-EPA-BN112221.M} Quant \ \mbox{Methods\SFAM-EPA-BN112221.M}$

Quant Title : SVOA CALIBRATION

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

Instrument : BNA_N ClientSampleId : DBLN9MS

Manual IntegrationsAPPROVED

	Compound	R.T.	QIon	Response	Conc Units Dev	(Min)
Inte	rnal Standards					
1)	1,4-Dichlorobenzene-d4	7.840	152	306076	20.000 ng/ul	0.00
20)	Naphthalene-d8	10.640	136		20.000 ng/ul	0.00
38)	Acenaphthene-d10	14.475	164		20.000 ng/ul	0.00
64)	Phenanthrene-d10	17.222	188	1580444	20.000 ng/ul	0.00
79)	Chrysene-d12	21.410	240	1111761	20.000 ng/ul	0.00
88)	Perylene-d12	23.774	264		20.000 ng/ul	0.00
Syst	em Monitoring Compounds					
3)	1,4-Dioxane-d8	3.311	96	26369	3.079 ng/uL	0.00
4)	Pyridine-d5	3.711	84	236434	10.521 ng/ul	0.00
7)	Phenol-d5	7.005	99	501015	17.981 ng/ul	0.00
9)	Bis-(2-Chloroethyl)eth	7.169	67	316134	18.105 ng/ul	0.00
11)	2-Chlorophenol-d4	7.375	132	390146	18.492 ng/ul	0.00
15)	4-Methylphenol-d8	8.546	113	367321	16.759 ng/ul	0.00
	Nitrobenzene-d5	8.993	128	192896	18.700 ng/ul	0.00
24)	2-Nitrophenol-d4	9.716	143	199058	19.504 ng/ul	0.00
28)	2,4-Dichlorophenol-d3	10.257	165	390020	19.411 ng/ul	0.00
31)	4-Chloroaniline-d4	10.769	131	476453	16.514 ng/ul	0.00
46)	Dimethylphthalate-d6	13.887	166	1218234	20.169 ng/ul	0.00
49)	Acenaphthylene-d8	14.169	160	1507388	19.437 ng/ul	0.00
	4-Nitrophenol-d4	14.663	143	198898	19.459 ng/ul	0.00
	Fluorene-d10	15.469	176	1012728	19.879 ng/ul	0.00
65)	4,6-Dinitro-2-methylph	15.586	200	161907	19.724 ng/ul	0.00
	Anthracene-d10	17.322	188	1504861	20.486 ng/ul	0.00
81)	Pyrene-d10	19.616	212	1606781	22.767 ng/ul	0.00
92)	Benzo(a)pyrene-d12	23.621	264	1071251	20.769 ng/ul	0.00
Targe	et Compounds				Ova	lue
2)	1,4-Dioxane	3.340	88	59885	7.082 ng/uL	96
5)	Pyridine	3.728	79	335132	14.818 ng/ul	92
6)	Benzaldehyde	6.975	77	332420	22.489 ng/ul	89
	Phenol	7.028	94	632636	22.417 ng/ul	96
10)	Bis(2-Chloroethyl)ether	7.263	93	473513	20.464 ng/ul	100
	2-Chlorophenol	7.405	128	456399	20.993 ng/ul	97
13)	2-Methylphenol	8.281	108	441907	20.790 ng/ul	100
	2,2'-oxybis(1-Chloropr	8.381	45	640170	20.521 ng/ul	96
	Acetophenone	8.657	105	722867	22.054 ng/ul	95
17)	N-Nitroso-di-n-propyla	8.652	70	375305	21.969 ng/ul	93
	4-Methylphenol	8.610	108	487583	21.413 ng/ul	97
	Hexachloroethane	8.928	117	189205	20.597 ng/ul	98
22)	Nitrobenzene	9.040	77	545316	20.918 ng/ul	96
	Isophorone	9.563	82	1072192	21.974 ng/ul	96
25)	2-Nitrophenol	9.751	139	248302	22.056 ng/ul#	89
	2,4-Dimethylphenol	9.810	107	380106	14.600 ng/ul	94
	Bis(2-Chloroethoxy)met	10.051	93	650511	20.483 ng/ul	99
	2,4-Dichlorophenol	10.287	162	437717	21.604 ng/ul	98
	Naphthalene	10.687	128	1494175	20.512 ng/ul	98
•	4-Chloroaniline	10.793	127	571622	19.526 ng/ul	99
	Hexachlorobutadiene	10.987	225	268536	20.150 ng/ul	99
	Caprolactam	11.540	113	149472m >	٠.	11/30/21 J
	4-Chloro-3-methylphenol	11.922	107	518043	22.727 ng/ul	96 111 20121 J

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

Data File : BN017566.D

Acq On : 23 Nov 2021 14:57 Operator : CG/JU

Sample : M4702-03MS

Misc

ALS Vial : 7 Sample Multiplier: 1

Quant Time: Nov 23 17:12:32 2021

 $\label{thm:local_power_power_power} \textbf{Quant Methods: Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN112221.M}$

Quant Title : SVOA CALIBRATION

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

Instrument : BNA_N

ClientSampleId :

DBLN9MS

Manual IntegrationsAPPROVED

Compound	R.T.	QIon	Response	Conc Units Dev	(Min)
36) 2-Methylnaphthalene	12.298	142	1039730	21.386 ng/ul	100
37) 1-Methylnaphthalene	12.522		1043686	21.202 ng/ul	98
39) 1,2,4,5-Tetrachloroben	12.675		522500	21.205 ng/ul	98
40) Hexachlorocyclopentadiene	12.657	237		13.256 ng/ul	97
41) 2,4,6-Trichlorophenol	12.910	196	347386	23.077 ng/ul	99
42) 2,4,5-Trichlorophenol	12.975		382501	23.303 ng/ul	99
43) 1,1'-Biphenyl	13.310	154	1384036	21.124 ng/ul	99
44) 2-Chloronaphthalene	13.351	162	1052684	21.129 ng/ul	97
45) 2-Nitroaniline	13.551	65	331151	24.015 ng/ul	96
47) Dimethylphthalate	13.934	163	1337817	22.345 ng/ul	99
48) 2,6-Dinitrotoluene	14.051	165	270351	24.061 ng/ul	92
50) Acenaphthylene	14.198	152	1714827	21.697 ng/ul	99
51) 3-Nitroaniline	14.375	138	269820	22.682 ng/ul	89
52) Acenaphthene	14.539	153	1120073	21.980 ng/ul	97
53) 2,4-Dinitrophenol	14.581	184	120535	21.340 ng/ul	96
55) 4-Nitrophenol	14.681	109	207868	23.317 ng/ul	84
56) Dibenzofuran	14.875	168	1586161	21.682 ng/ul	97
57) 2,4-Dinitrotoluene	14.834	165	396415	24.897 ng/ul	98
58) 2,3,4,6-Tetrachlorophenol	15.104	232	280997	23.137 ng/ul	98
59) Diethylphthalate	15.304	149	1400669	23.360 ng/ul	99
61) Fluorene	15.528	166	1195652	21.773 ng/ul	100
62) 4-Chlorophenyl-phenyle	15.522	204	573494	21.400 ng/ul	92
63) 4-Nitroaniline	15.534	138	259025	25.825 ng/ul	88
66) 4,6-Dinitro-2-methylph	15.598	198	193371	23.464 ng/ul#	88
67) N-Nitrosodiphenylamine	15.733	169	1089200	22.863 ng/ul	98
68) 4-Bromophenyl-phenylether	16.416	248	360992	23.134 ng/ul	92
69) Hexachlorobenzene	16.533	284	378519	22.761 ng/ul#	89
70) Atrazine	16.686	200	407878	24.243 ng/ul	95
71) Pentachlorophenol	16.875	266	183273	19.381 ng/ul	97
72) Phenanthrene	17.263	178	2020896	23.436 ng/ul	98
74) Anthracene	17.357	178	1922013	22.476 ng/ul	99
75) 1,2,3,4-Tetrachloroben	13.281	216	518560	20.927 ng/uL	97
76) Pentachlorobenzene 77) Carbazole	14.798	250	473024	20.993 ng/uL	95
78) Di-n-butylphthalate	17.622	167	1783652	24.303 ng/ul	99
80) Fluoranthene	18.198	149	2417082	27.101 ng/ul	99
82) Pyrene	19.280	202	2287518	26.945 ng/ul	99
83) Butylbenzylphthalate	19.645	202 149	2212790	26.035 ng/ul	96
84) 3,3'-Dichlorobenzidine	20.551 21.327	252	959885	30.936 ng/ul	95
85) Benzo(a)anthracene	21.327	228	358959 1740293	16.973 ng/ul	92
86) Bis(2-ethylhexyl)phtha	21.332	149		23.631 ng/ul	98
87) Chrysene	21.445	228	1366815 1728071	30.436 ng/ul#	97
89) Di-n-octyl phthalate	22.251	149	2076503	23.880 ng/ul	99
90) Benzo(b)fluoranthene	23.057			27.868 ng/ul	100
91) Benzo(k)fluoranthene	23.104	252 252	1665600 1472691	24.553 ng/ul 22.991 ng/ul	97 99
93) Benzo(a)pyrene	23.668	252	1514986	23.522 ng/ul	99 100
94) Indeno(1,2,3-cd)pyrene	26.215	276	1700585	_	100
95) Dibenzo(a,h)anthracene	26.239	278	1419534	25.428 ng/ul 24.990 ng/ul	98 00
96) Benzo(g,h,i)perylene	26.968	276	1338546	23.417 ng/ul	98 96
, 23.00 (8).17-7-0-1-0-1-0-1-0-1-0-1-0-1-0-1-0-1-0-1				•	

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